

FORMER MUGLER SHORING INC. SITE
SITE No. C203052
2401 THIRD AVENUE, BRONX, NY
Block 2319 Lot 2

REMEDIAL ACTION WORK PLAN

May 2016
Revised August 2016

Prepared for:
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CERTIFICATIONS

I _____certify that I am currently a NYS registered professional engineer and that this Remedial Action Work Plan 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).

NYS Professional Engineer #

Date

Signature

It is a violation of Article 145 of New York State Education Law for any person to alter this document in any way without the express written verification of adoption by any New York State licensed engineer in accordance with Section 7209(2), Article 145, New York State Education Law.

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LIST OF ACRONYMS

| Acronym | Definition |
|----------------|---|
| AMC | AMC Engineering |
| AWQS | Ambient Water Quality Standards |
| BCA | Brownfield Cleanup Agreement |
| BCP | Brownfield Cleanup Program |
| BTEX | Benzene, Toluene, Ethylbenzene and Xylene |
| CQMP | Construction Quality Management Plan |
| DUSR | Data Usability Statement Report |
| EBC | Environmental Business Consultants |
| FER | Final Engineering Report |
| HDPE | High Density Polyethylene |
| IRM | Interim Remedial Measure |
| NYC | New York City |
| NYCDEP | New York City Department of Environmental Protection |
| NYSDEC | New York State Department of Environmental Conservation |
| NYSDOH | New York State Department of Health |
| PS | Public School |
| PVC | Polyvinyl Chloride |
| RAO | Remedial Action Objectives |
| RAWP | Remedial Action Work Plan |
| RI | Remedial Investigation |
| RSCOs | Recommended Site Cleanup Objectives |
| SCG | Standards, Criteria, and Guidelines |
| SMMP | Soil/Materials Management Plan |
| SMP | Site Management Plan |
| SSDS | Sub-slab Depressurization System |
| SWPPP | Stormwater Pollution Prevention Plan |
| SVOCs | Semi-Volatile Organic Compounds |
| USEPA | United States Environmental Protection Agency |
| UST | Underground Storage Tank |
| VOCs | Volatile Organic Compounds |

EXECUTIVE SUMMARY

Site Description/Physical Setting/Site History

This Remedial Action Work Plan (RAWP) was prepared on behalf 2401 3rd Ave Associates Property LLC for the property known as the Former Mugler Shoring Inc. Site, located at 2401 Third Avenue, Bronx, New York. An application for acceptance into the New York State Brownfield Cleanup Program (BCP) was submitted to the New York State Department of Environmental Conservation (NYSDEC) on May 6, 2015. The application was determined to be complete on May 15, 2015. On June 29, 2015, the DEC notified 2401 3rd Ave Associates Property LLC that the Site had been accepted to the BCP (Site No. C203052) with the applicant defined as a Volunteer. The Brownfield Cleanup Agreement was executed on September 9, 2015.

A Restricted Residential use is proposed for the property. When completed, the Site will be redeveloped with 2 new 25-story residential buildings. Refer to the Brownfield Cleanup Program (BCP) application for additional details.

The street address for the Site is 2401 Third Avenue, Bronx, NY (**Figure 1**). The Site is located in the City of New York in the Port Morris neighborhood of the Borough of the Bronx. The Site is comprised of a single tax parcel covering 61,034.98 (1.4 acres). The subject property is located in the City of New York and Borough of the Bronx (Bronx County). The lot is located on the west side of Third Avenue and is identified as Block 2319 Lot 2 on the NY City tax map. The property has 159 feet of street frontage on Third Avenue and approximately 346 ft of frontage along the north side of the Harlem River (**Figure 2**).

The Site is currently owned by the Volunteer, 2401 3rd Ave Associates Property LLC. The property is currently vacant but was must recently used by a shoring company The Volunteer purchased the property in January 2015.

Summary of the Remedial Investigation

A Remedial Investigation was completed at the Site in November 2015 through December 2015 and documented in a Remedial Investigation Report dated February 2016. The goals of the Remedial Investigation were to define the nature and extent of contamination in soil,

groundwater and any other impacted media; to identify the source(s) of the contamination; to assess the impact of the contamination on public health and/or the environment; and to provide information to support the development of a Remedial Work Plan to address the contamination.

Activities completed under the RI:

- Sampling for non-petroleum contaminants such as pesticides, PCBs and metals in soil and groundwater including the analysis of soil and groundwater samples
- Soil sampling and analysis for petroleum compounds in soil samples from 23 soil boring locations;
- The installation of 12 groundwater monitoring wells
- The collection and analysis of groundwater samples for petroleum compounds;
- The collection of analysis of soil gas samples for VOCs from 8 soil gas sampling locations and 5 subslab vapor sampling locations.

The results of sampling performed during the RI, identified petroleum related contaminants in soil to a depth of 8 feet in the vicinity of a suspect 550 gallon underground storage tank (UST) that was previously abandoned in place.

Historic fill material has been identified across the Site to depths as great as 7 feet below grade. The historic fill material contains semi-volatile organic compounds and metals including arsenic, barium, chromium, copper, lead, mercury, nickel, silver and zinc above unrestricted, restricted residential or restricted commercial use SCOs.

Groundwater impacts from petroleum were not reported during the RI but were reported previously in the vicinity of the suspect underground gasoline tank. SVOC detections above groundwater standards were limited to those polynuclear aromatic hydrocarbons (PAHs) with a 2 per trillion standard. Exceedances of the part per trillion standards for PAHs were reported in well MW9. SVOCs reported in the parts per trillion range are a function of the laboratories ability to achieve extremely low detection limits and general background conditions. The PCB Arochlor 1254 was also reported above standards in MW9.

Several dissolved metals were detected above standards including sodium, iron and manganese were detected above standards throughout the Site. The metals detected in groundwater are related to brackish conditions and are associated with the Site's proximity to the Harlem River.

With the exception of lighter compounds such as heptane and hexane in several samples, total petroleum-related VOCs were generally low in soil vapor samples across the Site. There did not appear to be any correlation in concentration or distribution of petroleum VOCs in soil vapor with the source area.

CVOCs were also generally low and were reported in all of the soil vapor samples with concentrations ranging from 0.28 µg/m³ to 26.06 µg/m³.

Qualitative Human Health Exposure Assessment

The qualitative exposure assessment identified potential completed routes of exposure to construction workers and remediation workers through inhalation, ingestion and dermal contact of petroleum VOCs, SVOCs, CVOCs, pesticides and heavy metals during excavation activities. The Health and Safety Plan prepared for the site identifies such exposures and provides instructions for on-site workers to minimize potential exposure.

Potential completed routes of exposure were also identified for future occupants of the new building including commercial retail workers, residents and visitors to SVOCs and heavy metals in soil through ingestion and dermal contact if these contaminants were to remain in exposed soils at the Site.

Off-Site residents could also be exposed to dust or vapors during the excavation of impacted soil. A site specific Community Air Monitoring Plan has been developed to identify and minimize the potential for off-site exposure to residents through continuous air monitoring during excavation activity.

Potential environmental impacts through the groundwater to surface water discharge were considered unlikely based on the concentrations of VOCs in groundwater and the absence of any evidence of migration toward the River.

Summary of the Remedy

The remedy recommended for the Site is a Track 4 alternative (Alternative 2) which consists of the excavation of petroleum impacted soil to a depth of 9 feet within the UST source area (USTs 1 & 2), excavation and disposal of historic fill within the planned landscape areas with soil above Restricted Residential SCOs and capping the entire Site with the building foundations, concrete walkways / driveways or 2 ft of soil meeting Restricted Residential SCOs. The remedy will include the following items:

1. Removal of underground storage tanks (USTs);
2. Excavation of soil/fill exceeding groundwater protection SCOs for those VOC parameters in groundwater above standards as listed in **Table 1** to depths as great as 9 feet below grade, or as needed to remove impacted soil in the UST source area (USTs 1 & 2);
3. Screening for indications of contamination (by visual means, odor, and monitoring with PID) of all excavated soil during any intrusive Site work;
4. Excavation and disposal of historic fill materials with parameters above RRSCOs from the buildings basement level foundations;
5. Capping areas of the Site in which soil is present with parameters above Restricted Residential SCOs with the concrete building slab, concrete driveways, concrete pavers or 2 feet of clean soil meeting RRSCOs;
6. Collection and analysis of end-point samples to evaluate the performance of the remedy with respect to attainment of groundwater protection SCOs for VOCs present in groundwater above standards;
7. Appropriate off-Site disposal of all material removed from the Site in accordance with all Federal, State and local rules and regulations for handling, transport, and disposal;
8. Import of materials to be used for backfill and cover in compliance with: (1) chemical limits and other specifications included in **Table 1**, (2) all Federal, State and local rules and regulations for handling and transport of material.
9. Implementation of a Site Management Plan (SMP) for long term maintenance of the Engineering Controls. An Environmental Easement will be filed against the Site to ensure implementation of the SMP.

REMEDIAL ACTION WORK PLAN

1.0 INTRODUCTION

This Remedial Action Work Plan (RAWP) was prepared on behalf 2401 3rd Ave Associates Property LLC for the property known as the Former Mugler Shoring Inc. Site, located at 2401 Third Avenue, Bronx, New York. An application for acceptance into the New York State Brownfield Cleanup Program (BCP) was submitted to the New York State Department of Environmental Conservation (NYSDEC) on May 6, 2015. The application was determined to be complete on May 15, 2015. On June 29, 2015, the DEC notified 2401 3rd Ave Associates Property LLC that the Site had been accepted to the BCP (Site No. C203052) with the applicant defined as a Volunteer. The Brownfield Cleanup Agreement was executed on September 9, 2015.

This RAWP summarizes the nature and extent of contamination as determined from data gathered during the Remedial Investigation (RI), performed between November 20, 2015 and December 2, 2015. It provides an evaluation of a Track 1 cleanup and other applicable Remedial Action alternatives, their associated costs, and the recommended and preferred remedy. The remedy described in this document is consistent with the procedures defined in DER-10 and complies with all applicable standards, criteria and guidance. The remedy described in this document also complies with all applicable Federal, State and local laws, regulations and requirements. The NYSDEC and New York State Department of Health (NYSDOH) have determined that this Site does not pose a significant threat to human health and the environment. The RI for this Site did not identify fish and wildlife resources.

A formal Remedial Design document will not be prepared.

1.1 SITE LOCATION AND DESCRIPTION

The street address for the Site is 2401 Third Avenue, Bronx, NY (**Figure 1**). The Site is located in the City of New York in the Port Morris neighborhood of the Borough of the Bronx. The Site is comprised of a single tax parcel covering 61,034.98 (1.4 acres). The subject property is located in the City of New York and Borough of the Bronx (Bronx County). The lot is located on the

west side of Third Avenue and is identified as Block 2319 Lot 2 on the NY City tax map. The property has 159 feet of street frontage on Third Avenue and approximately 346 ft of frontage along the north side of the Harlem River (**Figure 2**).

The lot is developed with a one-story 19,450 sf commercial building which according to the NYC Department of Buildings was constructed in 1931. The Property was assigned “E” designations (E-143) for air, noise, and hazardous materials as a result of the Port Morris/Bruckner Boulevard Rezoning action completed by the City in March 2005 (CEQR # 05DCP005X).

The property has an elevation of approximately 2 feet above the National Geodetic Vertical Datum (NGVD). Based upon regional groundwater contour maps, the depth to groundwater beneath the site is approximately 5-10 feet below existing grade and flows southwest toward the Harlem River.

A boundary map will be attached to the BCA as required by Environmental Conservation Law (ECL) Title 14 Section 27-1419. The 0.43-acre property is fully described in **Attachment A – Metes and Bounds**.

1.2 CONTEMPLATED REDEVELOPMENT PLAN

The redevelopment project consists of the construction of two new 25-story residential tower buildings identified as the east and the west towers. The two towers flank an entry court which is open to the river and Manhattan views to the south. The East Tower is free-standing, serving as a “portal” on Third Avenue, whereas the West Tower has a 7-story base with common amenities and parking for 200 cars. The residential component consists of 475 rental units in an area of 399,798 gross square feet. Commercial (retail) and community space will be included in the west tower base.

1.3 DESCRIPTION OF SURROUNDING PROPERTY

The area surrounding the property (**Figure 3**) is highly urbanized and predominantly consists of heavy commercial / industrial / warehouse properties to the north along a corridor adjacent to the

Harlem River. Multi-use residential / commercial (retail) properties are present to the east along Bruckner Boulevard and a large housing project is located to the northeast.

2.0 DESCRIPTION OF REMEDIAL INVESTIGATION FINDINGS

The field work portion of the Remedial Investigation was conducted by EBC in November and December 2015. The investigation is summarized in the sections below. Further details are provided in the Remedial Investigation Report (EBC February 2016).

2.1 SUMMARY OF REMEDIAL INVESTIGATIONS PERFORMED

2.1.1 Soil Borings

A total of twenty-Three soil borings (15B1-15B23) were advanced on November 20, 23 and 24, 2015 to identify source areas and to obtain general soil quality information present at the Site (**Figure 4**).

At each soil boring location soil samples were collected continuously in 5-foot intervals from grade to a depth of 15 feet below grade using a Geoprobe™ 6720DT, probe drilling machine. The Geoprobe™ system uses a direct push hydraulic percussion system to drive and retrieve core samplers. Soil samples were retrieved using a 1.25-inch diameter, 5-foot long dual-tube sampler with disposable acetate liners. Soil recovered from each soil boring was field screened by an environmental professional for the presence of VOCs with a photo-ionization detector (PID) and visually inspected for evidence of contamination. Soil samples were retained for laboratory analysis from all borings in accordance with the RIWP with the exception of B16 due to an error in completing the chain of custody form..

Retained samples were submitted for laboratory analysis of one or more of the following analyses: volatile organic compounds (VOCs) by EPA Method 8260, semi-volatile organic compounds (SVOCs) by EPA Method 8270, TAL Metals, pesticides and PCBs by EPA Method 8081/8082. Soil boring locations are identified in **Figure 4**.

2.1.2 Monitoring Wells

Twelve groundwater monitoring wells, MW1 - MW5 and MW7 - MW13D, were installed at the Site on November 20, November 24 and December 2, 2015. All of the wells were installed with a track mounted probe drilling machine. Wells MW1-MW10 were installed to a depth of

approximately 15 feet below grade with 10 feet of 1-inch diameter 0.010-inch slotted PVC well screen and 5 feet of PVC riser. Wells MW11D-MW13D were installed to a depth of 30 ft with 10 ft of well screen.

A No. 00 morie filter-pack sand filled the annulus surrounding the screen within two feet above the top of the screen. A one-foot hydrated bentonite seal was then placed on top of the filter sand and the remainder of the borehole was backfilled to grade. Following installation, each of the wells were surveyed to determine relative casing elevation to the nearest 0.01 ft and horizontal position to the nearest 0.1 ft. Monitoring well locations are identified in **Figure 5**. Well completion reports detailing monitoring well construction are provided in **Appendix B**.

Prior to sampling, a synoptic round of depth-to-groundwater (DTW) measurements was obtained from the wells on December 1, 2015 to determine the water table elevation and to calculate the volume of standing water in the well.

2.1.3 Samples Collected

A summary of the sampling performed during the RI is provided in **Table 2**.

2.1.3.1 Soil Samples

A total of thirty-four soil samples were collected from 23 soil borings for laboratory analysis of one or more of the following parameters: VOCs (EPA Method 8260), SVOCs (EPA Method 8270), TAL metals and pesticides/PCBs (EPA Method 8081/8082).

2.1.3.2 Groundwater Samples

A total of twelve groundwater samples were collected from the groundwater monitoring wells for laboratory analysis of VOCs (EPA Method 8260), SVOCs (EPA Method 8270), TAL metals and pesticides/PCBs (EPA Method 8081/8082).

2.1.3.3 Soil Vapor Samples

To assess the presence of VOCs in soil vapor beneath the site, eight soil vapor implants and five subsurface vapor points were installed at the Site and sampled on December 1 and 2, 2015. The

vapor implants (Geoprobe™ Model AT86 series), were constructed of a 6-inch length of double woven stainless steel wire and installed to a depth of 3.5 ft below grade using Geoprobe™ equipment. Five subslab samples installed by drilling a 1/2 inch hole through the concrete slab with a handheld drill and inserting 1/4 inch polyethylene tubing to no more than 2 inches below the base of the slab. The tubing was then sealed at the surface with hydrated granular bentonite.

All soil vapor samples were collected over a 2 hr sampling period.

Soil vapor samples were collected in accordance with the procedures as described in section 2.4 of the approved RIR and the *Guidance for Evaluating Soil Vapor Intrusion in the State of New York (NYSDOH 10/06)*.

2.1.4 Chemical Analytical Work Performed

Each soil and groundwater sample was placed in pre-cleaned laboratory supplied glassware, and placed in a cooler packed with ice for transport to the laboratory. Laboratory services for soil and groundwater sample analysis were provided by Phoenix Environmental Laboratories of Manchester, CT, a New York State ELAP certified environmental laboratory (ELAP Certification No. 11301).

Retained soil samples were submitted for laboratory analysis of one or more of the following analyses: volatile organic compounds (VOCs) by EPA Method 8260, semi-volatile organic compounds (SVOCs) by EPA Method 8270, TAL Metals, pesticides and PCBs by EPA Method 8081/8082.

All groundwater samples from the monitoring wells were analyzed for VOCs / SVOCs by EPA method 8260 / 8270, target analyte list (TAL) metals by EPA method 6010 and Pesticides/PCBs by method 8081/8082. Soil gas samples were analyzed for VOCs by USEPA Method TO-15.

2.1.5 Documentation

A map showing the locations of the soil borings is provided in **Figure 4**. The locations of the monitoring wells and soil gas sample collection points are provided in **Figures 5** and **6**. The

results of sample soil, groundwater and soil gas samples collected during the RI are summarized in **Tables 3** through **14**. Below is a summary of the RI findings.

The results of sampling performed during the RI, identified petroleum related contaminants in soil to a depth of 8 feet in the vicinity of a suspect 550 gallon underground storage tank (UST) that was previously abandoned in place.

Historic fill material has been identified across the Site to depths as great as 7 feet below grade. The historic fill material contains semi-volatile organic compounds and metals including arsenic, barium, chromium, copper, lead, mercury, nickel, silver and zinc above unrestricted, restricted residential or restricted commercial use SCOs.

Groundwater impacts from petroleum were not reported during the RI but were reported previously in the vicinity of the suspect underground gasoline tank. SVOC detections above groundwater standards were limited to those polynuclear aromatic hydrocarbons (PAHs) with a 2 per trillion standard. Exceedances of the part per trillion standards for PAHs were reported in well MW9. SVOCs reported in the parts per trillion range are a function of the laboratories ability to achieve extremely low detection limits and general background conditions.

The PCB Arochlor 1254 was also reported above standards in MW9.

Several dissolved metals were detected above standards including sodium, iron and manganese were detected above standards throughout the Site. The metals detected in groundwater are related to brackish conditions and are associated with the Site's proximity to the Harlem River.

With the exception of lighter compounds such as heptane and hexane in several samples, total petroleum-related VOCs were generally low in soil vapor samples across the Site. There did not appear to be any correlation in concentration or distribution of petroleum VOCs in soil vapor with the source area.

CVOCs were reported in all of the soil vapor and subslab samples with total CVOCs ranging from 0.28 $\mu\text{g}/\text{m}^3$ (SG1) to 26.06 $\mu\text{g}/\text{m}^3$ (SG9) in the soil gas samples and from 0.69 $\mu\text{g}/\text{m}^3$ (SS4) to 5.15 $\mu\text{g}/\text{m}^3$ (SS2) in the subslab samples.

2.2 SIGNIFICANT THREAT

The NYSDEC and NYSDOH will review the RI Report and will determine whether the Site does or does not pose a significant threat to human health and the environment. Notice of that determination will be provided during the public comment period, through fact sheet No. 2 and the Proposed Decision Document.

2.3 SITE HISTORY

2.3.1 Past Uses and Ownership

The Site is currently owned by the Requestor, 2401 3rd Ave Associates Property LLC. The property is currently vacant but was most recently used by a shoring company and has historically been used as an iron works, steel fabrication, gear manufacturing, gasket manufacturing and shoring equipment fabrication. The Requestor purchased the property in January 2015.

2401 3rd Avenue Associates Property LLC (the Requestor) is the current owner of the property and has owned the property since January 20, 2015. The property has been owned by various corporations associated with the Mugler family since at least 1986. The Muglers have operated a shoring equipment business on the premises since at least 1965. There appear to have been multiple tenants on the property through the 1960's and 1970's with the Mugler operation. The property has been underutilized for years and was proposed for inclusion in a newly designated BOA as a strategic property by the South Bronx Overall Economic Development Corp.

Previous Owners

| Dates | Name | Comments | Contact Info |
|------------------------------|--|-------------------------|---|
| From 1953 to 2/1/1986 | Richard C. Mugler Mary Mugler | Deed Owner Knowledge | Red Rock Road Austerlitz, NY 12017 |
| From 2/1/1986 to 11/25/1997 | Richard C. Mugler, Jr. | Deed | 177 Hudson Avenue Tenafly, New Jersey 07670 |
| From 11/25/1997 to 3/20/2000 | KAI Development Corp. | Deed | 732 W. 58 th Street, New York, NY 10019 |
| From 3/20/2000 to 1/20/2015 | 2401 3 rd Avenue, LLC | Deed | 177 Hudson Avenue Tenafly, New Jersey 07670 |
| From 1/20/2015 to present | 2401 3 rd Ave Associates Property LLC | Deed | 512 7 th Avenue 15 th Floor, New York, NY 10018 |

Previous Operators

| Dates | Name | Comments | Contact Info |
|---|---|--------------------------------|---|
| From sometime prior to 1891 to sometime between 1908 & 1922 | J.L. Mott Iron Works | Sanborn Maps | Unknown 2401 Third Avenue, Bronx, NY 10451 |
| From sometime between 1908 & 1922 to sometime between 1928 & 1935 | Hydraulic Steel Co. | Sanborn Maps | Unknown 2401 Third Avenue, Bronx, NY 10451 |
| From sometime between 1928 and 1935 to 1956. | General Building Supply Inc. (1949, 1956) Brill Equip. Co. (1949, 1956) | Sanborn Maps City Directory | Unknown 2401 Third Avenue, Bronx, NY 10451 |
| From 1965-1971 | US Gear Man. (1965, 1971) Ohio Gasket Co. (1971) | City Directory | 2401 Third Avenue, Bronx, NY 10451 |
| From 1965-present | Mugler Shoring Co. | City Directory | 2401 Third Avenue, Bronx, NY 10451 |

2.3.2 Summary of Previous Reports

Environmental investigations performed at the Site include the following:

- Phase I Environmental Site Assessment - Langan Engineering, Environmental, Surveying and Landscape Architecture, D.P.C. October 28, 2014
- Phase II Environmental Site Investigation Report - Langan Engineering, Environmental, Surveying and Landscape Architecture, D.P.C. October 28, 2014

October 2014 - Phase I Environmental Site Assessment Report (Langan)

A Phase I Environmental Site Assessment Report was prepared by Langan Engineering, Environmental, Surveying and Landscape Architecture, D.P.C. (Langan) on October 28, 2014.

Based upon reconnaissance of the subject and surrounding properties, interviews and review of historical records and regulatory agency databases, Langan identified the following recognized environmental conditions:

- Current and Historical Manufacturing and Industrial Use:

The Subject Property was historically used for manufacturing purposes since at least 1891. Historical operators include; J.L. Mott Iron Works (1891–1922), Hydraulic Steel Company (1922–1935), General Builders Supply Corporation (1935–1968), Brill Equipment Company (1949–1956), US Gear Manufacturing Company (1965–1971), Ohio Gasket Manufacturing Corporation (1971), and Mugler Inc. (1965–present). Inadvertent releases of solvents, petroleum products, metals, polychlorinated biphenyls (PCB) and/or other chemicals used during manufacturing operations may have adversely impacted soil, groundwater, building components and/or soil vapor. The Subject Property is presently operated by Mugler, Inc. for shoring equipment fabrication, storage, truck loading/unloading, and equipment repairs. As such, there are typical tools, maintenance/repair materials, and miscellaneous equipment used and stored throughout. Discoloration and staining were apparent throughout the interior of the building, suggesting incidental releases of petroleum products during truck maintenance, and the concrete slab was compromised in several areas. Because fractures in the slab provide a

conduit for spilled motor oils and/or petroleum products to impact subsurface conditions, current use by Mugler, Inc. constitutes a REC.

- On-Site Petroleum Bulk Storage:

The following historical underground storage tanks (USTs) were identified:

- One 550-gallon gasoline UST, located outside the southeast corner of the building, was reportedly closed-in place circa 1999.
- An area of patched concrete and an apparent abandoned fill port southeast of the building, which is suspected to be associated with another decommissioned UST, was identified during the site reconnaissance.
- One 1,550-gallon gasoline UST was identified on the 1935 through 1946 Sanborn maps.

Based on the known and suspected presence of historic tanks and lack of any documentation of tank closure, the historical USTs are a REC.

- Current and Historical Use of Surrounding Properties:

Include a private garage (1946–1951), lubricating oils storage (1935–1983), two chemical corporations (1935–1947), a paint manufacturer (1951), a coal yard (1935–1951), an auto building (1908–1947), an auto house (1935–1944), a private garage (1935–1944), a printing facility (1935–1947), and several auto repair shops (2005–2012). In addition, an active New York State Department of Environmental Conservation (NYSDEC) Brownfield Cleanup Program (BCP) site was identified approximately 660 feet northeast (up-gradient) of the Subject Property. Based on investigations conducted to date, the primary contaminants of concern in soil and groundwater are petroleum related compounds. No information was provided with respect to the off-Site migration of contaminants; however, the Environmental Database Report (EDR) report indicates that the potential exists for off-Site migration of site-related contaminants in soil vapor. Potential petroleum and solvent releases associated with the historical surrounding

property uses may have adversely impacted groundwater and/or soil vapor at the Subject Property and is considered a REC.

October 2014 – Phase II Environmental Site Investigation Report (Langan)

A remedial investigation was performed by Langan for the property in July and August of 2014. The Phase II included the following scope of work:

1. Completion of a geophysical investigation;
2. Completion of seven soil borings and collection of 14 soil samples;
3. Installation and sampling of six temporary monitoring wells; and
4. Installation and sampling of six soil vapor points. Installed five soil vapor probes around Site perimeter and collected five samples for chemical analysis.

The findings of the Phase II are summarized as follows:

- The geophysical survey identified two major anomaly areas along the southeast corner of the Subject Property Building. Based on their reflection rates and their proximities, these anomalies are indicative of USTs located approximately 2 to 3 ft bgs. Another anomaly area consistent with disturbed soil was located along the western edge of the Subject Property.
- A layer of historic fill was identified throughout the Subject Property with thicknesses ranging from approximately 5 to 13 feet. This fill layer contained SCO exceedances (SVOC and metal) throughout the majority of the Site.
- The suite of metals and SVOCs and their concentrations detected in soil and groundwater across the site are attributed to historic fill material.
- One of the seven soil borings (SB-3) was installed next to a suspect 550-gallon gasoline UST that was reportedly decommissioned in 1999. Petroleum-like odors and PID readings (up to 700 ppm) were observed in the soil boring at depths of approximately 8 to 9 ft bgs. Groundwater analytical sampling results from this location confirmed the

presence of a petroleum release. Article 12 of the New York State Navigation Law requires that anyone with knowledge of a petroleum release must report the discovery to the New York State Spill Hotline as soon as possible. Therefore, on 13 August, 2014, Langan called in a spill and Spill Case No. 1405230 was opened by NYSDEC.

- Soil vapor concentrations are below NYSDOH AGV values; however, based on the presence of VOCs in the soil vapor, remedial measures should include installation of a soil vapor mitigation system.
- Due to the E-designation of the site, additional investigation including soil, groundwater and soil vapor sampling will be required in order to comply with NYCOER guidelines and obtain permits for construction and occupancy of proposed building structures.
- If site development is planned, a Remedial Action Plan (RAP) and Construction Health and Safety Plan (CHASP) should be prepared and implemented. In accordance with the requirements of the E-Designation program, these documents must be submitted to OER for approval prior to the start of Site work. OER will require additional site investigation to support preparation of the RAP. The RAP and CHASP may need to address the following:
 - Handling, transport and disposal of excess material including fill and debris as required for development with all appropriate requirements.
 - A composite cover system for impacted material that may be left in-place after development. The cover system will prevent direct contact with or inhalation of contaminated material.
 - Odor and soil vapor intrusion should be addressed if odorous or VOC-impacted material is left in-place as part of the development.
 - Registration, removal and closure of known petroleum storage tanks.
 - Permitting and treatment, if required, of dewatering effluent.
 - Health and safety requirements addressing the contaminants of concern including a Community Air Monitoring Plan (CAMP) to monitor odor and dust.
 - A contingency plan for removal of currently unknown underground storage tanks or other subsurface structures

2.4 GEOLOGICAL CONDITIONS

The bedrock geology at the property and in the immediate vicinity consists of Inwood Marble of Lower Ordovician to Lower Cambrian age with steep westerly dip of its upper surface. The depth to bedrock is anticipated at approximately 20 to 30 feet below land surface (ft-bls). Bedrock is overlain by an unconsolidated overburden of an unsorted heterogeneous mix of Pleistocene and Recent glacial material (i.e., glacial till) including clay, silt, sands, gravel, cobbles, and boulders. This overburden is overlain by historic urban fill.

Subsurface soils at the Site consist of historic fill materials in and silty-sand matrix to a depth of approximately 5 feet below grade. A silty clay was present below this layer to a depth of 15 feet in those borings located closer to the shore. Further inland a medium to coarse sand was encountered beneath the fill to a depth of 15 feet (**Figure 7**).

Groundwater at the Site is present under water table conditions at a depth of 5.17 to 6.75 feet below grade. Based upon on-site measurements, groundwater flow is to the southwest (**Figure 8**).

Considering the poor quality of groundwater in the area, including high levels of sodium and magnesium associated with saltwater intrusion, there is no anticipated future groundwater use.

2.5 CONTAMINATION CONDITIONS

2.5.1 Conceptual Model of Site Contamination

VOC contamination at the Site consists of petroleum related contaminants in soil to a depth of 8 feet in the vicinity of a suspect 550 gallon underground storage tank (UST) that was previously abandoned in place. Released gasoline migrated through the soil column to the water table. Some gasoline constituents dissolved into the groundwater in the immediate vicinity of the tank. There does not appear to be any migration of dissolved VOCs from the source area.

Lighter end VOC components such as heptane and hexane have transferred to the vapor phase from impacted soil and / or groundwater.

2.5.2 Description of Areas of Concern

The source area identified during the RI includes petroleum VOCs in the north-central area of the Site in the vicinity of a suspect UST.

2.5.3 Soil/Fill Contamination

VOC contamination at the Site consists of petroleum related contaminants in soil to a depth of 8 feet in the north-central area of the Site in the vicinity of a suspect UST.

Historic fill material has been identified across the Site to depths as great as 7 feet below grade. The historic fill material contains semi-volatile organic compounds and metals including arsenic, barium, chromium, copper, lead, mercury, nickel, silver and zinc above unrestricted, restricted residential or restricted commercial use SCOs.

2.5.3.1 Summary of Soil/Fill Data

Soil sample results from the RI are summarized in **Tables 3-6**. Further information on soil sample collection, handling and analysis can be found in the RI Report (EBC 2/16).

2.5.3.2 Comparison of Soil/Fill with SCGs

Table 7 shows sample results above Track 1 Unrestricted SCOs for all overburden soil at the Site. **Figure 9** is spider map which shows soil sampling locations and summarizes shallow and deep sample results above Track 1 Unrestricted SCOs for all overburden soil.

2.5.4 On-Site and Off-Site Groundwater Contamination

Groundwater impacts from petroleum were limited to methyl tertiary butyl ether (MTBE) at two locations during the RI but were not near the underground gasoline tank. However petroleum impacts were reported previously during a Phase II investigation (Langan 10/2014) in the vicinity of this tank. Chloromethane was also reported above its groundwater standard at one location.

SVOC detections above groundwater standards were limited to those polynuclear aromatic hydrocarbons (PAHs) with a 2 per trillion standard. Exceedances of the part per trillion standards

for PAHs were reported in well MW9. SVOCs reported in the parts per trillion range are a function of the laboratories ability to achieve extremely low detection limits and general background conditions.

The PCB Arochlor 1254 was also reported above standards in MW9.

Several dissolved metals were detected above standards including sodium, iron and manganese were detected above standards throughout the Site. The metals detected in groundwater are related to brackish conditions and are associated with the Site's proximity to the Harlem River.

2.5.4.1 Summary of Groundwater Data

The results of groundwater samples collected during the RI are summarized in **Tables 8-12**. Further information on groundwater sample collection, handling and analysis can be found in the RI Report (EBC 2/16).

2.5.4.2 Comparison of Groundwater with SCGs

Sample results above GA groundwater standards in monitor wells prior to the remedy are shown in **Table 13**. Spider maps which show groundwater sampling locations and summarize results above GA groundwater standards prior to the remedy are shown in **Figure 10**.

2.5.5 On-Site and Off-Site Soil Vapor Contamination

With the exception of lighter compounds such as heptane and hexane in several samples, total petroleum-related VOCs were generally low in soil vapor samples across the Site. There did not appear to be any correlation in concentration or distribution of petroleum VOCs in soil vapor with the source area.

CVOCs were also generally low and were reported in all of the soil vapor samples with concentrations ranging from 0.28 $\mu\text{g}/\text{m}^3$ to 26.06 $\mu\text{g}/\text{m}^3$.

2.5.5.1 Summary of Soil Vapor Data

A table of soil vapor data collected prior to the remedy is shown in **Table 14**. Further information on soil gas sample collection, handling and analysis can be found in the RI Report (EBC 2/16). Soil vapor results are posted on **Figure 11**.

2.6 ENVIRONMENTAL AND PUBLIC HEALTH ASSESSMENTS

2.6.1 Qualitative Human Health Exposure Assessment

The objective of the qualitative exposure assessment under the Brownfields Cleanup Program (BCP) is to identify potential receptors to the contaminants of concern (COC) that are present at, or migrating from, the Site. The identification of exposure pathways describes the route that the COC takes to travel from the source to the receptor. An identified pathway indicates that the potential for exposure exists; it does not imply that exposures actually occur. An exposure pathway has five elements; a contaminant source, release and transport mechanisms, point of exposure, route of exposure and a receptor population.

The potential exposure pathways identified below, represent both current and future exposure scenarios.

Contaminant Source

Source areas of the Site include petroleum VOCs in the north-central area of the Site in the vicinity of a suspect UST.

Elevated levels of metals and PAHs are also present in fill materials throughout the Site to depths as great as 7 ft below grade.

Contaminant Release and Transport Mechanism

Petroleum contamination is present in subsurface soil in the north-central area of the Site in the vicinity of a suspect gasoline storage tank. The contamination extends vertically to the groundwater interface. There appears to be minimal transfer of petroleum VOCs to the groundwater despite the fact that impacted soil is in contact with the groundwater. There does

not appear to be any significant transfer of petroleum VOCs to the vapor phase or concern for off-site vapor migration.

Point of Exposure, Route of Exposure and Potentially Exposed Populations

Potential On-Site Exposures: Remediation workers and construction workers engaged in the excavation of impacted and non-impacted soil at the site may be exposed to petroleum VOCs / SVOCs, and heavy metals through several routes including , ingestion and dermal contact. A site specific Health and Safety Plan has been developed to identify and minimize the potential hazards to on-site workers. Site trespassers could also be exposed to impacted soil during excavation, however security measures including an 8 ft high construction fence and 24 hr security will minimize potential exposure through this route.

Future occupants of the new building including commercial retail workers, residents and visitors could be exposed to SVOCs and heavy metals in soil through ingestion and dermal contact if these contaminants were to remain in exposed soils at the Site.

Potential Off-Site Exposures: Off-Site residents could also be exposed to dust or vapors during the excavation of impacted soil. A site specific Community Air Monitoring Plan has been developed to identify and minimize the potential for off-site exposure to residents through continuous air monitoring during excavation activity.

The entire area is serviced by the New York City Water System which distributes water from the Croton Reservoir system. Since there are no public or private potable supply wells in the area, exposure from contact with tap water is not a concern.

Potential Off-Site Environmental Impacts: Since petroleum VOCs in groundwater were reported in the vicinity of the suspect gasoline tank during a previous investigation at the Site, the groundwater to surface water discharge pathway was evaluated. The nearest body of water is the Harlem River which borders the Site to the south. Based on the low concentrations of contaminants reported and the absence of VOCs in groundwater during the RI downgradient of

the tank area, there are no expected impacts to surface water environments from contaminants migrating from the Site.

2.6.2 Fish & Wildlife Remedial Impact Analysis

Based on the nature and location of the contamination at the site, with limited impact to subsurface soils and groundwater and no evidence of migration / discharge of contaminants to the River, there are no expected adverse impacts to River. However, the presence of elevated levels of SVOCs, metals and pesticides in surficial soil at the Site would be a concern if these contaminants were to be discharged to the river in sediments from surface runoff. Therefore, capping or otherwise stabilizing surficial soil should be included in the Remedial Plan developed for the Site.

2.7 REMEDIAL ACTION OBJECTIVES

Based on the results of the Remedial Investigation, the following Remedial Action Objectives (RAOs) have been identified for this Site.

2.7.1 Groundwater

RAOs for Public Health Protection

- Prevent ingestion of groundwater containing contaminant levels exceeding drinking water standards.
- Prevent contact with, or inhalation of, volatiles emanating from contaminated groundwater.

RAOs for Environmental Protection

- Restore ground water aquifer, to the extent practicable, to pre-disposal/pre-release conditions.
- Prevent the discharge of contaminants to surface water.

2.7.2 Soil

RAOs for Public Health Protection

- Prevent ingestion/direct contact with contaminated soil.

RAOs for Environmental Protection

- Prevent migration of contaminants that would result in groundwater or surface water contamination.

2.7.3 Soil Vapor

- Mitigate impacts to public health resulting from existing, or the potential for, soil vapor intrusion into buildings at a site.

3.0 DESCRIPTION OF REMEDIAL ACTION PLAN

3.1 EVALUATION OF REMEDIAL ALTERNATIVES

The goal of the remedy selection process under the BCP is to select a remedy that is protective of human health and the environment taking into consideration the current, intended and reasonably anticipated future use of the property. The remedy selection process begins by establishing RAOs for media in which chemical constituents were found in exceedance of NYSDEC standards, criteria and guidance values (SCGs). A remedy is then developed based on the following nine criteria:

- Protection of human health and the environment;
- Compliance with SCGs;
- Short-term effectiveness and impacts;
- Long-term effectiveness and permanence;
- Reduction of toxicity, mobility, or volume of contaminated material;
- Implementability;
- Cost effectiveness;
- Community Acceptance; and
- Land use.

The first two criteria are threshold criteria and must be satisfied in order for an alternative to be considered for selection. The remaining seven criteria are balancing criteria which are used to compare the positive and negative aspects of each of the remedial alternatives, provided the alternative satisfies the threshold criteria.

3.2 STANDARDS, CRITERIA AND GUIDANCE (SCG)

A criterion for remedy selection is evaluation for conformance with SCGs that are applicable, relevant and appropriate. Principal SCGs that are applicable, relevant and appropriate for evaluating the alternatives for remediation of this BCP site include the following:

- 29 CFR Part 1910.120 - Hazardous Waste Operations and Emergency Response

- 10 NYCRR Part 67 – Lead
- 6 NYCRR Part 371 - Identification and Listing of Hazardous Wastes (November 1998)
- 6 NYCRR Part 372 - Hazardous Waste Manifest System and Related Standards for Generators, Transporters and Facilities (November 1998)
- 6 NYCRR Subpart 374-1 - Standards for the Management of Specific Hazardous Wastes and Specific Types of Hazardous Waste Management Facilities (November 1998)
- 6 NYCRR Part 375 - 6 NYCRR Part 375 Environmental Remediation Programs Subparts 375-1, 375-3 and 375-6 (December 2006)
- 6 NYCRR Part 376 - Land Disposal Restrictions
- 6 NYCRR Part 608 - Use and Protection of Waters
- 6 NYCRR Parts 700-706 - Water Quality Standards (June 1998)
- 6 NYCRR Part 750 through 758 - Implementation of NPDES Program in NYS (“SPDES Regulations”)
- 6 NYCRR Part 375-6 Soil Cleanup Objectives
- New York State Groundwater Quality Standards – 6 NYCRR Part 703;
- NYSDEC Ambient Water Quality Standards and Guidance Values – TOGS 1.1.1;
- NYSDEC DER-10 Technical Guidance for Site Investigation and Remediation - May 2010;
- NYSDEC Draft Brownfield Cleanup Program Guide – May 2004;
- New York State Department of Health (NYSDOH) Generic Community Air Monitoring Plan
- NYS Waste Transporter Permits – 6 NYCRR Part 364;
- NYS Solid Waste Management Requirements – 6 NYCRR Part 360 and Part 364.
- TAGM 4059 - Making Changes To Selected Remedies (May 1998)
- STARS #1 - Petroleum-Contaminated Soil Guidance Policy
- TAGM 3028 - "Contained In" Criteria for Environmental Media: Soil Action Levels (August 1997)
- DER-10, Technical Guidance for Site Investigation and Remediation, May 2010
- DER-23 / Citizen Participation Handbook for Remedial Programs, January 2010

- OSWER Directive 9200.4-17 - Use of Monitored Natural Attenuation at Superfund, RCRA Corrective Action, and Underground Storage Tank Sites (November 1997)

Additional regulations and guidance are applicable, relevant, and appropriate to the remedial alternatives and will be complied in connection with implementation of the remedial program; however, the list above is intended to represent the principal SCGs which should be considered in evaluating the remedial alternatives for the BCP site.

Conformance with the appropriate standards for remediation of contaminated soil is an important criterion in evaluating the remedial alternatives for the BCP site. Presently, in New York State 6 NYCRR Part 375 establishes the primary SCGs associated with remediation of contaminated soil at sites which are in the BCP. If proposing remediation pursuant to a Track other than Track 1 (Unrestricted Use), 6 NYCRR Part 375 requires evaluation of at least one remedial alternative pursuant to Track I (Unrestricted Use) and one other alternative developed by the applicant for the proposed use of the BCP site. The remedial alternatives presented in Section 3.3 of this work plan have been prepared in conformance with this requirement.

3.3 ALTERNATIVES ANALYSIS

The goal of the remedy selection process under the BCP is to select a remedy that is protective of human health and the environment taking into consideration the current, intended and reasonably anticipated future use of the property. The remedy selection process begins by establishing RAOs for media in which chemical constituents were found in exceedance of NYSDEC standards, criteria and guidance values (SCGs). A remedy is then developed based on the following nine criteria:

- Protection of human health and the environment;
- Compliance with SCGs;
- Short-term effectiveness and impacts;
- Long-term effectiveness and permanence;
- Reduction of toxicity, mobility, or volume of contaminated material;
- Implementability;

- Cost effectiveness;
- Community Acceptance; and
- Land use.

The following is a detailed description of the alternatives analysis and remedy selection to address impacted media at the Site. This analysis was prepared in accordance with 6 NYCRR Part 375-1.8(f) and Part 375-3.8(f) and Section 4.3(c) of NYSDEC DER-10. As required, a minimum of two remedial alternatives (including a Track 1 scenario) are evaluated, as follows:

- Alternative 1 - Track 1, remediation of all soils above bedrock to unrestricted use criteria. Excavation to a minimum depth of 8 feet across the Site with the excavation of the petroleum impacted area to a depth of approximately 9 feet. This alternative does not allow the use of long-term institutional /engineering controls to address impacted media or prevent exposures which may be required beneath the new building.
- Alternative 2 - Track 4, removal / replacement of the top 2 feet of soil in all planned exposed soil areas, and excavation / removal of all petroleum impacted soils with VOCs above groundwater protection SCOs which are also present above groundwater standards. This alternative would require a lesser degree of excavation than Alternative 1 consisting of the excavation of petroleum impacted soil in the UST source area to an approximate depth of 9 feet. An engineered cap consisting of the concrete building slab, paved driveways and walkways and removal /replacement of soil within landscaped areas with 2 feet of clean fill / soil meeting Restricted Residential use SCOs. Long-term institutional /engineering controls are not necessary but would be allowed to address or prevent exposures from soil vapor, if required.

3.4 REMEDIAL ALTERNATIVE 1

The following sections provide an evaluation of Alternative 1 based on the nine evaluation criteria as previously discussed.

3.4.1 Overall Protection of Human Health and the Environment

Alternative 1 will be protective of human health and the environment by eliminating constituents in soil related to petroleum and historic fill. The potential for human and environmental exposure to these constituents on-site will be eliminated by excavation of all petroleum contaminated and historic fill soils with parameters in excess of unrestricted criteria, disposing of excavated materials off-site and backfilling as needed with certified clean fill, virgin mined materials or recycled concrete materials from a NYSDEC permitted recycling facility.

Potential post-remediation exposures to on-site residents from soil vapors are not expected to require the operation of SSD systems, though groundwater will be restricted at the Site even though such a restriction is unnecessary due to brackish conditions and NYC prohibitions on groundwater usage.

During remedial and construction activity workers and area residents may be exposed to impacted soil and vapors. Worker exposure to soil and vapors will be minimized through implementation of a Health and Safety Plan. Exposures to area residents from dust and/or vapors will be minimized through the use of engineering controls and through implementation of a Community Air Monitoring Plan (CAMP).

3.4.2 Compliance with Remedial Goals, SCGs and RAOs

Alternative 1 will achieve compliance with the remedial goals, SCGs and RAOs for soil through source removal to Track 1 unrestricted cleanup levels. SCGs for groundwater may not be achieved as the groundwater is of poor quality affected by brackish conditions from the adjacent river. Compliance with SCGs for soil vapor is expected following completion of the remedial action.

3.4.3 Long-Term Effectiveness and Permanence

Alternative 1 achieves long term effectiveness and permanence by permanently removing and/or remediating all soils affected by Site contaminants or historic fill materials and by remediating groundwater. Under this Alternative, risk from soil impacts and groundwater will be eliminated.

Alternative 1 will continue to meet RAOs for soil, groundwater and soil vapor in the future, providing a permanent long-term solution for the Site.

3.4.4 Reduction in Toxicity, Mobility or Volume through Treatment

Alternative 1 will permanently eliminate the toxicity, mobility, and volume of contaminants from on-site soil by meeting unrestricted objectives.

3.4.5 Short-Term Effectiveness

The potential for short-term adverse impacts and risks to the workers, the community, and the environment during the implementation of Alternative 1 is minimal.

Short-term exposure to on-site workers during excavation and loading activities will be addressed with a HASP and mitigated through the use of personal protective equipment, monitoring and engineering controls. Potential short-term exposure to the surrounding community will be addressed through the use of odor and dust-suppression techniques and through the implementation of a CAMP which will require air monitoring activities during all excavation and soil disturbance activities.

Other potential impacts to the community such as construction-related noise, vibrations and traffic, will be controlled and regulated under the terms of the NYS Department of Buildings issued building permit which can place a Stop Work Order on the property for unsafe conditions, community impacts or violation of the terms and conditions of the permit. Decontamination procedures of equipment, including trucks transporting soil to off-site disposal facilities, will minimize the potential for impacted soil to be dispersed beyond the Site boundary. A truck traffic plan has also been prepared to minimize disturbance to the local roads and community.

3.4.6 Implementability

The techniques, materials and equipment to implement Alternative 1 are readily available and have been proven effective in remediating the contaminants associated with the Site. Excavation for the remediation of soils is both a "low tech" and reliable method which has a long and proven track record on the remediation of hazardous waste and petroleum spill sites.

However, excavation to a depth of 8 ft or more over a 1.4 acre site located adjacent to a river represents considerable technical challenges which in this case limit its feasibility. Sloping and shoring requirements and dewatering would present significant construction challenges and would impede and interfere with the construction of the new buildings.

3.4.7 Cost

Costs associated with Alternative 1 are estimated at approximately \$ 4,155,490. This cost estimate includes the following elements and assumptions:

- Removal of existing USTs;
- Excavate a minimum of 8 feet across entire Site. Over-excavate as necessary to remediate hotspot areas and remove all historic fill at the Site;
- Excavate petroleum impacted soil area to a depth of 9 ft in the UST source area;
- Shoring to allow excavation to 8 ft depth at property lines;
- Dewatering in southern third of the Site to allow excavation to 8 ft depth;
- Disposal of approximately 150 tons of petroleum impacted non-hazardous soil from the former fueling area;
- Disposal of approximately 27,126 tons of historic fill soil as non-hazardous;
- Backfilling excavations with certified clean fill meeting unrestricted SCOs (approximately 15,861 cy);
- HASP and CAMP monitoring for the duration of the remedial activities.

3.4.8 Compatibility with Land Use

The proposed redevelopment of the Site is compatible with its current M1-2 / R6A zoning. Following remediation, the Site will meet unrestricted use objectives which will exceed the objectives for its planned commercial-retail and residential use. A groundwater use restriction may be required to prevent future exposure to affected groundwater.

3.4.9 Community Acceptance

No questions regarding the Site have been raised regarding remedial options to date. This RAWP will be subject to a 45-day public comment period to determine if the community had comments

on the presented remedial alternatives and selected remedy. If no comments are received regarding Alternative 1, it will be considered to be acceptable to the community.

3.5 REMEDIAL ALTERNATIVE 2

The following sections provide an evaluation of Alternative 2 based on the nine evaluation criteria as previously discussed.

3.5.1 Overall Protection of Human Health and the Environment

Alternative 2 will be protective of human health and the environment by eliminating constituents related to petroleum in soil. The potential for human and environmental exposure to these constituents on-site will be eliminated by excavation of the petroleum impacted areas and then capping exposed soil areas with a combination of the concrete building slab, concrete walkways / driveways and 2 feet of soil / fill which meets restricted residential SCOs.

Potential post-remediation exposures to on-site residents from soil vapors are not expected to require the operation of SSD systems, though groundwater use will be restricted at the Site due to brackish conditions.

During remedial and construction activity, workers and area residents may be exposed to impacted soil and vapors. Worker exposure to soil and vapors will be minimized through implementation of a HASP. Exposures to area residents from dust and or vapors will be minimized through the use of engineering controls and through implementation of a CAMP.

3.5.2 Compliance with Remedial Goals, SCGs and RAOs

Alternative 2 will achieve compliance with the remedial goals, SCGs and RAOs for soil through petroleum source and then capping the remainder of the Site with concrete building slab, concrete walkways / driveways and 2 feet of soil / fill which meets restricted residential SCOs. SCGs for groundwater may not be achieved as the groundwater is affected by brackish conditions from the adjacent river. Compliance with SCGs for soil vapor is expected following completion of the remedial action.

3.5.3 Long-term Effectiveness and Permanence

Alternative 2 achieves long term effectiveness and permanence by permanently removing and/or remediating soils affected by Site contaminants above site specific objectives. Under this Alternative risk from soil impacts is eliminated for on-site residents. Alternative 2 will continue to meet RAOs for soil in the future, providing a permanent long-term solution for the Site.

3.5.4 Reduction in Toxicity, Mobility or Volume through Treatment

Alternative 2 will permanently eliminate the toxicity, mobility, and volume of contaminants from on-site soil by removing petroleum impacted soil and by meeting groundwater protection SCOs for VOCs present above standards in groundwater and restricted residential objectives in the upper 2 feet of exposed soil.

3.5.5 Short-term Effectiveness

The potential for short-term adverse impacts and risks to the workers, the community, and the environment during the implementation of Alternative 2 is minimal. Short-term exposure to on-site workers during excavation and loading activities will be addressed with a HASP and mitigated through the use of personal protective equipment, monitoring and engineering controls. Potential short-term exposure to the surrounding community will be addressed through the use of odor and dust-suppression techniques and through the implementation of a CAMP which will require air monitoring activities during all excavation and soil disturbance activities.

Other potential impacts to the community such as construction-related noise, vibrations and traffic will be controlled and regulated under the terms of the NYS Department of Buildings issued building permit which can place a Stop Work Order on the property for unsafe conditions, community impacts or violation of the terms and conditions of the permit. Decontamination procedures of equipment, including trucks transporting soil to off-site disposal facilities will minimize the potential for impacted soil to be dispersed beyond the Site boundary. A truck traffic plan will also be prepared to minimize disturbance to the local roads and community.

3.5.6 Implementability

The techniques, materials and equipment to implement Alternative 2 are readily available and have been proven effective in remediating the contaminants associated with the Site. Excavation for the remediation of soils is both a "low tech" and reliable method which has a long and proven track record on the remediation of hazardous waste and petroleum spill sites.

3.5.7 Cost

Costs associated with Alternative 2 are estimated at approximately \$ 916,038. This cost estimate includes the following elements and assumptions:

- Removal and close - out of existing USTs;
- Excavate petroleum impacted soil area to a depth of 9 ft in the UST source area;
- Disposal of approximately 150 tons of petroleum impacted non-hazardous soil from the UST source area;
- Disposal of approximately 7,641 tons of non-hazardous, historic fill soil with parameters above RRSCOs from the building excavation;
- Backfilling excavations and exposed soil areas with certified clean fill meeting Restricted Residential Use and Groundwater Protection SCOs (approximately 1,000 cy);
- HASP and CAMP monitoring for the duration of the remedial activities;
- Implementation of a Site Management Plan (SMP) for long term maintenance of the Engineering Controls;
- Filing an Environmental Easement to ensure implementation of the SMP.

3.5.8 Compatibility with Land Use

The proposed redevelopment of the Site is compatible with its current M1-3 / R8 zoning. Following remediation, the Site will meet restricted-residential use objectives in the top 2 feet of soil which will meet objectives for its planned mixed commercial-retail and residential use. A groundwater use restriction may be required to prevent future exposure to brackish groundwater.

3.5.9 Community Acceptance

No questions regarding the Site have been raised regarding remedial options to date. This RAWP will be subject to a 45-day public comment period to determine if the community has any comments on the presented remedial alternatives and selected remedy. If no comments are received, it will be considered to be acceptable to the community.

3.6 SELECTION OF THE PREFERRED REMEDY

The remedy recommended for the site is a Track 4 alternative (Alternative 2) which consists of the removal of existing USTs, excavation of petroleum impacted soil to a depth of approximately 9 feet within the UST source area (USTs 1&2), excavation and disposal of historic fill soil from the building foundation areas and capping areas of the Site where soil remains above RRSCOs with the building foundations, concrete walkways / driveways or 2 ft of soil meeting Restricted Residential SCOs.

Any backfill materials used at the site will either consist of clean native soil excavated from other areas of the site, virgin mined materials, recycled materials or certified fill which meets Restricted Residential SCOs.

3.6.1 Preferred Remedy Land Use Factor Evaluation

As required by Article 27, Title 14 of the Environmental Conservation Law 27-1415, the following land use factor evaluation examines whether the preferred alternative is acceptable based on the 14 criteria presented in the following subsections.

Zoning

The proposed redevelopment project, which includes the construction of two new 25-story residential tower buildings, is in compliance with the M1-3 / R8 zoning. Therefore the project will be constructed as-of-right regardless of the remedy implemented. The preferred remedy will comply with current zoning.

Applicable Comprehensive Community Master Plans or Land Use Plans

The proposed redevelopment project and selected remedy are consistent with comprehensive master and land use plans, specifically the Port Morris / Bruckner Boulevard rezoning action (CEQR No. 05DCP005X), and the Port Morris Harlem Riverfront Brownfield Opportunity Area (BOA).

The rezoning action, covering an eleven block area of Port Morris, Bronx Community District 1, was completed by the New York City Department of City Planning and adopted by the City Council in March 2005. The BOA, which was designated in April 2015, identified the Site as a strategic Site in the nomination document, referring to it as a sizeable and underutilized waterfront site strategically located on the Harlem River waterfront.

The preferred remedy will be in full compliance with these applicable land use plans.

Surrounding Property Uses

The area surrounding the property is highly urbanized and predominantly consists of heavy commercial / industrial / warehouse properties to the north along a corridor adjacent to the Harlem River. Multi-use residential / commercial (retail) properties are present to the east along Bruckner Boulevard and a large housing project is located to the northeast.

The Lot is currently zoned M1-3/R8 residential / commercial. The M1-3/R8 is a Special Mixed Use District established to encourage investment in, and enhance the vitality of, existing neighborhoods with mixed residential and industrial uses in close proximity and create expanded opportunities for new mixed use communities. New residential and non-residential uses (commercial, community facility and light industrial) can be developed as-of-right and be located side-by-side or within the same building. Pairing an M1 district with an R3 through R10 district ensures a balanced variety of uses.

The proposed project is compatible with the surrounding land use and will be in compliance with the current zoning. The proposed remedy will not interfere with surrounding property uses and considers the short term affects to neighboring properties.

Citizen Participation

Citizen participation for implementation of the preferred alternative will be performed in accordance with DER 23 and NYCRR Part 375-1.10 and Part 375-3.10. A Citizen Participation Plan has been prepared and is available for public review at the identified document repositories (NYSDEC Region 2 Office, Mott Haven Branch of the New York Public Library).

Environmental Justice Concerns

The Site is not located within a potential environmental justice area. The NYSDEC defines a potential environmental justice area as a "minority or low-income community that may bear a disproportionate share of the negative environmental consequences resulting from industrial, municipal, and commercial operations or the execution of federal, state, local, and tribal programs and policies.

Environmental justice means the fair treatment and meaningful involvement of all people regardless of race, color, or income with respect to the development, implementation, and enforcement of environmental laws, regulations, and policies. Fair treatment means that no group of people, including a racial, ethnic, or socioeconomic group, should bear a disproportionate share of the negative environmental consequences resulting from industrial, municipal, and commercial operations or the execution of federal, state, local, and tribal programs and policies.

Since the goal of the remedy will achieve the highest level of cleanup and will remove contaminated materials from the community, the remedy poses no environmental justice concerns.

Land use designations

The proposed remedy is consistent with land-use designations.

Population growth patterns

Population growth patterns support the proposed use for the Site. The preferred remedy will not negatively affect on population growth patterns.

Accessibility to existing infrastructure

The Site is accessible to existing infrastructure. The close proximity of the Site to the Major Deegan Expressway (I-87) will assist soil transportation and contractor access to the Site. The Site is also accessible to mass transit and is within walking distance to the 6 line of the NYC Subway system with a stop located on 138th Street at Third Avenue (2 blocks to the north). The preferred remedy will not alter accessibility to existing infrastructure.

Proximity to cultural resources

The proposed remedy will not negatively impact cultural resources.

Proximity to natural resources

The proposed remedy will improve the local environment and will not negatively impact affect natural resources.

Off-Site groundwater impacts

The proposed remedy will improve potential off-site groundwater impacts by removing petroleum impacted soil from the Site and capping exposed fill. The proposed remedy will not affect natural resources other than to potentially improve the quality of groundwater on a local basis and eliminate potentially impacted surface runoff and sediment into the Harlem River.

Proximity to floodplains

The entire Site is located within a designated high risk flood zone area. Capping of the Site will reduce the potential impacts from flooding.

Geography and geology of the Site

The selected remedy will excavate petroleum impacted soils from the source area and historic fill materials to a depth of 6 feet within the proposed building footprint areas and to a depth of 2 feet in the proposed landscaped areas. The selected alternative and development of the site have considered the geography and geology of the Site.

Current Institutional Controls

The Site was assigned an E-designation for hazardous materials as part of the rezoning action completed by the City. The compliance with the E-designation for hazardous materials will require the approval of the NYC Office of Environmental Remediation (NYCOER) of this RAWP. NYCOER must approve this RAWP in the form of a Notice to Proceed (NTP) letter before building permits will be released by the NYC Department of Buildings (DOB). Documentation in the form of a Final Engineering Report (FER) for site remediation must be approved by NYCOER in the form of a Notice of Satisfaction (NOS) before the NYCDOB will issue permanent Certificates of Occupancy for the new buildings.

3.7 SUMMARY OF SELECTED REMEDIAL ACTIONS

Summary of the Remedy

The remedy recommended for the Site is a Track 4 alternative (Alternative 2) which consists of the excavation of petroleum impacted soil to a depth of 9 feet within the UST source area (USTs 1 & 2), excavation and disposal of historic fill within the planned landscape areas with soil above Restricted Residential SCOs and capping the entire Site with the building foundations, concrete walkways / driveways or 2 ft of soil meeting Restricted Residential SCOs. The remedy will include the following items:

1. Removal of underground storage tanks (USTs);
2. Excavation of soil/fill exceeding groundwater protection SCOs for those VOC parameters in groundwater above standards as listed in **Table 1** to depths as great as 9 feet below grade, or as needed to remove impacted soil in the UST source area;
3. Screening for indications of contamination (by visual means, odor, and monitoring with PID) of all excavated soil during any intrusive Site work;
4. Excavation and disposal of historic fill materials with parameters above RRSCOs from the buildings basement level foundations;
5. Capping areas of the Site in which soil is present with parameters above Restricted Residential SCOs with the concrete building slab, concrete driveways, concrete pavers or 2 feet of clean soil meeting RRSCOs;

6. Collection and analysis of end-point samples to evaluate the performance of the remedy with respect to attainment of groundwater protection SCOs for VOCs present in groundwater above standards;
7. Appropriate off-Site disposal of all material removed from the Site in accordance with all Federal, State and local rules and regulations for handling, transport, and disposal;
8. Import of materials to be used for backfill and cover in compliance with: (1) chemical limits and other specifications included in **Table 1**, (2) all Federal, State and local rules and regulations for handling and transport of material.
9. Implementation of a Site Management Plan (SMP) for long term maintenance of the Engineering Controls. An Environmental Easement will be filed against the Site to ensure implementation of the SMP.

All responsibilities associated with the Remedial Action, including permitting requirements and pretreatment requirements, will be addressed in accordance with all applicable Federal, State and local rules and regulations.

Remedial activities will be performed at the Site in accordance with this NYSDEC-approved RAWP. Any anticipated deviations to the RAWP shall be submitted to the NYSDEC for review.

4.0 REMEDIAL ACTION PROGRAM

The objective of this section of the Remedial Action Work Plan, is to present a scope of work which will be approved by NYSDEC and when completely implemented will ready the BCP site for development under the Contemplated Use consistent with the requirements of the Brownfield Cleanup Program.

4.1 GOVERNING DOCUMENTS

Governing documents and procedures included in the Remedial Work Plan include a Site-specific Health and Safety Plan (HASP), a Community Air Monitoring Plan (CAMP), a Citizen Participation Plan, a Soil Management Plan (SoMP), a Quality Assurance Project Plan (QAPP), fluid management procedures, and contractors' site operations and quality control procedures. Highlights of these documents and procedures are provided in the following sections.

4.1.1 Health & Safety Plan (HASP)

Contractors and subcontractors will have the option of adopting this HASP or developing their own site-specific document. If a contractor or subcontractor chooses to prepare their own HASP, the Remedial Engineer will insure that it meets the minimum requirements as detailed in the site-specific HASP prepared for the Site.

Activities performed under the HASP will comply with applicable parts of OSHA Regulations, primarily 29 CFR Parts 1910 and 1926. Modifications to the HASP may be made with the approval of the Remedial Engineer (RE), Site Safety Manager (SSM) and/or Project Manager (PM).

All remedial work performed under this plan will be in full compliance with governmental requirements, including Site and worker safety requirements mandated by Federal OSHA.

The Volunteer and associated parties preparing the remedial documents submitted to the State and those performing the construction work, are completely responsible for the preparation of an

appropriate Health and Safety Plan and for the appropriate performance of work according to that plan and applicable laws.

The Health and Safety Plan (HASP) and requirements defined in this Remedial Action Work Plan pertain to all remedial and invasive work performed at the Site until the issuance of a Certificate of Completion.

The Site Safety Coordinator will be Ms. Chawinie Miller. Her resume is provided in **Attachment F**. Confined space entry will comply with all OSHA requirements to address the potential risk posed by combustible and toxic gasses. A copy of the Site Specific Health and Safety Plan is provided in **Attachment B**.

4.1.2 Quality Assurance Project Plan (QAPP)

The fundamental QA objective with respect to accuracy, precision, and sensitivity of analysis for laboratory analytical data is to achieve the QC acceptance of the analytical protocol. The accuracy, precision and completeness requirements will be addressed by the laboratory for all data generated.

Collected samples will be appropriately packaged, placed in coolers and shipped via overnight courier or delivered directly to the analytical laboratory by field personnel. Samples will be containerized in appropriate laboratory provided glassware and shipped in plastic coolers. Samples will be preserved through the use of ice or a cold-pak(s) to maintain a temperature of 4°C.

Dedicated disposable sampling materials will be used for both soil and groundwater samples (if collected), eliminating the need to prepare field equipment (rinsate) blanks. However, if non-disposable equipment is used, (stainless steel scoop, etc.) field rinsate blanks will be prepared at the rate of 1 for every eight samples collected.

Decontamination of non-dedicated sampling equipment will consist of the following:

- Gently tap or scrape to remove adhered soil
- Rinse with tap water
- Wash withalconox® detergent solution and scrub
- Rinse with tap water
- Rinse with distilled or deionized water

Prepare field blanks by pouring distilled or deionized water over decontaminated equipment and collecting the water in laboratory provided containers. Trip blanks will accompany samples each time they are transported to the laboratory. Matrix spike and matrix spike duplicates (MS/MSD) will be collected at the rate of one per 20 samples submitted to the laboratory. Laboratory reports will be upgradeable to ASP category B deliverables for use in the preparation of a data usability report (DUSR). The QAPP for the Site is provided in **Attachment C**.

4.1.3 Construction Quality Assurance Plan (CQAP)

All construction work related to the remedy (i.e. soil excavation) will be monitored by EBC / AMC field personnel under the direct supervision of the Remedial Engineer. Monitoring during soil excavation will be performed to protect the health of site workers and the surrounding community. A Health and Safety Plan (HASP) and Community Air Monitoring Plan (CAMP) have been specifically developed for this project. These plans specify the monitoring procedures, action levels, and contingency measures that are required to protect public health.

All intrusive and soil disturbance activities will be monitored by an environmental professional (EP) under the direct supervision of the Remedial Engineer who will record observations in the site field book and complete a photographic log of the daily activities. The EP will provide daily updates to the Project Manager and Remedial Engineer who will both make periodic visits to the site as needed to assure construction quality. Daily updates will also be submitted to the NYSDEC. See section 4.4.1 Daily Reports.

4.1.4 Soil/Materials Management Plan (SoMP)

A SoMP has been prepared for excavation, handling, storage, transport and disposal of all soils/materials that are disturbed / excavated at the Site. The SoMP includes all of the controls that will be applied to these efforts to assure effective, nuisance-free performance in compliance with all applicable Federal, State and local laws and regulations. The SoMP is presented in Section 5.4.

4.1.5 Erosion and Sediment Control Plan (ESCP)

Erosion and sediment controls will be performed in conformance with requirements presented in the New York State Guidelines for Urban Erosion and Sediment Control. Typical measures that will be utilized at various stages of the project to limit the potential for erosion and migration of soil include the use of hay bales, temporary stabilized construction entrances/exits, placement of silt fencing and/or hay bales around soil stockpiles, and dust control measures.

4.1.6 Community Air Monitoring Plan (CAMP)

The CAMP provides measures for protection for on-site workers and the downwind community (i.e., off-site receptors including residences, businesses, and on-site workers not directly involved in the remedial work) from potential airborne contaminant releases resulting from remedial activities.

The action levels specified require increased monitoring, corrective actions to abate emissions, and/or work shutdown. Additionally, the CAMP helps to confirm that the remedial work did not spread contamination off-site through the air. The primary concerns for this site are vapors, nuisance odors and dust particulates.

The primary concerns for this site are vapors, nuisance odors and dust particulates. The CAMP prepared for implementation of the RAWP is provided in **Attachment D**.

4.1.7 Contractors Site Operations Plan (SOP)

The Remedial Engineer has reviewed all plans and submittals for this remedial project (including those listed above and contractor and sub-contractor document submittals) and confirms that they

are in compliance with this RAWP. The Remedial Engineer is responsible to ensure that all later document submittals for this remedial project, including contractor and sub-contractor document submittals, are in compliance with this RAWP. All remedial documents will be submitted to NYSDEC and NYSDOH in a timely manner and prior to the start of work.

4.1.8 Citizen Participation Plan (CPP)

The Citizen Participation Plan prepared for this project is provided in **Attachment E**. The public will be informed of key project documents and events through the distribution of fact sheets through the Department's List Serv. The public was initially informed of the Site and the opportunity to join the List Serv through an ad placed in the local newspaper and mailed fact sheets.

No changes will be made to approved Fact Sheets authorized for release by NYSDEC without written consent of the NYSDEC. No other information, such as brochures and flyers, will be included with the Fact Sheet mailing.

Document repositories have been established at the following locations and contain all applicable project documents:

New York Public Library - Mott Haven Branch

321 East 140th Street
Bronx, NY 10454
(718) 665-4878

Hours:

Mon 10:00 AM - 6:00 PM
Tue 10:00 AM - 7:00 PM
Wed 10:00 AM - 7:00 PM
Thu 10:00 AM - 6:00 PM
Fri 10:00 AM - 5:00 PM
Sat 10:00 AM - 5:00 PM
Sun closed

Bronx 1 Community District Information
Chairperson: George Rodriguez
District Manager: Mr. Cedric Loftin
Address: 3024 Third Avenue, Bronx, NY, 10455
Phone: 718-585-7117

4.2 GENERAL REMEDIAL ACTION INFORMATION

4.2.1 Project Organization

The Project Manager for the Remedial Activity will be Mr. Robert Bennett. Overall responsibility for the BCP project will be Mr. Charles B. Sosik, P.G., P.HG. The Remedial Engineer for this project is Mr. Ariel Czemerinski, P.E. Resumes of key personnel involved in the Remedial Action are included in **Attachment F**.

4.2.2 Remedial Engineer

The Remedial Engineer for this project will be Mr. Ariel Czemerinski, P.E. The Remedial Engineer is a registered professional engineer licensed by the State of New York. The Remedial Engineer will have primary direct responsibility for implementation of the remedial program for the Site. The Remedial Engineer will certify in the Final Engineering Report that the remedial activities were observed by qualified environmental professionals under his supervision and that the remediation requirements set forth in the Remedial Action Work Plan and any other relevant provisions of ECL 27-1419 have been achieved in full conformance with that Plan. Other Remedial Engineer certification requirements are listed later in this RAWP.

The Remedial Engineer will review all pre-remedial plans submitted by contractors and subcontractors involved in all aspects of remedial construction, including soil excavation, stockpiling, characterization, removal and disposal, air monitoring, emergency spill response services, import of back fill material, and management of waste transport and disposal, and will certify compliance in the Final Remediation Report. The Remedial Engineer will provide the certifications listed in Section 10.1 in the Final Engineering Report.

4.2.3 Remedial Action Schedule

The remedial action will begin with mobilization of equipment and material to the Site, which will begin following RAWP approval and 10 days after the distribution of the remedial construction Fact Sheet. A pre-construction meeting will be held among NYSDEC, the Remedial Engineer, and the selected remedial contractor prior to site mobilization. Mobilization will be

followed by soil removal and disposal and confirmation sampling. The work is expected to take 6 months as part of the construction excavation and foundation installation.

4.2.4 Work Hours

The hours for operation of remedial construction will conform to the New York City Department of Buildings construction code requirements or according to specific variances issued by that agency. DEC will be notified by the Applicant of any variances issued by the Department of Buildings. NYSDEC reserves the right to deny alternate remedial construction hours.

4.2.5 Site Security

A construction fence will be erected around the entire property as required by the NYC Department of Buildings. The fence will be maintained as required and secured at the end of each work day.

4.2.6 Traffic Control

The Volunteer's construction management personnel will direct the arrival or departure of construction vehicles, and provide flag services as needed to maintain safe travel exiting and entering the Site from Third Avenue. All traffic will enter and leave the Site via existing gates on Third Avenue. The excavation contractor will direct the arrival or departure of construction vehicles, and provide flag services as needed to maintain safe travel exiting and entering the Site from the Third Avenue exit / entrance. Traffic related to the Work Plan activity will not require the staging of trucks or equipment along the streets. All trucks and equipment will remain on-site within the construction fence. The local transport route will be as follows:

- **ENTERING SITE** - from the Major Deegan Expressway (I87) heading south; take the Willis Ave / Third Ave Exit (Exit 2) and turn right heading north on Bruckner Blvd. to Third Avenue. Turn left, heading west on Third Avenue 1 block to Site entrance on the left.
- **EXITING SITE** – Turn left onto Third Avenue heading northeast to 138th Street. Turn right onto 138th Street heading southeast to Willis Avenue. Turn right onto Willis

Avenue heading south to E. 135th Street. Turn right on E.135th Street and continue to the on-ramp (bearing left) for the Major Deegan Expressway (I-87).

- This route was designed to minimize or eliminate the time trucks will be on local streets. Equipment and trucks needed to install the grade beam will remain within the Site boundaries and not on residential Streets. Site personnel will be required to park on Site or in legal all-day on-street parking spaces, near the Site or in an off-street parking lot/garage.

A map showing the truck routes is included as **Figure 12**.

4.2.7 Worker Training and Monitoring

An excavation contractor with appropriate experience, personnel and training (40 hr OSHA) is required to perform the removal of the UST and VOC impacted soil. After this material is removed the contractor will remove historic fill as needed for the foundation installation and landscaped areas. The excavation contractor's on-site personnel engaged in historic fill and / or native soil removal (if present) will have a minimum of 24 hour Hazardous Waste Operations and Emergency Response Operations training.

All field personnel involved in remedial activities will participate in training, if required under 29 CFR 1910.120, including 24 and 40-hour hazardous waste operator training and annual 8-hour refresher training. The Site Safety Officer will be responsible for maintaining workers training records.

Personnel entering any exclusion zone will be trained in the provisions of the HASP and be required to sign a HASP acknowledgment.

All on-site personnel engaged in remedial or sampling activities must receive adequate site-specific training in the form of an on-site Health and Safety briefing prior to participating in field work with emphasis on the following:

- Protection of the adjacent community from hazardous vapors and / or dust which may be released during intrusive activities.
- Identification of chemicals known or suspected to be present on-site and the health effects and hazards of those substances.
- The need for vigilance in personnel protection, and the importance of attention to proper use, fit and care of personnel protective equipment.
- Decontamination procedures.
- Site control including work zones, access and security.
- Hazards and protection against heat or cold.
- The proper observance of daily health and safety practices, such as entry and exit of work zones and site. Proper hygiene during lunch, break, etc.
- Emergency procedures to be followed in case of fire, explosion and sudden release of hazardous gases.

4.2.8 Agency Approvals

The Applicant has addressed all SEQRA requirements for this Site. All permits or government approvals required for remedial construction have been, obtained prior to the start of remedial construction.

The planned end use for the Site is in conformance with the current zoning for the property as determined by New York City Department of Planning. A Certificate of Completion will not be issued for the project unless conformance with zoning designation is demonstrated.

A complete list of all local, regional and national governmental permits, certificates or other approvals or authorizations required to perform the remedial and development work is attached in **Table 15**. This list includes a citation of the law, statute or code to be complied with, the originating agency, and a contact name and phone number in that agency. This list will be updated in the Final Remediation Report.

4.2.9 NYSDEC BCP Signage

A project sign will be erected at the main entrance to the Site prior to the start of any remedial activities. The sign will indicate that the project is being performed under the New York State Brownfield Cleanup Program. The sign will meet the detailed specifications provided by the NYSDEC Project Manager and contained in **Attachment G**.

4.2.10 Pre-Construction Meeting with NYSDEC

A pre-construction meeting or conference call with the Project Manager, Remedial Engineer, Construction Manager, Owner's Representative and the NYSDEC will take place prior to the start of major construction activities.

4.2.11 Emergency Contact Information

An emergency contact sheet with names and phone numbers is included in **Table 16**. That document will define the specific project contacts for use by NYSDEC and NYSDOH in the case of a day or night emergency.

4.2.12 Remedial Action Costs

The total estimated cost of the Remedial Action is \$ 1,857,574. An itemized and detailed summary of estimated costs for all remedial activity is provided as **Attachment H**. This will be revised based on actual costs and submitted as an Appendix to the Final Remediation Report.

4.3 SITE PREPARATION

4.3.1 Mobilization

Mobilization will include the delivery of construction equipment and materials to the site. All construction personnel will receive site orientation and training in accordance with the site specific HASP, CAMP and established policies and procedures to be followed during the implementation of the RAWP. The remediation contractor, construction manager and all associated subcontractors will each receive a copy of the RAWP and the site specific HASP and will be briefed on their contents.

4.3.2 Erosion and Sedimentation Controls

Soil erosion and sediment control measures for management of storm water will be installed in accordance with the New York Guidelines for Urban Erosion and Sediment Control. Haybales and/or silt fence will be placed by the remedial contractor at locations surrounding excavation areas and within the perimeter fencing as needed, to control stormwater runoff and surface water from exiting the excavation. These control measures will be installed prior to initiating the soil excavation.

4.3.3 Stabilized Construction Entrance(s)

Stabilized construction entrances will be installed at all points of vehicle ingress and egress to the Site. The construction entrance will be stabilized with crushed stone, gravel, plywood sheeting, existing asphalt or existing concrete as necessary to provide a safe egress and ingress to the Site. The stabilized entrances will be inspected on a daily basis during soil loading activities and reinforced as needed with additional stone/concrete material to prevent the accumulation of ruts, mud or soil.

4.3.4 Utility Marker and Easements Layout

The Applicant and its contractors are solely responsible for the identification of utilities that might be affected by work under the RAWP and implementation of all required, appropriate, or necessary health and safety measures during performance of work under this RAWP. The Applicant and its contractors are solely responsible for safe execution of all invasive and other work performed under this RAWP. The Applicant and its contractors must obtain any local, State or Federal permits or approvals pertinent to such work that may be required to perform work under this RAWP. Approval of this RAWP by NYSDEC does not constitute satisfaction of these requirements.

The presence of utilities and easements on the Site has been investigated by the Remedial Engineer. It has been determined that no risk or impediment to the planned work under this Remedial Action Work Plan is posed by utilities or easements on the Site.

4.3.5 Sheeting and Shoring

Appropriate management of structural stability of on-Site or off-Site structures during on-Site activities including excavation is the sole responsibility of the Applicant and its contractors. The Applicant and its contractors are solely responsible for safe execution of all invasive and other work performed under this Plan. The Applicant and its contractors must obtain any local, State or Federal permits or approvals that may be required to perform work under this Plan. Further, the Applicant and its contractors are solely responsible for the implementation of all required, appropriate, or necessary health and safety measures during performance of work under the approved Plan.

4.3.6 Equipment and Material Staging

Equipment used for excavation work will be staged on Site within the fenced perimeter. All equipment and work materials will be staged on-Site in areas as designated by the General Contractor, and / or Construction Site Superintendent.

4.3.7 Decontamination Area

All materials and equipment (except disposable items) will be decontaminated on specially constructed “pads”, an existing asphalt surface or an existing concrete surface located at the exit point from the Site. At a minimum, the constructed pad (if used) will consist of a layer of crushed stone underlain by an impervious plastic liner that has been graded to drain to the interior of the Site. If an existing surface is used it will be surrounded by hay bales or other silt controlling materials. The pad will be sized to accommodate the largest piece of equipment used on the project. Where effective, the equipment will be “dry” decontaminated using a broom and/or brushes. If significant amounts of soil or other contaminants remain after the dry decontamination, the equipment will also be pressure washed before leaving the Site. Disposable items will be containerized within the site and transported for appropriate off- site disposal.

4.3.8 Site Fencing

The lot currently has an 8 ft high chain link fence and gate along Third Avenue (south side). It also has a 6 ft high chain link fence along the north and a portion of the east property lines. The remainder of the east property line is bordered by an adjacent building. The west side of the

property is bordered by the Harlem River. The fences and gates will be maintained during the pile driving activity and properly secured at the end of the day.

4.3.9 Demobilization

Demobilization will consist of the restoration of material staging areas and the disposal of materials and/or general refuse in accordance with acceptable rules and regulations. Materials used in remedial activities will be removed and disposed properly. All equipment will be decontaminated prior to leaving the Site.

4.4 REPORTING

All daily and monthly Reports will be included in the Final Engineering Report.

4.4.1 Daily Reports

Daily reports will be submitted to NYSDEC and NYSDOH Project Managers by the end of each day in which remedial activity takes place. Daily reports will include:

- An update of progress made during the reporting day;
- A summary of any and all complaints with relevant details (names, phone numbers);
- A summary of CAMP finding, including excursions;
- An explanation of notable Site conditions.

Daily reports are not intended to be the mode of communication for notification to the NYSDEC of emergencies (accident, spill), requests for changes to the RAWP or other sensitive or time critical information. However, such conditions must also be included in the daily reports. Emergency conditions and changes to the RAWP will be addressed directly to NYSDEC Project Manager via personal communication.

These reports will include a summary of air sampling results, odor and dust problems and corrective actions, and all complaints received from the public.

4.4.2 Monthly Reports

Monthly reports will be submitted to NYSDEC and NYSDOH Project Managers within one week following the end of the month of the reporting period and will include:

- Activities relative to the Site during the previous reporting period and those anticipated for the next reporting period, including a quantitative presentation of work performed (i.e. tons of material exported and imported, etc.);
- Description of approved activity modifications, including changes of work scope and/or schedule;
- Sampling results received following internal data review and validation, as applicable; and,
- An update of the remedial schedule including the percentage of project completion, unresolved delays encountered or anticipated that may affect the future schedule, and efforts made to mitigate such delays.

4.4.3 Other Reporting

Photographs will be taken of all remedial activities and submitted to NYSDEC in digital (JPEG, PDF) format. Photos will illustrate all remedial program elements and will be of acceptable quality. Representative photos of the Site prior to any Remedial Actions will be provided. Representative photos will be provided of each contaminant source, source area and Site structures before, during and after remediation. Photos will be included in the daily reports as needed, and a comprehensive collection of photos will be included in the Final Engineering Report.

Job-site record keeping for all remedial work will be appropriately documented. These records will be maintained on-Site at all times during the project and be available for inspection by NYSDEC and NYSDOH staff.

4.4.4 Complaint Management Plan

Complaints from the public regarding nuisance or other Site conditions including noise, odor, truck traffic etc., will be recorded in the Site field book and reported to the NYSDEC via email on the same day as the complaint is received.

4.4.5 Deviations from the Remedial Action Work Plan

Minor deviations from the RAWP will be identified in the daily update report and will be noted in the Final Engineering Report. When deviations are reported, a brief discussion will be provided which will state the following:

- Reasons for deviating from the approved RAWP;
- Effect of the deviations on overall remedy.

Major changes to the scope of work must be discussed with the NYSDEC and the NYSDOH prior to implementation. If the changes are considered to be significant enough, an addendum to the RAWP Work Plan will be prepared and submitted to NYSDEC / NYSDOH for review.

5.0 REMEDIAL ACTION: MATERIAL REMOVAL FROM SITE

The removal of materials from the Site includes the following; the removal of existing USTs, the excavation and off-Site disposal of petroleum impacted soil to a depth of 9 feet within the UST source area (150 cy) and the excavation and offsite disposal of historic fill soil to a depth of 4 feet within the building footprint area (5,094 cy).

Soil excavation and loading will be performed using conventional equipment such as track-mounted excavators, backhoes and loaders.

All excavation work will be performed in accordance with the Site-specific HASP and CAMP. Removal of the petroleum impacted soil and historic fill soil will be performed by a qualified contractor and trained personnel (Minimum 24HR OSHA HAZWOPER). Removal of underground storage tanks (UST) believed to be present will be removed by a qualified contractor and trained personnel (Minimum 40HR OSHA HAZWOPER).in accordance with DER-10, NYSDEC PBS regulations and NYC Fire Department regulations. It is anticipated that the excavation of petroleum impacted and historic fill materials will be performed by the excavation contractor for the construction project.

An excavation plan showing the excavation depths to achieve the Track 4 remedy is provided in **Figure 13**. Some dewatering may be needed for the excavation of contaminated areas but is not anticipated for foundation construction.

5.1 UST REMOVAL METHODS

USTs, which are believed to be present at the Site, will be removed in accordance with the procedures described under the NYSDEC Memorandum for the Permanent Abandonment of Petroleum Storage Tanks and Section 5.5 of Draft DER-10 as follows:

- Remove all product to its lowest draw-off point
- Drain and flush piping into the tank
- Vacuum out the tank bottom consisting of water product and sludge
- Dig down to the top of the tank and expose the upper half of the tank

- Remove the fill tube and disconnect the fill, gauge, product and vent lines and pumps. Cap and plug open ends of lines
- Temporarily plug all tank openings, complete the excavation, remove the tank and place it in a secure location
- Render the tank safe and check the tank atmosphere to ensure that petroleum vapors have been satisfactorily purged from the tank
- Clean tank or remove to a storage yard for cleaning
- If the tank is to be moved it must be transported by licensed waste transporter. Plug and cap all holes prior to transport leaving a 1/8 inch vent hole located at the top of the tank during transport
- After cleaning the tank must be made acceptable for disposal at a scrap yard cleaning the tank interior with a high pressure rinse and cutting the tank in several pieces.

During the tank and pipe line removal the following field observations should be made and recorded:

- A description and photographic documentation of the tank and pipe line condition (pitting, holes, staining, leak points, evidence of repairs, etc.)
- Examination of the excavation floor and sidewalls for physical evidence of contamination (odor, staining, sheen, etc.)
- Periodic field screening (through bucket return) of the floor and sidewalls of the excavation with a calibrated photoionization detector (PID).

5.2 SOIL CLEANUP OBJECTIVES

The Soil Cleanup Objectives for this Site are listed in **Table 1**. **Table 7** summarizes all soil samples that exceed unrestricted SCOs at the Site. Spider maps showing all soil samples that exceed the unrestricted SCO are shown in **Figure 9**.

5.3 REMEDIAL PERFORMANCE EVALUATION (POST EXCAVATION END-POINT SAMPLING)

Post excavation (endpoint) soil samples will be collected from across the Site to verify that remedial goals have been achieved. Endpoint soil samples will be collected from the Site as follows:

- (1) Endpoint soil samples will be collected following the removal of soil from the petroleum areas to verify that remedial goals have been achieved (**Figure 14**). The endpoint soil samples from the petroleum area hotspot will be analyzed for VOCs and SVOCs. Sidewall endpoint soil samples will also be collected from each of the petroleum impacted and SVOC hotspot excavation areas.

5.3.1 End-Point Sampling Frequency

Endpoint sampling frequency will be in accordance with DER-10 section 5.4 which recommends the collection of one bottom sample per 900 sf of bottom area and one sidewall sample per 30 liner feet.

5.3.2 Methodology

Collected samples be placed in glass jars supplied by the analytical laboratory and stored in a cooler with ice to maintain a temperature of 4 degrees C. Samples will either be picked up at the Site by a laboratory dispatched courier at the end of the day or transported back to the EBC /AMC office where they will be picked up the following day by the laboratory courier. All samples will be analyzed by a NYSDOH ELAP certified environmental laboratory

All post excavation (endpoint) soil samples will be analyzed for VOCs by EPA Method 8260B, SVOCs by EPA method 8270 PAHs.

5.3.3 Reporting of Results

Sample analysis will be provided by a New York State certified environmental laboratory. Laboratory reports will include ASP category B deliverables for use in the preparation of a data usability summary report (DUSR). All results will be provided in accordance with the NYSDEC Environmental Information Management System (EIMS) electronic data deliverable (EDD) format.

5.3.4 QA/QC

The fundamental QA objective with respect to accuracy, precision, and sensitivity of analysis for laboratory analytical data is to achieve the QC acceptance of the analytical protocol. The accuracy, precision and completeness requirements will be addressed by the laboratory for all data generated.

Collected samples will be appropriately packaged, placed in coolers and shipped via overnight courier or delivered directly to the analytical laboratory by field personnel. Samples will be containerized in appropriate laboratory provided glassware and shipped in plastic coolers. Samples will be preserved through the use of ice or cold-pak(s) to maintain a temperature of 4°C.

Dedicated disposable sampling materials will be used for soil samples, eliminating the need to prepare field equipment (rinsate) blanks. However, if non-disposable equipment is used, (stainless steel scoop, etc.) field rinsate blanks will be prepared at the rate of 1 for every eight samples collected. Field blanks will be prepared by pouring distilled or deionized water over decontaminated equipment and collecting the water in laboratory provided containers.

Trip blanks will accompany samples each time they are transported to the laboratory. Matrix spike and matrix spike duplicates (MS/MSD) will be collected at the rate of one per 20 samples submitted to the laboratory.

5.3.5 DUSR

The DUSR provides a thorough evaluation of analytical data without third party data validation. The primary objective of a DUSR is to determine whether or not the data, as presented, meets the

site/project specific criteria for data quality and data use. Verification and/or performance monitoring samples collected under this RAWP will be reviewed and evaluated in accordance with the Guidance for the Development of Data Usability Summary Reports as presented in Appendix 2B of DER-10. The completed DUSR for verification/performance samples collected during implementation of this RAWP will be included in the final Engineering Report.

5.3.6 Reporting of End-Point Data in FER

All endpoint data collected as part of this remedial action will be summarized and presented in the Final Engineering Report. The summary tables will include comparison of results to unrestricted SCOs to verify attainment of Track 1. Laboratory reports and the DUSR will be included as an appendix in the FER.

5.4 ESTIMATED MATERIAL REMOVAL QUANTITIES

It is expected that 150 cubic yards (225 tons) of petroleum impacted soil will be generated by excavating the UST source area to a depth of 9 ft. Historic fill material was also documented throughout the Site to depths as great as 5 feet below grade. Therefore, an estimated 5,094 cubic yards (7,641 tons) of historic fill material will be generated by excavating the basement foundations of the new buildings.

5.5 SOIL/MATERIALS MANAGEMENT PLAN

Excavated soil will be secured and temporarily stored on-site until arrangements can be made for off-site disposal. As an alternative, pre-characterization samples may be collected to allow the soil to be loaded directly on to trucks for transport to the disposal facility. Excavated soils including petroleum contaminated soil and historic fill soil are expected to be classified as non-hazardous.

The final determination on classification will be based on the results of waste characterization analysis and the NYSDEC.

Soil excavation will be performed in accordance with the procedures described under Section 5.5 of DER-10 as follows:

- A description and photographic documentation of the excavation.
- Examination of the excavation floor and sidewalls for physical evidence of contamination (odor, staining, sheen, etc.).
- Periodic field screening (through bucket return) of the floor and sidewalls of the excavation with a calibrated photoionization detector (PID).

Final excavation depth, length, and width will be determined by the Remedial Engineer or his designee, and will depend on the horizontal and vertical extent of contaminated soils as identified through physical examination (PID response, odor, staining, etc.). Expansion of the excavation beyond the planned hotspot area is anticipated and can easily be accommodated.

The following procedure will be used for the excavation of impacted soil (as necessary and appropriate):

- Wear appropriate health and safety equipment as outlined in the HASP;
- Prior to excavation, ensure that the area is clear of utility lines or other obstructions. Lay plastic sheeting on the ground next to the area to be excavated;
- Using a rubber-tired backhoe or track mounted excavator, remove overburden soils and stockpile or dispose of separate from the impacted soil;
- If USTs are discovered, the NYSDEC will be notified and the best course of action to remove the structure should be determined in the field. This may involve the continued removal of overburden to access the top of the structure or continued trenching around the perimeter to minimize its disturbance;
- If physically contaminated soil is present (e.g., staining, odors, sheen, PID response, etc), an attempt will be made to remove it to the extent not limited by the site boundaries. If possible, physically impacted soil will be removed using the backhoe or excavator,

segregated from clean soils and overburden, and staged on separate dedicated plastic sheeting or live loaded into trucks from the disposal facility. Removal of the impacted soils will continue until visibly clean material is encountered and monitoring instruments indicate that no contaminants are present;

- Excavated soils which are temporarily stockpiled on-site will be covered with 6-mil polyethylene sheeting while disposal options are determined. Sheeting will be checked on a daily basis and replaced, repaired or adjusted as needed to provide full coverage. The sheeting will be shaped and secured in such a manner as to drain runoff and direct it toward the interior of the property;
- Once the Remedial Engineer is satisfied with the removal effort, verification or confirmatory samples will be collected from the excavation as described in **Section 6.2** of this document.

5.5.1 Excavation of Petroleum Contaminated Soil

Petroleum impacted soil is known to be present in the vicinity of a suspect UST located in the north-central area of the Site. The vertical extent is approximately 9 feet below grade. However, soil screening will be performed to determine the limits of the excavation with verification sampling performed to confirm removal of all petroleum impacted soil. The excavation of the petroleum soil will be performed by a qualified contractor and trained (24 hr HAZWOPER) personnel.

5.5.2 Excavation of Historic Fill Soil

Historic fill material is present beneath the Site to depths which vary from 5 to 8 feet below grade. The historic fill material contains SVOCs, metals and pesticides above unrestricted and restricted-residential use SCOs. Historic fill material will be segregated from non-contaminated native soils and disposed of off-Site at a permitted disposal facility.

Excavated historic fill materials will either be pre-characterized and live-loaded into trucks for transport to the disposal facility or temporarily stored on-Site until arrangements can be made for

off-Site disposal. It is anticipated that the historic fill material will be classified as non-hazardous material. If this material is classified as non-hazardous, then the excavation of historic fill material will be performed by the excavation contractor for the construction project using trained personnel (24 hr HAZWOPER). If this material is classified as hazardous, then 40 hr HAZWOPER trained personnel will be needed to perform the excavation of this material.

5.5.3 Excavation of Native Soils

Native soils are present directly below the fill materials and may be encountered during excavation for basement areas and foundation components during construction of the new buildings. Since excavation of the basement areas will begin following removal of petroleum / CVOC contaminated soil and historic fill, it is expected that native soils will not be contaminated. However, if evidence of contamination is discovered during the excavation of basement areas, the contamination will be removed to the extent possible and segregated from clean native soils for proper disposal. Clean native soils will be stockpiled on-site and characterized for reuse on-site in areas over-excavated to remove historic fill or petroleum. Any excess soil will be disposed of off-site as a beneficial re-use material upon approval by the NYSDEC Region 2's Division of Materials Management. Clean native soils utilized on-site will be subject to a testing program to verify that they meet restricted residential SCOs prior to use.

It is anticipated that the excavation of native soil materials will be performed by the excavation contractor for the construction project.

5.5.4 Soil Screening Methods

Visual, olfactory and PID soil screening and assessment will be performed by an environmental professional during all remedial and development excavations into known or potentially contaminated material (Residual Contamination Zone). Soil screening will be performed regardless of when the invasive work is done and will include all excavation and invasive work performed during the remedy and during development phase, such as excavations for foundations and utility work, prior to issuance of the COC.

All primary contaminant sources (including but not limited to tanks and hotspots) identified during Site Characterization, Remedial Investigation, and Remedial Action will be surveyed by a surveyor licensed to practice in the State of New York. This information will be provided on maps in the Final Engineering Report.

Screening will be performed by qualified environmental professionals. Resumes will be provided for all personnel responsible for field screening (i.e. those representing the Remedial Engineer) of invasive work for unknown contaminant sources during remediation and development work.

5.5.5 Stockpile Methods

Stockpiles will be inspected at a minimum once each week and after every storm event. Results of inspections will be recorded in a logbook and maintained at the Site and available for inspection by NYSDEC.

Stockpiles will be kept covered at all times with appropriately anchored tarps. Stockpiles will be routinely inspected and damaged tarp covers will be promptly replaced. Soils which exhibit strong odors will be completely sealed with heavy tarps or vapor suppressant foam.

5.5.6 Materials Excavation and Load Out

The Remedial Engineer or an EP under his/her supervision will oversee all invasive work and the excavation and load-out of all excavated material. The Volunteer and its contractors are solely responsible for safe execution of all invasive and other work performed under this Plan.

Loaded vehicles leaving the Site will be appropriately lined, tarped, securely covered, manifested, and placarded in accordance with appropriate Federal, State, local, and NYSDOT requirements (and all other applicable transportation requirements).

Where effective, the equipment will be “dry” decontaminated using a broom and/or brushes. If significant amounts of soil or other contaminants remain after the dry decontamination, the equipment will also be pressure washed before leaving the Site. The EP will be responsible for ensuring that all outbound trucks are dry-brushed or washed on the truck wash/equipment pad before leaving the Site until the remedial construction is complete. Locations where vehicles

enter or exit the Site shall be inspected daily for evidence of off-Site sediment tracking. The EP will be responsible for ensuring that all egress points for truck and equipment transport from the Site will be clean of dirt and other materials derived from the Site during Site remediation and development. Cleaning of the adjacent streets will be performed as needed to maintain a clean condition with respect to Site derived materials.

The Volunteer and associated parties preparing the remedial documents submitted to the State, and parties performing this work, are completely responsible for the safe performance of all invasive work, the structural integrity of excavations, and for structures that may be affected by excavations (such as building foundations and bridge footings).

The Remedial Engineer will ensure that Site development activities will not interfere with, or otherwise impair or compromise, remedial activities proposed in this Remedial Action Work Plan.

Development-related grading cuts and fills will not interfere with, or otherwise impair or compromise, the performance of remediation required by this plan.

Mechanical processing of historical fill material and contaminated soil on-Site is prohibited. All primary contaminant sources (including but not limited to tanks and hotspots) identified during Site Characterization, Remedial Investigation, and Remedial Action will be located and shown on maps to be reported in the Final Engineering Report.

5.5.7 Materials Transport Off-Site

All transport of materials will be performed by licensed haulers in accordance with appropriate local, State, and Federal regulations, including 6 NYCRR Part 364. Haulers will be appropriately licensed and trucks properly placarded.

All traffic will enter and leave the Site via existing gates on Third Avenue. The excavation contractor will direct the arrival or departure of construction vehicles, and provide flag services as needed to maintain safe travel exiting and entering the Site from the Third Avenue exit / entrance. The local transport route will be as follows:

- ENTERING SITE - from the Major Deegan Expressway (I87) heading south; take the Willis Ave / Third Ave Exit (Exit 2) and turn right heading north on Bruckner Blvd. to Third Avenue. Turn left, heading west on Third Avenue 1 block to Site entrance on the left.
- EXITING SITE – Turn left onto Third Avenue heading northeast to 138th Street. Turn right onto 138th Street heading southeast to Willis Avenue. Turn right onto Willis Avenue heading south to E. 135th Street. Turn right on E.135th Street and continue to the on-ramp (bearing left) for the Major Deegan Expressway (I-87).
- This route was designed to minimize or eliminate the time trucks will be on local streets. Equipment and trucks needed to install the grade beam will remain within the Site boundaries and not on residential Streets. Site personnel will be required to park on Site or in legal all-day on-street parking spaces, near the Site or in an off-street parking lot/garage.

These routes are shown in **Figure 12**.

Trucks will be prohibited from stopping and idling in the neighborhood outside the project Site. Egress points for truck and equipment transport from the Site will be kept clean of dirt and other materials during Site remediation and development. Wet loads are not anticipated since the entire site will be dewatered prior to excavating soils. However, if wet soils are excavated they will be stockpiled within the excavation to dry or blended with dry soils. No loads of material capable of generating free liquid will be allowed to leave the Site. All trucks will be inspected, dry-brushed and / or washed, as needed, before leaving the site.

5.5.8 Materials Disposal Off-Site

Multiple disposal facility designations may be employed for the materials removed from the Site. Once final arrangements have been made, the disposal location(s) will be reported to the NYSDEC Project Manager. It is anticipated that the soil will be disposed of at up to 3 different facilities, based on the following classification:

- Non Hazardous - Contaminated (petroleum) Low Lead < 1,000 mg/kg
- Non Hazardous - Contaminated (historic fill) Low Lead < 1,000 mg/kg
- Uncontaminated Native Soil - meets NJDSC Criteria for beneficial Reuse

The total quantity of material expected to be disposed off-Site is 5,244 cubic yards, including 150 cubic yards of petroleum impacted soil and 5,094 cubic yards of historic fill material from the cellar level and foundation excavations.

Hazardous Soil Disposal and Transport

Although not anticipated to be present, soil classified as hazardous will be shipped under a hazardous waste manifest system. All hazardous waste transported and disposed of must have a USEPA ID Number and waste code and must be distributed in accordance with the regulatory requirements.

The multi-part manifest will be filled out for each load of soil shipped off of the Site. At a minimum, the following information will be recorded on each manifest:

- 1) Generator's Name, Address, and Phone Number
- 2) Destination Facility Name, Address and Phone Number
- 3) EPA ID Number
- 4) Waste classification code
- 5) Transporter Name, Address, Phone Number, License Plate Number, Driver Name, and SW Haulers Permit #
- 6) Signatures – Generator or an authorized agent for the generator shall print, sign, and date each non-hazardous material manifest after each truck is loaded. The transporter shall then sign and date noting time material was picked up at the site. Both the transporter and a representative of the disposal facility will sign the non-hazardous material manifest when the material has been delivered to disposal facility.

Non-Hazardous Soil Disposal and Transport

Non-hazardous historic fill material and petroleum contaminated soil classified as non-hazardous, will be handled, at a minimum, as a Municipal Solid Waste per 6NYCRR Part 360-1.2. Historical fill material and contaminated soils from the Site are prohibited from being disposed at Part 360-16 Registration Facilities (also known as Soil Recycling Facilities).

Soils that are contaminated but non-hazardous and are being removed from the Site are considered by the Division of Materials Management (DMM) in NYSDEC to be Construction and Demolition (C/D) materials with contamination not typical of virgin soils. These soils may be sent to a permitted Part 360 landfill. They may be sent to a permitted C/D processing facility without permit modifications only upon prior notification of NYSDEC Region 2 DSHM. This material is prohibited from being sent or redirected to a Part 360-16 Registration Facility. In this case, as dictated by DMM, special procedures will include, at a minimum, a letter to the C/D facility that provides a detailed explanation that the material is derived from a DER remediation Site, that the soil material is contaminated and that it must not be redirected to on-Site or off-Site Soil Recycling Facilities. The letter will provide the project identity and the name and phone number of the Remedial Engineer. The letter will include as an attachment a summary of all chemical data for the material being transported.

Soil classified as non-hazardous fill will be transported under a non-hazardous waste manifest obtained from the selected disposal facility. The multi-part manifest will be filled out for each load of soil shipped off of the Site. At a minimum, the following information will be recorded on each manifest:

- 1) Generator's Name, Address, and Phone Number
- 2) Destination Facility Name, Address and Phone Number
- 3) Transporter Name, Address, Phone Number, License Plate Number, Driver Name, and SW Haulers Permit #
- 4) Signatures – Generator or an authorized agent for the generator shall print, sign, and date each non-hazardous material manifest after each truck is loaded. The transporter shall then sign and date noting time material was picked up at the site. Both the transporter and

a representative of the disposal facility will sign the non-hazardous material manifest when the material has been delivered to disposal facility.

A copy of the manifest will be retained by on-Site personnel for each shipment. Final signed manifests will be forwarded by the disposal facility to the generator. Copies of the final manifests will be presented in the FER.

Clean Soil Disposal

Clean native soil removed from the Site for development purposes (i.e. basement levels) will be handled as unregulated or beneficial use disposal. This soil will undergo a testing program to confirm that it meets Unrestricted Use SCOs or Residential / Groundwater Protection SCOs prior to unregulated disposal or meets Restricted Residential Use SCOs prior to reuse on-Site. Confirmation testing of clean soils will be in accordance with DER-10 Section 5.4(e)(10) as follows:

| Contaminant | VOCs | | SVOCs, Inorganics & PCBs/Pesticides | |
|--------------------|--|-------------------------|--|-----------------------------------|
| | Soil Quantity (cubic yards) | Discrete Samples | Composite | Discrete Samples/Composite |
| 0-50 | 1 | 1 | Each composite sample for analysis is created from 3-5 discrete samples from representative locations in the fill. | |
| 50-100 | 2 | 1 | | |
| 100-200 | 3 | 1 | | |
| 200-300 | 4 | 1 | | |
| 300-400 | 4 | 2 | | |
| 400-500 | 5 | 2 | | |
| 500-800 | 6 | 2 | | |
| 800-1000 | 7 | 2 | | |
| 1000 | Add an additional 2 VOC and 1 composite for each additional 1000 Cubic yards or consult with DER | | | |

Uncontaminated native soil confirmed by the above testing program and removed from the site, will be disposed of as C&D material (if approved) or sent to a beneficial re-use facility. The final destination of soils whether classified as contaminated or uncontaminated must be approved by the NYSDEC.

C&D and Scrap Metal Disposal

Concrete demolition material generated on the Site from building slabs, parking areas and other structures will be segregated, sized and shipped to a concrete recycling facility. Concrete

crushing or processing on-Site is prohibited. Asphalt removed from the parking areas will be sent to a separate recycling facility.

Additionally, it is common to encounter scrap metals and large boulders (greater than one foot in diameter) during excavation which may not be accepted by either the licensed disposal facility or the C&D facility. These materials will be segregated and subsequently recycled at local facilities. Uncontaminated metal objects will be taken to a local scrap metal facility.

Bricks and other C&D material are also not accepted by most soil disposal facilities if present at greater than 5% by volume. This material, if encountered, will be sent to a C&D landfill or other C&D processing facility if approved by the DEC. C&D material of this type is most often encountered on sites in which former basement structures have been filled in with material from demolishing a former building. There was no evidence of former basement areas identified during previous investigations performed at the Site.

Scale Tickets

All trucks to be utilized for transport of hazardous or non-hazardous contaminated soil shall be weighed before and after unloading at the disposal facility. Disposal facilities must provide truck scales capable of generating load tickets measured in tons. The tonnage transported and disposed will be determined by the disposal facility and reported on a certified scale ticket which will be attached to each returned manifest. Weights will be reported on the certified scale ticket as Tare and Gross weights.

C&D Transport Tickets / Bills of Lading

Bill of Lading system or equivalent will be used for the disposal of C&D and related materials. Documentation for materials disposed of at recycling facilities (such as metal, concrete, asphalt) and as non-regulated C&D will include transport tickets for each load stating the origin of the material, the destination of the material and the quantity transported. This information will be reported in the Final Engineering Report.

Disposal Facility Documentation

The following documentation will be obtained and reported by the Remedial Engineer for each disposal location used in this project to fully demonstrate and document that the disposal of material derived from the Site conforms with all applicable laws: (1) a letter from the Remedial Engineer or BCP Applicant to the receiving facility describing the material to be disposed and requesting formal written acceptance of the material. This letter will state that material to be disposed is contaminated material generated at an environmental remediation Site in New York State. The letter will provide the project identity and the name and phone number of the Remedial Engineer. The letter will include as an attachment a summary of all chemical data for the material being transported (including Site Characterization data); and (2) a letter from all receiving facilities stating it is in receipt of the correspondence (above) and is approved to accept the material. These documents will be included in the FER.

The Final Engineering Report will include an accounting of the destination of all material removed from the Site during this Remedial Action, including excavated soil, contaminated soil, historic fill, solid waste, and hazardous waste, non-regulated material, and fluids. Documentation associated with disposal of all material must also include records and approvals for receipt of the material. This information will also be presented in a tabular form in the FER.

5.5.9 Materials Reuse On-Site

Re-use of on-Site soil will only be allowed if the material is found to meet Restricted Residential Use SCOs through the verification testing program detailed above. The Remedial Engineer will ensure that procedures defined for materials reuse in this RAWP are followed and that unacceptable material will not remain on-Site.

Acceptable demolition material proposed for reuse on-Site, if any, will be sampled for asbestos. Concrete crushing or processing on-Site is prohibited. Contaminated on-Site material, including historic fill material and contaminated soil, removed for grading or other purposes will not be reused within a cover soil layer, within landscaping berms, or as backfill for subsurface utility lines.

5.5.10 Fluids Management

As the depth to groundwater at the site is approximately 6 to 7 feet below grade in the planned basement excavation areas, some dewatering operations may be employed during construction. Dewatering fluids will be handled, transported and disposed in accordance with applicable local, State, and Federal regulations. Liquids discharged into the New York City sewer system will be addressed through approval by the NYCDEP.

Dewatered fluids will not be recharged back to the land surface or subsurface of the Site. Dewatering fluids will be managed off-Site. Discharge of water generated during remedial construction to surface waters (i.e. a local pond, stream or river) is prohibited without a SPDES permit.

5.5.11 Backfill from Off-Site Sources

Off-site fill material may be needed to stabilize the entrance - exit areas of the Site, for temporary driveways for loading trucks and as an underlayment to structural components of the new buildings including slabs and footings. Recycled Concrete Aggregate (RCA) derived from recognizable and uncontaminated concrete and supplied by facilities permitted by, and in full compliance with Part 360-16 and DSNY regulations, is an acceptable form of backfill material. The Remedial Engineer is responsible for ensuring that the facility is compliant with the registration and permitting requirements of 6 NYCRR Part 360 and DSNY regulations at the time the RCA is acquired. RCA imported from compliant facilities does not require additional testing unless required by NYS DEC and DSNY under its terms of operations for the facility. Documentation of part 360-16 and DSNY compliance must be provided to the Remedial Engineer before the RCA is transported to the Site.

Fill material may also consist of virgin mined sand, gravel or stone products. Materials from a virgin mined source may be imported to the Site without testing provided that that the material meets the specifications of the geotechnical engineer, Remedial Engineer, and Redevelopment Construction Documents and that the source of the material is approved by the Remediation Engineer and the NYSDEC Project Manager.

The source approval process will require a review of the following information:

- The origin of the material;
- The address of the facility which mines/processes the material;
- A letter from the facility stating that the material to be delivered to the site is a virgin mined material and that it has not been co-mingled with other materials during processing or stockpiling.

All materials proposed for import onto the Site will be approved by the Remedial Engineer and will be in compliance with provisions in this RAWP prior to receipt at the Site. Material from industrial sites, spill sites or other potentially contaminated sites will not be imported to the Site.

The Final Engineering Report will include the following certification by the Remedial Engineer: “I certify that all import of soils from off-Site, including source evaluation, approval and sampling, has been performed in a manner that is consistent with the methodology defined in the Remedial Action Work Plan”.

Under no circumstances will fill materials be imported to the site without prior approval from the NYSDEC Project Manager. If sufficient documentation is not obtained, fill materials will be tested at a frequency consistent with that as specified in Table 4 of NYSDEC CP-51 Soil Cleanup Guidance Policy. Soils that meet ‘exempt’ fill requirements under 6 NYCRR Part 360, but do not meet backfill or cover soil objectives for this Site, will not be imported onto the Site without prior approval by NYSDEC. Solid waste will not be imported onto the Site.

5.5.12 Stormwater Pollution Prevention

Barriers and hay bale checks will be installed and inspected once a week and after every storm event. Results of inspections will be recorded in a logbook and maintained at the Site and available for inspection by NYSDEC. All necessary repairs shall be made immediately. Accumulated sediments will be removed as required to keep the barrier and hay bale check functional. All undercutting or erosion of the silt fence toe anchor shall be repaired immediately with appropriate backfill materials. Manufacturer's recommendations will be followed for

replacing silt fencing damaged due to weathering. Erosion and sediment control measures identified in the RAWP shall be observed to ensure that they are operating correctly. Where discharge locations or points are accessible, they shall be inspected to ascertain whether erosion control measures are effective in preventing significant impacts to receiving waters. Silt fencing or hay bales will be installed around the entire perimeter of the remedial construction area.

5.5.13 Contingency Plan

If underground tanks or other previously unidentified contaminant sources are found during on-site remedial excavation or development related construction, sampling will be performed on product, sediment and surrounding soils, etc. Chemical analytical work will be for full scan parameters (TAL metals; TCL volatiles and semi-volatiles, TCL pesticides and PCBs) if the use/contents of the tank cannot be verified as petroleum. These analyses will not be limited to STARS parameters where tanks are identified without prior approval by NYSDEC.

Identification of unknown or unexpected contaminated media identified by screening during invasive Site work will be promptly communicated by phone to NYSDEC's Project Manager. These findings will be also included in daily and periodic electronic media reports.

5.5.14 Community Air Monitoring Plan

The Community Air Monitoring Plan (CAMP) provides measures for protection for on-site workers and the downwind community (i.e., off-site receptors including residences, businesses, and on-site workers not directly involved in the remedial work) from potential airborne contaminant releases resulting from remedial activities at construction sites.

The action levels specified herein require increased monitoring, corrective actions to abate emissions, and/or work shutdown. Additionally, the CAMP helps to confirm that the remedial work did not spread contamination off-site through the air. The primary concerns for this site are odors associated with groundwater purging and sampling.

Exceedances observed in the CAMP will be reported to NYSDEC and NYSDOH Project Managers and included in the Daily Report. The complete CAMP developed for this site is included in **Attachment D** or this Work Plan.

5.5.15 Odor, Dust and Nuisance Control Plan

The Final Engineering Report will include the following certification by the Remedial Engineer: “I certify that all invasive work during the remediation and all invasive development work were conducted in accordance with dust and odor suppression methodology defined in the Remedial Action Work Plan.”

5.5.15.1 Odor Control Plan

This odor control plan is capable of controlling emissions of nuisance odors off-Site and on-Site. If nuisance odors are identified, work will be halted and the source of odors will be identified and corrected. Work will not resume until all nuisance odors have been abated. NYSDEC and NYSDOH will be notified of all odor events and of all other complaints about the project. Implementation of all odor controls, including the halt of work, will be the responsibility of the Applicant’s Remediation Engineer, who is responsible for certifying the Final Engineering Report.

All necessary means will be employed to prevent on and off-Site nuisances. At a minimum, procedures will include: (a) limiting the area of open excavations; (b) shrouding open excavations with tarps and other covers; and (c) using foams to cover exposed odorous soils. If odors develop and cannot be otherwise controlled, additional means to eliminate odor nuisances will include: (d) direct load-out of soils to trucks for off-Site disposal; (e) use of chemical odorants in spray or misting systems; and, (f) use of staff to monitor odors in surrounding neighborhoods.

Where odor nuisances have developed during remedial work and cannot be corrected, or where the release of nuisance odors cannot otherwise be avoided due to on-Site conditions or close proximity to sensitive receptors, odor control will be achieved by sheltering excavation and

handling areas under tented containment structures equipped with appropriate air venting/filtering systems.

5.5.15.2 Dust Control Plan

A dust suppression plan that addresses dust management during invasive on-Site work, will include, at a minimum, the items listed below:

- Dust suppression will be achieved though spraying water directly onto off-road areas including excavations and stockpiles.
- Gravel will be used on roadways to provide a clean and dust-free road surface.
- On-Site roads will be limited in total area to minimize the area required for water application.

5.5.15.3 Nuisance Control Plan

A plan for rodent control will be developed and utilized by the contractor prior to and during Site clearing and Site grubbing, and during all remedial work. A plan will be developed and utilized by the contractor for all remedial work and conforms, to NYCDEP noise control standards.

6.0 RESIDUAL CONTAMINATION TO REMAIN ON-SITE

Since soil with parameters above restricted residential SCOs will remain at depth after the remedy is complete, an Institutional Control (IC) is required to protect human health and the environment. The IC is described hereafter. Long-term management of the IC will be executed under a deed restriction recorded with the NYC Department of Finance, Office of the City Register.

ECs will be implemented to protect public health and the environment by appropriately managing residual contamination. The Controlled Property (the Site) will have the following EC systems:

1. Site Cover will be required to allow for residential use of the Site. The cover will consist of the new building foundation, concrete capped rear parking area, and a demarcation barrier and 2 feet of certified clean soil/top soil in landscaped areas.

The FER will report residual contamination on the Site in tabular and map form.

7.0 ENGINEERING CONTROLS

A site cover (**Figure 14**) will be required to allow for restricted residential use of the site. The cover will consist of the structures such as buildings, pavement, sidewalks comprising the site development and / or a soil cover in areas where the upper one foot (for commercial use) or two feet (for restricted residential use) of exposed surface soil will exceed the applicable soil cleanup objectives (SCOs). Where the soil cover is required it will be a minimum of one foot (for commercial use) or two feet (for restricted residential use) of soil, meeting the SCOs for cover material as set forth in 6 NYCRR Part 375-6.7(d). The soil cover will be placed over a demarcation layer, with the upper six inches of the soil of sufficient quality to maintain a vegetation layer. Any fill material brought to the site will meet the requirements for the identified site use as set forth in 6 NYCRR Part 375-6.7(d).

A Soil Management Plan will be included in the Site Management Plan and will outline the procedures to be followed in the event that the soil cover system and underlying residual contamination are disturbed after the Remedial Action is complete. Maintenance of the soil cover system will be described in the Site Management Plan in the FER.

8.0 INSTITUTIONAL CONTROLS

Institutional Controls (ICs) will be incorporated into the remedy to render the overall Site remedy protective of public health and the environmental. Two elements have been designed to ensure continual and proper management of residual contamination in perpetuity: an Environmental Easement and a Site Management Plan (SMP).

All as-build drawings, diagram, calculation and manufacturer documentation for treatment will be presented in the FER. A Site-Specific Environmental Easement will be recorded with the City of New York to provide an enforceable means of ensuring the continual and proper management of residual contamination and protection of public health and the environment in perpetuity or until released in writing by NYSDEC. It requires that the grantor of the Environmental Easement and the grantor's successors and assigns adhere to all Engineering and Institutional Controls (ECs/ICs) placed on the Site by this NYSDEC-approved remedy. ICs provide restrictions on Site usage and mandate operation, maintenance, monitoring and reporting measures for all ECs and ICs.

The SMP describes appropriate methods and procedures to ensure compliance with all ECs and ICs that are required by the Environmental Easement. Once the SMP has been approved by the NYSDEC, compliance with the SMP is required by the grantor of the Environmental Easement and grantor's successors and assigns.

8.1 ENVIRONMENTAL EASEMENT

An Environmental Easement, as defined in Article 71 Title 36 of the Environmental Conservation Law, is required when residual contamination is left on-Site after the Remedial Action is complete. If the Site will have residual contamination after completion of all Remedial Actions than an Environmental Easement is required. If an Environmental Easement is needed following completion of the remedy an Environmental Easement approved by NYSDEC will be filed and recorded with the City of New York. The Environmental Easement (if needed) will be submitted as part of the Final Remediation Report.

The Environmental Easement renders the Site a Controlled Property. The Environmental Easement must be recorded with the City of New York before the Certificate of Completion can be issued by NYSDEC. These Institutional Controls are requirements or restrictions placed on the Site that are listed in, and required by, the Environmental Easement. Institutional Controls can, generally, be subdivided between controls that support Engineering Controls, and those that place general restrictions on Site usage or other requirements. Institutional Controls in both of these groups are closely integrated with the Site Management Plan (SMP), which provides all of the methods and procedures to be followed to comply with this remedy.

The Institutional Controls which will be needed to support Engineering Controls are:

- Use of groundwater underlying the Controlled Property is prohibited without treatment rendering it safe for intended purpose;
- Compliance with the Environmental Easement by the Grantee and the Grantee's successor's is required;
- Grantor agrees to submit to NYSDEC a written statement that certifies, under penalty of perjury, that: (1) controls employed at the Controlled Property are unchanged from the previous certification or that any changes to the controls were approved by the NYSDEC; and, (2) nothing has occurred that impairs the ability of the controls to protect public health and environment or that constitute a violation or failure to comply with the Controls;
- NYSDEC retains the right to access such Controlled Property at any time in order to evaluate the continued maintenance of any and all controls. This certification shall be submitted annually, or an alternate period of time that NYSDEC may allow. This annual statement must be certified by an expert that the NYSDEC finds acceptable;

8.2 SITE MANAGEMENT PLAN

Site Management is the last phase of remediation and begins with the approval of the Final Engineering Report and issuance of the Certificate of Completion (COC) for the Remedial

Action. The Site Management Plan is submitted as a separate and independent document from the FER. Site Management continues in perpetuity or until released in writing by NYSDEC. The property owner is responsible to ensure that all Site Management responsibilities defined in the Environmental Easement and the Site Management Plan are performed.

The SMP is intended to provide a detailed description of the procedures required to manage residual contamination left in place at the Site following completion of the Remedial Action in accordance with the BCA with the NYSDEC. This includes: (1) development, implementation, and management of all Engineering and Institutional Controls; (2) development and implementation of monitoring systems and a Monitoring Plan; (3) development of a plan to operate and maintain any treatment, collection, containment, or recovery systems (including, where appropriate, preparation of an Operation and Maintenance Manual); (4) submittal of Site Management Reports, performance of inspections and certification of results, and demonstration of proper communication of Site information to NYSDEC; and (5) defining criteria for termination of treatment system operation.

To address these needs, this SMP will include four plans: (1) an Engineering and Institutional Control Plan for implementation and management of EC/ICs; (2) a Monitoring Plan for implementation of Site Monitoring; (3) an Operation and Maintenance Plan for implementation of remedial collection, containment, treatment, and recovery systems; and (4) a Site Management Reporting Plan for submittal of data, information, recommendations, and certifications to NYSDEC. The SMP will be prepared in accordance with the requirements in NYSDEC Draft DER-10 Technical Guidance for Site Investigation and Remediation, dated [month, year], and the guidelines provided by NYSDEC.

Site management activities, reporting, and EC/IC certification will be scheduled on a certification period basis. The certification period will be annually. The Site Management Plan will be based on a calendar year and will be due for submission to NYSDEC by March 1 of the year following the reporting period.

No exclusions for handling of residual contaminated soils will be provided in the Site Management Plan (SMP). All handling of residual contaminated material will be subject to provisions contained in the SMP.

9.0 FINAL ENGINEERING REPORT

A Final Engineering Report (FER) and Certificate of Completion (COC) will be submitted to NYSDEC following implementation of the Remedial Action defined in this RAWP. The FER provides the documentation that the remedial work required under this RAWP has been completed and has been performed in compliance with this plan. The FER will provide a comprehensive account of the locations and characteristics of all material removed from the Site including the surveyed map(s) of all sources. The Final Engineering Report will include as-built drawings for all constructed elements, certifications, manifests, bills of lading as well as the complete Site Management Plan (formerly the Operation and Maintenance Plan). The FER will provide a description of the changes in the Remedial Action from the elements provided in the RAWP and associated design documents. The FER will provide a tabular summary of all performance evaluation sampling results and all material characterization results and other sampling and chemical analysis performed as part of the Remedial Action. The FER will provide test results demonstrating that all mitigation and remedial systems are functioning properly. The FER will be prepared in conformance with DER-10.

Where determined to be necessary by NYSDEC, a Financial Assurance Plan will be required to ensure the sufficiency of revenue to perform long-term operations, maintenance and monitoring tasks defined in the Site Management Plan and Environmental Easement. This determination will be made by NYSDEC in the context of the Final Engineering Report review.

The Final Engineering Report will include written and photographic documentation of all remedial work performed under this remedy. The FER will include an itemized tabular description of actual costs incurred during all aspects of the Remedial Action.

The FER will provide a thorough summary of all residual contamination left on the Site after the remedy is complete. Residual contamination includes all contamination that exceeds the Track 1 Unrestricted Use SCO in 6NYCRR Part 375-6. A table that shows exceedances from Track 1 Unrestricted SCOs for all soil/fill remaining at the Site after the Remedial Action and a map that

shows the location and summarizes exceedances from Track 1 Unrestricted SCOs for all soil/fill remaining at the Site after the Remedial Action will be included in the FER.

The FER will provide a thorough summary of all residual contamination that exceeds the SCOs defined for the Site in the RAWP and must provide an explanation for why the material was not removed as part of the Remedial Action. A table that shows residual contamination in excess of Site SCOs and a map that shows residual contamination in excess of Site SCOs will be included in the FER.

The Final Engineering Report will include an accounting of the destination of all material removed from the Site, including excavated contaminated soil, historic fill, solid waste, hazardous waste, non-regulated material, and fluids. Documentation associated with disposal of all material must also include records and approvals for receipt of the material. It will provide an accounting of the origin and chemical quality of all material imported onto the Site.

Before approval of a FER and issuance of a Certificate of Completion, all project reports must be submitted in digital form on electronic media (PDF).

9.1 CERTIFICATIONS

The following certification will appear in front of the Executive Summary of the Final Engineering Report. The certification will be signed by the Remedial Engineer who is a Professional Engineer registered in New York State. This certification will be appropriately signed and stamped. The certification will include the following statements:

I _____ certify that I am currently a NYS registered professional engineer, I had primary direct responsibility for the implementation of the subject construction program, and I certify that the Remedial Work Plan (or Remedial Design or Plans and Specifications) was implemented and that all construction activities were completed in substantial conformance with the DER-approved Remedial Work Plan (or Remedial Design or Plans and Specifications).

If the Remedial Action Work Plan (or Remedial Design or Plans and Specifications) identifies time frames to be achieved by the remedial program, the certification must include:

The data submitted to DER demonstrates that the remediation requirements set forth in the Remedial Work Plan (or Remedial Design or Plans and Specifications) and all applicable statutes and regulations have been or will be achieved in accordance with the time frames, if any, established in the work plan (or Remedial Design or Plans and Specifications).

If the remedial program requires ICs or ECs, the certification will include:

All use restrictions, institutional controls, engineering controls and/or any operation and maintenance requirements applicable to the site are contained in an environmental easement created and recorded pursuant to ECL 71-3605 and that any affected local governments, as defined in ECL 71-3603, have been notified that such easement has been recorded.

If the remedial program requires applicable SMP, the certification will include:

A Site Management Plan has been submitted for the continual and proper operation, maintenance, and monitoring of any engineering controls employed at the site including the proper maintenance of any remaining monitoring wells, and that such plan has been approved by DER.

If the remedial program requires financial assurance, the certification will include:

Any financial assurance mechanisms required by DEC pursuant to Environmental Conservation Law have been executed.

10.0 SCHEDULE

The remedial action will begin with mobilization of equipment and material to the Site which will begin approximately 4 weeks following RAWP approval and within 10 days of the distribution of the Construction Fact Sheet. Mobilization will be followed by removal and disposal of the USTs (if present) followed by excavation and disposal of petroleum impacted soil, historic fill materials and native soil, followed by confirmation sampling. The work is expected to take approximately 6 months as part of the construction excavation and foundation installation. The schedule of tasks completed under this RAWP is as follows:

| | |
|---|--|
| Conduct pre-construction meeting with NYSDEC | Within 2 weeks of RAWP approval (Oct 2016) |
| Mobilize equipment to the site and construct truck pad and other designated areas | Within 2 weeks following the pre-construction meeting and issuance of Pre-Construction Fact Sheet (Oct 2016) |
| Begin excavation of USTs (if present) | Immediately following mobilization (Oct 2016) |
| Mobilize Excavation Contractor and equipment to the Site | Immediately following excavation of UST Week (Oct 2016) |
| Complete excavation and disposal of historic fill material and clean native soil. | Within 6 months of mobilization (May 2017) |
| Perform endpoint verification of UST areas | Performed in sequence as final depth of each excavated area is complete (Oct. 2016-May2017) |
| Submit SMP (as a contingency) if Track 1 Cleanup is not achieved | By August 15 th of the year in which the COC is sought (Dec 2017) or as required by DEC. |
| Submit FER | By September 15 th of the year in which the COC is sought (Dec 2017) or as required by DEC. |

TABLES

TABLE 1
Soil Cleanup Objectives
Soil Import Criteria

| Contaminant | CAS Number | Top 2 ft in Exposed Soil Areas |
|-----------------------------------|------------|--------------------------------------|
| | | Restricted- Residential |
| METALS | | |
| Arsenic | 7440-38 -2 | 16f |
| Barium | 7440-39 -3 | 400 |
| Beryllium | 7440-41 -7 | 72 |
| Cadmium | 7440-43 -9 | 4.3 |
| Chromium, hexavalent ^h | 18540-29-9 | 110 |
| Chromium, trivalent ^h | 16065-83-1 | 180 |
| Copper | 7440-50 -8 | 270 |
| Total Cyanide ^h | | 27 |
| Lead | 7439-92 -1 | 400 |
| Manganese | 7439-96 -5 | 2,000f |
| Total Mercury | | 0.81j |
| Nickel | 7440-02 -0 | 310 |
| Selenium | 7782-49 -2 | 180 |
| Silver | 7440-22 -4 | 180 |
| Zinc | 7440-66 -6 | 10,000 d |
| PESTICIDES / PCBs | | |
| 2,4,5-TP Acid (Silvex) | 93-72-1 | 100a |
| 4,4'-DDE | 72-55-9 | 8.9 |
| 4,4'-DDT | 50-29-3 | 7.9 |
| 4,4'-DDD | 72-54-8 | 13 |
| Aldrin | 309-00-2 | 0.097 |
| alpha-BHC | 319-84-6 | 0.48 |
| beta-BHC | 319-85-7 | 0.36 |
| Chlordane (alpha) | 5103-71 -9 | 4.2 |
| delta-BHC | 319-86-8 | 100a |
| Dibenzofuran | 132-64-9 | 59 |
| Dieldrin | 60-57-1 | 0.2 |
| Endosulfan I | 959-98-8 | 24i |
| Endosulfan II | 33213-65-9 | 24i |
| Endosulfan sulfate | 1031-07 -8 | 24i |
| Endrin | 72-20-8 | 11 |
| Heptachlor | 76-44-8 | 2.1 |
| Lindane | 58-89-9 | 1.3 |
| Polychlorinated biphenyls | 1336-36 -3 | 1 |

| Contaminant | CAS Number | Top 2 ft in Exposed Soil Areas |
|--------------------------|------------|--------------------------------------|
| | | Restricted- Residential |
| VOLATILES | | |
| 1,1,1-Trichloroethane | 71-55-6 | 100a |
| 1,1-Dichloroethane | 75-34-3 | 26 |
| 1,1-Dichloroethene | 75-35-4 | 100a |
| 1,2-Dichlorobenzene | 95-50-1 | 100a |
| 1,2-Dichloroethane | 107-06-2 | 3.1 |
| cis-1,2-Dichloroethene | 156-59-2 | 100a |
| trans-1,2-Dichloroethene | 156-60-5 | 100a |
| 1,3-Dichlorobenzene | 541-73-1 | 49 |
| 1,4-Dichlorobenzene | 106-46-7 | 13 |
| 1,4-Dioxane | 123-91-1 | 13 |
| Acetone | 67-64-1 | 100b |
| Benzene | 71-43-2 | 4.8 |
| Butylbenzene | 104-51-8 | 100a |
| Carbon tetrachloride | 56-23-5 | 2.4 |
| Chlorobenzene | 108-90-7 | 100a |
| Chloroform | 67-66-3 | 49 |
| Ethylbenzene | 100-41-4 | 41 |
| Hexachlorobenzene | 118-74-1 | 1.2 |
| Methyl ethyl ketone | 78-93-3 | 100a |
| Methyl tert-butyl ether | 1634-04 -4 | 100a |
| Methylene chloride | 75-09-2 | 100a |
| n-Propylbenzene | 103-65-1 | 100a |
| sec-Butylbenzene | 135-98-8 | 100a |
| tert-Butylbenzene | 98-06-6 | 100a |
| Tetrachloroethene | 127-18-4 | 19 |
| Toluene | 108-88-3 | 100a |
| Trichloroethene | 79-01-6 | 21 |
| 1,2,4-Trimethylbenzene | 95-63-6 | 52 |
| 1,3,5-Trimethylbenzene | 108-67-8 | 52 |
| Vinyl chloride | 75-01-4 | 0.9 |
| Xylene (mixed) | 1330-20 -7 | 100a |

| Contaminant | CAS Number | Top 2 ft in Exposed Soil Areas |
|------------------------|------------|--------------------------------------|
| | | Restricted- Residential |
| SEMI-VOLATILES | | |
| Acenaphthene | 83-32-9 | 100a |
| Acenaphthylene | 208-96-8 | 100a |
| Anthracene | 120-12-7 | 100a |
| Benzo(a)anthracene | 56-55-3 | 1f |
| Benzo(a)pyrene | 50-32-8 | 1f |
| Benzo(b)fluoranthene | 205-99-2 | 1f |
| Benzo(g,h,i) perylene | 191-24-2 | 100a |
| Benzo(k)fluoranthene | 207-08-9 | 3.9 |
| Chrysene | 218-01-9 | 3.9 |
| Dibenz(a,h)anthracene | 53-70-3 | 0.33e |
| Fluoranthene | 206-44-0 | 100a |
| Fluorene | 86-73-7 | 100a |
| Indeno(1,2,3-cd)pyrene | 193-39-5 | 0.5f |
| m-Cresol | 108-39-4 | 100a |
| Naphthalene | 91-20-3 | 100a |
| o-Cresol | 95-48-7 | 100a |
| p-Cresol | 106-44-5 | 100a |
| Pentachlorophenol | 87-86-5 | 6.7 |
| Phenanthrene | 85-01-8 | 100a |
| Phenol | 108-95-2 | 100a |
| Pyrene | 129-00-0 | 100a |

All soil cleanup objectives (SCOs) are in parts per million (ppm).

TABLE 2
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Remedial Investigation Sampling Summary

| Matrix | Location | Number of Samples | Rationale for Sampling | Laboratory Analysis |
|---|---|-------------------|---|---|
| Total (Soils) | | 34 | | |
| Subsurface soil (0 to 5 feet) | 11 soil borings throughout the Site. Samples collected at 0-2 ft, 2-4 ft, and 3-5 ft intervals. | 11 | To evaluate the extent of soil impact and obtain information on soil quality at the Site. | VOCs EPA Method 8260B, SVOCs EPA Method 8270, Pesticides / PCBs EPA Method 8081/8082, TAL Metals |
| Subsurface soil (5 to 10 feet) | 10 soil borings throughout the Site. Samples collected at 5-7 ft and 6-8 ft intervals. | 10 | To evaluate the extent of soil impact and obtain information on soil quality at the Site. | VOCs EPA Method 8260B, SVOCs EPA Method 8270, Pesticides / PCBs EPA Method 8081/8082, TAL Metals |
| Subsurface soil (5 to 10 feet) | 10 soil borings throughout the Site. Samples collected at 5-7 ft, 6-8 ft and 8-10 ft intervals. | 10 | To evaluate the extent of soil impact and obtain information on soil quality at the Site. | VOCs EPA Method 8260B, SVOCs EPA Method 8270 |
| Subsurface soil (10 to 15 feet) | 1 soil boring throughout the site. Samples collected at the 10-12 ft interval. | 1 | To evaluate the extent of soil impact and obtain information on soil quality at the Site. | VOCs EPA Method 8260B, SVOCs EPA Method 8270, Pesticides / PCBs EPA Method 8081/8082, TAL Metals |
| Subsurface soil (10 to 15 feet) | 2 soil boring throughout the site. Samples collected at 10-12 ft and 13-15 ft intervals. | 2 | To evaluate the extent of soil impact and obtain information on soil quality at the Site. | VOCs EPA Method 8260B and SVOCs EPA Method 8270 |
| Total (Groundwater) | | 12 | | |
| Groundwater (below water table) | From 12 monitoring wells across the Site. | 12 | To assess groundwater quality at the Site. | VOCs EPA Method 8260B, SVOCs EPA Method 8270, Pesticide / PCBs EPA Method 8081/8082, Dissolved Metals |
| Total (Soil Gas) | | 13 | | |
| Soil Gas (3 ft below grade) | 8 soil gas implant locations installed 3 ft beneath the first floor level across the Site. | 8 | Evaluate soil gas across the Site. | VOCs EPA Method TO15 |
| Sub-Slab Soil Gas (6 in. below basement slab) | 5 soil gas implant locations installed beneath the basement level slab-on-grade foundation throughout the Site. | 5 | Evaluate soil gas across the Site. | VOCs EPA Method TO15 |
| Total (QA / QC Samples) | | 4 | | |
| Trip Blanks (Groundwater) | Two laboratory prepared trip blanks to accompany samples each time they are delivered to the laboratory. | 2 | To meet requirements of QA / QC program. | VOCs EPA Method 8260B |
| Duplicates (Soil) | Created a duplicate of another soil sample in the field for the purpose of comparison and quality control. | 2 | To meet requirements of QA / QC program. | VOCs EPA Method 8260B, SVOCs EPA Method 8270, Pesticides / PCBs EPA Method 8081/8082, TAL Metals |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B1 | | | | | | | | 15B2 | | | |
|----------------------------------|---|---|-----------------|-------------|------|------|-----------------|-----------|------|------|-----------------|-------------|------|------|
| | | | 11/24/2015 | | | | | | | | 11/24/2015 | | | |
| | | | (0-2') µg/Kg | | | | (5-7') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 1100 | 1,100 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 21 | 21 | U | 1.0 |
| 1,1,2,2-Tetrachloroethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,1,2-Trichloroethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,1-Dichloroethane | 270 | 26,000 | < 140 | 140 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,1-Dichloroethene | 330 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,1-Dichloropropene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2,3-Trichlorobenzene | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,2,3-Trichloropropane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2,4-Trichlorobenzene | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | 100 | 270 | J | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2-Dibromo-3-chloropropane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,2-Dibromomethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2-Dichloroethane | 20 | 3,100 | < 27 | 27 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,2-Dichloropropane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | 62 | 270 | J | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,3-Dichloropropane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 1,4-dioxane | | | < 2200 | 2,200 | U | 2200 | < 100 | 100 | U | 42 | < 100 | 100 | U | 41 |
| 2,2-Dichloropropane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 2-Chlorotoluene | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 1400 | 1,400 | U | 270 | < 27 | 27 | U | 5.3 | < 26 | 26 | U | 5.1 |
| 2-Isopropyltoluene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 4-Chlorotoluene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| 4-Methyl-2-Pentanone | | | < 1400 | 1,400 | U | 270 | < 27 | 27 | U | 5.3 | < 26 | 26 | U | 5.1 |
| Acetone | 50 | 100,000 | < 270 | 270 | U | 270 | 10 | 50 | JS | 5.3 | < 50 | 50 | U | 5.1 |
| Acrolein | | | < 1100 | 1,100 | U | 140 | < 21 | 21 | U | 2.7 | < 21 | 21 | U | 2.6 |
| Acrylonitrile | | | < 1100 | 1,100 | U | 27 | < 11 | 11 | U | 1.1 | < 10 | 10 | U | 1.0 |
| Benzene | 60 | 4,800 | < 55 | 55 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Bromobenzene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Bromochloromethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Bromodichloromethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Bromoform | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Bromomethane | | | < 270 | 270 | U | 110 | < 5.3 | 5.3 | U | 2.1 | < 5.1 | 5.1 | U | 2.1 |
| Carbon Disulfide | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | 2.8 | 5.1 | J | 1.0 |
| Carbon tetrachloride | 760 | 2,400 | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Chlorobenzene | 1,100 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Chloroethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Chloroform | 370 | 49,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Chloromethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 140 | 140 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| cis-1,3-Dichloropropene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Dibromochloromethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Dibromomethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Dichlorodifluoromethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Ethylbenzene | 1,000 | 41,000 | 60 | 270 | J | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Hexachlorobutadiene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Isopropylbenzene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| m&p-Xylenes | 260 | 100,000 | 250 | 270 | J | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 270 | 270 | U | 270 | < 32 | 32 | U | 5.3 | < 31 | 31 | U | 5.1 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 550 | 550 | U | 55 | < 11 | 11 | U | 1.1 | < 10 | 10 | U | 1.0 |
| Methylene chloride | 50 | 100,000 | < 270 | 270 | U | 270 | < 5.3 | 5.3 | U | 5.3 | < 5.1 | 5.1 | U | 5.1 |
| Naphthalene | 12,000 | 100,000 | 2,900 | 270 | | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| n-Butylbenzene | 12,000 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| n-Propylbenzene | 3,900 | 100,000 | < 270 | 270 | U | 49 | < 5.3 | 5.3 | U | 0.95 | < 5.1 | 5.1 | U | 0.93 |
| o-Xylene | 260 | 100,000 | 55 | 270 | J | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| p-Isopropyltoluene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| sec-Butylbenzene | 11,000 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Styrene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Tert-butyl alcohol | | | < 5500 | 5,500 | U | 1100 | < 110 | 110 | U | 21 | 35 | 100 | J | 21 |
| tert-Butylbenzene | 5,900 | 100,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Tetrachloroethene | 1,300 | 19,000 | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Tetrahydrofuran (THF) | | | < 550 | 550 | U | 140 | < 11 | 11 | U | 2.7 | < 10 | 10 | U | 2.6 |
| Toluene | 700 | 100,000 | 400 | 270 | | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 140 | 140 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| trans-1,3-Dichloropropene | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| trans-1,4-dichloro-2-butene | | | < 550 | 550 | U | 140 | < 11 | 11 | U | 2.7 | < 10 | 10 | U | 2.6 |
| Trichloroethene | 470 | 21,000 | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Trichlorofluoromethane | | | < 270 | 270 | U | 55 | < 5.3 | 5.3 | U | 1.1 | < 5.1 | 5.1 | U | 1.0 |
| Trichlorotrifluoroethane | | | < 270 | 270 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Vinyl Chloride | 20 | 900 | < 27 | 27 | U | 27 | < 5.3 | 5.3 | U | 0.53 | < 5.1 | 5.1 | U | 0.51 |
| Total BTEX Concentration | | | | 765 | | | | 0 | | | | 0 | | |
| Total VOCs Concentration | | | | 3827 | | | | 10 | | | | 37.8 | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives
RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B3 | | | | | | | | | | | |
|----------------------------------|---|---|------------|-------|------|------------|------------|------------|------|------|------------|-----|------------|------|
| | | | 11/24/2015 | | | | | | | | | | | |
| | | | (0-2') | | | | (5-7') | | | | (13-15') | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 720 | 720 | U | 36 | < 920 | 920 | U | 46 | < 14 | 14 | U | 0.70 |
| 1,1,2,2-Tetrachloroethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,1,2-Trichloroethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,1-Dichloroethane | 270 | 26,000 | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,1-Dichloroethene | 330 | 100,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,1-Dichloropropene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,2,3-Trichlorobenzene | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,2,3-Trichloropropane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,2,4-Trichlorobenzene | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | 63 | 180 | J | 18 | 110 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,2-Dibromo-3-chloropropane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,2-Dibromomethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,2-Dichloroethane | 20 | 3,100 | < 18 | 18 | U | 18 | < 20 | 20 | U | 20 | < 3.5 | 3.5 | U | 0.35 |
| 1,2-Dichloropropane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | 50 | 180 | J | 18 | 42 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,3-Dichloropropane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 1,4-dioxane | | | < 1400 | 1,400 | U | 1400 | < 1800 | 1,800 | U | 1800 | < 70 | 70 | U | 28 |
| 2,2-Dichloropropane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 2-Chlorotoluene | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 900 | 900 | U | 180 | < 1200 | 1,200 | U | 230 | < 18 | 18 | U | 3.5 |
| 2-Isopropyltoluene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 4-Chlorotoluene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| 4-Methyl-2-Pentanone | | | < 900 | 900 | U | 180 | < 1200 | 1,200 | U | 230 | < 18 | 18 | U | 3.5 |
| Acetone | 50 | 100,000 | < 180 | 180 | U | 180 | < 230 | 230 | U | 230 | 3.8 | 35 | JS | 3.5 |
| Acrolein | | | < 720 | 720 | U | 90 | < 920 | 920 | U | 120 | < 14 | 14 | U | 1.8 |
| Acrylonitrile | | | < 720 | 720 | U | 18 | < 920 | 920 | U | 23 | < 7.0 | 7.0 | U | 0.70 |
| Benzene | 60 | 4,800 | 23 | 58 | J | 18 | 65 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| Bromobenzene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Bromochloromethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Bromodichloromethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Bromoform | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Bromomethane | | | < 180 | 180 | U | 72 | < 230 | 230 | U | 92 | < 3.5 | 3.5 | U | 1.4 |
| Carbon Disulfide | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Carbon tetrachloride | 760 | 2,400 | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Chlorobenzene | 1,100 | 100,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Chloroethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Chloroform | 370 | 49,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Chloromethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| cis-1,3-Dichloropropene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Dibromochloromethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Dibromomethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Dichlorodifluoromethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Ethylbenzene | 1,000 | 41,000 | 39 | 180 | J | 18 | 65 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| Hexachlorobutadiene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Isopropylbenzene | | | < 180 | 180 | U | 18 | 72 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| m&p-Xylenes | 260 | 100,000 | 60 | 180 | J | 36 | 140 | 230 | J | 46 | < 3.5 | 3.5 | U | 0.70 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 180 | 180 | U | 180 | < 230 | 230 | U | 230 | < 21 | 21 | U | 3.5 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 360 | 360 | U | 36 | < 460 | 460 | U | 46 | < 7.0 | 7.0 | U | 0.70 |
| Methylene chloride | 50 | 100,000 | < 180 | 180 | U | 180 | < 230 | 230 | U | 230 | < 3.5 | 3.5 | U | 3.5 |
| Naphthalene | 12,000 | 100,000 | 56 | 180 | J | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| n-Butylbenzene | 12,000 | 100,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| n-Propylbenzene | 3,900 | 100,000 | < 180 | 180 | U | 36 | 68 | 230 | J | 42 | < 3.5 | 3.5 | U | 0.63 |
| o-Xylene | 260 | 100,000 | < 180 | 180 | U | 32 | 58 | 230 | J | 46 | < 3.5 | 3.5 | U | 0.70 |
| p-Isopropyltoluene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| sec-Butylbenzene | 11,000 | 100,000 | < 180 | 180 | U | 18 | 75 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| Styrene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Tert-butyl alcohol | | | < 3600 | 3,600 | U | 720 | < 4600 | 4,600 | U | 920 | < 70 | 70 | U | 14 |
| tert-Butylbenzene | 5,900 | 100,000 | < 180 | 180 | U | 18 | 41 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| Tetrachloroethene | 1,300 | 19,000 | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Tetrahydrofuran (THF) | | | < 360 | 360 | U | 90 | < 460 | 460 | U | 120 | < 7.0 | 7.0 | U | 1.8 |
| Toluene | 700 | 100,000 | 54 | 180 | J | 18 | 99 | 230 | J | 23 | < 3.5 | 3.5 | U | 0.35 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 180 | 180 | U | 18 | < 120 | 120 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| trans-1,3-Dichloropropene | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| trans-1,4-dichloro-2-butene | | | < 360 | 360 | U | 90 | < 460 | 460 | U | 120 | < 7.0 | 7.0 | U | 1.8 |
| Trichloroethene | 470 | 21,000 | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Trichlorofluoromethane | | | < 180 | 180 | U | 36 | < 230 | 230 | U | 46 | < 3.5 | 3.5 | U | 0.70 |
| Trichlorotrifluoroethane | | | < 180 | 180 | U | 18 | < 230 | 230 | U | 23 | < 3.5 | 3.5 | U | 0.35 |
| Vinyl Chloride | 20 | 900 | < 18 | 18 | U | 18 | < 20 | 20 | U | 20 | < 3.5 | 3.5 | U | 0.35 |
| Total BTEX Concentration | | | | | | 176 | | 427 | | | | | 0 | |
| Total VOCs Concentration | | | | | | 345 | | 835 | | | | | 3.8 | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives
 RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank The value is estimated. This flag is used |
| J | a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B4 | | | | | | | | | | | | 15B5 | | | |
|----------------------------------|---|---|------------|-----------|-------|------|------------|------------|-------|------|----------|--------|-------|----------|------------|-----|-------|------|
| | | | 11/24/2015 | | | | | | | | | | | | | | | |
| | | | (0-2') | | | | (6-8') | | | | (10-12') | | | | (6-8') | | | |
| | | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,1,1,2-Tetrachloroethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,1,2,2-Tetrachloroethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,1,2-Trichloroethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,1-Dichloroethane | 270 | 26,000 | < 220 | 220 | U | 44 | < 170 | 170 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,1-Dichloroethene | 330 | 100,000 | < 220 | 220 | U | 22 | < 170 | 170 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,1-Dichloropropene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,2,3-Trichlorobenzene | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,2,3-Trichloropropane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,2,4-Trichlorobenzene | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 220 | 220 | U | 22 | 91 | 340 | J | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,2-Dibromo-3-chloropropane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,2-Dibromomethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,2-Dichloroethane | 20 | 3,100 | < 20 | 20 | U | 20 | < 34 | 34 | U | 34 | < 20 | 20 | U | 20 | < 5.9 | 5.9 | U | 0.59 |
| 1,2-Dichloropropane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,3-Dichloropropane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 1,4-dioxane | | | < 1800 | 1,800 | U | 1800 | < 2700 | 2,700 | U | 2700 | < 4500 | 4,500 | U | 1800 | < 94 | 94 | U | 47 |
| 2,2-Dichloropropane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 2-Chlorotoluene | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 1100 | 1,100 | U | 220 | < 1700 | 1,700 | U | 340 | < 1100 | 1,100 | U | 220 | < 30 | 30 | U | 5.9 |
| 2-Isopropyltoluene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 4-Chlorotoluene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| 4-Methyl-2-Pentanone | | | < 1100 | 1,100 | U | 220 | < 1700 | 1,700 | U | 340 | < 1100 | 1,100 | U | 220 | < 30 | 30 | U | 5.9 |
| Acetone | 50 | 100,000 | < 220 | 220 | U | 220 | < 340 | 340 | U | 340 | < 220 | 220 | U | 220 | 29 | 47 | JS | 5.9 |
| Acrolein | | | < 880 | 880 | U | 110 | < 1300 | 1,300 | U | 170 | < 890 | 890 | U | 110 | < 24 | 24 | U | 3.0 |
| Acrylonitrile | | | < 880 | 880 | U | 22 | < 670 | 670 | U | 67 | < 450 | 450 | U | 45 | < 12 | 12 | U | 1.2 |
| Benzene | 60 | 4,800 | 43 | 57 | J | 22 | 140 | 340 | J | 34 | < 58 | 58 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Bromobenzene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Bromochloromethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Bromodichloromethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Bromoform | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Bromomethane | | | < 220 | 220 | U | 88 | < 340 | 340 | U | 130 | < 220 | 220 | U | 89 | < 5.9 | 5.9 | U | 2.4 |
| Carbon Disulfide | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Carbon tetrachloride | 760 | 2,400 | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Chlorobenzene | 1,100 | 100,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Chloroethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Chloroform | 370 | 49,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Chloromethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 220 | 220 | U | 22 | < 170 | 170 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| cis-1,3-Dichloropropene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Dibromochloromethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Dibromomethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Dichlorodifluoromethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Ethylbenzene | 1,000 | 41,000 | < 220 | 220 | U | 22 | 61 | 340 | J | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Hexachlorobutadiene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Isopropylbenzene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| m&p-Xylenes | 260 | 100,000 | < 220 | 220 | U | 44 | 140 | 340 | J | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 220 | 220 | U | 220 | < 340 | 340 | U | 340 | < 220 | 220 | U | 220 | < 35 | 35 | U | 5.9 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 440 | 440 | U | 44 | < 670 | 670 | U | 67 | < 450 | 450 | U | 45 | 2.1 | 12 | J | 1.2 |
| Methylene chloride | 50 | 100,000 | < 220 | 220 | U | 220 | < 340 | 340 | U | 340 | < 220 | 220 | U | 220 | < 5.9 | 5.9 | U | 5.9 |
| Naphthalene | 12,000 | 100,000 | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| n-Butylbenzene | 12,000 | 100,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| n-Propylbenzene | 3,900 | 100,000 | < 220 | 220 | U | 40 | < 340 | 340 | U | 61 | < 220 | 220 | U | 40 | < 5.9 | 5.9 | U | 1.1 |
| o-Xylene | 260 | 100,000 | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| p-Isopropyltoluene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| sec-Butylbenzene | 11,000 | 100,000 | < 220 | 220 | U | 22 | 44 | 340 | J | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Styrene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Tert-butyl alcohol | | | < 4400 | 4,400 | U | 880 | < 6700 | 6,700 | U | 1300 | < 4500 | 4,500 | U | 890 | < 120 | 120 | U | 24 |
| tert-Butylbenzene | 5,900 | 100,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Tetrachloroethene | 1,300 | 19,000 | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Tetrahydrofuran (THF) | | | < 440 | 440 | U | 110 | < 670 | 670 | U | 170 | < 450 | 450 | U | 110 | < 12 | 12 | U | 3.0 |
| Toluene | 700 | 100,000 | 55 | 220 | J | 22 | 87 | 340 | J | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 110 | 110 | U | 22 | < 170 | 170 | U | 34 | < 110 | 110 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| trans-1,3-Dichloropropene | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| trans-1,4-dichloro-2-butene | | | < 440 | 440 | U | 110 | < 670 | 670 | U | 170 | < 450 | 450 | U | 110 | < 12 | 12 | U | 3.0 |
| Trichloroethene | 470 | 21,000 | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Trichlorofluoromethane | | | < 220 | 220 | U | 44 | < 340 | 340 | U | 67 | < 220 | 220 | U | 45 | < 5.9 | 5.9 | U | 1.2 |
| Trichlorotrifluoroethane | | | < 220 | 220 | U | 22 | < 340 | 340 | U | 34 | < 220 | 220 | U | 22 | < 5.9 | 5.9 | U | 0.59 |
| Vinyl Chloride | 20 | 900 | < 20 | 20 | U | 20 | < 34 | 34 | U | 34 | < 20 | 20 | U | 20 | < 5.9 | 5.9 | U | 0.59 |
| Total BTEX Concentration | | | | 98 | | | | 428 | | | 0 | | | 0 | | | </ | |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B6 | | | | | | | | 15SB7 | | | | 15B8 | | | |
|----------------------------------|---|---|----------------|-----|------|------|------------------|-----|------|------|----------------|-----|------|------|----------------|-----|------|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (3-5) µg/Kg | | | | (10-12) µg/Kg | | | | (6-8) µg/Kg | | | | (6-8) µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,1,2,2-Tetrachloroethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,1,2-Trichloroethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,1-Dichloroethane | 270 | 26,000 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,1-Dichloroethene | 330 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,1-Dichloropropene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2,3-Trichlorobenzene | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,2,3-Trichloropropane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2,4-Trichlorobenzene | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2-Dibromo-3-chloropropane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,2-Dibromomethane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2-Dichloroethane | 20 | 3,100 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,2-Dichloropropane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,3-Dichloropropane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 1,4-dioxane | | | < 100 | 100 | U | 42 | < 100 | 100 | U | 71 | < 94 | 94 | U | 38 | < 100 | 100 | U | 61 |
| 2,2-Dichloropropane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 2-Chlorotoluene | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 26 | 26 | U | 5.3 | < 44 | 44 | U | 8.9 | < 23 | 23 | U | 4.7 | < 38 | 38 | U | 7.6 |
| 2-Isopropyltoluene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 4-Chlorotoluene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| 4-Methyl-2-Pentanone | | | < 26 | 26 | U | 5.3 | < 44 | 44 | U | 8.9 | < 23 | 23 | U | 4.7 | < 38 | 38 | U | 7.6 |
| Acetone | 50 | 100,000 | 22 | 50 | JS | 5.3 | 39 | 50 | JS | 8.9 | 11 | 47 | JS | 4.7 | 110 | 50 | S | 7.6 |
| Acrolein | | | < 21 | 21 | U | 2.6 | < 35 | 35 | U | 4.4 | < 19 | 19 | U | 2.3 | < 31 | 31 | U | 3.8 |
| Acrylonitrile | | | < 11 | 11 | U | 1.1 | < 18 | 18 | U | 1.8 | < 9.4 | 9.4 | U | 0.94 | < 15 | 15 | U | 1.5 |
| Benzene | 60 | 4,800 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | 0.91 | 7.6 | J | 0.76 |
| Bromobenzene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Bromochloromethane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Bromodichloromethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Bromoform | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Bromomethane | | | < 5.3 | 5.3 | U | 2.1 | < 8.9 | 8.9 | U | 3.5 | < 4.7 | 4.7 | U | 1.9 | < 7.6 | 7.6 | U | 3.1 |
| Carbon Disulfide | | | < 5.3 | 5.3 | U | 1.1 | 8.3 | 8.9 | J | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Carbon tetrachloride | 760 | 2,400 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Chlorobenzene | 1,100 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Chloroethane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Chloroform | 370 | 49,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Chloromethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| cis-1,3-Dichloropropene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Dibromochloromethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Dibromomethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Dichlorodifluoromethane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Ethylbenzene | 1,000 | 41,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Hexachlorobutadiene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Isopropylbenzene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| m&p-Xylenes | 260 | 100,000 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 32 | 32 | U | 5.3 | < 53 | 53 | U | 8.9 | < 28 | 28 | U | 4.7 | 33 | 46 | J | 7.6 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 11 | 11 | U | 1.1 | < 18 | 18 | U | 1.8 | < 9.4 | 9.4 | U | 0.94 | < 15 | 15 | U | 1.5 |
| Methylene chloride | 50 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Naphthalene | 12,000 | 100,000 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| n-Butylbenzene | 12,000 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| n-Propylbenzene | 3,900 | 100,000 | < 5.3 | 5.3 | U | 0.95 | < 8.9 | 8.9 | U | 1.6 | < 4.7 | 4.7 | U | 0.84 | < 7.6 | 7.6 | U | 1.4 |
| o-Xylene | 260 | 100,000 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| p-Isopropyltoluene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| sec-Butylbenzene | 11,000 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Styrene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Tert-butyl alcohol | | | < 110 | 110 | U | 21 | < 180 | 180 | U | 35 | < 94 | 94 | U | 19 | < 150 | 150 | U | 31 |
| tert-Butylbenzene | 5,900 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Tetrachloroethene | 1,300 | 19,000 | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Tetrahydrofuran (THF) | | | < 11 | 11 | U | 2.6 | < 18 | 18 | U | 4.4 | < 9.4 | 9.4 | U | 2.3 | < 15 | 15 | U | 3.8 |
| Toluene | 700 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| trans-1,3-Dichloropropene | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| trans-1,4-dichloro-2-butene | | | < 11 | 11 | U | 2.6 | < 18 | 18 | U | 4.4 | < 9.4 | 9.4 | U | 2.3 | < 15 | 15 | U | 3.8 |
| Trichloroethene | 470 | 21,000 | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Trichlorofluoromethane | | | < 5.3 | 5.3 | U | 1.1 | < 8.9 | 8.9 | U | 1.8 | < 4.7 | 4.7 | U | 0.94 | < 7.6 | 7.6 | U | 1.5 |
| Trichlorotrifluoroethane | | | < 5.3 | 5.3 | U | 0.53 | < 8.9 | 8.9 | U | 0.89 | < 4.7 | 4.7 | U | 0.47 | < 7.6 | 7.6 | U | 0.76 |
| Vinyl Chloride | 20 | 900 | < 5.3 | 5.3 | | | | | | | | | | | | | | |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B9 | | | | | | | | 15B10 | | | |
|----------------------------------|---|---|------------|-----|------|------|------------|-----|------|------|------------|-----|------|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | |
| | | | (0-2') | | | | (6-8') | | | | (6-8') | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,1,2,2-Tetrachloroethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,1,2-Trichloroethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,1-Dichloroethane | 270 | 26,000 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,1-Dichloroethene | 330 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,1-Dichloropropene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2,3-Trichlorobenzene | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,2,3-Trichloropropane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2,4-Trichlorobenzene | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2-Dibromo-3-chloropropane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,2-Dibromomethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2-Dichloroethane | 20 | 3,100 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,2-Dichloropropane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,3-Dichloropropane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 1,4-dioxane | | | < 85 | 85 | U | 34 | < 89 | 89 | U | 35 | < 94 | 94 | U | 38 |
| 2,2-Dichloropropane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 2-Chlorotoluene | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 21 | 21 | U | 4.2 | < 22 | 22 | U | 4.4 | < 24 | 24 | U | 4.7 |
| 2-Isopropyltoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 4-Chlorotoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| 4-Methyl-2-Pentanone | | | < 21 | 21 | U | 4.2 | < 22 | 22 | U | 4.4 | < 24 | 24 | U | 4.7 |
| Acetone | 50 | 100,000 | 5 | 42 | JS | 4.2 | 6.2 | 44 | JS | 4.4 | 17 | 47 | JS | 4.7 |
| Acrolein | | | < 17 | 17 | U | 2.1 | < 18 | 18 | U | 2.2 | < 19 | 19 | U | 2.4 |
| Acrylonitrile | | | < 8.5 | 8.5 | U | 0.85 | < 8.9 | 8.9 | U | 0.89 | < 9.4 | 9.4 | U | 0.94 |
| Benzene | 60 | 4,800 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Bromobenzene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Bromochloromethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Bromodichloromethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Bromoform | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Bromomethane | | | < 4.2 | 4.2 | U | 1.7 | < 4.4 | 4.4 | U | 1.8 | < 4.7 | 4.7 | U | 1.9 |
| Carbon Disulfide | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Carbon tetrachloride | 760 | 2,400 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Chlorobenzene | 1,100 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Chloroethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Chloroform | 370 | 49,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Chloromethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| cis-1,3-Dichloropropene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Dibromochloromethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Dibromomethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Dichlorodifluoromethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Ethylbenzene | 1,000 | 41,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Hexachlorobutadiene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Isopropylbenzene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| m&p-Xylenes | 260 | 100,000 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 25 | 25 | U | 4.2 | < 27 | 27 | U | 4.4 | < 28 | 28 | U | 4.7 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 8.5 | 8.5 | U | 0.85 | < 8.9 | 8.9 | U | 0.89 | < 9.4 | 9.4 | U | 0.94 |
| Methylene chloride | 50 | 100,000 | < 4.2 | 4.2 | U | 4.2 | < 4.4 | 4.4 | U | 4.4 | < 4.7 | 4.7 | U | 4.7 |
| Naphthalene | 12,000 | 100,000 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| n-Butylbenzene | 12,000 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| n-Propylbenzene | 3,900 | 100,000 | < 4.2 | 4.2 | U | 0.76 | < 4.4 | 4.4 | U | 0.80 | < 4.7 | 4.7 | U | 0.85 |
| o-Xylene | 260 | 100,000 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| p-Isopropyltoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| sec-Butylbenzene | 11,000 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Styrene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Tert-butyl alcohol | | | < 85 | 85 | U | 17 | < 89 | 89 | U | 18 | < 94 | 94 | U | 19 |
| tert-Butylbenzene | 5,900 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Tetrachloroethene | 1,300 | 19,000 | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Tetrahydrofuran (THF) | | | < 8.5 | 8.5 | U | 2.1 | < 8.9 | 8.9 | U | 2.2 | < 9.4 | 9.4 | U | 2.4 |
| Toluene | 700 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| trans-1,3-Dichloropropene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| trans-1,4-dichloro-2-butene | | | < 8.5 | 8.5 | U | 2.1 | < 8.9 | 8.9 | U | 2.2 | < 9.4 | 9.4 | U | 2.4 |
| Trichloroethene | 470 | 21,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Trichlorofluoromethane | | | < 4.2 | 4.2 | U | 0.85 | < 4.4 | 4.4 | U | 0.89 | < 4.7 | 4.7 | U | 0.94 |
| Trichlorotrifluoroethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Vinyl Chloride | 20 | 900 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 4.7 | 4.7 | U | 0.47 |
| Total BTEX Concentration | | | 0 | | | | 0 | | | | 0 | | | |
| Total VOCs Concentration | | | 5 | | | | 6.2 | | | | 17 | | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B11 | | | | | | | | 15B12 | | | |
|----------------------------------|---|---|-----------------|-------------|------|------|-----------------|-----------|------|------|-----------------|--------------|------|------|
| | | | 11/23/2015 | | | | | | | | 11/23/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 30 | 30 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 18 | 18 | U | 0.89 |
| 1,1,2,2-Tetrachloroethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| 1,1,2-Trichloroethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| 1,1-Dichloroethane | 270 | 26,000 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| 1,1-Dichloroethene | 330 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 1,1-Dichloropropene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 1,2,3-Trichlorobenzene | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| 1,2,3-Trichloropropane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,2,4-Trichlorobenzene | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,2-Dibromo-3-chloropropane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| 1,2-Dibromomethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,2-Dichloroethane | 20 | 3,100 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 1,2-Dichloropropane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,3-Dichloropropane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 1,4-dioxane | | | < 100 | 100 | U | 60 | < 73 | 73 | U | 29 | < 89 | 89 | U | 36 |
| 2,2-Dichloropropane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| 2-Chlorotoluene | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 37 | 37 | U | 7.5 | < 18 | 18 | U | 3.7 | < 22 | 22 | U | 4.4 |
| 2-Isopropyltoluene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 4-Chlorotoluene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| 4-Methyl-2-Pentanone | | | < 37 | 37 | U | 7.5 | < 18 | 18 | U | 3.7 | < 22 | 22 | U | 4.4 |
| Acetone | 50 | 100,000 | < 37 | 37 | U | 7.5 | 19 | 37 | JS | 3.7 | 100 | 44 | S | 4.4 |
| Acrolein | | | < 30 | 30 | U | 3.7 | < 15 | 15 | U | 1.8 | < 18 | 18 | U | 2.2 |
| Acrylonitrile | | | < 15 | 15 | U | 1.5 | < 7.3 | 7.3 | U | 0.73 | < 8.9 | 8.9 | U | 0.89 |
| Benzene | 60 | 4,800 | 0.79 | 7.5 | J | 0.75 | < 3.7 | 3.7 | U | 0.37 | 61 | 60 | U | 27 |
| Bromobenzene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| Bromochloromethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Bromodichloromethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Bromoforn | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Bromomethane | | | < 7.5 | 7.5 | U | 3.0 | < 3.7 | 3.7 | U | 1.5 | < 4.4 | 4.4 | U | 1.8 |
| Carbon Disulfide | | | 2.8 | 7.5 | J | 1.5 | < 3.7 | 3.7 | U | 0.73 | 2.3 | 4.4 | J | 0.89 |
| Carbon tetrachloride | 760 | 2,400 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Chlorobenzene | 1,100 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Chloroethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Chloroform | 370 | 49,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Chloromethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| cis-1,3-Dichloropropene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Dibromochloromethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Dibromomethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Dichlorodifluoromethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Ethylbenzene | 1,000 | 41,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Hexachlorobutadiene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| Isopropylbenzene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| m&p-Xylenes | 260 | 100,000 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 45 | 45 | U | 7.5 | < 22 | 22 | U | 3.7 | < 27 | 27 | U | 4.4 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 15 | 15 | U | 1.5 | < 7.3 | 7.3 | U | 0.73 | < 8.9 | 8.9 | U | 0.89 |
| Methylene chloride | 50 | 100,000 | < 7.5 | 7.5 | U | 7.5 | < 3.7 | 3.7 | U | 3.7 | < 4.4 | 4.4 | U | 4.4 |
| Naphthalene | 12,000 | 100,000 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 270 | 270 | U | 54 |
| n-Butylbenzene | 12,000 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| n-Propylbenzene | 3,900 | 100,000 | < 7.5 | 7.5 | U | 1.3 | < 3.7 | 3.7 | U | 0.66 | < 270 | 270 | U | 49 |
| o-Xylene | 260 | 100,000 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| p-Isopropyltoluene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| sec-Butylbenzene | 11,000 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| Styrene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Tert-butyl alcohol | | | < 150 | 150 | U | 30 | < 73 | 73 | U | 15 | < 89 | 89 | U | 18 |
| tert-Butylbenzene | 5,900 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 270 | 270 | U | 27 |
| Tetrachloroethene | 1,300 | 19,000 | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Tetrahydrofuran (THF) | | | < 15 | 15 | U | 3.7 | < 7.3 | 7.3 | U | 1.8 | < 8.9 | 8.9 | U | 2.2 |
| Toluene | 700 | 100,000 | 1.3 | 7.5 | J | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| trans-1,3-Dichloropropene | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| trans-1,4-dichloro-2-butene | | | < 15 | 15 | U | 3.7 | < 7.3 | 7.3 | U | 1.8 | < 540 | 540 | U | 140 |
| Trichloroethene | 470 | 21,000 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Trichlorofluoromethane | | | < 7.5 | 7.5 | U | 1.5 | < 3.7 | 3.7 | U | 0.73 | < 4.4 | 4.4 | U | 0.89 |
| Trichlorotrifluoroethane | | | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Vinyl Chloride | 20 | 900 | < 7.5 | 7.5 | U | 0.75 | < 3.7 | 3.7 | U | 0.37 | < 4.4 | 4.4 | U | 0.44 |
| Total BTEX Concentration | | | | 2.09 | | | | 0 | | | | 61 | | |
| Total VOCs Concentration | | | | 4.89 | | | | 19 | | | | 163.3 | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B13 | | | | | | | | 15B14 | | | | 15B15 | | | |
|----------------------------------|---|---|-------------|--------|----|------|--------|--------|----|------|------------|--------|----|------|------------|-----|----|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (0-2') | | | | (6-8') | | | | (5-7') | | | | (6-8') | | | |
| | | | µg/Kg | | | | µg/Kg | | | | µg/Kg | | | | µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,1,1,2-Tetrachloroethane | | | < 5.9 | 5.9 | U | 1.2 | < 20 | 20 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.46 |
| 1,1,2,2-Tetrachloroethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,1,2-Trichloroethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,1-Dichloroethane | 270 | 26,000 | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,1-Dichloroethene | 330 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,1-Dichloropropene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2,3-Trichlorobenzene | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,2,3-Trichloropropane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2,4-Trichlorobenzene | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | 85 | 410 | J | 41 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2-Dibromo-3-chloropropane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,2-Dibromomethane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2-Dichloroethane | 20 | 3,100 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,2-Dichloropropane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,3-Dichloropropane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 1,4-dioxane | | | < 100 | 100 | U | 48 | < 100 | 100 | U | 40 | < 100 | 100 | U | 62 | < 93 | 93 | U | 37 |
| 2,2-Dichloropropane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 2-Chlorotoluene | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 30 | 30 | U | 5.9 | < 25 | 25 | U | 5.1 | < 39 | 39 | U | 7.7 | < 23 | 23 | U | 4.6 |
| 2-Isopropyltoluene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 4-Chlorotoluene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| 4-Methyl-2-Pentanone | | | < 30 | 30 | U | 5.9 | < 25 | 25 | U | 5.1 | < 39 | 39 | U | 7.7 | < 23 | 23 | U | 4.6 |
| Acetone | 50 | 100,000 | 34 | 50 | JS | 5.9 | < 50 | 50 | U | 5.1 | < 50 | 50 | U | 7.7 | 18 | 46 | JS | 4.6 |
| Acrolein | | | < 24 | 24 | U | 3.0 | < 20 | 20 | U | 2.5 | < 31 | 31 | U | 3.9 | < 19 | 19 | U | 2.3 |
| Acrylonitrile | | | < 24 | 24 | U | 0.59 | < 10 | 10 | U | 1.0 | < 15 | 15 | U | 1.5 | < 9.3 | 9.3 | U | 0.93 |
| Benzene | 60 | 4,800 | 0.73 | 5.9 | J | 0.59 | < 5.1 | 5.1 | U | 0.51 | 4.5 | 7.7 | J | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Bromobenzene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Bromochloromethane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Bromodichloromethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Bromoform | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Bromomethane | | | < 5.9 | 5.9 | U | 2.4 | < 5.1 | 5.1 | U | 2.0 | < 7.7 | 7.7 | U | 3.1 | < 4.6 | 4.6 | U | 1.9 |
| Carbon Disulfide | | | 3.8 | 5.9 | J | 1.2 | < 5.1 | 5.1 | U | 1.0 | 3.7 | 7.7 | J | 1.5 | 4 | 4.6 | J | 0.93 |
| Carbon tetrachloride | 760 | 2,400 | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Chlorobenzene | 1,100 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Chloroethane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Chloroform | 370 | 49,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Chloromethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| cis-1,3-Dichloropropene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Dibromochloromethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Dibromomethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Dichlorodifluoromethane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Ethylbenzene | 1,000 | 41,000 | 45 | 410 | J | 41 | < 5.1 | 5.1 | U | 0.51 | 1 | 7.7 | J | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Hexachlorobutadiene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Isopropylbenzene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| m&p-Xylenes | 260 | 100,000 | 210 | 410 | J | 82 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 36 | 36 | U | 5.9 | < 30 | 30 | U | 5.1 | < 46 | 46 | U | 7.7 | < 28 | 28 | U | 4.6 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 12 | 12 | U | 1.2 | < 10 | 10 | U | 1.0 | < 15 | 15 | U | 1.5 | < 9.3 | 9.3 | U | 0.93 |
| Methylene chloride | 50 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Naphthalene | 12,000 | 100,000 | 170 | 410 | J | 82 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| n-Butylbenzene | 12,000 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| n-Propylbenzene | 3,900 | 100,000 | < 5.9 | 5.9 | U | 1.1 | < 5.1 | 5.1 | U | 0.91 | < 7.7 | 7.7 | U | 1.4 | < 4.6 | 4.6 | U | 0.83 |
| o-Xylene | 260 | 100,000 | 150 | 410 | J | 82 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| p-Isopropyltoluene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| sec-Butylbenzene | 11,000 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Styrene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Tert-butyl alcohol | | | < 120 | 120 | U | 24 | < 100 | 100 | U | 20 | < 150 | 150 | U | 31 | < 93 | 93 | U | 19 |
| tert-Butylbenzene | 5,900 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Tetrachloroethene | 1,300 | 19,000 | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Tetrahydrofuran (THF) | | | < 12 | 12 | U | 3.0 | < 10 | 10 | U | 2.5 | < 15 | 15 | U | 3.9 | < 9.3 | 9.3 | U | 2.3 |
| Toluene | 700 | 100,000 | 96 | 410 | J | 41 | < 5.1 | 5.1 | U | 0.51 | 2.3 | 7.7 | J | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| trans-1,3-Dichloropropene | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| trans-1,4-dichloro-2-butene | | | < 12 | 12 | U | 3.0 | < 10 | 10 | U | 2.5 | < 15 | 15 | U | 3.9 | < 9.3 | 9.3 | U | 2.3 |
| Trichloroethene | 470 | 21,000 | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Trichlorofluoromethane | | | < 5.9 | 5.9 | U | 1.2 | < 5.1 | 5.1 | U | 1.0 | < 7.7 | 7.7 | U | 1.5 | < 4.6 | 4.6 | U | 0.93 |
| Trichlorotrifluoroethane | | | < 5.9 | 5.9 | U | 0.59 | < 5.1 | 5.1 | U | 0.51 | < 7.7 | 7.7 | U | 0.77 | < 4.6 | 4.6 | U | 0.46 |
| Vinyl Chloride | 20 | 900 | < 5.9</ | | | | | | | | | | | | | | | |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B17 | | | | 15B18 | | | | | | | |
|----------------------------------|---|---|------------|-----|------|------|-------------|-----|------|------|--------|-----|------|------|
| | | | 11/20/2015 | | | | 11/20/2015 | | | | | | | |
| | | | (6-8") | | | | (0-2") | | | | (6-8") | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,1,2,2-Tetrachloroethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,1,2-Trichloroethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,1-Dichloroethane | 270 | 26,000 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,1-Dichloroethene | 330 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,1-Dichloropropene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2,3-Trichlorobenzene | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,2,3-Trichloropropane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2,4-Trichlorobenzene | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2-Dibromo-3-chloropropane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,2-Dibromomethane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2-Dichloroethane | 20 | 3,100 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,2-Dichloropropane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,3-Dichloropropane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 1,4-dioxane | | | < 100 | 100 | U | 47 | < 69 | 69 | U | 28 | < 82 | 82 | U | 33 |
| 2,2-Dichloropropane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 2-Chlorotoluene | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 29 | 29 | U | 5.8 | < 17 | 17 | U | 3.5 | < 20 | 20 | U | 4.1 |
| 2-Isopropyltoluene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 4-Chlorotoluene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| 4-Methyl-2-Pentanone | | | < 29 | 29 | U | 5.8 | < 17 | 17 | U | 3.5 | < 20 | 20 | U | 4.1 |
| Acetone | 50 | 100,000 | < 50 | 50 | U | 5.8 | 6.1 | 35 | JS | 3.5 | < 4.1 | 4.1 | U | 4.1 |
| Acrolein | | | < 23 | 23 | U | 2.9 | < 14 | 14 | U | 1.7 | < 16 | 16 | U | 2.0 |
| Acrylonitrile | | | < 12 | 12 | U | 1.2 | < 6.9 | 6.9 | U | 0.69 | < 8.2 | 8.2 | U | 0.82 |
| Benzene | 60 | 4,800 | < 5.8 | 5.8 | U | 0.58 | 1.3 | 3.5 | J | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Bromobenzene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Bromochloromethane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Bromodichloromethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Bromoform | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Bromomethane | | | < 5.8 | 5.8 | U | 2.3 | < 3.5 | 3.5 | U | 1.4 | < 4.1 | 4.1 | U | 1.6 |
| Carbon Disulfide | | | 1.2 | 5.8 | J | 1.2 | 1.3 | 3.5 | J | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Carbon tetrachloride | 760 | 2,400 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Chlorobenzene | 1,100 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Chloroethane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Chloroform | 370 | 49,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Chloromethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| cis-1,3-Dichloropropene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Dibromochloromethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Dibromomethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Dichlorodifluoromethane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Ethylbenzene | 1,000 | 41,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Hexachlorobutadiene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Isopropylbenzene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| m&p-Xylenes | 260 | 100,000 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 35 | 35 | U | 5.8 | < 21 | 21 | U | 3.5 | < 25 | 25 | U | 4.1 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 12 | 12 | U | 1.2 | < 6.9 | 6.9 | U | 0.69 | < 8.2 | 8.2 | U | 0.82 |
| Methylene chloride | 50 | 100,000 | < 5.8 | 5.8 | U | 5.8 | < 3.5 | 3.5 | U | 3.5 | < 4.1 | 4.1 | U | 4.1 |
| Naphthalene | 12,000 | 100,000 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| n-Butylbenzene | 12,000 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| n-Propylbenzene | 3,900 | 100,000 | < 5.8 | 5.8 | U | 1.1 | < 3.5 | 3.5 | U | 0.62 | < 4.1 | 4.1 | U | 0.74 |
| o-Xylene | 260 | 100,000 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| p-Isopropyltoluene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| sec-Butylbenzene | 11,000 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Styrene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Tert-butyl alcohol | | | < 120 | 120 | U | 23 | < 69 | 69 | U | 14 | < 82 | 82 | U | 16 |
| tert-Butylbenzene | 5,900 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Tetrachloroethene | 1,300 | 19,000 | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Tetrahydrofuran (THF) | | | < 12 | 12 | U | 2.9 | < 6.9 | 6.9 | U | 1.7 | < 8.2 | 8.2 | U | 2.0 |
| Toluene | 700 | 100,000 | < 5.8 | 5.8 | U | 0.58 | 0.65 | 3.5 | J | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| trans-1,3-Dichloropropene | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| trans-1,4-dichloro-2-butene | | | < 12 | 12 | U | 2.9 | < 6.9 | 6.9 | U | 1.7 | < 8.2 | 8.2 | U | 2.0 |
| Trichloroethene | 470 | 21,000 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Trichlorofluoromethane | | | < 5.8 | 5.8 | U | 1.2 | < 3.5 | 3.5 | U | 0.69 | < 4.1 | 4.1 | U | 0.82 |
| Trichlorotrifluoroethane | | | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Vinyl Chloride | 20 | 900 | < 5.8 | 5.8 | U | 0.58 | < 3.5 | 3.5 | U | 0.35 | < 4.1 | 4.1 | U | 0.41 |
| Total BTEX Concentration | | | 0 | | | | 1.95 | | | | 0 | | | |
| Total VOCs Concentration | | | 1.2 | | | | 9.35 | | | | 0 | | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B19 | | | | | | | | 15B20 | | | | 15B21 | | | |
|----------------------------------|---|---|------------|--------|----|------|--------|--------|----|------|------------|--------|----|------|------------|-----|----|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (0-2') | | | | (6-8') | | | | (8-10') | | | | (6-8') | | | |
| | | | µg/Kg | | | | µg/Kg | | | | µg/Kg | | | | µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,1,1,2-Tetrachloroethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 30 | 30 | U | 1.5 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,1,2,2-Tetrachloroethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,1,2-Trichloroethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,1-Dichloroethane | 270 | 26,000 | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,1-Dichloroethene | 330 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,1-Dichloropropene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2,3-Trichlorobenzene | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,2,3-Trichloropropane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2,4-Trichlorobenzene | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | 3.7 | 4.2 | J | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2-Dibromo-3-chloropropane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,2-Dibromomethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2-Dichloroethane | 20 | 3,100 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,2-Dichloropropane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | 1.7 | 4.2 | J | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,3-Dichloropropane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 1,4-dioxane | | | < 84 | 84 | U | 34 | < 88 | 88 | U | 35 | < 70 | 70 | U | 28 | < 100 | 100 | U | 59 |
| 2,2-Dichloropropane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 2-Chlorotoluene | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 21 | 21 | U | 4.2 | < 22 | 22 | U | 4.4 | < 17 | 17 | U | 3.5 | < 37 | 37 | U | 7.4 |
| 2-Isopropyltoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 4-Chlorotoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| 4-Methyl-2-Pentanone | | | < 21 | 21 | U | 4.2 | < 22 | 22 | U | 4.4 | < 17 | 17 | U | 3.5 | < 37 | 37 | U | 7.4 |
| Acetone | 50 | 100,000 | 6.8 | 4.2 | JS | 4.2 | < 44 | 44 | U | 4.4 | 10 | 35 | JS | 3.5 | 72 | 74 | JS | 7.4 |
| Acrolein | | | < 17 | 17 | U | 2.1 | < 18 | 18 | U | 2.2 | < 14 | 14 | U | 1.7 | < 30 | 30 | U | 3.7 |
| Acrylonitrile | | | < 8.4 | 8.4 | U | 0.84 | < 8.8 | 8.8 | U | 0.88 | < 7.0 | 7.0 | U | 0.70 | < 30 | 30 | U | 0.74 |
| Benzene | 60 | 4,800 | 0.5 | 4.2 | J | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Bromobenzene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Bromochloromethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Bromodichloromethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Bromoform | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Bromomethane | | | < 4.2 | 4.2 | U | 1.7 | < 4.4 | 4.4 | U | 1.8 | < 3.5 | 3.5 | U | 1.4 | < 7.4 | 7.4 | U | 3.0 |
| Carbon Disulfide | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | 5.4 | 3.5 | U | 0.70 | 9.3 | 7.4 | U | 1.5 |
| Carbon tetrachloride | 760 | 2,400 | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Chlorobenzene | 1,100 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Chloroethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Chloroform | 370 | 49,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Chloromethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| cis-1,3-Dichloropropene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Dibromochloromethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Dibromomethane | | | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Dichlorodifluoromethane | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Ethylbenzene | 1,000 | 41,000 | 0.61 | 4.2 | J | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Hexachlorobutadiene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Isopropylbenzene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| m&p-Xylenes | 260 | 100,000 | 2.2 | 4.2 | J | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 25 | 25 | U | 4.2 | < 26 | 26 | U | 4.4 | < 21 | 21 | U | 3.5 | 21 | 45 | J | 7.4 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 8.4 | 8.4 | U | 0.84 | < 8.8 | 8.8 | U | 0.88 | < 7.0 | 7.0 | U | 0.70 | 1.9 | 15 | J | 1.5 |
| Methylene chloride | 50 | 100,000 | < 4.2 | 4.2 | U | 4.2 | < 4.4 | 4.4 | U | 4.4 | < 3.5 | 3.5 | U | 3.5 | < 7.4 | 7.4 | U | 7.4 |
| Naphthalene | 12,000 | 100,000 | 2 | 4.2 | J | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| n-Butylbenzene | 12,000 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| n-Propylbenzene | 3,900 | 100,000 | < 4.2 | 4.2 | U | 0.76 | < 4.4 | 4.4 | U | 0.79 | < 3.5 | 3.5 | U | 0.63 | < 7.4 | 7.4 | U | 1.3 |
| o-Xylene | 260 | 100,000 | 1.2 | 4.2 | J | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |
| p-Isopropyltoluene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| sec-Butylbenzene | 11,000 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Styrene | | | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Tert-butyl alcohol | | | < 84 | 84 | U | 17 | < 88 | 88 | U | 18 | < 70 | 70 | U | 14 | < 150 | 150 | U | 30 |
| tert-Butylbenzene | 5,900 | 100,000 | < 4.2 | 4.2 | U | 0.42 | < 4.4 | 4.4 | U | 0.44 | < 3.5 | 3.5 | U | 0.35 | < 7.4 | 7.4 | U | 0.74 |
| Tetrachloroethene | 1,300 | 19,000 | < 4.2 | 4.2 | U | 0.84 | < 4.4 | 4.4 | U | 0.88 | < 3.5 | 3.5 | U | 0.70 | < 7.4 | 7.4 | U | 1.5 |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B22 | | | | | | | | 15B23 | | | | | | | |
|----------------------------------|---|---|-------------|-----|--------|------|--------------|-------|--------|------|------------|-------|--------|------------|------------|-----|--------|------|
| | | | 11/20/2015 | | | | 11/20/2015 | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (0-2') | | (6-8') | | (2-4') | | (6-8') | | (2-4') | | (6-8') | | (2-4') | | (6-8') | |
| | | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | | µg/Kg | |
| | | | | | | | | | | | | | | | | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,1,2,2-Tetrachloroethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,1,2-Trichloroethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,1-Dichloroethane | 270 | 26,000 | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,1-Dichloroethene | 330 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,1-Dichloropropene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,2,3-Trichlorobenzene | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,2,3-Trichloropropane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,2,4-Trichlorobenzene | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 5.0 | 5.0 | U | 0.50 | 34 | 330 | J | 33 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,2-Dibromo-3-chloropropane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,2-Dibromomethane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,2-Dichloroethane | 20 | 3,100 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 20 | 20 | U | 20 | < 7.3 | 7.3 | U | 0.73 |
| 1,2-Dichloropropane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,3-Dichloropropane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 1,4-dioxane | | | < 100 | 100 | U | 40 | < 100 | 100 | U | 51 | < 1800 | 1,800 | U | 1800 | < 100 | 100 | U | 58 |
| 2,2-Dichloropropane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 2-Chlorotoluene | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 25 | 25 | U | 5.0 | < 32 | 32 | U | 6.3 | < 1100 | 1,100 | U | 230 | < 36 | 36 | U | 7.3 |
| 2-Isopropyltoluene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 4-Chlorotoluene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| 4-Methyl-2-Pentanone | | | < 25 | 25 | U | 5.0 | < 32 | 32 | U | 6.3 | < 1100 | 1,100 | U | 230 | < 36 | 36 | U | 7.3 |
| Acetone | 50 | 100,000 | 19 | 50 | JS | 5.0 | 1,500 | 3,300 | JS | 330 | 510 | 2,300 | JS | 230 | 23 | 50 | JS | 7.3 |
| Acrolein | | | < 20 | 20 | U | 2.5 | < 25 | 25 | U | 3.2 | < 910 | 910 | U | 110 | < 29 | 29 | U | 3.6 |
| Acrylonitrile | | | < 20 | 20 | U | 0.50 | < 13 | 13 | U | 1.3 | < 450 | 450 | U | 45 | < 15 | 15 | U | 1.5 |
| Benzene | 60 | 4,800 | 0.94 | 5.0 | J | 0.50 | < 6.3 | 6.3 | U | 0.63 | 29 | 59 | J | 23 | < 7.3 | 7.3 | U | 0.73 |
| Bromobenzene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Bromochloromethane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Bromodichloromethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Bromoform | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Bromomethane | | | < 5.0 | 5.0 | U | 2.0 | < 6.3 | 6.3 | U | 2.5 | < 230 | 230 | U | 91 | < 7.3 | 7.3 | U | 2.9 |
| Carbon Disulfide | | | < 5.0 | 5.0 | U | 1.0 | 180 | 330 | J | 65 | < 230 | 230 | U | 45 | 6.2 | 7.3 | J | 1.5 |
| Carbon tetrachloride | 760 | 2,400 | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Chlorobenzene | 1,100 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Chloroethane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Chloroform | 370 | 49,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Chloromethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | 120 | 230 | J | 45 | < 7.3 | 7.3 | U | 1.5 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| cis-1,3-Dichloropropene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Dibromochloromethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Dibromomethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Dichlorodifluoromethane | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Ethylbenzene | 1,000 | 41,000 | < 5.0 | 5.0 | U | 0.50 | 45 | 330 | J | 33 | 67 | 230 | J | 23 | < 7.3 | 7.3 | U | 0.73 |
| Hexachlorobutadiene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Isopropylbenzene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| m&p-Xylenes | 260 | 100,000 | < 5.0 | 5.0 | U | 1.0 | 89 | 330 | J | 65 | 58 | 230 | J | 45 | < 7.3 | 7.3 | U | 1.5 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 30 | 30 | U | 5.0 | < 38 | 38 | U | 6.3 | < 230 | 230 | U | 230 | < 44 | 44 | U | 7.3 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 10 | 10 | U | 1.0 | < 13 | 13 | U | 1.3 | < 450 | 450 | U | 45 | < 15 | 15 | U | 1.5 |
| Methylene chloride | 50 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 230 | < 7.3 | 7.3 | U | 7.3 |
| Naphthalene | 12,000 | 100,000 | < 5.0 | 5.0 | U | 1.0 | 400 | 330 | | 65 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| n-Butylbenzene | 12,000 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | 29 | 230 | J | 23 | < 7.3 | 7.3 | U | 0.73 |
| n-Propylbenzene | 3,900 | 100,000 | < 5.0 | 5.0 | U | 0.90 | < 6.3 | 6.3 | U | 1.1 | < 230 | 230 | U | 41 | < 7.3 | 7.3 | U | 1.3 |
| o-Xylene | 260 | 100,000 | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| p-Isopropyltoluene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| sec-Butylbenzene | 11,000 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Styrene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Tert-butyl alcohol | | | < 100 | 100 | U | 20 | < 130 | 130 | U | 25 | < 4500 | 4,500 | U | 910 | < 150 | 150 | U | 29 |
| tert-Butylbenzene | 5,900 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Tetrachloroethene | 1,300 | 19,000 | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Tetrahydrofuran (THF) | | | < 10 | 10 | U | 2.5 | < 13 | 13 | U | 3.2 | < 450 | 450 | U | 110 | < 15 | 15 | U | 3.6 |
| Toluene | 700 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | 60 | 230 | J | 23 | < 7.3 | 7.3 | U | 0.73 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 190 | 190 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| trans-1,3-Dichloropropene | | | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| trans-1,4-dichloro-2-butene | | | < 10 | 10 | U | 2.5 | < 13 | 13 | U | 3.2 | < 450 | 450 | U | 110 | < 15 | 15 | U | 3.6 |
| Trichloroethene | 470 | 21,000 | < 5.0 | 5.0 | U | 0.50 | < 6.3 | 6.3 | U | 0.63 | < 230 | 230 | U | 23 | < 7.3 | 7.3 | U | 0.73 |
| Trichlorofluoromethane | | | < 5.0 | 5.0 | U | 1.0 | < 6.3 | 6.3 | U | 1.3 | < 230 | 230 | U | 45 | < 7.3 | 7.3 | U | 1.5 |
| Trichlorotrifluoroethane | | | < 5.0 | 5.0 | U | 0.50 | < | | | | | | | | | | | |

TABLE 3
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Volatile Organic Compounds

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | Duplicate | | | | Duplicate | | | |
|----------------------------------|---|---|-------------|--------|----|------|------------|-------|---|------|
| | | | 11/20/2015 | | | | 11/23/2015 | | | |
| | | | µg/Kg | | | | µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,1,1,2-Tetrachloroethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 34 |
| 1,1,1-Trichloroethane | 680 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 69 |
| 1,1,2,2-Tetrachloroethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,1,2-Trichloroethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,1-Dichloroethane | 270 | 26,000 | < 9.7 | 9.7 | U | 1.9 | < 250 | 250 | U | 69 |
| 1,1-Dichloroethene | 330 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 330 | 330 | U | 34 |
| 1,1-Dichloropropene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2,3-Trichlorobenzene | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,2,3-Trichloropropane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2,4-Trichlorobenzene | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,2,4-Trimethylbenzene | 3,600 | 52,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2-Dibromo-3-chloropropane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,2-Dibromomethane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2-Dichlorobenzene | 1,100 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2-Dichloroethane | 20 | 3,100 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,2-Dichloropropane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,3,5-Trimethylbenzene | 8,400 | 52,000 | 1.6 | 9.7 | J | 0.97 | < 340 | 340 | U | 34 |
| 1,3-Dichlorobenzene | 2,400 | 4,900 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,3-Dichloropropane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 1,4-Dichlorobenzene | 1,800 | 13,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 1,4-dioxane | | | < 100 | 100 | U | 77 | < 6900 | 6,900 | U | 2800 |
| 2,2-Dichloropropane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 2-Chlorotoluene | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| 2-Hexanone (Methyl Butyl Ketone) | | | < 48 | 48 | U | 9.7 | < 1700 | 1,700 | U | 340 |
| 2-Isopropyltoluene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 4-Chlorotoluene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| 4-Methyl-2-Pentanone | | | < 48 | 48 | U | 9.7 | < 1700 | 1,700 | U | 340 |
| Acetone | 50 | 100,000 | 18 | 50 | JS | 9.7 | < 340 | 340 | U | 340 |
| Acrolein | | | < 39 | 39 | U | 4.8 | < 1400 | 1,400 | U | 170 |
| Acrylonitrile | | | < 39 | 39 | U | 0.97 | < 690 | 690 | U | 69 |
| Benzene | 60 | 4,800 | < 9.7 | 9.7 | U | 0.97 | 57 | 60 | J | 34 |
| Bromobenzene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Bromochloromethane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Bromodichloromethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Bromoform | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Bromomethane | | | < 9.7 | 9.7 | U | 3.9 | < 340 | 340 | U | 140 |
| Carbon Disulfide | | | 3.3 | 9.7 | J | 1.9 | < 340 | 340 | U | 69 |
| Carbon tetrachloride | 760 | 2,400 | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Chlorobenzene | 1,100 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Chloroethane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Chloroform | 370 | 49,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Chloromethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| cis-1,2-Dichloroethene | 250 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 250 | 250 | U | 34 |
| cis-1,3-Dichloropropene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Dibromochloromethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Dibromomethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Dichlorodifluoromethane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Ethylbenzene | 1,000 | 41,000 | < 9.7 | 9.7 | U | 0.97 | 44 | 340 | J | 34 |
| Hexachlorobutadiene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Isopropylbenzene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| m&p-Xylenes | 260 | 100,000 | < 9.7 | 9.7 | U | 1.9 | 73 | 340 | J | 69 |
| Methyl Ethyl Ketone (2-Butanone) | 120 | 100,000 | < 58 | 58 | U | 9.7 | < 340 | 340 | U | 340 |
| Methyl t-butyl ether (MTBE) | 930 | 100,000 | < 19 | 19 | U | 1.9 | < 690 | 690 | U | 69 |
| Methylene chloride | 50 | 100,000 | < 9.7 | 9.7 | U | 9.7 | < 340 | 340 | U | 340 |
| Naphthalene | 12,000 | 100,000 | < 9.7 | 9.7 | U | 1.9 | 160 | 340 | J | 69 |
| n-Butylbenzene | 12,000 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| n-Propylbenzene | 3,900 | 100,000 | < 9.7 | 9.7 | U | 1.7 | < 340 | 340 | U | 62 |
| o-Xylene | 260 | 100,000 | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| p-Isopropyltoluene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| sec-Butylbenzene | 11,000 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Styrene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Tert-butyl alcohol | | | < 190 | 190 | U | 39 | < 6900 | 6,900 | U | 1400 |
| tert-Butylbenzene | 5,900 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Tetrachloroethene | 1,300 | 19,000 | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Tetrahydrofuran (THF) | | | < 19 | 19 | U | 4.8 | < 690 | 690 | U | 170 |
| Toluene | 700 | 100,000 | < 9.7 | 9.7 | U | 0.97 | 75 | 340 | J | 34 |
| trans-1,2-Dichloroethene | 190 | 100,000 | < 9.7 | 9.7 | U | 0.97 | < 150 | 150 | U | 34 |
| trans-1,3-Dichloropropene | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| trans-1,4-dichloro-2-butene | | | < 19 | 19 | U | 4.8 | < 690 | 690 | U | 170 |
| Trichloroethene | 470 | 21,000 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Trichlorofluoromethane | | | < 9.7 | 9.7 | U | 1.9 | < 340 | 340 | U | 69 |
| Trichlorotrifluoroethane | | | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Vinyl Chloride | 20 | 900 | < 9.7 | 9.7 | U | 0.97 | < 340 | 340 | U | 34 |
| Total BTEX Concentration | | | 0 | | | | 249 | | | |
| Total VOCs Concentration | | | 22.9 | | | | 409 | | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B1 | | | | | | | | 15B2 | | | |
|-------------------------------|--|--|----------------|--------|----|------|----------------|--------|----|------|----------------|-------|----|-----|
| | | | 11/24/2015 | | | | | | | | 11/24/2015 | | | |
| | | | (0-2) µg/Kg | | | | (5-7) µg/Kg | | | | (6-8) µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,2,4,5-Tetrachlorobenzene | | | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| 1,2,4-Trichlorobenzene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| 1,2-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| 1,2-Diphenylhydrazine | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| 1,3-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| 1,4-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| 2,4,5-Trichlorophenol | | | < 260 | 260 | U | 210 | < 280 | 280 | U | 220 | < 320 | 320 | U | 250 |
| 2,4,6-Trichlorophenol | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 140 |
| 2,4-Dichlorophenol | | | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| 2,4-Dimethylphenol | | | < 260 | 260 | U | 94 | < 280 | 280 | U | 98 | < 320 | 320 | U | 110 |
| 2,4-Dinitrophenol | | | < 760 | 760 | U | 260 | < 790 | 790 | U | 280 | < 900 | 900 | U | 320 |
| 2,4-Dinitrotoluene | | | < 260 | 260 | U | 150 | < 280 | 280 | U | 160 | < 320 | 320 | U | 180 |
| 2,6-Dinitrotoluene | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| 2-Chloronaphthalene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| 2-Chlorophenol | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| 2-Methylnaphthalene | | | 1,900 | 260 | | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 260 | 260 | U | 180 | < 280 | 280 | U | 190 | < 320 | 320 | U | 210 |
| 2-Nitroaniline | | | < 760 | 760 | U | 380 | < 790 | 790 | U | 400 | < 900 | 900 | U | 460 |
| 2-Nitrophenol | | | < 260 | 260 | U | 240 | < 280 | 280 | U | 250 | < 320 | 320 | U | 290 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 260 | 260 | U | 150 | < 280 | 280 | U | 160 | < 320 | 320 | U | 180 |
| 3,3'-Dichlorobenzidine | | | < 760 | 760 | U | 180 | < 790 | 790 | U | 190 | < 900 | 900 | U | 210 |
| 3-Nitroaniline | | | < 760 | 760 | U | 760 | < 790 | 790 | U | 790 | < 900 | 900 | U | 900 |
| 4,6-Dinitro-2-methylphenol | | | < 1900 | 1,900 | U | 410 | < 2000 | 2,000 | U | 420 | < 2300 | 2,300 | U | 490 |
| 4-Bromophenyl phenyl ether | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| 4-Chloro-3-methylphenol | | | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| 4-Chloroaniline | | | < 300 | 300 | U | 180 | < 320 | 320 | U | 180 | < 360 | 360 | U | 210 |
| 4-Chlorophenyl phenyl ether | | | < 260 | 260 | U | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| 4-Nitroaniline | | | < 760 | 760 | U | 130 | < 790 | 790 | U | 130 | < 900 | 900 | U | 150 |
| 4-Nitrophenol | | | < 380 | 380 | U | 170 | < 390 | 390 | U | 180 | < 450 | 450 | U | 200 |
| Acenaphthene | 20,000 | 100,000 | 2,200 | 260 | | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Acenaphthylene | 100,000 | 100,000 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Acetophenone | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Aniline | | | < 300 | 300 | U | 300 | < 320 | 320 | U | 320 | < 360 | 360 | U | 360 |
| Anthracene | 100,000 | 100,000 | 1,600 | 260 | | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benz(a)anthracene | 1,000 | 1,000 | 630 | 260 | | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benzbidine | | | < 760 | 760 | U | 220 | < 790 | 790 | U | 230 | < 900 | 900 | U | 270 |
| Benzo(a)pyrene | 1,000 | 1,000 | 300 | 260 | | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 330 | 260 | | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 800 | 260 | | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benzo(k)fluoranthene | 800 | 3,900 | 180 | 260 | J | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Benzoic acid | | | < 1900 | 1,900 | U | 760 | < 2000 | 2,000 | U | 790 | < 2300 | 2,300 | U | 900 |
| Benzyl butyl phthalate | | | < 260 | 260 | U | 97 | < 280 | 280 | U | 100 | < 320 | 320 | U | 120 |
| Bis(2-chloroethoxy)methane | | | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 | < 320 | 320 | U | 120 |
| Bis(2-chloroethyl)ether | | | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 | < 320 | 320 | U | 120 |
| Bis(2-chloroisopropyl)ether | | | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Bis(2-ethylhexyl)phthalate | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Carbazole | | | < 1900 | 1,900 | U | 290 | < 2000 | 2,000 | U | 300 | < 2300 | 2,300 | U | 340 |
| Chrysene | 1,000 | 3,900 | 710 | 260 | | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Dibenz(a,h)anthracene | 330 | 330 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Dibenzofuran | 7,000 | 59,000 | 990 | 260 | | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| Diethyl phthalate | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Dimethylphthalate | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Di-n-butylphthalate | | | < 260 | 260 | U | 100 | < 280 | 280 | U | 100 | 130 | 320 | JB | 120 |
| Di-n-octylphthalate | | | < 260 | 260 | U | 97 | < 280 | 280 | U | 100 | < 320 | 320 | U | 120 |
| Fluoranthene | 100,000 | 100,000 | 3,200 | 260 | | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Fluorene | 30,000 | 100,000 | 1,400 | 260 | | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Hexachlorobenzene | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 130 |
| Hexachlorobutadiene | | | < 260 | 260 | U | 140 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| Hexachlorocyclopentadiene | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Hexachloroethane | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 | < 320 | 320 | U | 140 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 680 | 260 | | 130 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| Isophorone | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Naphthalene | 12,000 | 100,000 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Nitrobenzene | | | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| N-Nitrosodimethylamine | | | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| N-Nitrosodi-n-propylamine | | | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 150 |
| N-Nitrosodiphenylamine | | | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 | < 320 | 320 | U | 170 |
| Pentachloronitrobenzene | | | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 | < 320 | 320 | U | 170 |
| Pentachlorophenol | 800 | 6,700 | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 | < 320 | 320 | U | 170 |
| Phenanthrene | 100,000 | 100,000 | 5,400 | 260 | | 110 | < 280 | 280 | U | 110 | < 320 | 320 | U | 130 |
| Phenol | 330 | 100,000 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 | < 320 | 320 | U | 140 |
| Pyrene | 100,000 | 100,000 | 2,800 | 260 | | 130 | < 280 | 280 | U | 140 | < 320 | 320 | U | 160 |
| Pyridine | | | < 260 | 260 | U | 93 | < 280 | 280 | U | 97 | < 320 | 320 | U | 110 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B3 | | | | | | | | | | | |
|-------------------------------|--|--|----------------|-------|------|-----|----------------|-------|------|-----|------------------|-------|------|-----|
| | | | 11/24/2015 | | | | | | | | | | | |
| | | | (0-2) µg/Kg | | | | (5-7) µg/Kg | | | | (13-15) µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| 1,2,4-Trichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| 1,2-Dichlorobenzene | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| 1,2-Diphenylhydrazine | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| 1,3-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| 1,4-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| 2,4,5-Trichlorophenol | | | < 250 | 250 | U | 200 | < 280 | 280 | U | 220 | < 260 | 260 | U | 200 |
| 2,4,6-Trichlorophenol | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| 2,4-Dichlorophenol | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| 2,4-Dimethylphenol | | | < 250 | 250 | U | 89 | < 280 | 280 | U | 98 | < 260 | 260 | U | 91 |
| 2,4-Dinitrophenol | | | < 720 | 720 | U | 250 | < 790 | 790 | U | 280 | < 740 | 740 | U | 260 |
| 2,4-Dinitrotoluene | | | < 250 | 250 | U | 140 | < 280 | 280 | U | 160 | < 260 | 260 | U | 140 |
| 2,6-Dinitrotoluene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| 2-Chloronaphthalene | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| 2-Chlorophenol | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| 2-Methylnaphthalene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 250 | 250 | U | 170 | < 280 | 280 | U | 190 | < 260 | 260 | U | 170 |
| 2-Nitroaniline | | | < 720 | 720 | U | 360 | < 790 | 790 | U | 400 | < 740 | 740 | U | 370 |
| 2-Nitrophenol | | | < 250 | 250 | U | 230 | < 280 | 280 | U | 250 | < 260 | 260 | U | 230 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 250 | 250 | U | 140 | < 280 | 280 | U | 160 | < 260 | 260 | U | 140 |
| 3,3'-Dichlorobenzidine | | | < 720 | 720 | U | 170 | < 790 | 790 | U | 190 | < 740 | 740 | U | 170 |
| 3-Nitroaniline | | | < 720 | 720 | U | 720 | < 790 | 790 | U | 790 | < 740 | 740 | U | 740 |
| 4,6-Dinitro-2-methylphenol | | | < 1800 | 1,800 | U | 380 | < 2000 | 2,000 | U | 430 | < 1800 | 1,800 | U | 400 |
| 4-Bromophenyl phenyl ether | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| 4-Chloro-3-methylphenol | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| 4-Chloroaniline | | | < 290 | 290 | U | 170 | < 320 | 320 | U | 180 | < 290 | 290 | U | 170 |
| 4-Chlorophenyl phenyl ether | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| 4-Nitroaniline | | | < 720 | 720 | U | 120 | < 790 | 790 | U | 130 | < 740 | 740 | U | 120 |
| 4-Nitrophenol | | | < 360 | 360 | U | 160 | < 400 | 400 | U | 180 | < 370 | 370 | U | 170 |
| Acenaphthene | 20,000 | 100,000 | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Acenaphthylene | 100,000 | 100,000 | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| Acetophenone | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Aniline | | | < 290 | 290 | U | 290 | < 320 | 320 | U | 320 | < 290 | 290 | U | 290 |
| Anthracene | 100,000 | 100,000 | 220 | 250 | J | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Benz(a)anthracene | 1,000 | 1,000 | 480 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Benzbzidine | | | < 720 | 720 | U | 210 | < 790 | 790 | U | 230 | < 740 | 740 | U | 220 |
| Benzo(a)pyrene | 1,000 | 1,000 | 400 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 330 | 250 | | 120 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 220 | 250 | J | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Benzo(k)fluoranthene | 800 | 3,900 | 330 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Benzoic acid | | | < 1800 | 1,800 | U | 720 | < 2000 | 2,000 | U | 790 | < 1800 | 1,800 | U | 740 |
| Benzyl butyl phthalate | | | < 250 | 250 | U | 92 | < 280 | 280 | U | 100 | < 260 | 260 | U | 95 |
| Bis(2-chloroethoxy)methane | | | < 250 | 250 | U | 99 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-chloroethyl)ether | | | < 250 | 250 | U | 97 | < 280 | 280 | U | 110 | < 260 | 260 | U | 99 |
| Bis(2-chloroisopropyl)ether | | | < 250 | 250 | U | 99 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-ethylhexyl)phthalate | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 |
| Carbazole | | | < 1800 | 1,800 | U | 270 | < 2000 | 2,000 | U | 300 | < 1800 | 1,800 | U | 280 |
| Chrysene | 1,000 | 3,900 | 470 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Dibenz(a,h)anthracene | 330 | 330 | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Dibenzofuran | 7,000 | 59,000 | < 250 | 250 | U | 100 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Diethyl phthalate | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Dimethylphthalate | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Di-n-butylphthalate | | | 110 | 250 | JB | 95 | 110 | 280 | JB | 110 | < 260 | 260 | U | 98 |
| Di-n-octylphthalate | | | < 250 | 250 | U | 92 | < 280 | 280 | U | 100 | < 260 | 260 | U | 95 |
| Fluoranthene | 100,000 | 100,000 | 1,100 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Fluorene | 30,000 | 100,000 | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Hexachlorobenzene | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Hexachlorobutadiene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| Hexachlorocyclopentadiene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Hexachloroethane | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 230 | 250 | J | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Isophorone | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| Naphthalene | 12,000 | 100,000 | < 250 | 250 | U | 100 | 190 | 280 | J | 110 | < 260 | 260 | U | 110 |
| Nitrobenzene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| N-Nitrosodimethylamine | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 |
| N-Nitrosodi-n-propylamine | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| N-Nitrosodiphenylamine | | | < 250 | 250 | U | 140 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 |
| Pentachloronitrobenzene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 |
| Pentachlorophenol | 800 | 6,700 | < 250 | 250 | U | 140 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 |
| Phenanthrene | 100,000 | 100,000 | 790 | 250 | | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 |
| Phenol | 330 | 100,000 | < 250 | 250 | U | 110 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 |
| Pyrene | 100,000 | 100,000 | 970 | 250 | | 120 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 |
| Pyridine | | | < 250 | 250 | U | 88 | < 280 | 280 | U | 97 | < 260 | 260 | U | 91 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B4 | | | | | | | | | | | | 15B5 | | | |
|-------------------------------|--|--|-----------------|--------|----|------|-----------------|--------|----|------|-------------------|--------|----|------|-----------------|-------|---|-----|
| | | | 11/24/2015 | | | | | | | | | | | | 11/24/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (10-12') µg/Kg | | | | (6-8') µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,2,4,5-Tetrachlorobenzene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| 1,2,4-Trichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| 1,2-Dichlorobenzene | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 |
| 1,2-Diphenylhydrazine | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| 1,3-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| 1,4-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| 2,4,5-Trichlorophenol | | | < 250 | 250 | U | 200 | < 280 | 280 | U | 220 | < 260 | 260 | U | 200 | < 280 | 280 | U | 220 |
| 2,4,6-Trichlorophenol | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| 2,4-Dichlorophenol | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| 2,4-Dimethylphenol | | | < 250 | 250 | U | 90 | < 280 | 280 | U | 98 | < 260 | 260 | U | 93 | < 280 | 280 | U | 99 |
| 2,4-Dinitrophenol | | | < 730 | 730 | U | 250 | < 790 | 790 | U | 280 | < 750 | 750 | U | 260 | < 800 | 800 | U | 280 |
| 2,4-Dinitrotoluene | | | < 250 | 250 | U | 140 | < 280 | 280 | U | 160 | < 260 | 260 | U | 150 | < 280 | 280 | U | 160 |
| 2,6-Dinitrotoluene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| 2-Chloronaphthalene | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 |
| 2-Chlorophenol | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 |
| 2-Methylnaphthalene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 250 | 250 | U | 170 | < 280 | 280 | U | 180 | < 260 | 260 | U | 180 | < 280 | 280 | U | 190 |
| 2-Nitroaniline | | | < 730 | 730 | U | 370 | < 790 | 790 | U | 400 | < 750 | 750 | U | 380 | < 800 | 800 | U | 400 |
| 2-Nitrophenol | | | < 250 | 250 | U | 230 | < 280 | 280 | U | 250 | < 260 | 260 | U | 240 | < 280 | 280 | U | 250 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 250 | 250 | U | 140 | < 280 | 280 | U | 160 | < 260 | 260 | U | 150 | < 280 | 280 | U | 160 |
| 3,3'-Dichlorobenzidine | | | < 730 | 730 | U | 170 | < 790 | 790 | U | 190 | < 750 | 750 | U | 180 | < 800 | 800 | U | 190 |
| 3-Nitroaniline | | | < 730 | 730 | U | 730 | < 790 | 790 | U | 790 | < 750 | 750 | U | 750 | < 800 | 800 | U | 800 |
| 4,6-Dinitro-2-methylphenol | | | < 1800 | 1,800 | U | 390 | < 2000 | 2,000 | U | 420 | < 1900 | 1,900 | U | 400 | < 2000 | 2,000 | U | 430 |
| 4-Bromophenyl phenyl ether | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| 4-Chloro-3-methylphenol | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| 4-Chloroaniline | | | < 290 | 290 | U | 170 | < 310 | 310 | U | 180 | < 300 | 300 | U | 170 | < 320 | 320 | U | 190 |
| 4-Chlorophenyl phenyl ether | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 130 | < 280 | 280 | U | 130 |
| 4-Nitroaniline | | | < 730 | 730 | U | 120 | < 790 | 790 | U | 130 | < 750 | 750 | U | 120 | < 800 | 800 | U | 130 |
| 4-Nitrophenol | | | < 360 | 360 | U | 160 | < 390 | 390 | U | 180 | < 370 | 370 | U | 170 | < 400 | 400 | U | 180 |
| Acenaphthene | 20,000 | 100,000 | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Acenaphthylene | 100,000 | 100,000 | 110 | 250 | J | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 |
| Acetophenone | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 |
| Aniline | | | < 290 | 290 | U | 290 | < 310 | 310 | U | 310 | < 300 | 300 | U | 300 | < 320 | 320 | U | 320 |
| Anthracene | 100,000 | 100,000 | 450 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Benz(a)anthracene | 1,000 | 1,000 | 1,600 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 130 | < 280 | 280 | U | 130 |
| Benzidine | | | < 730 | 730 | U | 210 | < 790 | 790 | U | 230 | < 750 | 750 | U | 220 | < 800 | 800 | U | 240 |
| Benzo(a)pyrene | 1,000 | 1,000 | 1,300 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 1,100 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 700 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Benzo(k)fluoranthene | 800 | 3,900 | 1,100 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Benzoic acid | | | < 1800 | 1,800 | U | 730 | < 2000 | 2,000 | U | 790 | < 1900 | 1,900 | U | 750 | < 2000 | 2,000 | U | 800 |
| Benzyl butyl phthalate | | | < 250 | 250 | U | 94 | < 280 | 280 | U | 100 | < 260 | 260 | U | 96 | < 280 | 280 | U | 100 |
| Bis(2-chloroethoxy)methane | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 |
| Bis(2-chloroethyl)ether | | | < 250 | 250 | U | 98 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 |
| Bis(2-chloroisopropyl)ether | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 |
| Bis(2-ethylhexyl)phthalate | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Carbazole | | | < 1800 | 1,800 | U | 280 | < 2000 | 2,000 | U | 300 | < 1900 | 1,900 | U | 280 | < 2000 | 2,000 | U | 300 |
| Chrysene | 1,000 | 3,900 | 1,500 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 130 | < 280 | 280 | U | 130 |
| Dibenz(a,h)anthracene | 330 | 330 | 180 | 250 | J | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Dibenzofuran | 7,000 | 59,000 | < 250 | 250 | U | 110 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Diethyl phthalate | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Dimethylphthalate | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 120 | < 280 | 280 | U | 120 |
| Di-n-butylphthalate | | | < 250 | 250 | U | 97 | < 280 | 280 | U | 100 | 110 | 260 | JB | 99 | < 280 | 280 | U | 110 |
| Di-n-octylphthalate | | | < 250 | 250 | U | 94 | < 280 | 280 | U | 100 | < 260 | 260 | U | 96 | < 280 | 280 | U | 100 |
| Fluoranthene | 100,000 | 100,000 | 3,100 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Fluorene | 30,000 | 100,000 | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Hexachlorobenzene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Hexachlorobutadiene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 140 | < 280 | 280 | U | 140 |
| Hexachlorocyclopentadiene | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Hexachloroethane | | | < 250 | 250 | U | 110 | < 280 | 280 | U | 120 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 820 | 250 | | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Isophorone | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 100 | < 280 | 280 | U | 110 |
| Naphthalene | 12,000 | 100,000 | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 120 |
| Nitrobenzene | | | < 250 | 250 | U | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| N-Nitrosodimethylamine | | | < 250 | 250 | U | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 |
| N-Nitrosodi-n-propylamine | | | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| N-Nitrosodiphenylamine | | | < 250 | 250 | U | 140 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 |
| Pentachloronitrobenzene | | | < 250 | 250 | U | 140 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 |
| Pentachlorophenol | 800 | 6,700 | < 250 | 250 | U | 140 | < 280 | 280 | U | 150 | < 260 | 260 | U | 140 | < 280 | 280 | U | 150 |
| Phenanthrene | 100,000 | 100,000 | 1,400 | 250 | | 100 | < 280 | 280 | U | 110 | < 260 | 260 | U | 110 | < 280 | 280 | U | 110 |
| Phenol | 330 | 100,000 | < 250 | 250 | U | 120 | < 280 | 280 | U | 130 | < 260 | 260 | U | 120 | < 280 | 280 | U | 130 |
| Pyrene | 100,000 | 100,000 | 2,700 | 250 | | 130 | < 280 | 280 | U | 140 | < 260 | 260 | U | 130 | < 280 | 280 | U | 140 |
| Pyridine | | | < 250 | 250 | U | 90 | < 280 | 280 | U | 97 | < 260 | 260 | U | 92 | < 280 | 280 | U | 99 |

Notes:

RL - Laboratory Reporting Limit, MDL-Minimum Detection Limit
RL - Laboratory Reporting Limit

Bold/highlighted - Indicated exceedance of the NYSDEC UUSCO Guidance Value

**TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results**

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B6 | | | | | | | | 15SB7 | | | | 15B8 | | | |
|-------------------------------|--|--|-----------------|-------|------|-----|-------------------|-------|------|------|-----------------|-------|------|-----|-----------------|-------|------|-----|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (3-5') µg/Kg | | | | (10-12') µg/Kg | | | | (6-8') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 270 | 270 | U | 130 | < 370 | 370 | U | 190 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| 1,2,4-Trichlorobenzene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 150 |
| 1,2-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 140 |
| 1,2-Diphenylhydrazine | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| 1,3-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 1,4-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 2,4,5-Trichlorophenol | | | < 270 | 270 | U | 210 | < 370 | 370 | U | 290 | < 290 | 290 | U | 220 | < 340 | 340 | U | 270 |
| 2,4,6-Trichlorophenol | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| 2,4-Dichlorophenol | | | < 270 | 270 | U | 130 | < 370 | 370 | U | 190 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| 2,4-Dimethylphenol | | | < 270 | 270 | U | 94 | < 370 | 370 | U | 130 | < 290 | 290 | U | 100 | < 340 | 340 | U | 120 |
| 2,4-Dinitrophenol | | | < 760 | 760 | U | 270 | < 1100 | 1,100 | U | 370 | < 810 | 810 | U | 290 | < 970 | 970 | U | 340 |
| 2,4-Dinitrotoluene | | | < 270 | 270 | U | 150 | < 370 | 370 | U | 210 | < 290 | 290 | U | 160 | < 340 | 340 | U | 190 |
| 2,6-Dinitrotoluene | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 150 |
| 2-Chloronaphthalene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 2-Chlorophenol | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 2-Methylnaphthalene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 270 | 270 | U | 180 | < 330 | 330 | U | 250 | < 290 | 290 | U | 190 | < 330 | 330 | U | 230 |
| 2-Nitroaniline | | | < 760 | 760 | U | 380 | < 1100 | 1,100 | U | 530 | < 810 | 810 | U | 410 | < 970 | 970 | U | 490 |
| 2-Nitrophenol | | | < 270 | 270 | U | 240 | < 370 | 370 | U | 340 | < 290 | 290 | U | 260 | < 340 | 340 | U | 310 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 270 | 270 | U | 150 | < 370 | 370 | U | 210 | < 290 | 290 | U | 160 | < 340 | 340 | U | 190 |
| 3,3'-Dichlorobenzidine | | | < 760 | 760 | U | 180 | < 1100 | 1,100 | U | 250 | < 810 | 810 | U | 190 | < 970 | 970 | U | 230 |
| 3-Nitroaniline | | | < 760 | 760 | U | 760 | < 1100 | 1,100 | U | 1100 | < 810 | 810 | U | 810 | < 970 | 970 | U | 970 |
| 4,6-Dinitro-2-methylphenol | | | < 1900 | 1,900 | U | 410 | < 2600 | 2,600 | U | 570 | < 2000 | 2,000 | U | 440 | < 2400 | 2,400 | U | 520 |
| 4-Bromophenyl phenyl ether | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| 4-Chloro-3-methylphenol | | | < 270 | 270 | U | 130 | < 370 | 370 | U | 190 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| 4-Chloroaniline | | | < 300 | 300 | U | 180 | < 420 | 420 | U | 250 | < 330 | 330 | U | 190 | < 390 | 390 | U | 230 |
| 4-Chlorophenyl phenyl ether | | | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 160 |
| 4-Nitroaniline | | | < 760 | 760 | U | 130 | < 1100 | 1,100 | U | 180 | < 810 | 810 | U | 140 | < 970 | 970 | U | 160 |
| 4-Nitrophenol | | | < 380 | 380 | U | 170 | < 530 | 530 | U | 240 | < 410 | 410 | U | 180 | < 490 | 490 | U | 220 |
| Acenaphthene | 20,000 | 100,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 150 |
| Acenaphthylene | 100,000 | 100,000 | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 140 |
| Acetophenone | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 150 |
| Aniline | | | < 300 | 300 | U | 300 | < 420 | 420 | U | 420 | < 330 | 330 | U | 330 | < 390 | 390 | U | 390 |
| Anthracene | 100,000 | 100,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| Benz(a)anthracene | 1,000 | 1,000 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 160 |
| Benzidine | | | < 760 | 760 | U | 220 | < 1100 | 1,100 | U | 310 | < 810 | 810 | U | 240 | < 970 | 970 | U | 290 |
| Benzo(a)pyrene | 1,000 | 1,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| Benzo(ghi)perylene | 100,000 | 100,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| Benzo(k)fluoranthene | 800 | 3,900 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 160 |
| Benzoic acid | | | < 1900 | 1,900 | U | 760 | < 2600 | 2,600 | U | 1100 | < 2000 | 2,000 | U | 810 | < 2400 | 2,400 | U | 970 |
| Benzyl butyl phthalate | | | < 270 | 270 | U | 98 | < 370 | 370 | U | 140 | < 290 | 290 | U | 110 | < 340 | 340 | U | 130 |
| Bis(2-chloroethoxy)methane | | | < 270 | 270 | U | 100 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 130 |
| Bis(2-chloroethyl)ether | | | < 270 | 270 | U | 100 | < 370 | 370 | U | 140 | < 290 | 290 | U | 110 | < 340 | 340 | U | 130 |
| Bis(2-chloroisopropyl)ether | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 140 |
| Bis(2-ethylhexyl)phthalate | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| Carbazole | | | < 1900 | 1,900 | U | 290 | < 2600 | 2,600 | U | 400 | < 2000 | 2,000 | U | 310 | < 2400 | 2,400 | U | 370 |
| Chrysene | 1,000 | 3,900 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 160 |
| Dibenz(a,h)anthracene | 330 | 330 | < 270 | 270 | U | 120 | < 330 | 330 | U | 170 | < 290 | 290 | U | 130 | < 330 | 330 | U | 160 |
| Dibenzofuran | 7,000 | 59,000 | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| Diethyl phthalate | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 150 |
| Dimethylphthalate | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 160 | < 290 | 290 | U | 130 | < 340 | 340 | U | 150 |
| Di-n-butylphthalate | | | < 270 | 270 | U | 100 | < 370 | 370 | U | 140 | < 290 | 290 | U | 110 | < 340 | 340 | U | 130 |
| Di-n-octylphthalate | | | < 270 | 270 | U | 98 | < 370 | 370 | U | 140 | < 290 | 290 | U | 110 | < 340 | 340 | U | 130 |
| Fluoranthene | 100,000 | 100,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| Fluorene | 30,000 | 100,000 | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| Hexachlorobenzene | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| Hexachlorobutadiene | | | < 270 | 270 | U | 140 | < 370 | 370 | U | 190 | < 290 | 290 | U | 150 | < 340 | 340 | U | 180 |
| Hexachlorocyclopentadiene | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 150 |
| Hexachloroethane | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 160 | < 290 | 290 | U | 120 | < 340 | 340 | U | 150 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 160 |
| Isophorone | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 140 |
| Naphthalene | 12,000 | 100,000 | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| Nitrobenzene | | | < 270 | 270 | U | 130 | < 370 | 370 | U | 190 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| N-Nitrosodimethylamine | | | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 110 | < 340 | 340 | U | 140 |
| N-Nitrosodi-n-propylamine | | | < 270 | 270 | U | 120 | < 370 | 370 | U | 170 | < 290 | 290 | U | 130 | < 340 | 340 | U | 160 |
| N-Nitrosodiphenylamine | | | < 270 | 270 | U | 150 | < 370 | 370 | U | 200 | < 290 | 290 | U | 160 | < 340 | 340 | U | 190 |
| Pentachloronitrobenzene | | | < 270 | 270 | U | 140 | < 370 | 370 | U | 200 | < 290 | 290 | U | 150 | < 340 | 340 | U | 180 |
| Pentachlorophenol | 800 | 6,700 | < 270 | 270 | U | 140 | < 370 | 370 | U | 200 | < 290 | 290 | U | 150 | < 340 | 340 | U | 180 |
| Phenanthrene | 100,000 | 100,000 | < 270 | 270 | U | 110 | < 370 | 370 | U | 150 | < 290 | 290 | U | 120 | < 340 | 340 | U | 140 |
| Phenol | 330 | 100,000 | < 270 | 270 | U | 120 | < 330 | 330 | U | 170 | < 290 | 290 | U | 130 | < 330 | 330 | U | 160 |
| Pyrene | 100,000 | 100,000 | < 270 | 270 | U | 130 | < 370 | 370 | U | 180 | < 290 | 290 | U | 140 | < 340 | 340 | U | 170 |
| Pyridine | | | < 270 | 270 | U | 93 | < 370 | 370 | U | 130 | < 290 | 290 | U | 100 | < 340 | 340 | U | 120 |

Notes:

RL - Laboratory Reporting Limit

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B9 | | | | | | | | 15B10 | | | |
|-------------------------------|--|--|-----------------|--------|----|------|-----------------|--------|----|------|-----------------|-------|---|-----|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (6-8') µg/Kg | | | |
| Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 1,2,4,5-Tetrachlorobenzene | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 130 |
| 1,2,4-Trichlorobenzene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 110 |
| 1,2-Dichlorobenzene | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 1,2-Diphenylhydrazine | | | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| 1,3-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 1,4-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 2,4,5-Trichlorophenol | | | < 250 | 250 | U | 200 | < 290 | 290 | U | 230 | < 260 | 260 | U | 200 |
| 2,4,6-Trichlorophenol | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| 2,4-Dichlorophenol | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 130 |
| 2,4-Dimethylphenol | | | < 250 | 250 | U | 89 | < 290 | 290 | U | 100 | < 260 | 260 | U | 92 |
| 2,4-Dinitrophenol | | | < 710 | 710 | U | 250 | < 830 | 830 | U | 290 | < 750 | 750 | U | 260 |
| 2,4-Dinitrotoluene | | | < 250 | 250 | U | 140 | < 290 | 290 | U | 160 | < 260 | 260 | U | 150 |
| 2,6-Dinitrotoluene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| 2-Chloronaphthalene | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 2-Chlorophenol | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 2-Methylnaphthalene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 250 | 250 | U | 170 | < 290 | 290 | U | 200 | < 260 | 260 | U | 180 |
| 2-Nitroaniline | | | < 710 | 710 | U | 360 | < 830 | 830 | U | 420 | < 750 | 750 | U | 380 |
| 2-Nitrophenol | | | < 250 | 250 | U | 230 | < 290 | 290 | U | 260 | < 260 | 260 | U | 240 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 250 | 250 | U | 140 | < 290 | 290 | U | 160 | < 260 | 260 | U | 150 |
| 3,3'-Dichlorobenzidine | | | < 710 | 710 | U | 170 | < 830 | 830 | U | 200 | < 750 | 750 | U | 180 |
| 3-Nitroaniline | | | < 710 | 710 | U | 710 | < 830 | 830 | U | 830 | < 750 | 750 | U | 750 |
| 4,6-Dinitro-2-methylphenol | | | < 1800 | 1,800 | U | 380 | < 2100 | 2,100 | U | 450 | < 1900 | 1,900 | U | 400 |
| 4-Bromophenyl phenyl ether | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| 4-Chloro-3-methylphenol | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 130 |
| 4-Chloroaniline | | | < 290 | 290 | U | 170 | < 330 | 330 | U | 190 | < 300 | 300 | U | 170 |
| 4-Chlorophenyl phenyl ether | | | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 130 |
| 4-Nitroaniline | | | < 710 | 710 | U | 120 | < 830 | 830 | U | 140 | < 750 | 750 | U | 120 |
| 4-Nitrophenol | | | < 360 | 360 | U | 160 | < 420 | 420 | U | 190 | < 370 | 370 | U | 170 |
| Acenaphthene | 20,000 | 100,000 | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 110 |
| Acenaphthylene | 100,000 | 100,000 | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 100 |
| Acetophenone | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Aniline | | | < 290 | 290 | U | 290 | < 330 | 330 | U | 330 | < 300 | 300 | U | 300 |
| Anthracene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| Benz(a)anthracene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 130 |
| Benidine | | | < 710 | 710 | U | 210 | < 830 | 830 | U | 240 | < 750 | 750 | U | 220 |
| Benzo(a)pyrene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 130 |
| Benzo(ghi)perylene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Benzo(k)fluoranthene | 800 | 3,900 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| Benzoic acid | | | < 1800 | 1,800 | U | 710 | < 2100 | 2,100 | U | 830 | < 1900 | 1,900 | U | 750 |
| Benzyl butyl phthalate | | | < 250 | 250 | U | 92 | < 290 | 290 | U | 110 | < 260 | 260 | U | 96 |
| Bis(2-chloroethoxy)methane | | | < 250 | 250 | U | 99 | < 290 | 290 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-chloroethyl)ether | | | < 250 | 250 | U | 96 | < 290 | 290 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-chloroisopropyl)ether | | | < 250 | 250 | U | 99 | < 290 | 290 | U | 120 | < 260 | 260 | U | 100 |
| Bis(2-ethylhexyl)phthalate | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Carbazole | | | < 1800 | 1,800 | U | 270 | < 2100 | 2,100 | U | 320 | < 1900 | 1,900 | U | 280 |
| Chrysene | 1,000 | 3,900 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 130 |
| Dibenz(a,h)anthracene | 330 | 330 | < 250 | 250 | U | 120 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Dibenzofuran | 7,000 | 59,000 | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Diethyl phthalate | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Dimethylphthalate | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Di-n-butylphthalate | | | < 250 | 250 | U | 95 | < 290 | 290 | U | 110 | < 260 | 260 | U | 99 |
| Di-n-octylphthalate | | | < 250 | 250 | U | 92 | < 290 | 290 | U | 110 | < 260 | 260 | U | 96 |
| Fluoranthene | 100,000 | 100,000 | 130 | 250 | J | 120 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Fluorene | 30,000 | 100,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| Hexachlorobenzene | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Hexachlorobutadiene | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 130 |
| Hexachlorocyclopentadiene | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 110 |
| Hexachloroethane | | | < 250 | 250 | U | 110 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 120 |
| Isophorone | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 100 |
| Naphthalene | 12,000 | 100,000 | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Nitrobenzene | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 130 |
| N-Nitrosodimethylamine | | | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| N-Nitrosodi-n-propylamine | | | < 250 | 250 | U | 120 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| N-Nitrosodiphenylamine | | | < 250 | 250 | U | 140 | < 290 | 290 | U | 160 | < 260 | 260 | U | 140 |
| Pentachloronitrobenzene | | | < 250 | 250 | U | 130 | < 290 | 290 | U | 150 | < 260 | 260 | U | 140 |
| Pentachlorophenol | 800 | 6,700 | < 250 | 250 | U | 140 | < 290 | 290 | U | 160 | < 260 | 260 | U | 140 |
| Phenanthrene | 100,000 | 100,000 | < 250 | 250 | U | 100 | < 290 | 290 | U | 120 | < 260 | 260 | U | 110 |
| Phenol | 330 | 100,000 | < 250 | 250 | U | 110 | < 290 | 290 | U | 130 | < 260 | 260 | U | 120 |
| Pyrene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 290 | 290 | U | 140 | < 260 | 260 | U | 130 |
| Pyridine | | | < 250 | 250 | U | 88 | < 290 | 290 | U | 100 | < 260 | 260 | U | 92 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B11 | | | | | | | | 15B12 | | | |
|-------------------------------|--|--|-----------------|-------|------|-----|-----------------|-------|------|-----|-----------------|-------|------|-----|
| | | | 11/23/2015 | | | | | | | | 11/23/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 260 | 260 | U | 130 | < 270 | 270 | U | 140 | < 260 | 260 | U | 130 |
| 1,2,4-Trichlorobenzene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| 1,2-Dichlorobenzene | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| 1,2-Diphenylhydrazine | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| 1,3-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| 1,4-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| 2,4,5-Trichlorophenol | | | < 260 | 260 | U | 200 | < 270 | 270 | U | 210 | < 260 | 260 | U | 200 |
| 2,4,6-Trichlorophenol | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| 2,4-Dichlorophenol | | | < 260 | 260 | U | 130 | < 270 | 270 | U | 140 | < 260 | 260 | U | 130 |
| 2,4-Dimethylphenol | | | < 260 | 260 | U | 92 | < 270 | 270 | U | 97 | < 260 | 260 | U | 92 |
| 2,4-Dinitrophenol | | | < 740 | 740 | U | 260 | < 780 | 780 | U | 270 | < 740 | 740 | U | 260 |
| 2,4-Dinitrotoluene | | | < 260 | 260 | U | 150 | < 270 | 270 | U | 150 | < 260 | 260 | U | 150 |
| 2,6-Dinitrotoluene | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| 2-Chloronaphthalene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| 2-Chlorophenol | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| 2-Methylnaphthalene | | | 160 | 260 | J | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 260 | 260 | U | 170 | < 270 | 270 | U | 180 | < 260 | 260 | U | 170 |
| 2-Nitroaniline | | | < 740 | 740 | U | 370 | < 780 | 780 | U | 390 | < 740 | 740 | U | 380 |
| 2-Nitrophenol | | | < 260 | 260 | U | 230 | < 270 | 270 | U | 250 | < 260 | 260 | U | 240 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 260 | 260 | U | 150 | < 270 | 270 | U | 150 | < 260 | 260 | U | 150 |
| 3,3'-Dichlorobenzidine | | | < 740 | 740 | U | 170 | < 780 | 780 | U | 180 | < 740 | 740 | U | 180 |
| 3-Nitroaniline | | | < 740 | 740 | U | 740 | < 780 | 780 | U | 780 | < 740 | 740 | U | 740 |
| 4,6-Dinitro-2-methylphenol | | | < 1900 | 1,900 | U | 400 | < 1900 | 1,900 | U | 420 | < 1900 | 1,900 | U | 400 |
| 4-Bromophenyl phenyl ether | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| 4-Chloro-3-methylphenol | | | < 260 | 260 | U | 130 | < 270 | 270 | U | 140 | < 260 | 260 | U | 130 |
| 4-Chloroaniline | | | < 300 | 300 | U | 170 | < 310 | 310 | U | 180 | < 300 | 300 | U | 170 |
| 4-Chlorophenyl phenyl ether | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 130 |
| 4-Nitroaniline | | | < 740 | 740 | U | 120 | < 780 | 780 | U | 130 | < 740 | 740 | U | 120 |
| 4-Nitrophenol | | | < 370 | 370 | U | 170 | < 390 | 390 | U | 180 | < 370 | 370 | U | 170 |
| Acenaphthene | 20,000 | 100,000 | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| Acenaphthylene | 100,000 | 100,000 | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| Acetophenone | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| Aniline | | | < 300 | 300 | U | 300 | < 310 | 310 | U | 310 | < 300 | 300 | U | 300 |
| Anthracene | 100,000 | 100,000 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Benz(a)anthracene | 1,000 | 1,000 | 140 | 260 | J | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 130 |
| Benztidine | | | < 740 | 740 | U | 220 | < 780 | 780 | U | 230 | < 740 | 740 | U | 220 |
| Benzo(a)pyrene | 1,000 | 1,000 | 130 | 260 | J | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 130 | 260 | J | 130 | < 270 | 270 | U | 130 | < 260 | 260 | U | 130 |
| Benzo(ghi)perylene | 100,000 | 100,000 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Benzo(k)fluoranthene | 800 | 3,900 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Benzoic acid | | | < 1900 | 1,900 | U | 740 | < 1900 | 1,900 | U | 780 | < 1900 | 1,900 | U | 740 |
| Benzyl butyl phthalate | | | < 260 | 260 | U | 96 | < 270 | 270 | U | 100 | < 260 | 260 | U | 96 |
| Bis(2-chloroethoxy)methane | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-chloroethyl)ether | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-chloroisopropyl)ether | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| Bis(2-ethylhexyl)phthalate | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| Carbazole | | | < 1900 | 1,900 | U | 280 | < 1900 | 1,900 | U | 300 | < 1900 | 1,900 | U | 280 |
| Chrysene | 1,000 | 3,900 | 180 | 260 | J | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 130 |
| Dibenz(a,h)anthracene | 330 | 330 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Dibenzofuran | 7,000 | 59,000 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| Diethyl phthalate | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| Dimethylphthalate | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| Di-n-butylphthalate | | | < 260 | 260 | U | 99 | < 270 | 270 | U | 100 | < 260 | 260 | U | 99 |
| Di-n-octylphthalate | | | < 260 | 260 | U | 96 | < 270 | 270 | U | 100 | < 260 | 260 | U | 96 |
| Fluoranthene | 100,000 | 100,000 | 280 | 260 | | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Fluorene | 30,000 | 100,000 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Hexachlorobenzene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| Hexachlorobutadiene | | | < 260 | 260 | U | 130 | < 270 | 270 | U | 140 | < 260 | 260 | U | 130 |
| Hexachlorocyclopentadiene | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| Hexachloroethane | | | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 260 | 260 | U | 110 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| Isophorone | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| Naphthalene | 12,000 | 100,000 | 170 | 260 | J | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| Nitrobenzene | | | < 260 | 260 | U | 130 | < 270 | 270 | U | 140 | < 260 | 260 | U | 130 |
| N-Nitrosodimethylamine | | | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 260 | 260 | U | 100 |
| N-Nitrosodi-n-propylamine | | | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 260 | 260 | U | 120 |
| N-Nitrosodiphenylamine | | | < 260 | 260 | U | 140 | < 270 | 270 | U | 150 | < 260 | 260 | U | 140 |
| Pentachloronitrobenzene | | | < 260 | 260 | U | 140 | < 270 | 270 | U | 150 | < 260 | 260 | U | 140 |
| Pentachlorophenol | 800 | 6,700 | < 260 | 260 | U | 140 | < 270 | 270 | U | 150 | < 260 | 260 | U | 140 |
| Phenanthrene | 100,000 | 100,000 | 220 | 260 | J | 110 | < 270 | 270 | U | 110 | < 260 | 260 | U | 110 |
| Phenol | 330 | 100,000 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 260 | 260 | U | 120 |
| Pyrene | 100,000 | 100,000 | 220 | 260 | J | 130 | < 270 | 270 | U | 130 | < 260 | 260 | U | 130 |
| Pyridine | | | < 260 | 260 | U | 91 | < 270 | 270 | U | 96 | < 260 | 260 | U | 92 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B13 | | | | | | | | 15B14 | | | | 15B15 | | | |
|-------------------------------|--|--|-----------------|-------|------|-----|-----------------|-------|------|-----|-----------------|-------|------|-----|-----------------|-------|------|-----|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (5-7') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 140 | < 280 | 280 | U | 140 | < 280 | 280 | U | 140 |
| 1,2,4-Trichlorobenzene | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| 1,2-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| 1,2-Diphenylhydrazine | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| 1,3-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| 1,4-Dichlorobenzene | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| 2,4,5-Trichlorophenol | | | < 270 | 270 | U | 210 | < 270 | 270 | U | 210 | < 280 | 280 | U | 220 | < 280 | 280 | U | 220 |
| 2,4,6-Trichlorophenol | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| 2,4-Dichlorophenol | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 140 | < 280 | 280 | U | 140 | < 280 | 280 | U | 140 |
| 2,4-Dimethylphenol | | | < 270 | 270 | U | 95 | < 270 | 270 | U | 97 | < 280 | 280 | U | 98 | < 280 | 280 | U | 99 |
| 2,4-Dinitrophenol | | | < 770 | 770 | U | 270 | < 780 | 780 | U | 270 | < 790 | 790 | U | 280 | < 800 | 800 | U | 280 |
| 2,4-Dinitrotoluene | | | < 270 | 270 | U | 150 | < 270 | 270 | U | 150 | < 280 | 280 | U | 150 | < 280 | 280 | U | 160 |
| 2,6-Dinitrotoluene | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 130 |
| 2-Chloronaphthalene | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| 2-Chlorophenol | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| 2-Methylnaphthalene | | | 360 | 270 | | 110 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 270 | 270 | U | 180 | < 270 | 270 | U | 180 | < 280 | 280 | U | 180 | < 280 | 280 | U | 190 |
| 2-Nitroaniline | | | < 770 | 770 | U | 390 | < 780 | 780 | U | 390 | < 790 | 790 | U | 400 | < 800 | 800 | U | 400 |
| 2-Nitrophenol | | | < 270 | 270 | U | 240 | < 270 | 270 | U | 250 | < 280 | 280 | U | 250 | < 280 | 280 | U | 250 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 270 | 270 | U | 150 | < 270 | 270 | U | 150 | < 280 | 280 | U | 150 | < 280 | 280 | U | 160 |
| 3,3'-Dichlorobenzidine | | | < 770 | 770 | U | 180 | < 780 | 780 | U | 180 | < 790 | 790 | U | 190 | < 800 | 800 | U | 190 |
| 3-Nitroaniline | | | < 770 | 770 | U | 770 | < 780 | 780 | U | 780 | < 790 | 790 | U | 790 | < 800 | 800 | U | 800 |
| 4,6-Dinitro-2-methylphenol | | | < 1900 | 1,900 | U | 410 | < 2000 | 2,000 | U | 420 | < 2000 | 2,000 | U | 420 | < 2000 | 2,000 | U | 430 |
| 4-Bromophenyl phenyl ether | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| 4-Chloro-3-methylphenol | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 140 | < 280 | 280 | U | 140 | < 280 | 280 | U | 140 |
| 4-Chloroaniline | | | < 310 | 310 | U | 180 | < 310 | 310 | U | 180 | < 310 | 310 | U | 180 | < 320 | 320 | U | 190 |
| 4-Chlorophenyl phenyl ether | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| 4-Nitroaniline | | | < 770 | 770 | U | 130 | < 780 | 780 | U | 130 | < 790 | 790 | U | 130 | < 800 | 800 | U | 130 |
| 4-Nitrophenol | | | < 380 | 380 | U | 170 | < 390 | 390 | U | 180 | < 390 | 390 | U | 180 | < 400 | 400 | U | 180 |
| Acenaphthene | 20,000 | 100,000 | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Acenaphthylene | 100,000 | 100,000 | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Acetophenone | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Aniline | | | < 310 | 310 | U | 310 | < 310 | 310 | U | 310 | < 310 | 310 | U | 310 | < 320 | 320 | U | 320 |
| Anthracene | 100,000 | 100,000 | 200 | 270 | J | 130 | < 270 | 270 | U | 130 | 200 | 280 | J | 130 | < 280 | 280 | U | 130 |
| Benz(a)anthracene | 1,000 | 1,000 | 240 | 270 | J | 130 | < 270 | 270 | U | 130 | 580 | 280 | | 130 | < 280 | 280 | U | 130 |
| Benzidine | | | < 770 | 770 | U | 230 | < 780 | 780 | U | 230 | < 790 | 790 | U | 230 | < 800 | 800 | U | 230 |
| Benzo(a)pyrene | 1,000 | 1,000 | 250 | 270 | J | 130 | < 270 | 270 | U | 130 | 480 | 280 | | 130 | < 280 | 280 | U | 130 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 280 | 270 | | 130 | < 270 | 270 | U | 130 | 380 | 280 | | 130 | < 280 | 280 | U | 140 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 300 | 270 | | 120 | < 270 | 270 | U | 130 | 240 | 280 | J | 130 | < 280 | 280 | U | 130 |
| Benzo(k)fluoranthene | 800 | 3,900 | 230 | 270 | J | 130 | < 270 | 270 | U | 130 | 500 | 280 | | 130 | < 280 | 280 | U | 130 |
| Benzoic acid | | | < 1900 | 1,900 | U | 770 | < 2000 | 2,000 | U | 780 | < 2000 | 2,000 | U | 790 | < 2000 | 2,000 | U | 800 |
| Benzyl butyl phthalate | | | < 270 | 270 | U | 99 | < 270 | 270 | U | 100 | < 280 | 280 | U | 100 | < 280 | 280 | U | 100 |
| Bis(2-chloroethoxy)methane | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Bis(2-chloroethyl)ether | | | < 270 | 270 | U | 100 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Bis(2-chloroisopropyl)ether | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Bis(2-ethylhexyl)phthalate | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Carbazole | | | < 1900 | 1,900 | U | 290 | < 2000 | 2,000 | U | 300 | < 2000 | 2,000 | U | 300 | < 2000 | 2,000 | U | 300 |
| Chrysene | 1,000 | 3,900 | 320 | 270 | | 130 | < 270 | 270 | U | 130 | 580 | 280 | | 130 | < 280 | 280 | U | 130 |
| Dibenz(a,h)anthracene | 330 | 330 | < 270 | 270 | U | 120 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| Dibenzofuran | 7,000 | 59,000 | 120 | 270 | J | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 120 |
| Diethyl phthalate | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Dimethylphthalate | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Di-n-butylphthalate | | | < 270 | 270 | U | 100 | < 270 | 270 | U | 100 | < 280 | 280 | U | 100 | < 280 | 280 | U | 110 |
| Di-n-octylphthalate | | | < 270 | 270 | U | 99 | < 270 | 270 | U | 100 | < 280 | 280 | U | 100 | < 280 | 280 | U | 100 |
| Fluoranthene | 100,000 | 100,000 | 510 | 270 | | 120 | < 270 | 270 | U | 130 | 1,100 | 280 | | 130 | < 280 | 280 | U | 130 |
| Fluorene | 30,000 | 100,000 | < 270 | 270 | U | 130 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| Hexachlorobenzene | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 120 |
| Hexachlorobutadiene | | | < 270 | 270 | U | 140 | < 270 | 270 | U | 140 | < 280 | 280 | U | 140 | < 280 | 280 | U | 140 |
| Hexachlorocyclopentadiene | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Hexachloroethane | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 120 | < 280 | 280 | U | 120 | < 280 | 280 | U | 120 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 230 | 270 | J | 130 | < 270 | 270 | U | 130 | 260 | 280 | J | 130 | < 280 | 280 | U | 130 |
| Isophorone | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Naphthalene | 12,000 | 100,000 | 310 | 270 | | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| Nitrobenzene | | | < 270 | 270 | U | 130 | < 270 | 270 | U | 140 | < 280 | 280 | U | 140 | < 280 | 280 | U | 140 |
| N-Nitrosodimethylamine | | | < 270 | 270 | U | 110 | < 270 | 270 | U | 110 | < 280 | 280 | U | 110 | < 280 | 280 | U | 110 |
| N-Nitrosodi-n-propylamine | | | < 270 | 270 | U | 120 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| N-Nitrosodiphenylamine | | | < 270 | 270 | U | 150 | < 270 | 270 | U | 150 | < 280 | 280 | U | 150 | < 280 | 280 | U | 150 |
| Pentachloronitrobenzene | | | < 270 | 270 | U | 140 | < 270 | 270 | U | 150 | < 280 | 280 | U | 150 | < 280 | 280 | U | 150 |
| Pentachlorophenol | 800 | 6,700 | < 270 | 270 | U | 140 | < 270 | 270 | U | 150 | < 280 | 280 | U | 150 | < 280 | 280 | U | 150 |
| Phenanthrene | 100,000 | 100,000 | 600 | 270 | | 110 | < 270 | 270 | U | 110 | 740 | 280 | | 110 | < 280 | 280 | U | 110 |
| Phenol | 330 | 100,000 | < 270 | 270 | U | 120 | < 270 | 270 | U | 130 | < 280 | 280 | U | 130 | < 280 | 280 | U | 130 |
| Pyrene | 100,000 | 100,000 | 570 | 270 | | 130 | < 270 | 270 | U | 130 | 1,000 | 280 | | 140 | < 280 | 280 | U | 140 |
| Pyridine | | | < 270 | 270 | U | 94 | < 270 | 270 | U | 96 | < 280 | 280 | U | 97 | < 280 | 280 | U | 98 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 RL - Laboratory Reporting

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B17 | | | | 15B18 | | | | | | | |
|-------------------------------|--|--|------------|-------|------|-----|----------------|--------|----|------|------------|----------------|--------|-----|
| | | | 11/20/2015 | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | Result | RL | Qual | MDL | (6-8) µg/Kg | Result | RL | Qual | MDL | (0-2) µg/Kg | Result | RL |
| 1,2,4,5-Tetrachlorobenzene | | | < 290 | 290 | U | 150 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| 1,2,4-Trichlorobenzene | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 |
| 1,2-Dichlorobenzene | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 99 | < 260 | 260 | U | 100 |
| 1,2-Diphenylhydrazine | | | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| 1,3-Dichlorobenzene | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| 1,4-Dichlorobenzene | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| 2,4,5-Trichlorophenol | | | < 290 | 290 | U | 230 | < 250 | 250 | U | 190 | < 260 | 260 | U | 200 |
| 2,4,6-Trichlorophenol | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| 2,4-Dichlorophenol | | | < 290 | 290 | U | 150 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| 2,4-Dimethylphenol | | | < 290 | 290 | U | 100 | < 250 | 250 | U | 87 | < 260 | 260 | U | 92 |
| 2,4-Dinitrophenol | | | < 840 | 840 | U | 290 | < 700 | 700 | U | 250 | < 740 | 740 | U | 260 |
| 2,4-Dinitrotoluene | | | < 290 | 290 | U | 170 | < 250 | 250 | U | 140 | < 260 | 260 | U | 150 |
| 2,6-Dinitrotoluene | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| 2-Chloronaphthalene | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 |
| 2-Chlorophenol | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 |
| 2-Methylnaphthalene | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 290 | 290 | U | 200 | < 250 | 250 | U | 170 | < 260 | 260 | U | 170 |
| 2-Nitroaniline | | | < 840 | 840 | U | 420 | < 700 | 700 | U | 360 | < 740 | 740 | U | 370 |
| 2-Nitrophenol | | | < 290 | 290 | U | 270 | < 250 | 250 | U | 220 | < 260 | 260 | U | 230 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 290 | 290 | U | 170 | < 250 | 250 | U | 140 | < 260 | 260 | U | 150 |
| 3,3'-Dichlorobenzidine | | | < 840 | 840 | U | 200 | < 700 | 700 | U | 170 | < 740 | 740 | U | 170 |
| 3-Nitroaniline | | | < 840 | 840 | U | 840 | < 700 | 700 | U | 700 | < 740 | 740 | U | 740 |
| 4,6-Dinitro-2-methylphenol | | | < 2100 | 2,100 | U | 450 | < 1800 | 1,800 | U | 380 | < 1800 | 1,800 | U | 400 |
| 4-Bromophenyl phenyl ether | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| 4-Chloro-3-methylphenol | | | < 290 | 290 | U | 150 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| 4-Chloroaniline | | | < 340 | 340 | U | 200 | < 280 | 280 | U | 160 | < 300 | 300 | U | 170 |
| 4-Chlorophenyl phenyl ether | | | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| 4-Nitroaniline | | | < 840 | 840 | U | 140 | < 700 | 700 | U | 120 | < 740 | 740 | U | 120 |
| 4-Nitrophenol | | | < 420 | 420 | U | 190 | < 350 | 350 | U | 160 | < 370 | 370 | U | 170 |
| Acenaphthene | 20,000 | 100,000 | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 |
| Acenaphthylene | 100,000 | 100,000 | < 290 | 290 | U | 120 | < 250 | 250 | U | 99 | < 260 | 260 | U | 100 |
| Acetophenone | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Aniline | | | < 340 | 340 | U | 340 | < 280 | 280 | U | 280 | < 300 | 300 | U | 300 |
| Anthracene | 100,000 | 100,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Benz(a)anthracene | 1,000 | 1,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Benidine | | | < 840 | 840 | U | 250 | < 700 | 700 | U | 210 | < 740 | 740 | U | 220 |
| Benzo(a)pyrene | 1,000 | 1,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| Benzo(ghi)perylene | 100,000 | 100,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Benzo(k)fluoranthene | 800 | 3,900 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Benzoic acid | | | < 2100 | 2,100 | U | 840 | < 1800 | 1,800 | U | 700 | < 1800 | 1,800 | U | 740 |
| Benzyl butyl phthalate | | | < 290 | 290 | U | 110 | < 250 | 250 | U | 91 | < 260 | 260 | U | 95 |
| Bis(2-chloroethoxy)methane | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 97 | < 260 | 260 | U | 100 |
| Bis(2-chloroethyl)ether | | | < 290 | 290 | U | 110 | < 250 | 250 | U | 95 | < 260 | 260 | U | 100 |
| Bis(2-chloroisopropyl)ether | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 98 | < 260 | 260 | U | 100 |
| Bis(2-ethylhexyl)phthalate | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| Carbazole | | | < 2100 | 2,100 | U | 320 | < 1800 | 1,800 | U | 270 | < 1800 | 1,800 | U | 280 |
| Chrysene | 1,000 | 3,900 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Dibenz(a,h)anthracene | 330 | 330 | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Dibenzofuran | 7,000 | 59,000 | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| Diethyl phthalate | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Dimethylphthalate | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 |
| Di-n-butylphthalate | | | < 290 | 290 | U | 110 | < 250 | 250 | U | 94 | < 260 | 260 | U | 98 |
| Di-n-octylphthalate | | | < 290 | 290 | U | 110 | < 250 | 250 | U | 91 | < 260 | 260 | U | 95 |
| Fluoranthene | 100,000 | 100,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Fluorene | 30,000 | 100,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Hexachlorobenzene | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| Hexachlorobutadiene | | | < 290 | 290 | U | 150 | < 250 | 250 | U | 130 | < 260 | 260 | U | 130 |
| Hexachlorocyclopentadiene | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 |
| Hexachloroethane | | | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 |
| Isophorone | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 99 | < 260 | 260 | U | 100 |
| Naphthalene | 12,000 | 100,000 | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| Nitrobenzene | | | < 290 | 290 | U | 150 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| N-Nitrosodimethylamine | | | < 290 | 290 | U | 120 | < 250 | 250 | U | 99 | < 260 | 260 | U | 100 |
| N-Nitrosodi-n-propylamine | | | < 290 | 290 | U | 140 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| N-Nitrosodiphenylamine | | | < 290 | 290 | U | 160 | < 250 | 250 | U | 140 | < 260 | 260 | U | 140 |
| Pentachloronitrobenzene | | | < 290 | 290 | U | 160 | < 250 | 250 | U | 130 | < 260 | 260 | U | 140 |
| Pentachlorophenol | 800 | 6,700 | < 290 | 290 | U | 160 | < 250 | 250 | U | 130 | < 260 | 260 | U | 140 |
| Phenanthrene | 100,000 | 100,000 | < 290 | 290 | U | 120 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 |
| Phenol | 330 | 100,000 | < 290 | 290 | U | 130 | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 |
| Pyrene | 100,000 | 100,000 | < 290 | 290 | U | 140 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 |
| Pyridine | | | < 290 | 290 | U | 100 | < 250 | 250 | U | 87 | < 260 | 260 | U | 91 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B19 | | | | | | | | | | | | 15B20 | | | | 15B21 | | | |
|-------------------------------|--|--|-----------------|-------|------|-----|-----------------|-------|------|-----|------------------|-------|------|-----|-----------------|-------|------|------|------------|--|--|--|
| | | | 11/20/2015 | | | | | | | | | | | | 11/20/2015 | | | | 11/20/2015 | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (8-10') µg/Kg | | | | (6-8') µg/Kg | | | | | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | | |
| 1,2,4,5-Tetrachlorobenzene | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| 1,2,4-Trichlorobenzene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 380 | 380 | U | 160 | | | | |
| 1,2-Dichlorobenzene | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| 1,2-Diphenylhydrazine | | | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 180 | | | | |
| 1,3-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| 1,4-Dichlorobenzene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| 2,4,5-Trichlorophenol | | | < 250 | 250 | U | 200 | < 260 | 260 | U | 200 | < 270 | 270 | U | 210 | < 380 | 380 | U | 300 | | | | |
| 2,4,6-Trichlorophenol | | | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| 2,4-Dichlorophenol | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| 2,4-Dimethylphenol | | | < 250 | 250 | U | 89 | < 260 | 260 | U | 93 | < 270 | 270 | U | 95 | < 380 | 380 | U | 130 | | | | |
| 2,4-Dinitrophenol | | | < 720 | 720 | U | 250 | < 750 | 750 | U | 260 | < 770 | 770 | U | 270 | < 1100 | 1,100 | U | 380 | | | | |
| 2,4-Dinitrotoluene | | | < 250 | 250 | U | 140 | < 260 | 260 | U | 150 | < 270 | 270 | U | 150 | < 380 | 380 | U | 210 | | | | |
| 2,6-Dinitrotoluene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| 2-Chloronaphthalene | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| 2-Chlorophenol | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| 2-Methylnaphthalene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 250 | 250 | U | 170 | < 260 | 260 | U | 180 | < 270 | 270 | U | 180 | < 110 | 110 | U | 260 | | | | |
| 2-Nitroaniline | | | < 720 | 720 | U | 360 | < 750 | 750 | U | 380 | < 770 | 770 | U | 390 | < 1100 | 1,100 | U | 550 | | | | |
| 2-Nitrophenol | | | < 250 | 250 | U | 230 | < 260 | 260 | U | 240 | < 270 | 270 | U | 240 | < 380 | 380 | U | 350 | | | | |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 250 | 250 | U | 140 | < 260 | 260 | U | 150 | < 270 | 270 | U | 150 | < 380 | 380 | U | 210 | | | | |
| 3,3'-Dichlorobenzidine | | | < 720 | 720 | U | 170 | < 750 | 750 | U | 180 | < 770 | 770 | U | 180 | < 1100 | 1,100 | U | 260 | | | | |
| 3-Nitroaniline | | | < 720 | 720 | U | 720 | < 750 | 750 | U | 750 | < 770 | 770 | U | 770 | < 1100 | 1,100 | U | 1100 | | | | |
| 4,6-Dinitro-2-methylphenol | | | < 1800 | 1,800 | U | 390 | < 1900 | 1,900 | U | 400 | < 1900 | 1,900 | U | 410 | < 2700 | 2,700 | U | 590 | | | | |
| 4-Bromophenyl phenyl ether | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| 4-Chloro-3-methylphenol | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| 4-Chloroaniline | | | < 290 | 290 | U | 170 | < 300 | 300 | U | 170 | < 310 | 310 | U | 180 | < 440 | 440 | U | 250 | | | | |
| 4-Chlorophenyl phenyl ether | | | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| 4-Nitroaniline | | | < 720 | 720 | U | 120 | < 750 | 750 | U | 120 | < 770 | 770 | U | 130 | < 1100 | 1,100 | U | 180 | | | | |
| 4-Nitrophenol | | | < 360 | 360 | U | 160 | < 370 | 370 | U | 170 | < 380 | 380 | U | 170 | < 540 | 540 | U | 250 | | | | |
| Acenaphthene | 20,000 | 100,000 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| Acenaphthylene | 100,000 | 100,000 | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| Acetophenone | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| Aniline | | | < 290 | 290 | U | 290 | < 300 | 300 | U | 300 | < 310 | 310 | U | 310 | < 440 | 440 | U | 440 | | | | |
| Anthracene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Benz(a)anthracene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Benzidine | | | < 720 | 720 | U | 210 | < 750 | 750 | U | 220 | < 770 | 770 | U | 230 | < 1100 | 1,100 | U | 320 | | | | |
| Benzo(a)pyrene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 180 | | | | |
| Benzo(b)fluoranthene | 1,000 | 1,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| Benzo(ghi)perylene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 180 | | | | |
| Benzo(k)fluoranthene | 800 | 3,900 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Benzoic acid | | | < 1800 | 1,800 | U | 720 | < 1900 | 1,900 | U | 750 | < 1900 | 1,900 | U | 770 | < 2700 | 2,700 | U | 1100 | | | | |
| Benzyl butyl phthalate | | | < 250 | 250 | U | 93 | < 260 | 260 | U | 96 | < 270 | 270 | U | 99 | < 380 | 380 | U | 140 | | | | |
| Bis(2-chloroethoxy)methane | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| Bis(2-chloroethyl)ether | | | < 250 | 250 | U | 97 | < 260 | 260 | U | 100 | < 270 | 270 | U | 100 | < 380 | 380 | U | 150 | | | | |
| Bis(2-chloroisopropyl)ether | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| Bis(2-ethylhexyl)phthalate | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Carbazole | | | < 1800 | 1,800 | U | 270 | < 1900 | 1,900 | U | 280 | < 1900 | 1,900 | U | 290 | < 2700 | 2,700 | U | 410 | | | | |
| Chrysene | 1,000 | 3,900 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Dibenz(a,h)anthracene | 330 | 330 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 110 | 110 | U | 180 | | | | |
| Dibenzofuran | 7,000 | 59,000 | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Diethyl phthalate | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| Dimethylphthalate | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| Di-n-butylphthalate | | | < 250 | 250 | U | 96 | < 260 | 260 | U | 99 | < 270 | 270 | U | 100 | < 380 | 380 | U | 140 | | | | |
| Di-n-octylphthalate | | | < 250 | 250 | U | 93 | < 260 | 260 | U | 96 | < 270 | 270 | U | 99 | < 380 | 380 | U | 140 | | | | |
| Fluoranthene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 180 | | | | |
| Fluorene | 30,000 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Hexachlorobenzene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Hexachlorobutadiene | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 140 | < 270 | 270 | U | 140 | < 380 | 380 | U | 200 | | | | |
| Hexachlorocyclopentadiene | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 120 | < 380 | 380 | U | 170 | | | | |
| Hexachloroethane | | | < 250 | 250 | U | 110 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 130 | < 380 | 380 | U | 180 | | | | |
| Isophorone | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 100 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| Naphthalene | 12,000 | 100,000 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Nitrobenzene | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| N-Nitrosodimethylamine | | | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 150 | | | | |
| N-Nitrosodi-n-propylamine | | | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 380 | 380 | U | 180 | | | | |
| N-Nitrosodiphenylamine | | | < 250 | 250 | U | 140 | < 260 | 260 | U | 140 | < 270 | 270 | U | 150 | < 380 | 380 | U | 210 | | | | |
| Pentachloronitrobenzene | | | < 250 | 250 | U | 130 | < 260 | 260 | U | 140 | < 270 | 270 | U | 140 | < 380 | 380 | U | 200 | | | | |
| Pentachlorophenol | 800 | 6,700 | < 250 | 250 | U | 140 | < 260 | 260 | U | 140 | < 270 | 270 | U | 140 | < 380 | 380 | U | 210 | | | | |
| Phenanthrene | 100,000 | 100,000 | < 250 | 250 | U | 100 | < 260 | 260 | U | 110 | < 270 | 270 | U | 110 | < 380 | 380 | U | 160 | | | | |
| Phenol | 330 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 120 | < 270 | 270 | U | 120 | < 110 | 110 | U | 170 | | | | |
| Pyrene | 100,000 | 100,000 | < 250 | 250 | U | 120 | < 260 | 260 | U | 130 | < 270 | 270 | U | 130 | < 380 | 380 | U | 190 | | | | |
| Pyridine | | | < 250 | 250 | U | 89 | < 260 | 260 | U | 92 | < 270 | 270 | U | 94 | < 380 | 380 | U | 130 | | | | |

Notes:
RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
RL - Laboratory Reporting Limit

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TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B22 | | | | | | | | 15B23 | | | | | | | |
|-------------------------------|--|--|-----------------|-------|------|-----|-----------------|-------|------|------|-----------------|--------|------|------|-----------------|-------|------|-----|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | |
| | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (2-4') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 320 | 320 | U | 160 | < 360 | 360 | U | 180 | < 2500 | 2,500 | U | 1300 | < 350 | 350 | U | 170 |
| 1,2,4-Trichlorobenzene | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| 1,2-Dichlorobenzene | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| 1,2-Diphenylhydrazine | | | < 320 | 320 | U | 150 | < 360 | 360 | U | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 160 |
| 1,3-Dichlorobenzene | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| 1,4-Dichlorobenzene | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| 2,4,5-Trichlorophenol | | | < 320 | 320 | U | 250 | < 360 | 360 | U | 280 | < 2500 | 2,500 | U | 2000 | < 350 | 350 | U | 270 |
| 2,4,6-Trichlorophenol | | | < 320 | 320 | U | 150 | < 360 | 360 | U | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 160 |
| 2,4-Dichlorophenol | | | < 320 | 320 | U | 160 | < 360 | 360 | U | 180 | < 2500 | 2,500 | U | 1300 | < 350 | 350 | U | 170 |
| 2,4-Dimethylphenol | | | < 320 | 320 | U | 110 | < 360 | 360 | U | 130 | < 2500 | 2,500 | U | 890 | < 350 | 350 | U | 120 |
| 2,4-Dinitrophenol | | | < 920 | 920 | U | 320 | < 1000 | 1,000 | U | 360 | < 7200 | 7,200 | U | 2500 | < 990 | 990 | U | 350 |
| 2,4-Dinitrotoluene | | | < 320 | 320 | U | 180 | < 360 | 360 | U | 200 | < 2500 | 2,500 | U | 1400 | < 350 | 350 | U | 200 |
| 2,6-Dinitrotoluene | | | < 320 | 320 | U | 150 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 160 |
| 2-Chloronaphthalene | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| 2-Chlorophenol | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| 2-Methylnaphthalene | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 320 | 320 | U | 220 | < 330 | 330 | U | 240 | < 720 | 720 | U | 720 | < 330 | 330 | U | 230 |
| 2-Nitroaniline | | | < 920 | 920 | U | 470 | < 1000 | 1,000 | U | 520 | < 7200 | 7,200 | U | 3600 | < 990 | 990 | U | 500 |
| 2-Nitrophenol | | | < 320 | 320 | U | 290 | < 360 | 360 | U | 330 | < 2500 | 2,500 | U | 2300 | < 350 | 350 | U | 310 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 320 | 320 | U | 180 | < 360 | 360 | U | 200 | < 2500 | 2,500 | U | 1400 | < 350 | 350 | U | 200 |
| 3,3'-Dichlorobenzidine | | | < 920 | 920 | U | 220 | < 1000 | 1,000 | U | 240 | < 7200 | 7,200 | U | 1700 | < 990 | 990 | U | 230 |
| 3-Nitroaniline | | | < 920 | 920 | U | 920 | < 1000 | 1,000 | U | 1000 | < 7200 | 7,200 | U | 7200 | < 990 | 990 | U | 990 |
| 4,6-Dinitro-2-methylphenol | | | < 2300 | 2,300 | U | 500 | < 2600 | 2,600 | U | 560 | < 18000 | 18,000 | U | 3900 | < 2500 | 2,500 | U | 530 |
| 4-Bromophenyl phenyl ether | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| 4-Chloro-3-methylphenol | | | < 320 | 320 | U | 160 | < 360 | 360 | U | 180 | < 2500 | 2,500 | U | 1300 | < 350 | 350 | U | 170 |
| 4-Chloroaniline | | | < 370 | 370 | U | 220 | < 410 | 410 | U | 240 | < 2900 | 2,900 | U | 1700 | < 400 | 400 | U | 230 |
| 4-Chlorophenyl phenyl ether | | | < 320 | 320 | U | 160 | < 360 | 360 | U | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 170 |
| 4-Nitroaniline | | | < 920 | 920 | U | 150 | < 1000 | 1,000 | U | 170 | < 7200 | 7,200 | U | 1200 | < 990 | 990 | U | 170 |
| 4-Nitrophenol | | | < 460 | 460 | U | 210 | < 520 | 520 | U | 230 | < 3600 | 3,600 | U | 1600 | < 500 | 500 | U | 220 |
| Acenaphthene | 20,000 | 100,000 | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| Acenaphthylene | 100,000 | 100,000 | < 320 | 320 | U | 130 | 190 | 360 | J | 140 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| Acetophenone | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| Aniline | | | < 370 | 370 | U | 370 | < 410 | 410 | U | 410 | < 2900 | 2,900 | U | 2900 | < 400 | 400 | U | 400 |
| Anthracene | 100,000 | 100,000 | 190 | 320 | J | 150 | 210 | 360 | J | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 160 |
| Benz(a)anthracene | 1,000 | 1,000 | 560 | 320 | | 160 | 870 | 360 | | 170 | 1,300 | 2,500 | J | 1200 | < 350 | 350 | U | 170 |
| Benzidine | | | < 920 | 920 | U | 270 | < 1000 | 1,000 | U | 300 | < 7200 | 7,200 | U | 2100 | < 990 | 990 | U | 290 |
| Benzo(a)pyrene | 1,000 | 1,000 | 520 | 320 | | 150 | 1,200 | 360 | | 170 | < 1000 | 1,000 | U | 1000 | < 350 | 350 | U | 160 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 540 | 320 | | 160 | 690 | 360 | | 180 | < 1000 | 1,000 | U | 1000 | < 350 | 350 | U | 170 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 430 | 320 | | 150 | 590 | 360 | | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 160 |
| Benzo(k)fluoranthene | 800 | 3,900 | 470 | 320 | | 150 | 790 | 360 | | 170 | 730 | 720 | | 720 | < 350 | 350 | U | 160 |
| Benzoic acid | | | < 2300 | 2,300 | U | 920 | < 2600 | 2,600 | U | 1000 | < 18000 | 18,000 | U | 7200 | < 2500 | 2,500 | U | 990 |
| Benzyl butyl phthalate | | | < 320 | 320 | U | 120 | < 360 | 360 | U | 130 | < 2500 | 2,500 | U | 930 | < 350 | 350 | U | 130 |
| Bis(2-chloroethoxy)methane | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 140 | < 2500 | 2,500 | U | 990 | < 350 | 350 | U | 140 |
| Bis(2-chloroethyl)ether | | | < 320 | 320 | U | 120 | < 360 | 360 | U | 140 | < 2500 | 2,500 | U | 970 | < 350 | 350 | U | 130 |
| Bis(2-chloroisopropyl)ether | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 140 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| Bis(2-ethylhexyl)phthalate | | | 380 | 320 | | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| Carbazole | | | < 2300 | 2,300 | U | 350 | < 2600 | 2,600 | U | 390 | < 18000 | 18,000 | U | 2700 | < 2500 | 2,500 | U | 380 |
| Chrysene | 1,000 | 3,900 | 680 | 320 | | 160 | 860 | 360 | | 170 | 1,200 | 1,100 | | 1100 | < 350 | 350 | U | 170 |
| Dibenz(a,h)anthracene | 330 | 330 | < 320 | 320 | U | 150 | < 330 | 330 | U | 170 | < 720 | 720 | U | 720 | < 330 | 330 | U | 160 |
| Dibenzofuran | 7,000 | 59,000 | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 140 |
| Diethyl phthalate | | | < 320 | 320 | U | 150 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 160 |
| Dimethylphthalate | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| Di-n-butylphthalate | | | < 320 | 320 | U | 120 | < 360 | 360 | U | 140 | < 2500 | 2,500 | U | 960 | < 350 | 350 | U | 130 |
| Di-n-octylphthalate | | | < 320 | 320 | U | 120 | < 360 | 360 | U | 130 | < 2500 | 2,500 | U | 930 | < 350 | 350 | U | 130 |
| Fluoranthene | 100,000 | 100,000 | 960 | 320 | | 150 | 710 | 360 | | 170 | 2,200 | 2,500 | J | 1200 | < 350 | 350 | U | 160 |
| Fluorene | 30,000 | 100,000 | < 320 | 320 | U | 150 | < 360 | 360 | U | 170 | < 2500 | 2,500 | U | 1200 | < 350 | 350 | U | 160 |
| Hexachlorobenzene | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 140 |
| Hexachlorobutadiene | | | < 320 | 320 | U | 170 | < 360 | 360 | U | 190 | < 2500 | 2,500 | U | 1300 | < 350 | 350 | U | 180 |
| Hexachlorocyclopentadiene | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| Hexachloroethane | | | < 320 | 320 | U | 140 | < 360 | 360 | U | 160 | < 2500 | 2,500 | U | 1100 | < 350 | 350 | U | 150 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 350 | 320 | | 150 | 500 | 360 | | 170 | < 720 | 720 | U | 720 | < 350 | 350 | U | 160 |
| Isophorone | | | < 320 | 320 | U | 130 | < 360 | 360 | U | 140 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| Naphthalene | 12,000 | 100,000 | < 320 | 320 | U | 130 | < 360 | 360 | U | 150 | < 2500 | 2,500 | U | 1000 | < 350 | 350 | U | 140 |
| Nitrobenzene | | | < 320 | 320 | U | 160 | < 360 | 360 | U | 180 | < 2500 | 2,500 | U | 1300 | < 350 | 3 | | |

TABLE 4
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | Duplicate 11/20/2015 | | | | Duplicate 11/23/2015 | | | |
|-------------------------------|--|--|-------------------------|-------|------|-----|-------------------------|-------|------|-----|
| | | | µg/Kg | | | | µg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| 1,2,4,5-Tetrachlorobenzene | | | < 260 | 260 | U | 130 | < 250 | 250 | U | 130 |
| 1,2,4-Trichlorobenzene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 110 |
| 1,2-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| 1,2-Diphenylhydrazine | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 120 |
| 1,3-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 110 |
| 1,4-Dichlorobenzene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 110 |
| 2,4,5-Trichlorophenol | | | < 260 | 260 | U | 210 | < 250 | 250 | U | 190 |
| 2,4,6-Trichlorophenol | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| 2,4-Dichlorophenol | | | < 260 | 260 | U | 130 | < 250 | 250 | U | 130 |
| 2,4-Dimethylphenol | | | < 260 | 260 | U | 93 | < 250 | 250 | U | 88 |
| 2,4-Dinitrophenol | | | < 750 | 750 | U | 260 | < 710 | 710 | U | 250 |
| 2,4-Dinitrotoluene | | | < 260 | 260 | U | 150 | < 250 | 250 | U | 140 |
| 2,6-Dinitrotoluene | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| 2-Chloronaphthalene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| 2-Chlorophenol | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| 2-Methylnaphthalene | | | 510 | 260 | | 110 | < 250 | 250 | U | 110 |
| 2-Methylphenol (o-cresol) | 330 | 100,000 | < 260 | 260 | U | 180 | < 250 | 250 | U | 170 |
| 2-Nitroaniline | | | < 750 | 750 | U | 380 | < 710 | 710 | U | 360 |
| 2-Nitrophenol | | | < 260 | 260 | U | 240 | < 250 | 250 | U | 230 |
| 3&4-Methylphenol (m&p-cresol) | 330 | 100,000 | < 260 | 260 | U | 150 | < 250 | 250 | U | 140 |
| 3,3'-Dichlorobenzidine | | | < 750 | 750 | U | 180 | < 710 | 710 | U | 170 |
| 3-Nitroaniline | | | < 750 | 750 | U | 750 | < 710 | 710 | U | 710 |
| 4,6-Dinitro-2-methylphenol | | | < 1900 | 1,900 | U | 400 | < 1800 | 1,800 | U | 380 |
| 4-Bromophenyl phenyl ether | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| 4-Chloro-3-methylphenol | | | < 260 | 260 | U | 130 | < 250 | 250 | U | 130 |
| 4-Chloroaniline | | | < 300 | 300 | U | 180 | < 280 | 280 | U | 170 |
| 4-Chlorophenyl phenyl ether | | | < 260 | 260 | U | 130 | < 250 | 250 | U | 120 |
| 4-Nitroaniline | | | < 750 | 750 | U | 130 | < 710 | 710 | U | 120 |
| 4-Nitrophenol | | | < 380 | 380 | U | 170 | < 360 | 360 | U | 160 |
| Acenaphthene | 20,000 | 100,000 | 2,800 | 260 | | 110 | < 250 | 250 | U | 110 |
| Acenaphthylene | 100,000 | 100,000 | 430 | 260 | | 110 | < 250 | 250 | U | 99 |
| Acetophenone | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| Aniline | | | < 300 | 300 | U | 300 | < 280 | 280 | U | 280 |
| Anthracene | 100,000 | 100,000 | 2,300 | 260 | | 120 | < 250 | 250 | U | 120 |
| Benz(a)anthracene | 1,000 | 1,000 | 1,300 | 260 | | 130 | 120 | 250 | J | 120 |
| Benzidine | | | < 750 | 750 | U | 220 | < 710 | 710 | U | 210 |
| Benzo(a)pyrene | 1,000 | 1,000 | 1,100 | 260 | | 120 | < 250 | 250 | U | 120 |
| Benzo(b)fluoranthene | 1,000 | 1,000 | 1,400 | 260 | | 130 | < 250 | 250 | U | 120 |
| Benzo(ghi)perylene | 100,000 | 100,000 | 910 | 260 | | 120 | < 250 | 250 | U | 120 |
| Benzo(k)fluoranthene | 800 | 3,900 | 760 | 260 | | 120 | < 250 | 250 | U | 120 |
| Benzoic acid | | | < 1900 | 1,900 | U | 750 | < 1800 | 1,800 | U | 710 |
| Benzyl butyl phthalate | | | < 260 | 260 | U | 97 | < 250 | 250 | U | 92 |
| Bis(2-chloroethoxy)methane | | | < 260 | 260 | U | 100 | < 250 | 250 | U | 98 |
| Bis(2-chloroethyl)ether | | | < 260 | 260 | U | 100 | < 250 | 250 | U | 96 |
| Bis(2-chloroisopropyl)ether | | | < 260 | 260 | U | 100 | < 250 | 250 | U | 99 |
| Bis(2-ethylhexyl)phthalate | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| Carbazole | | | 520 | 1,900 | J | 290 | < 1800 | 1,800 | U | 270 |
| Chrysene | 1,000 | 3,900 | 1,600 | 260 | | 130 | 140 | 250 | J | 120 |
| Dibenz(a,h)anthracene | 330 | 330 | < 260 | 260 | U | 120 | < 250 | 250 | U | 120 |
| Dibenzofuran | 7,000 | 59,000 | 550 | 260 | | 110 | < 250 | 250 | U | 100 |
| Diethyl phthalate | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| Dimethylphthalate | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| Di-n-butylphthalate | | | < 260 | 260 | U | 100 | < 250 | 250 | U | 94 |
| Di-n-octylphthalate | | | < 260 | 260 | U | 97 | < 250 | 250 | U | 92 |
| Fluoranthene | 100,000 | 100,000 | 5,400 | 260 | | 120 | 240 | 250 | J | 120 |
| Fluorene | 30,000 | 100,000 | 980 | 260 | | 120 | < 250 | 250 | U | 120 |
| Hexachlorobenzene | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| Hexachlorobutadiene | | | < 260 | 260 | U | 140 | < 250 | 250 | U | 130 |
| Hexachlorocyclopentadiene | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| Hexachloroethane | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 110 |
| Indeno(1,2,3-cd)pyrene | 500 | 500 | 640 | 260 | | 120 | < 250 | 250 | U | 120 |
| Isophorone | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 99 |
| Naphthalene | 12,000 | 100,000 | 340 | 260 | | 110 | < 250 | 250 | U | 100 |
| Nitrobenzene | | | < 260 | 260 | U | 130 | < 250 | 250 | U | 120 |
| N-Nitrosodimethylamine | | | < 260 | 260 | U | 110 | < 250 | 250 | U | 100 |
| N-Nitrosodi-n-propylamine | | | < 260 | 260 | U | 120 | < 250 | 250 | U | 120 |
| N-Nitrosodiphenylamine | | | < 260 | 260 | U | 140 | < 250 | 250 | U | 140 |
| Pentachloronitrobenzene | | | < 260 | 260 | U | 140 | < 250 | 250 | U | 130 |
| Pentachlorophenol | 800 | 6,700 | < 260 | 260 | U | 140 | < 250 | 250 | U | 130 |
| Phenanthrene | 100,000 | 100,000 | 4,100 | 260 | | 110 | 160 | 250 | J | 100 |
| Phenol | 330 | 100,000 | < 260 | 260 | U | 120 | < 250 | 250 | U | 110 |
| Pyrene | 100,000 | 100,000 | 6,900 | 260 | | 130 | 210 | 250 | J | 120 |
| Pyridine | | | < 260 | 260 | U | 93 | < 250 | 250 | U | 87 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|---|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank. The value is estimated. This flag is used |
| J | a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 5
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Pesticides and PCBs

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B1 | | | | | | | | | 15B3 | | | | | | | | | 15B4 | | | | | | | | |
|--------------------|---|--|-----------------|-----|------|-----------------|--------|-----|-----------------|-----|--------|-----------------|------|-----|-----------------|-----|------|-----------------|--------|-----|-----------------|-----|--------|-----------------|------|-----|--------|----|------|
| | | | 11/24/2015 | | | | | | | | | 11/24/2015 | | | | | | | | | 11/24/2015 | | | | | | | | |
| | | | (0-2') µg/Kg | | | (5-7') µg/Kg | | | (0-2') µg/Kg | | | (5-7') µg/Kg | | | (0-2') µg/Kg | | | (5-7') µg/Kg | | | (0-2') µg/Kg | | | (5-7') µg/Kg | | | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual |
| 4,4' -DDD | 3.3 | 13,000 | < 9.5 | 9.5 | U | 9.5 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.4 | 2.4 | U | 2.4 | < 2.2 | 2.2 | U | 2.2 | < 2.4 | 2.4 | U | 2.4 | | | |
| 4,4' -DDE | 3.3 | 8,900 | < 9.5 | 9.5 | U | 9.5 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.4 | 2.4 | U | 2.4 | < 2.2 | 2.2 | U | 2.2 | < 2.4 | 2.4 | U | 2.4 | | | |
| 4,4' -DDT | 3.3 | 7,900 | < 9.5 | 9.5 | U | 9.5 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.4 | 2.4 | U | 2.4 | < 2.2 | 2.2 | U | 2.2 | < 2.4 | 2.4 | U | 2.4 | | | |
| a-BHC | 20 | 480 | < 9.5 | 9.5 | U | 9.5 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| a-Chlordane | 94 | 4,200 | < 19 | 19 | U | 19 | < 4.0 | 4.0 | U | 4.0 | < 3.6 | 3.6 | U | 3.6 | < 3.9 | 3.9 | U | 3.9 | < 3.6 | 3.6 | U | 3.6 | < 4.0 | 4.0 | U | 4.0 | | | |
| Aldrin | 5 | 97 | < 4.9 | 4.9 | U | 4.9 | < 4.0 | 4.0 | U | 4.0 | < 3.6 | 3.6 | U | 3.6 | < 3.9 | 3.9 | U | 3.9 | < 3.6 | 3.6 | U | 3.6 | < 4.0 | 4.0 | U | 4.0 | | | |
| b-BHC | 36 | 360 | < 9.5 | 9.5 | U | 9.5 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Chlordane | | | < 190 | 190 | U | 190 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| d-BHC | 40 | 100,000 | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Dieldrin | 5 | 200 | < 4.9 | 4.9 | U | 4.9 | < 4.0 | 4.0 | U | 4.0 | < 3.6 | 3.6 | U | 3.6 | < 3.9 | 3.9 | U | 3.9 | < 3.6 | 3.6 | U | 3.6 | < 4.0 | 4.0 | U | 4.0 | | | |
| Endosulfan I | 2,400 | 24,000 | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Endosulfan II | 2,400 | 24,000 | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Endosulfan sulfate | 2,400 | 24,000 | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Endrin | 14 | 11,000 | < 9.5 | 9.5 | U | 9.5 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Endrin aldehyde | | | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Endrin ketone | | | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| g-BHC | | | < 7.6 | 7.6 | U | 7.6 | < 3.0 | 3.0 | U | 3.0 | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 | | | |
| g-Chlordane | | | < 19 | 19 | U | 19 | < 4.0 | 4.0 | U | 4.0 | < 3.6 | 3.6 | U | 3.6 | < 3.9 | 3.9 | U | 3.9 | < 3.6 | 3.6 | U | 3.6 | < 4.0 | 4.0 | U | 4.0 | | | |
| Heptachlor | 42 | 2,100 | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Heptachlor epoxide | | | < 38 | 38 | U | 38 | < 7.9 | 7.9 | U | 7.9 | < 7.1 | 7.1 | U | 7.1 | < 7.9 | 7.9 | U | 7.9 | < 7.2 | 7.2 | U | 7.2 | < 7.9 | 7.9 | U | 7.9 | | | |
| Methoxychlor | | | < 190 | 190 | U | 190 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| Toxaphene | | | < 760 | 760 | U | 760 | < 160 | 160 | U | 160 | < 140 | 140 | U | 140 | < 160 | 160 | U | 160 | < 140 | 140 | U | 140 | < 160 | 160 | U | 160 | | | |
| PCB-1016 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1221 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1232 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1242 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1248 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1254 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1260 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1262 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |
| PCB-1268 | 100 | 1,000 | < 38 | 38 | U | 38 | < 40 | 40 | U | 40 | < 36 | 36 | U | 36 | < 39 | 39 | U | 39 | < 36 | 36 | U | 36 | < 40 | 40 | U | 40 | | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 5
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Pesticides and PCBs

| | COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B6 | | | | | | | | 15B9 | | | | | | | | 15B10 | | | |
|--------------|--------------------|---|--|----------------|-----|------|-------|------------------|-----|------|-------|----------------|-----|------|-------|----------------|-----|------|-------|----------------|-----|------|-----|
| | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | |
| | | | | (3-5) µg/Kg | | | | (10-12) µg/Kg | | | | (0-2) µg/Kg | | | | (6-8) µg/Kg | | | | (6-8) µg/Kg | | | |
| | | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| Pesticides | 4,4' -DDD | 3.3 | 13,000 | < 2.3 | 2.3 | U | 2.3 | < 3.2 | 3.2 | U | 3.2 | < 2.1 | 2.1 | U | 2.1 | < 2.5 | 2.5 | U | 2.5 | < 2.3 | 2.3 | U | 2.3 |
| | 4,4' -DDE | 3.3 | 8,900 | < 2.3 | 2.3 | U | 2.3 | < 3.2 | 3.2 | U | 3.2 | < 2.1 | 2.1 | U | 2.1 | < 2.5 | 2.5 | U | 2.5 | < 2.3 | 2.3 | U | 2.3 |
| | 4,4' -DDT | 3.3 | 7,900 | < 2.3 | 2.3 | U | 2.3 | < 3.2 | 3.2 | U | 3.2 | < 2.1 | 2.1 | U | 2.1 | < 2.5 | 2.5 | U | 2.5 | < 2.3 | 2.3 | U | 2.3 |
| | a-BHC | 20 | 480 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | a-Chlordane | 94 | 4,200 | < 3.8 | 3.8 | U | 3.8 | < 5.3 | 5.3 | U | 5.3 | < 3.5 | 3.5 | U | 3.5 | < 4.2 | 4.2 | U | 4.2 | < 3.8 | 3.8 | U | 3.8 |
| | Aldrin | 5 | 97 | < 3.8 | 3.8 | U | 3.8 | < 2.6 | 2.6 | U | 2.6 | < 3.5 | 3.5 | U | 3.5 | < 4.2 | 4.2 | U | 4.2 | < 3.8 | 3.8 | U | 3.8 |
| | b-BHC | 36 | 360 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Chlordane | | | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | d-BHC | 40 | 100,000 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Dieldrin | 5 | 200 | < 3.8 | 3.8 | U | 3.8 | < 2.6 | 2.6 | U | 2.6 | < 3.5 | 3.5 | U | 3.5 | < 4.2 | 4.2 | U | 4.2 | < 3.8 | 3.8 | U | 3.8 |
| | Endosulfan I | 2,400 | 24,000 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Endosulfan II | 2,400 | 24,000 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Endosulfan sulfate | 2,400 | 24,000 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Endrin | 14 | 11,000 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Endrin aldehyde | | | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Endrin ketone | | | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | g-BHC | | | < 1.5 | 1.5 | U | 1.5 | < 2.1 | 2.1 | U | 2.1 | < 1.4 | 1.4 | U | 1.4 | < 1.7 | 1.7 | U | 1.7 | < 1.5 | 1.5 | U | 1.5 |
| | g-Chlordane | | | < 3.8 | 3.8 | U | 3.8 | < 5.3 | 5.3 | U | 5.3 | < 3.5 | 3.5 | U | 3.5 | < 4.2 | 4.2 | U | 4.2 | < 3.8 | 3.8 | U | 3.8 |
| | Heptachlor | 42 | 2,100 | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| | Heptachlor epoxide | | | < 7.6 | 7.6 | U | 7.6 | < 11 | 11 | U | 11 | < 7.0 | 7.0 | U | 7.0 | < 8.4 | 8.4 | U | 8.4 | < 7.6 | 7.6 | U | 7.6 |
| Methoxychlor | | | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 | |
| Toxaphene | | | < 150 | 150 | U | 150 | < 210 | 210 | U | 210 | < 140 | 140 | U | 140 | < 170 | 170 | U | 170 | < 150 | 150 | U | 150 | |
| PCBs | PCB-1016 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1221 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1232 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1242 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1248 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1254 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1260 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| | PCB-1262 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 |
| PCB-1268 | 100 | 1,000 | < 38 | 38 | U | 38 | < 53 | 53 | U | 53 | < 35 | 35 | U | 35 | < 42 | 42 | U | 42 | < 38 | 38 | U | 38 | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 5
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Pesticides and PCBs

| | COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B11 | | | | | | | | 15B13 | | | | | | | | 15B18 | | | | | | | |
|--------------|--------------------|---|--|-----------------|-----|------|-------|-----------------|-----|------|-------|-----------------|-----|------|-------|-----------------|-----|------|-------|-----------------|-----|------|-------|-----------------|-----|------|-----|
| | | | | 11/23/2015 | | | | | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | |
| | | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | | (0-2') µg/Kg | | | | (6-8') µg/Kg | | | |
| | | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| Pesticides | 4,4' -DDD | 3.3 | 13,000 | < 2.2 | 2.2 | U | 2.2 | < 2.4 | 2.4 | U | 2.4 | 3.1 | 2.3 | | 2.3 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.2 | 2.2 | U | 2.2 |
| | 4,4' -DDE | 3.3 | 8,900 | < 2.2 | 2.2 | U | 2.2 | < 2.4 | 2.4 | U | 2.4 | 7.9 | 2.3 | | 2.3 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.2 | 2.2 | U | 2.2 |
| | 4,4' -DDT | 3.3 | 7,900 | 4.2 | 2.2 | | 2.2 | < 2.4 | 2.4 | U | 2.4 | 9.4 | 2.3 | | 2.3 | < 2.4 | 2.4 | U | 2.4 | < 2.1 | 2.1 | U | 2.1 | < 2.2 | 2.2 | U | 2.2 |
| | a-BHC | 20 | 480 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | a-Chlordane | 94 | 4,200 | < 3.7 | 3.7 | U | 3.7 | < 4.0 | 4.0 | U | 4.0 | < 3.9 | 3.9 | U | 3.9 | < 4.0 | 4.0 | U | 4.0 | < 3.5 | 3.5 | U | 3.5 | < 3.7 | 3.7 | U | 3.7 |
| | Aldrin | 5 | 97 | < 3.7 | 3.7 | U | 3.7 | < 4.0 | 4.0 | U | 4.0 | < 3.9 | 3.9 | U | 3.9 | < 4.0 | 4.0 | U | 4.0 | < 3.5 | 3.5 | U | 3.5 | < 3.7 | 3.7 | U | 3.7 |
| | b-BHC | 36 | 360 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Chlordane | | | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | d-BHC | 40 | 100,000 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Dieldrin | 5 | 200 | < 3.7 | 3.7 | U | 3.7 | < 4.0 | 4.0 | U | 4.0 | < 3.9 | 3.9 | U | 3.9 | < 4.0 | 4.0 | U | 4.0 | < 3.5 | 3.5 | U | 3.5 | < 3.7 | 3.7 | U | 3.7 |
| | Endosulfan I | 2,400 | 24,000 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Endosulfan II | 2,400 | 24,000 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Endosulfan sulfate | 2,400 | 24,000 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Endrin | 14 | 11,000 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Endrin aldehyde | | | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Endrin ketone | | | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | g-BHC | | | < 1.5 | 1.5 | U | 1.5 | < 1.6 | 1.6 | U | 1.6 | < 2.0 | 2.0 | U | 2.0 | < 1.6 | 1.6 | U | 1.6 | < 1.4 | 1.4 | U | 1.4 | < 1.5 | 1.5 | U | 1.5 |
| | g-Chlordane | | | < 3.7 | 3.7 | U | 3.7 | < 4.0 | 4.0 | U | 4.0 | < 12 | 12 | U | 12 | < 4.0 | 4.0 | U | 4.0 | < 3.5 | 3.5 | U | 3.5 | < 3.7 | 3.7 | U | 3.7 |
| | Heptachlor | 42 | 2,100 | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| | Heptachlor epoxide | | | < 7.4 | 7.4 | U | 7.4 | < 8.0 | 8.0 | U | 8.0 | < 7.8 | 7.8 | U | 7.8 | < 7.9 | 7.9 | U | 7.9 | < 6.9 | 6.9 | U | 6.9 | < 7.5 | 7.5 | U | 7.5 |
| Methoxychlor | | | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 | |
| Toxaphene | | | < 150 | 150 | U | 150 | < 160 | 160 | U | 160 | < 160 | 160 | U | 160 | < 160 | 160 | U | 160 | < 140 | 140 | U | 140 | < 150 | 150 | U | 150 | |
| PCBs | PCB-1016 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1221 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1232 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1242 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1248 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1254 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1260 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| | PCB-1262 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 |
| PCB-1268 | 100 | 1,000 | < 37 | 37 | U | 37 | < 40 | 40 | U | 40 | < 39 | 39 | U | 39 | < 40 | 40 | U | 40 | < 35 | 35 | U | 35 | < 37 | 37 | U | 37 | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 5
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Pesticides and PCBs

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B19 | | | | | | | | | 15B22 | | | | | | | | | 15B23 | | | | | | | | |
|--------------------|---|--|-----------------|-----|------|-----------------|--------|-----|-----------------|-----|--------|-----------------|------|-----|-----------------|-----|------|-----------------|--------|-----|------------|-----|--------|-----|------|-----|--|--|--|
| | | | 11/20/2015 | | | | | | | | | 11/20/2015 | | | | | | | | | 11/20/2015 | | | | | | | | |
| | | | (0-2') µg/Kg | | | (6-8') µg/Kg | | | (0-2') µg/Kg | | | (6-8') µg/Kg | | | (2-4') µg/Kg | | | (6-8') µg/Kg | | | | | | | | | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | | | |
| 4,4' -DDD | 3.3 | 13,000 | < 2.2 | 2.2 | U | 2.2 | < 2.2 | 2.2 | U | 2.2 | < 14 | 14 | U | 14 | < 3.1 | 3.1 | U | 3.1 | < 2.2 | 2.2 | U | 2.2 | < 3.0 | 3.0 | U | 3.0 | | | |
| 4,4' -DDE | 3.3 | 8,900 | < 2.2 | 2.2 | U | 2.2 | < 2.2 | 2.2 | U | 2.2 | 130 | 14 | | 14 | < 3.1 | 3.1 | U | 3.1 | < 2.2 | 2.2 | U | 2.2 | < 3.0 | 3.0 | U | 3.0 | | | |
| 4,4' -DDT | 3.3 | 7,900 | < 2.2 | 2.2 | U | 2.2 | < 2.2 | 2.2 | U | 2.2 | 280 | 14 | | 14 | < 3.1 | 3.1 | U | 3.1 | < 2.2 | 2.2 | U | 2.2 | < 3.0 | 3.0 | U | 3.0 | | | |
| a-BHC | 20 | 480 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 12 | 12 | U | 12 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| a-Chlordane | 94 | 4,200 | < 3.6 | 3.6 | U | 3.6 | < 3.7 | 3.7 | U | 3.7 | 300 | 23 | | 23 | < 5.2 | 5.2 | U | 5.2 | < 3.7 | 3.7 | U | 3.7 | < 4.9 | 4.9 | U | 4.9 | | | |
| Aldrin | 5 | 97 | < 3.6 | 3.6 | U | 3.6 | < 3.7 | 3.7 | U | 3.7 | < 12 | 12 | U | 12 | < 2.6 | 2.6 | U | 2.6 | < 3.7 | 3.7 | U | 3.7 | < 4.9 | 4.9 | U | 4.9 | | | |
| b-BHC | 36 | 360 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 12 | 12 | U | 12 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Chlordane | | | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | 1,600 | 230 | | 230 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| d-BHC | 40 | 100,000 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 23 | 23 | U | 23 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Dieldrin | 5 | 200 | < 3.6 | 3.6 | U | 3.6 | < 3.7 | 3.7 | U | 3.7 | < 23 | 23 | U | 23 | < 1.6 | 1.6 | U | 1.6 | < 3.7 | 3.7 | U | 3.7 | < 4.9 | 4.9 | U | 4.9 | | | |
| Endosulfan I | 2,400 | 24,000 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Endosulfan II | 2,400 | 24,000 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Endosulfan sulfate | 2,400 | 24,000 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Endrin | 14 | 11,000 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 23 | 23 | U | 23 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Endrin aldehyde | | | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Endrin ketone | | | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| g-BHC | | | < 1.4 | 1.4 | U | 1.4 | < 1.5 | 1.5 | U | 1.5 | < 9.2 | 9.2 | U | 9.2 | < 2.1 | 2.1 | U | 2.1 | < 1.5 | 1.5 | U | 1.5 | < 2.0 | 2.0 | U | 2.0 | | | |
| g-Chlordane | | | < 3.6 | 3.6 | U | 3.6 | < 3.7 | 3.7 | U | 3.7 | 250 | 23 | | 23 | < 5.2 | 5.2 | U | 5.2 | < 3.7 | 3.7 | U | 3.7 | < 4.9 | 4.9 | U | 4.9 | | | |
| Heptachlor | 42 | 2,100 | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 23 | 23 | U | 23 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Heptachlor epoxide | | | < 7.2 | 7.2 | U | 7.2 | < 7.4 | 7.4 | U | 7.4 | < 46 | 46 | U | 46 | < 10 | 10 | U | 10 | < 7.3 | 7.3 | U | 7.3 | < 9.9 | 9.9 | U | 9.9 | | | |
| Methoxychlor | | | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 230 | 230 | U | 230 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| Toxaphene | | | < 140 | 140 | U | 140 | < 150 | 150 | U | 150 | < 920 | 920 | U | 920 | < 210 | 210 | U | 210 | < 150 | 150 | U | 150 | < 200 | 200 | U | 200 | | | |
| PCB-1016 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1221 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1232 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1242 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1248 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1254 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1260 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | 210 | 46 | | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1262 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |
| PCB-1268 | 100 | 1,000 | < 36 | 36 | U | 36 | < 37 | 37 | U | 37 | < 46 | 46 | U | 46 | < 52 | 52 | U | 52 | < 37 | 37 | U | 37 | < 49 | 49 | U | 49 | | | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 5
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
Pesticides and PCBs

| | COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | Duplicate 11/20/2015 | | | | Duplicate 11/23/2015 | | | |
|--------------------|----------|---|--|-------------------------|-----------|------|--------|-------------------------|-----|------|-----|
| | | | | µg/Kg | | | | µg/Kg | | | |
| | | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| | | | | Pesticides | 4,4' -DDD | 3.3 | 13,000 | 8 | 2.3 | | 2.3 |
| 4,4' -DDE | 3.3 | 8,900 | 22 | | 2.3 | | 2.3 | < 2.2 | 2.2 | U | 2.2 |
| 4,4' -DDT | 3.3 | 7,900 | 32 | | 2.3 | | 2.3 | 3.6 | 2.2 | | 2.2 |
| a-BHC | 20 | 480 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| a-Chlordane | 94 | 4,200 | < 11 | | 11 | U | 11 | < 3.6 | 3.6 | U | 3.6 |
| Aldrin | 5 | 97 | < 3.8 | | 3.8 | U | 3.8 | < 3.6 | 3.6 | U | 3.6 |
| b-BHC | 36 | 360 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Chlordane | | | < 38 | | 38 | U | 38 | < 36 | 36 | U | 36 |
| d-BHC | 40 | 100,000 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Dieldrin | 5 | 200 | < 3.8 | | 3.8 | U | 3.8 | < 3.6 | 3.6 | U | 3.6 |
| Endosulfan I | 2,400 | 24,000 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Endosulfan II | 2,400 | 24,000 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Endosulfan sulfate | 2,400 | 24,000 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Endrin | 14 | 11,000 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Endrin aldehyde | | | 220 | | 75 | | 75 | < 7.2 | 7.2 | U | 7.2 |
| Endrin ketone | | | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| g-BHC | | | < 1.5 | | 1.5 | U | 1.5 | < 1.4 | 1.4 | U | 1.4 |
| g-Chlordane | | | < 15 | | 15 | U | 15 | < 3.6 | 3.6 | U | 3.6 |
| Heptachlor | 42 | 2,100 | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Heptachlor epoxide | | | < 7.5 | | 7.5 | U | 7.5 | < 7.2 | 7.2 | U | 7.2 |
| Methoxychlor | | | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 | |
| Toxaphene | | | < 150 | 150 | U | 150 | < 140 | 140 | U | 140 | |
| PCBs | PCB-1016 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1221 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1232 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1242 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1248 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1254 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| | PCB-1260 | 100 | 1,000 | 110 | 38 | | 38 | < 36 | 36 | U | 36 |
| | PCB-1262 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 |
| PCB-1268 | 100 | 1,000 | < 38 | 38 | U | 38 | < 36 | 36 | U | 36 | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response to the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 6
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
TAL Metals

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B1 | | | | | | | | 15B3 | | | | | | | | 15B4 | | | | | | | |
|-----------|---|---|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|
| | | | 11/24/2015 | | | | | | | | 11/24/2015 | | | | | | | | 11/24/2015 | | | | | | | |
| | | | (0-2') mg/Kg | | | | (5-7') mg/Kg | | | | (0-2') mg/Kg | | | | (5-7') mg/Kg | | | | (0-2') mg/Kg | | | | (6-8') mg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| Aluminum | | | 4,390 | 360 | * | 72 | 10,300 | 39 | * | 7.8 | 5,030 | 34 | * | 6.9 | 3,260 | 41 | * | 8.2 | 6,980 | 35 | * | 7.0 | 8,910 | 41 | * | 8.1 |
| Antimony | | | < 1.8 | 1.8 | U | 1.8 | < 2.0 | 2.0 | U | 2.0 | < 1.7 | 1.7 | U | 1.7 | < 2.1 | 2.1 | U | 2.1 | < 1.8 | 1.8 | U | 1.8 | < 2.0 | 2.0 | U | 2.0 |
| Arsenic | 13 | 16 | 3.7 | 0.7 | | 0.72 | 7.7 | 0.8 | | 0.78 | 19.5 | 0.7 | | 0.69 | 33.5 | 0.8 | | 0.82 | 19.5 | 0.7 | | 0.70 | 2 | 0.8 | | 0.81 |
| Barium | 350 | 350 | 38 | 0.7 | | 0.36 | 33.9 | 0.8 | | 0.39 | 43.4 | 0.7 | | 0.34 | 21 | 0.8 | | 0.41 | 48.9 | 0.7 | | 0.35 | 33 | 0.8 | | 0.41 |
| Beryllium | 7.2 | 14 | 0.23 | 0.29 | B | 0.14 | 0.5 | 0.31 | | 0.16 | 0.18 | 0.28 | B | 0.14 | < 0.33 | 0.33 | U | 0.16 | 0.28 | 0.28 | B | 0.14 | | 0.33 | | 0.16 |
| Cadmium | 2.5 | 2.5 | 1.23 | 0.36 | | 0.14 | 1 | 0.39 | | 0.16 | 3.41 | 0.34 | | 0.14 | 4.01 | 0.41 | | 0.16 | 1.58 | 0.35 | | 0.14 | 1 | 0.41 | | 0.16 |
| Calcium | | | 20,500 | 360 | * | 330 | 26,800 | 39 | * | 36 | 3,000 | 3.4 | * | 3.2 | 5,910 | 4.1 | * | 3.8 | 7,990 | 3.5 | * | 3.2 | 10,400 | 4.1 | * | 3.7 |
| Chromium | 30 | 180 | 14.7 | 0.36 | | 0.36 | 22.5 | 0.39 | | 0.39 | 25.3 | 0.34 | | 0.34 | 69.6 | 0.41 | | 0.41 | 27.5 | 0.35 | | 0.35 | 20 | 0.41 | | 0.41 |
| Cobalt | | | 22.1 | 0.36 | | 0.36 | 34.3 | 0.39 | | 0.39 | 37.2 | 0.34 | | 0.34 | 92.7 | 0.41 | | 0.41 | 24.9 | 0.35 | | 0.35 | 19 | 0.41 | | 0.41 |
| Copper | 50 | 270 | 36.6 | 0.36 | | 0.36 | 33.8 | 0.39 | | 0.39 | 72.6 | 0.34 | | 0.34 | 194 | 4.1 | | 4.1 | 100 | 0.35 | | 0.35 | 21 | 0.41 | | 0.41 |
| Iron | | | 73,700 | 360 | * | 360 | 69,000 | 39 | * | 39 | 193,000 | 340 | * | 340 | 222,000 | 410 | * | 410 | 97,500 | 350 | * | 350 | 47,700 | 41 | * | 41 |
| Lead | 63 | 400 | 76 | 0.7 | * | 0.36 | 60.8 | 0.8 | * | 0.39 | 282 | 6.9 | * | 3.4 | 16.1 | 0.8 | * | 0.41 | 2,280 | 70 | * | 35 | 38 | 0.8 | * | 0.41 |
| Magnesium | | | 11,300 | 360 | * | 360 | 19,600 | 39 | * | 39 | 1,110 | 3.4 | * | 3.4 | 2,250 | 4.1 | * | 4.1 | 3,660 | 3.5 | * | 3.5 | 6,310 | 4.1 | * | 4.1 |
| Manganese | 1,600 | 2,000 | 522 | 36 | * | 36 | 970 | 3.9 | * | 3.9 | 1,030 | 3.4 | * | 3.4 | 2,750 | 4.1 | * | 4.1 | 1,020 | 3.5 | * | 3.5 | 741 | 4.1 | * | 4.1 |
| Mercury | 0.18 | 0.81 | 0.04 | 0.03 | | 0.02 | 0.04 | 0.03 | | 0.02 | 0.31 | 0.03 | | 0.02 | < 0.03 | 0.03 | U | 0.02 | 0.92 | 0.03 | | 0.02 | < 0.03 | 0.03 | U | 0.02 |
| Nickel | 30 | 140 | 26.5 | 0.36 | | 0.36 | 36 | 0.39 | | 0.39 | 36.8 | 0.34 | | 0.34 | 143 | 0.41 | | 0.41 | 35.7 | 0.35 | | 0.35 | 23 | 0.41 | | 0.41 |
| Potassium | | | 766 | 7 | | 2.8 | 3,160 | 8 | | 3.1 | 647 | 7 | | 2.7 | 376 | 8 | | 3.2 | 1,090 | 7 | | 2.7 | 1,750 | 8 | | 3.2 |
| Selenium | 3.9 | 36 | < 1.4 | 1.4 | U | 1.2 | < 1.6 | 1.6 | U | 1.3 | < 1.4 | 1.4 | U | 1.2 | < 1.6 | 1.6 | U | 1.4 | < 1.4 | 1.4 | U | 1.2 | < 1.6 | 1.6 | U | 1.4 |
| Silver | 2 | 36 | < 0.36 | 0.36 | U | 0.36 | < 0.39 | 0.39 | U | 0.39 | < 0.34 | 0.34 | U | 0.34 | < 0.41 | 0.41 | U | 0.41 | < 0.35 | 0.35 | U | 0.35 | < 0.41 | 0.41 | U | 0.41 |
| Sodium | | | 1,480 | 7 | | 3.1 | 3,150 | 8 | | 3.4 | 144 | 7 | | 3.0 | 889 | 8 | | 3.5 | 204 | 7 | | 3.0 | 810 | 8 | | 3.5 |
| Thallium | | | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 |
| Vanadium | | | 29.8 | 0.4 | | 0.36 | 36.9 | 0.4 | | 0.39 | 73.3 | 0.3 | | 0.34 | 92.8 | 0.4 | | 0.41 | 57.4 | 0.4 | | 0.35 | 26 | 0.4 | | 0.41 |
| Zinc | 109 | 2,200 | 29.6 | 0.7 | | 0.36 | 50.6 | 0.8 | | 0.39 | 89.1 | 0.7 | | 0.34 | < 8.2 | 8.2 | U | 4.1 | 95.5 | 0.7 | | 0.35 | 34.1 | 0.8 | | 0.41 |

Notes:

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 RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value
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| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
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| (*) | See report for comment. |

TABLE 6
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
TAL Metals

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B6 | | | | | | | | 15B9 | | | | | | | | 15B10 | | | | 15B11 | | | | | | | | |
|-----------|---|---|----------------|------|------|------|------------------|------|------|------|----------------|------|------|------|----------------|------|------|------|----------------|------|------|------|----------------|------|------|------|----------------|------|------|------|--------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/23/2015 | | | | | | | | |
| | | | (3-5) mg/Kg | | | | (10-12) mg/Kg | | | | (0-2) mg/Kg | | | | (6-8) mg/Kg | | | | (6-8) mg/Kg | | | | (0-2) mg/Kg | | | | (6-8) mg/Kg | | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result |
| Aluminum | | | 4,000 | 36 | | 7.2 | 15,500 | 51 | | 10 | 5,910 | 35 | | 7.0 | 9,780 | 42 | | 8.3 | 3,990 | 36 | | 7.3 | 6,420 | 36 | | 7.3 | 9,670 | 39 | | 7.7 | |
| Antimony | | | < 1.8 | 1.8 | U | 1.8 | < 2.6 | 2.6 | U | 2.6 | < 1.7 | 1.7 | U | 1.7 | < 2.1 | 2.1 | U | 2.1 | < 1.8 | 1.8 | U | 1.8 | < 1.8 | 1.8 | U | 1.8 | < 1.9 | 1.9 | U | 1.9 | |
| Arsenic | 13 | 16 | 11.9 | 0.7 | | 0.72 | 8 | 1.0 | | 1.0 | 9.8 | 0.7 | | 0.70 | 4 | 0.8 | | 0.83 | 22.8 | 0.7 | | 0.73 | 14.2 | 0.7 | | 0.73 | 2 | 0.8 | | 0.77 | |
| Barium | 350 | 350 | 53.3 | 0.7 | | 0.36 | 39 | 1.0 | | 0.51 | 44.7 | 0.7 | | 0.35 | 38 | 0.8 | | 0.42 | 30.4 | 0.7 | | 0.36 | 87.5 | 0.7 | | 0.36 | 27 | 0.8 | | 0.39 | |
| Beryllium | 7.2 | 14 | 0.18 | 0.29 | B | 0.14 | 1 | 0.41 | | 0.20 | 0.25 | 0.28 | B | 0.14 | 1 | 0.33 | | 0.17 | 0.19 | 0.29 | B | 0.15 | 0.4 | 0.29 | | 0.15 | | 0.31 | | 0.15 | |
| Cadmium | 2.5 | 2.5 | 3.03 | 0.36 | | 0.14 | 1 | 0.51 | | 0.20 | 2.98 | 0.35 | | 0.14 | | 0.42 | B | 0.17 | 3.55 | 0.36 | | 0.15 | 2.06 | 0.36 | | 0.15 | 2 | 0.39 | | 0.15 | |
| Calcium | | | 4,700 | 3.6 | | 3.3 | 2,070 | 5.1 | | 4.7 | 4,510 | 3.5 | | 3.2 | 15,800 | 42 | | 38 | 411 | 3.6 | | 3.3 | 13,500 | 36 | | 34 | 5,850 | 39 | | 36 | |
| Chromium | 30 | 180 | 28.2 | 0.36 | | 0.36 | 32.8 | 0.51 | | 0.51 | 25.7 | 0.35 | | 0.35 | 25 | 0.42 | | 0.42 | 45.9 | 0.36 | | 0.36 | 18.4 | 0.36 | | 0.36 | 23 | 0.39 | | 0.39 | |
| Cobalt | | | 48.7 | 0.36 | | 0.36 | 12 | 0.51 | | 0.51 | 47.7 | 0.35 | | 0.35 | 10 | 0.42 | | 0.42 | 54 | 0.36 | | 0.36 | 26.8 | 0.36 | | 0.36 | 17 | 0.39 | | 0.39 | |
| Copper | 50 | 270 | 64.6 | 0.36 | * | | 17 | 0.51 | * | 0.51 | 28 | 0.35 | * | 0.35 | 20 | 0.42 | * | 0.42 | 40.6 | 0.36 | * | 0.36 | 51.2 | 0.36 | | 0.36 | 31 | 0.39 | | 0.39 | |
| Iron | | | 139,000 | 360 | | 360 | 31,100 | 51 | | 51 | 148,000 | 350 | | 350 | 20,700 | 42 | | 42 | 164,000 | 360 | | 360 | 96,400 | 360 | | 360 | 82,500 | 39 | | 39 | |
| Lead | 63 | 400 | 42.2 | 0.7 | | 0.36 | 12 | 1.0 | | 0.51 | 24.5 | 0.7 | | 0.35 | 33 | 0.8 | | 0.42 | 10.3 | 0.7 | | 0.36 | 317 | 7.3 | N | 3.6 | 51 | 0.8 | N | 0.39 | |
| Magnesium | | | 518 | 3.6 | N | 3.6 | 7,430 | 5.1 | N | 5.1 | 2,150 | 3.5 | N | 3.5 | 11,400 | 42 | N | 42 | 774 | 3.6 | N | 3.6 | 1,460 | 36 | | 36 | 6,540 | 39 | | 39 | |
| Manganese | 1,600 | 2,000 | 1,660 | 36 | | 36 | 459 | 5.1 | | 5.1 | 1,150 | 3.5 | | 3.5 | 357 | 4.2 | | 4.2 | 484 | 3.6 | | 3.6 | 793 | 3.6 | | 3.6 | 2,450 | 39 | | 39 | |
| Mercury | 0.18 | 0.81 | 0.05 | 0.03 | * | 0.02 | < 0.04 | 0.04 | U* | 0.03 | 0.14 | 0.03 | * | 0.02 | < 0.03 | 0.03 | U* | 0.02 | < 0.03 | 0.03 | U* | 0.02 | 0.22 | 0.03 | * | 0.02 | < 0.03 | 0.03 | U* | 0.02 | |
| Nickel | 30 | 140 | 49.6 | 0.36 | | 0.36 | 26 | 0.51 | | 0.51 | 45.9 | 0.35 | | 0.35 | 20 | 0.42 | | 0.42 | 42.8 | 0.36 | | 0.36 | 33 | 0.36 | | 0.36 | 23 | 0.39 | | 0.39 | |
| Potassium | | | 334 | 7 | | 2.8 | 4,220 | 10 | | 4.0 | 613 | 7 | | 2.7 | 3,170 | 8 | | 3.2 | 537 | 7 | | 2.8 | 707 | 7 | | 2.8 | 916 | 77 | | 30 | |
| Selenium | 3.9 | 36 | < 1.4 | 1.4 | U | 1.2 | < 2.0 | 2.0 | U | 1.7 | < 1.4 | 1.4 | U | 1.2 | < 1.7 | 1.7 | U | 1.4 | < 1.5 | 1.5 | U | 1.2 | < 1.5 | 1.5 | U | 1.2 | < 1.5 | 1.5 | U | 1.3 | |
| Silver | 2 | 36 | < 0.36 | 0.36 | U | 0.36 | < 0.51 | 0.51 | U | 0.51 | < 0.35 | 0.35 | U | 0.35 | < 0.42 | 0.42 | U | 0.42 | < 0.36 | 0.36 | U | 0.36 | < 0.36 | 0.36 | U | 0.36 | < 0.39 | 0.39 | U | 0.39 | |
| Sodium | | | 149 | 7 | N | 3.1 | 587 | 10 | N | 4.4 | 214 | 7 | N | 3.0 | 2,730 | 8 | N | 3.6 | 1,010 | 7 | N | 3.1 | 238 | 7 | | 3.1 | 1,610 | 77 | | 33 | |
| Thallium | | | < 1.4 | 1.4 | U | 1.4 | < 2.0 | 2.0 | U | 2.0 | < 1.4 | 1.4 | U | 1.4 | < 1.7 | 1.7 | U | 1.7 | < 1.5 | 1.5 | U | 1.5 | < 1.5 | 1.5 | U | 1.5 | < 1.5 | 1.5 | U | 1.5 | |
| Vanadium | | | 71.6 | 0.4 | | 0.36 | 45 | 0.5 | | 0.51 | 80.8 | 0.3 | | 0.35 | 30 | 0.4 | | 0.42 | 127 | 0.4 | | 0.36 | 47.1 | 0.4 | | 0.36 | 34 | 0.4 | | 0.39 | |
| Zinc | 109 | 2,200 | < 7.2 | 7.2 | U | 3.6 | 67.6 | 1.0 | | 0.51 | 4.6 | 0.7 | | 0.35 | 45.7 | 0.8 | | 0.42 | < 7.3 | 7.3 | U | 3.6 | 123 | 7.3 | | 3.6 | 24.6 | 0.8 | | 0.39 | |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSO Guidance Value

| | |
|-----|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 6
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
TAL Metals

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B13 | | | | | | | | 15B18 | | | | | | | | 15B19 | | | | | | | |
|-----------|---|---|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | |
| | | | (0-2') mg/Kg | | | | (6-8') mg/Kg | | | | (0-2') mg/Kg | | | | (6-8') mg/Kg | | | | (0-2') mg/Kg | | | | (6-8') mg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| Aluminum | | | 5,130 | 36 | | 7.3 | 416 | 4.1 | | 0.83 | 5,800 | 35 | | 6.9 | 7,740 | 36 | | 7.1 | 11,500 | 39 | | 7.8 | 13,500 | 35 | | 6.9 |
| Antimony | | | < 1.8 | 1.8 | U | 1.8 | < 2.1 | 2.1 | U | 2.1 | < 1.7 | 1.7 | U | 1.7 | < 1.8 | 1.8 | U | 1.8 | < 2.0 | 2.0 | U | 2.0 | < 1.7 | 1.7 | U | 1.7 |
| Arsenic | 13 | 16 | 14.2 | 0.7 | | 0.73 | < 0.8 | 0.8 | U | 0.83 | 20.1 | 0.7 | | 0.69 | 3 | 0.7 | | 0.71 | 2.3 | 0.8 | | 0.78 | 1 | 0.7 | | 0.69 |
| Barium | 350 | 350 | 95.7 | 0.7 | | 0.36 | 3 | 0.8 | | 0.41 | 49.1 | 0.7 | | 0.35 | 15 | 0.7 | | 0.36 | 30.3 | 0.8 | | 0.39 | 10.7 | 0.7 | | 0.35 |
| Beryllium | 7.2 | 14 | 0.39 | 0.29 | | 0.15 | < 0.33 | 0.33 | U | 0.17 | 0.28 | 0.28 | B | 0.14 | | 0.28 | | 0.14 | 0.64 | 0.31 | | 0.16 | 0.53 | 0.28 | | 0.14 |
| Cadmium | 2.5 | 2.5 | 0.79 | 0.36 | | 0.15 | < 0.41 | 0.41 | U | 0.17 | 1.95 | 0.35 | | 0.14 | | 0.36 | B | 0.14 | 0.2 | 0.39 | B | 0.16 | < 0.35 | 0.35 | U | 0.14 |
| Calcium | | | 19,400 | 36 | | 33 | 1,680 | 4.1 | | 3.8 | 11,300 | 35 | | 32 | 7,920 | 3.6 | | 3.3 | 46,500 | 39 | | 36 | 1,450 | 3.5 | | 3.2 |
| Chromium | 30 | 180 | 15.3 | 0.36 | | 0.36 | 2 | 0.41 | | 0.41 | 17.2 | 0.35 | | 0.35 | 19 | 0.36 | | 0.36 | 19.7 | 0.39 | | 0.39 | 28.7 | 0.35 | | 0.35 |
| Cobalt | | | 8.28 | 0.36 | | 0.36 | 1 | 0.41 | | 0.41 | 16.3 | 0.35 | | 0.35 | 8 | 0.36 | | 0.36 | 9.58 | 0.39 | | 0.39 | 9.31 | 0.35 | | 0.35 |
| Copper | 50 | 270 | 44.8 | 0.36 | | 0.36 | 3 | 0.41 | | 0.41 | 37.5 | 0.35 | * | 0.35 | 21 | 0.36 | * | 0.36 | 16.5 | 0.39 | * | 0.39 | 18 | 0.35 | * | 0.35 |
| Iron | | | 29,000 | 36 | | 36 | 1,100 | 4.1 | | 4.1 | 79,900 | 35 | | 35 | 13,600 | 36 | | 36 | 18,600 | 39 | | 39 | 16,300 | 35 | | 35 |
| Lead | 63 | 400 | 89.3 | 0.7 | | 0.36 | 6 | 0.8 | | 0.41 | 65.3 | 0.7 | | 0.35 | 5 | 0.7 | | 0.36 | 5.2 | 0.8 | | 0.39 | 4.8 | 0.7 | | 0.35 |
| Magnesium | | | 5,670 | 36 | | 36 | 486 | 4.1 | | 4.1 | 1,590 | 3.5 | N | 3.5 | 9,240 | 36 | N | 36 | 8,190 | 39 | N | 39 | 14,200 | 35 | N | 35 |
| Manganese | 1,600 | 2,000 | 216 | 3.6 | N | 3.6 | 62 | 0.41 | N | 0.41 | 938 | 3.5 | | 3.5 | 279 | 3.6 | | 3.6 | 435 | 3.9 | | 3.9 | 244 | 3.5 | | 3.5 |
| Mercury | 0.18 | 0.81 | 0.06 | 0.03 | | 0.02 | < 0.03 | 0.03 | U | 0.02 | 0.02 | 0.03 | B* | 0.02 | < 0.03 | 0.03 | U* | 0.02 | < 0.03 | 0.03 | U* | 0.02 | < 0.03 | 0.03 | U* | 0.02 |
| Nickel | 30 | 140 | 16.3 | 0.36 | | 0.36 | 2 | 0.41 | | 0.41 | 22 | 0.35 | | 0.35 | 13 | 0.36 | | 0.36 | 16.1 | 0.39 | | 0.39 | 16.4 | 0.35 | | 0.35 |
| Potassium | | | 829 | 7 | N | 2.8 | 153 | 8 | N | 3.2 | 1,130 | 7 | | 2.7 | 2,070 | 7 | | 2.8 | 1,030 | 8 | | 3.1 | 1,870 | 7 | | 2.7 |
| Selenium | 3.9 | 36 | < 1.5 | 1.5 | U | 1.2 | < 1.7 | 1.7 | U | 1.4 | < 1.4 | 1.4 | U | 1.2 | < 1.4 | 1.4 | U | 1.2 | < 1.6 | 1.6 | U | 1.3 | < 1.4 | 1.4 | U | 1.2 |
| Silver | 2 | 36 | < 0.36 | 0.36 | U | 0.36 | < 0.41 | 0.41 | U | 0.41 | < 0.35 | 0.35 | U | 0.35 | < 0.36 | 0.36 | U | 0.36 | < 0.39 | 0.39 | U | 0.39 | < 0.35 | 0.35 | U | 0.35 |
| Sodium | | | 893 | 7 | | 3.1 | 702 | 8 | | 3.5 | 792 | 7 | N | 3.0 | 963 | 7 | N | 3.1 | 1,330 | 8 | N | 3.4 | 1,170 | 7 | N | 3.0 |
| Thallium | | | < 1.5 | 1.5 | U | 1.5 | < 1.7 | 1.7 | U | 1.7 | < 1.4 | 1.4 | U | 1.4 | < 1.4 | 1.4 | U | 1.4 | < 1.6 | 1.6 | U | 1.6 | < 1.4 | 1.4 | U | 1.4 |
| Vanadium | | | 29.7 | 0.4 | | 0.36 | 2 | 0.4 | | 0.41 | 50.3 | 0.3 | | 0.35 | 26 | 0.4 | | 0.36 | 25.7 | 0.4 | | 0.39 | 36.8 | 0.3 | | 0.35 |
| Zinc | 109 | 2,200 | 100 | 0.7 | | 0.36 | 4.4 | 0.8 | | 0.41 | 49.6 | 0.7 | | 0.35 | 41.1 | 0.7 | | 0.36 | 37.4 | 0.8 | | 0.39 | 51.4 | 0.7 | | 0.35 |

Notes:

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 RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value
Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

| | |
|-----|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank. |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 6
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Analytical Results
TAL Metals

| COMPOUND | NYSDEC Part 375.6 Unrestricted Use Soil Cleanup Objectives* | NYDEC Part 375.6 Restricted Residential Soil Cleanup Objectives* | 15B22 | | | | | | | | 15B23 | | | | | | | | Duplicate | | | | Duplicate | | | |
|-----------|---|---|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|-----------------|------|------|------|---------------|------|------|------|----------------|------|------|------|
| | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | | | | | 11/20/2015 | | | | 11/23/2015 | | | |
| | | | (0-2') mg/Kg | | | | (6-8') mg/Kg | | | | (2-4') mg/Kg | | | | (6-8') mg/Kg | | | | mg/Kg | | | | mg/Kg | | | |
| | | | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL | Result | RL | Qual | MDL |
| Aluminum | | | 6,890 | 46 | | 9.3 | 9,890 | 48 | | 9.7 | 5,950 | 38 | | 7.5 | 17,100 | 49 | | 9.7 | 4,720 | 35 | | 7.0 | 4,910 | 34 | | 6.8 |
| Antimony | | | < 2.3 | 2.3 | U | 2.3 | < 2.4 | 2.4 | U | 2.4 | < 1.9 | 1.9 | U | 1.9 | < 2.4 | 2.4 | U | 2.4 | < 1.8 | 1.8 | U | 1.8 | < 1.7 | 1.7 | U | 1.7 |
| Arsenic | 13 | 16 | 14.2 | 0.9 | | 0.93 | 34.9 | 1.0 | | 0.97 | 9.2 | 0.8 | | 0.75 | 8.7 | 1.0 | | 0.97 | 6.7 | 0.7 | | 0.70 | 18.6 | 0.7 | | 0.68 |
| Barium | 350 | 350 | 537 | 0.9 | | 0.46 | 121 | 1.0 | | 0.48 | 35.8 | 0.8 | | 0.38 | 57.5 | 1.0 | | 0.49 | 104 | 0.7 | | 0.35 | 84 | 0.7 | | 0.34 |
| Beryllium | 7.2 | 14 | 0.25 | 0.37 | B | 0.19 | 1 | 0.39 | | 0.19 | 0.27 | 0.30 | B | 0.15 | 0.79 | 0.39 | | 0.19 | 0.27 | 0.28 | B | 0.14 | 0.24 | 0.27 | B | 0.14 |
| Cadmium | 2.5 | 2.5 | 4.96 | 0.46 | | 0.19 | 1 | 0.48 | | 0.19 | 2.54 | 0.38 | | 0.15 | 0.35 | 0.49 | B | 0.19 | 1.04 | 0.35 | | 0.14 | 2.33 | 0.34 | | 0.14 |
| Calcium | | | 18,700 | 46 | | 43 | 71,800 | 48 | | 45 | 2,490 | 3.8 | | 3.5 | 6,830 | 4.9 | | 4.5 | 32,400 | 35 | | 32 | 9,820 | 34 | | 31 |
| Chromium | 30 | 180 | 128 | 0.46 | | 0.46 | 39.5 | 0.48 | | 0.48 | 22.7 | 0.38 | | 0.38 | 35.4 | 0.49 | | 0.49 | 21.7 | 0.35 | | 0.35 | 21.4 | 0.34 | | 0.34 |
| Cobalt | | | 15.9 | 0.46 | | 0.46 | 14 | 0.48 | | 0.48 | 21.1 | 0.38 | | 0.38 | 12 | 0.49 | | 0.49 | 7.1 | 0.35 | | 0.35 | 28.4 | 0.34 | | 0.34 |
| Copper | 50 | 270 | 1340 | 4.6 | | 4.6 | 145 | 0.48 | | 0.48 | 13.6 | 0.38 | * | 0.38 | 14.3 | 0.49 | * | 0.49 | 75.8 | 0.35 | | 0.35 | 35.5 | 0.34 | | 0.34 |
| Iron | | | 71,300 | 46 | | 46 | 30,700 | 48 | | 48 | 122,000 | 380 | | 380 | 33,000 | 49 | | 49 | 25,900 | 35 | | 35 | 116,000 | 340 | | 340 |
| Lead | 63 | 400 | 1550 | 9.3 | | 4.6 | 255 | 9.7 | | 4.8 | 155 | 7.5 | | 3.8 | 18 | 1.0 | | 0.49 | 497 | 7.0 | | 3.5 | 476 | 6.8 | N | 3.4 |
| Magnesium | | | 3,630 | 4.6 | | 4.6 | 9,500 | 48 | | 48 | 1,580 | 3.8 | N | 3.8 | 9,100 | 49 | N | 49 | 11,500 | 35 | | 35 | 1,280 | 34 | | 34 |
| Manganese | 1,600 | 2,000 | 992 | 4.6 | N | 4.6 | 382 | 4.8 | N | 4.8 | 853 | 3.8 | | 3.8 | 407 | 4.9 | | 4.9 | 284 | 3.5 | N | 3.5 | 1,950 | 34 | | 34 |
| Mercury | 0.18 | 0.81 | 0.34 | 0.04 | | 0.02 | 2.85 | 0.18 | | 0.11 | < 0.03 | 0.03 | U* | 0.02 | 0.03 | 0.04 | B* | 0.02 | 0.09 | 0.03 | | 0.02 | 0.21 | 0.03 | * | 0.02 |
| Nickel | 30 | 140 | 131 | 0.46 | | 0.46 | 28 | 0.48 | | 0.48 | 24.9 | 0.38 | | 0.38 | 23.8 | 0.49 | | 0.49 | 22.2 | 0.35 | | 0.35 | 38.6 | 0.34 | | 0.34 |
| Potassium | | | 1,640 | 9 | N | 3.6 | 4,470 | 10 | N | 3.8 | 674 | 8 | | 2.9 | 4,990 | 97 | | 38 | 993 | 7 | N | 2.7 | 518 | 7 | | 2.6 |
| Selenium | 3.9 | 36 | < 1.9 | 1.9 | U | 1.6 | < 1.9 | 1.9 | U | 1.6 | < 1.5 | 1.5 | U | 1.3 | < 1.9 | 1.9 | U | 1.7 | < 1.4 | 1.4 | U | 1.2 | < 1.4 | 1.4 | U | 1.2 |
| Silver | 2 | 36 | 3.09 | 0.46 | | 0.46 | 2 | 0.48 | | 0.48 | < 0.38 | 0.38 | U | 0.38 | < 0.49 | 0.49 | U | 0.49 | < 0.35 | 0.35 | U | 0.35 | < 0.34 | 0.34 | U | 0.34 |
| Sodium | | | 1,350 | 9 | | 4.0 | 4,140 | 10 | | 4.2 | 165 | 8 | N | 3.2 | 2,460 | 10 | N | 4.2 | 885 | 7 | | 3.0 | 274 | 7 | | 2.9 |
| Thallium | | | < 1.9 | 1.9 | U | 1.9 | < 1.9 | 1.9 | U | 1.9 | < 1.5 | 1.5 | U | 1.5 | < 1.9 | 1.9 | U | 1.9 | < 1.4 | 1.4 | U | 1.4 | < 1.4 | 1.4 | U | 1.4 |
| Vanadium | | | 53.3 | 0.5 | | 0.46 | 36 | 0.5 | | 0.48 | 66.1 | 0.4 | | 0.38 | 42.7 | 0.5 | | 0.49 | 31.6 | 0.4 | | 0.35 | 55 | 0.3 | | 0.34 |
| Zinc | 109 | 2,200 | 2840 | 93 | | 46 | 223 | 9.7 | | 4.8 | 22.9 | 0.8 | | 0.38 | 69.5 | 1.0 | | 0.49 | 225 | 7.0 | | 3.5 | 236 | 6.8 | | 3.4 |

Notes:

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Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

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| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 7
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Parameters Detected Above Track 1 Soil Cleanup Objectives

| COMPOUND | Range in Exceedances | Frequency of Detection | 15B1 | | 15B3 | | 15B4 | | 15B6 | | 15B8 | 15B9 | 15B10 |
|--------------------------------|----------------------|------------------------|------------|--------|------------|--------|------------|--------|------------|----------|------------|------------|------------|
| | | | 11/24/2015 | | 11/24/2015 | | 11/24/2015 | | 11/20/2015 | | 11/20/2015 | 11/20/2015 | 11/20/2015 |
| | | | (0-2') | (5-7') | (0-2') | (5-7') | (0-2') | (6-8') | (3-5') | (10-12') | (6-8') | (0-2') | (6-8') |
| <i>Sample Results in ug/kg</i> | | | | | | | | | | | | | |
| Acetone | 72-1500 | 5 | - | - | - | - | - | - | - | - | 110 | - | - |
| Benzene | 61-140 | 3 | - | - | - | 65 | - | 140 | - | - | - | - | - |
| <i>Sample Results in ug/kg</i> | | | | | | | | | | | | | |
| Benz(a)anthracene | 1300-1600 | 3 | - | - | - | - | 1,600 | - | - | - | - | - | - |
| Benzo(a)pyrene | 1100-1300 | 3 | - | - | - | - | 1,300 | - | - | - | - | - | - |
| Benzo(b)fluoranthene | 1100-1400 | 2 | - | - | - | - | 1,100 | - | - | - | - | - | - |
| Benzo(k)fluoranthene | 1100 | 1 | - | - | - | - | 1,100 | - | - | - | - | - | - |
| Chrysene | 1200-1600 | 3 | - | - | - | - | 1,500 | - | - | - | - | - | - |
| Indeno(1,2,3-cd)pyrene | 640-820 | 3 | 680 | - | - | - | 820 | - | - | - | - | - | - |
| <i>Sample Results in mg/kg</i> | | | | | | | | | | | | | |
| 4,4' -DDD | 8 | 1 | - | - | - | - | - | - | - | - | - | - | - |
| 4,4' -DDE | 7.9-130 | 3 | - | - | - | - | - | - | - | - | - | - | - |
| 4,4' -DDT | 3.6-280 | 5 | - | - | - | - | - | - | - | - | - | - | - |
| a-Chlordane | 300 | 1 | - | - | - | - | - | - | - | - | - | - | - |
| PCB-1260 | 110-210 | 2 | - | - | - | - | - | - | - | - | - | - | - |
| <i>Sample Results in mg/kg</i> | | | | | | | | | | | | | |
| Arsenic | 14.2-34.9 | 10 | - | - | 19.5 | 33.5 | 19.5 | - | - | - | - | - | 22.8 |
| Barium | 537 | 1 | - | - | - | - | - | - | - | - | - | - | - |
| Cadmium | 2.54-4.96 | 7 | - | - | 3.41 | 4.01 | - | - | 3.03 | - | - | 2.98 | 3.55 |
| Chromium | 32.8-128 | 6 | - | - | - | 69.6 | - | - | - | 32.8 | - | - | 45.9 |
| Copper | 51.2-1340 | 8 | - | - | 72.6 | 194 | 100 | - | 64.6 | - | - | - | - |
| Lead | 65.3-2280 | 11 | 76 | - | 282 | - | 2,280 | - | - | - | - | - | - |
| Manganese | 1660-2750 | 4 | - | - | - | 2,750 | - | - | 1,660 | - | - | - | - |
| Mercury | 0.21-2.85 | 6 | - | - | 0.31 | - | 0.92 | - | - | - | - | - | - |
| Nickel | 33-131 | 10 | - | 36 | 36.8 | 143 | 35.7 | - | 49.6 | - | - | 45.9 | 42.8 |
| Silver | 3.09 | 1 | - | - | - | - | - | - | - | - | - | - | - |
| Zinc | 123-2840 | 5 | - | - | - | - | - | - | - | - | - | - | - |

Notes:

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RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

TABLE 7
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Parameters Detected Above Track 1 Soil Cleanup Objectives

| COMPOUND | Range in Exceedances | Frequency of Detection | 15B11 | | 15B12 | 15B13 | 15B18 | 15B21 | 15B22 | | 15B23 | | Duplicate | Duplicate |
|--------------------------------|----------------------|------------------------|------------|--------|------------|------------|------------|------------|------------|--------|------------|--------|------------|------------|
| | | | 11/23/2015 | | 11/23/2015 | 11/20/2015 | 11/20/2015 | 11/20/2015 | 11/20/2015 | | 11/20/2015 | | 11/20/2015 | 11/23/2015 |
| | | | (0-2') | (6-8') | (6-8') | (0-2') | (0-2') | (6-8') | (0-2') | (6-8') | (2-4') | (6-8') | | |
| <i>Sample Results in ug/kg</i> | | | | | | | | | | | | | | |
| Acetone | 72-1500 | 5 | - | - | 100 | - | - | 72 | - | 1,500 | 510 | - | - | - |
| Benzene | 61-140 | 3 | - | - | 61 | - | - | - | - | - | - | - | - | - |
| <i>Sample Results in ug/kg</i> | | | | | | | | | | | | | | |
| Benz(a)anthracene | 1300-1600 | 3 | - | - | - | - | - | - | - | - | 1,300 | - | 1,300 | - |
| Benzo(a)pyrene | 1100-1300 | 3 | - | - | - | - | - | - | - | 1,200 | - | - | 1,100 | - |
| Benzo(b)fluoranthene | 1100-1400 | 2 | - | - | - | - | - | - | - | - | - | - | 1,400 | - |
| Benzo(k)fluoranthene | 1100 | 1 | - | - | - | - | - | - | - | - | - | - | - | - |
| Chrysene | 1200-1600 | 3 | - | - | - | - | - | - | - | - | 1,200 | - | 1,600 | - |
| Indeno(1,2,3-cd)pyrene | 640-820 | 3 | - | - | - | - | - | - | - | - | - | - | 640 | - |
| <i>Sample Results in ug/kg</i> | | | | | | | | | | | | | | |
| 4,4' -DDD | 8 | 1 | - | - | - | - | - | - | - | - | - | - | 8 | - |
| 4,4' -DDE | 7.9-130 | 3 | - | - | - | 7.9 | - | - | 130 | - | - | - | 22 | - |
| 4,4' -DDT | 3.6-280 | 5 | 4.2 | - | - | 9.4 | - | - | 280 | - | - | - | 32 | 3.6 |
| a-Chlordane | 300 | 1 | - | - | - | - | - | - | 300 | - | - | - | - | - |
| PCB-1260 | 110-210 | 2 | - | - | - | - | - | - | 210 | - | - | - | 110 | - |
| <i>Sample Results in mg/kg</i> | | | | | | | | | | | | | | |
| Arsenic | 14.2-34.9 | 10 | 14.2 | - | - | 14.2 | 20.1 | - | 14.2 | 34.9 | - | - | - | 18.6 |
| Barium | 537 | 1 | - | - | - | - | - | - | 537 | - | - | - | - | - |
| Cadmium | 2.54-4.96 | 7 | - | - | - | - | - | - | 4.96 | - | 2.54 | - | - | - |
| Chromium | 32.8-128 | 6 | - | - | - | - | - | - | 128 | 39.5 | - | 35.4 | - | - |
| Copper | 51.2-1340 | 8 | 51.2 | - | - | - | - | - | 1340 | 145 | - | - | 75.8 | - |
| Lead | 65.3-2280 | 11 | 317 | - | - | 89.3 | 65.3 | - | 1550 | 255 | 155 | - | 497 | 476 |
| Manganese | 1660-2750 | 4 | - | 2,450 | - | - | - | - | - | - | - | - | - | 1,950 |
| Mercury | 0.21-2.85 | 6 | 0.22 | - | - | - | - | - | 0.34 | 2.85 | - | - | - | 0.21 |
| Nickel | 33-131 | 10 | 33 | - | - | - | - | - | 131 | - | - | - | - | 38.6 |
| Silver | 3.09 | 1 | - | - | - | - | - | - | 3.09 | - | - | - | - | - |
| Zinc | 123-2840 | 5 | 123 | - | - | - | - | - | 2840 | 223 | - | - | 225 | 236 |

Notes:

* - 6 NYCRR Part 375-6 Remedial Program Soil Cleanup Objectives

RL - Laboratory Reporting Limit

Bold/highlighted- Indicated exceedance of the NYSDEC UUSCO Guidance Value

Bold/highlighted- Indicated exceedance of the NYSDEC RRSCO Guidance Value

TABLE 8
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Ground Water Analytical Results
Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW1 | | | | MW2 | | | | MW3 | | | |
|----------------------------------|--|-----------|------|------|------|------------|------|------|------|-------------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,1-Trichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1,2,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,2-Trichloroethane | 1 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloropropene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichloropropane | 0.04 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dibromo-3-chloropropane | 0.04 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 |
| 1,2-Dibromoethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichloroethane | 0.6 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 |
| 1,2-Dichloropropane | 0.94 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3,5-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,4-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2,2-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Hexanone (Methyl Butyl Ketone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| 2-Isopropyltoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Methyl-2-Pentanone | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Acetone | 50 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Acrolein | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Acrylonitrile | 5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Benzene | 1 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 |
| Bromobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromochloromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromodichloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromoform | | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Bromomethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | 1.7 | 5.0 | J | 0.25 |
| Carbon Disulfide | 60 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Carbon tetrachloride | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Chlorobenzene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroform | 7 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloromethane | 5 | < 5.0 | 5.0 | U | 0.25 | 1.1 | 5.0 | J | 0.25 | 39 | 10 | D | 0.50 |
| cis-1,2-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| cis-1,3-Dichloropropene | | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| Dibromochloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dibromomethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dichlorodifluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Ethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 |
| Isopropylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| m&p-Xylenes | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Methyl Ethyl Ketone (2-Butanone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Methyl t-butyl ether (MTBE) | 10 | < 1.0 | 1.0 | U | 0.25 | 1.3 | 1.0 | | 0.25 | 0.27 | 1.0 | J | 0.25 |
| Methylene chloride | 5 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 |
| Naphthalene | 10 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| n-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| n-Propylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| o-Xylene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| p-Isopropyltoluene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| sec-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Styrene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| tert-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrachloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrahydrofuran (THF) | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Toluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| trans-1,2-Dichloroethene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| trans-1,3-Dichloropropene | 0.4 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| trans-1,4-dichloro-2-butene | 5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Trichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorofluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorotrifluoroethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Vinyl Chloride | 2 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 8
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Ground Water Analytical Results
Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW4 | | | | MW5 | | | | MW7 | | | |
|----------------------------------|--|------------|------|------|------|-------------|------|------|------|-----------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,1,1-Trichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| 1,1,2,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,1,2-Trichloroethane | 1 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 |
| 1,1-Dichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| 1,1-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,1-Dichloropropene | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2,3-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2,3-Trichloropropane | 0.04 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2,4-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2,4-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2-Dibromo-3-chloropropane | 0.04 | < 1.0 | 1.0 | U | 0.50 | < 2.0 | 2.0 | U | 1.0 | < 2.0 | 2.0 | U | 1.0 |
| 1,2-Dibromoethane | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,2-Dichloroethane | 0.6 | < 0.60 | 0.60 | U | 0.25 | < 0.6 | 0.6 | U | 0.50 | < 0.6 | 0.6 | U | 0.50 |
| 1,2-Dichloropropane | 0.94 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 |
| 1,3,5-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,3-Dichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,3-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 1,4-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 2,2-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 2-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 2-Hexanone (Methyl Butyl Ketone) | | < 2.5 | 2.5 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 2-Isopropyltoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 4-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| 4-Methyl-2-Pentanone | | < 2.5 | 2.5 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| Acetone | 50 | < 5.0 | 5.0 | U | 2.5 | 9.7 | 10 | JS | 5.0 | < 10 | 10 | U | 5.0 |
| Acrolein | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| Acrylonitrile | 5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| Benzene | 1 | < 0.70 | 0.70 | U | 0.25 | < 0.7 | 0.7 | U | 0.50 | < 0.7 | 0.7 | U | 0.50 |
| Bromobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Bromochloromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Bromodichloromethane | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Bromoform | | < 5.0 | 5.0 | U | 0.25 | < 10 | 10 | U | 0.50 | < 10 | 10 | U | 0.50 |
| Bromomethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| Carbon Disulfide | 60 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Carbon tetrachloride | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Chlorobenzene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| Chloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| Chloroform | 7 | < 5.0 | 5.0 | U | 0.25 | < 7.0 | 7.0 | U | 0.50 | < 7.0 | 7.0 | U | 0.50 |
| Chloromethane | 5 | 4.3 | 5.0 | J | 0.25 | 8.3 | 5.0 | | 0.50 | 2 | 5.0 | J | 0.50 |
| cis-1,2-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| cis-1,3-Dichloropropene | | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.50 | < 0.40 | 0.40 | U | 0.50 |
| Dibromochloromethane | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Dibromomethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Dichlorodifluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Ethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.40 | < 0.50 | 0.50 | U | 0.40 |
| Isopropylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| m&p-Xylenes | 5 | < 1.0 | 1.0 | U | 0.25 | 0.98 | 2.0 | J | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Methyl Ethyl Ketone (2-Butanone) | | < 2.5 | 2.5 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| Methyl t-butyl ether (MTBE) | 10 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Methylene chloride | 5 | < 3.0 | 3.0 | U | 1.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 |
| Naphthalene | 10 | < 1.0 | 1.0 | U | 1.0 | < 2.0 | 2.0 | U | 2.0 | < 2.0 | 2.0 | U | 2.0 |
| n-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| n-Propylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| o-Xylene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| p-Isopropyltoluene | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| sec-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Styrene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| tert-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Tetrachloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Tetrahydrofuran (THF) | | < 5.0 | 5.0 | U | 2.5 | < 10 | 10 | U | 5.0 | < 10 | 10 | U | 5.0 |
| Toluene | 5 | < 1.0 | 1.0 | U | 0.25 | 2.3 | 2.0 | | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| trans-1,2-Dichloroethene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.50 | < 5.0 | 5.0 | U | 0.50 |
| trans-1,3-Dichloropropene | 0.4 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.50 | < 0.40 | 0.40 | U | 0.50 |
| trans-1,4-dichloro-2-butene | 5 | < 2.5 | 2.5 | U | 2.5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| Trichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Trichlorofluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Trichlorotrifluoroethane | | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |
| Vinyl Chloride | 2 | < 1.0 | 1.0 | U | 0.25 | < 2.0 | 2.0 | U | 0.50 | < 2.0 | 2.0 | U | 0.50 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 8
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Ground Water Analytical Results
Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW8 | | | | MW9 | | | | MW10 | | | |
|----------------------------------|--|------------|------|------|------|-------------|------|------|------|-------------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,1-Trichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1,2,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,2-Trichloroethane | 1 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloropropene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichloropropane | 0.04 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dibromo-3-chloropropane | 0.04 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 |
| 1,2-Dibromoethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichloroethane | 0.6 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 |
| 1,2-Dichloropropane | 0.94 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3,5-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,4-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2,2-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Hexanone (Methyl Butyl Ketone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| 2-Isopropyltoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Methyl-2-Pentanone | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Acetone | 50 | < 5.0 | 5.0 | U | 2.5 | 4.8 | 5.0 | JS | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Acrolein | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Acrylonitrile | 5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Benzene | 1 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 |
| Bromobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromochloromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromodichloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromoform | | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Bromomethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Carbon Disulfide | 60 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Carbon tetrachloride | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Chlorobenzene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroform | 7 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloromethane | 5 | 1.2 | 5.0 | J | 0.25 | 3 | 5.0 | J | 0.25 | 1.5 | 5.0 | J | 0.25 |
| cis-1,2-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| cis-1,3-Dichloropropene | | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| Dibromochloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dibromomethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dichlorodifluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Ethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 |
| Isopropylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| m&p-Xylenes | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Methyl Ethyl Ketone (2-Butanone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Methyl t-butyl ether (MTBE) | 10 | < 1.0 | 1.0 | U | 0.25 | 0.61 | 1.0 | J | 0.25 | 0.33 | 1.0 | J | 0.25 |
| Methylene chloride | 5 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 |
| Naphthalene | 10 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| n-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| n-Propylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| o-Xylene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| p-Isopropyltoluene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| sec-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Styrene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| tert-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrachloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrahydrofuran (THF) | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Toluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| trans-1,2-Dichloroethene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| trans-1,3-Dichloropropene | 0.4 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| trans-1,4-dichloro-2-butene | 5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Trichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorofluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorotrifluoroethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Vinyl Chloride | 2 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 8
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Ground Water Analytical Results
Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW11D | | | | MW12D | | | | MW13D | | | |
|----------------------------------|--|------------|------|------|------|------------|------|------|------|-------------|------|------|------|
| | | 12/2/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,1-Trichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1,2,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,2-Trichloroethane | 1 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| 1,1-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloropropene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichloropropane | 0.04 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dibromo-3-chloropropane | 0.04 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 | < 1.0 | 1.0 | U | 0.50 |
| 1,2-Dibromoethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichloroethane | 0.6 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 | < 0.60 | 0.60 | U | 0.25 |
| 1,2-Dichloropropane | 0.94 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3,5-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichlorobenzene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 1,4-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2,2-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 2-Hexanone (Methyl Butyl Ketone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| 2-Isopropyltoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| 4-Methyl-2-Pentanone | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Acetone | 50 | 3 | 5.0 | JS | 2.5 | < 5.0 | 5.0 | U | 2.5 | 3.3 | 5.0 | JS | 2.5 |
| Acrolein | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Acrylonitrile | 5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Benzene | 1 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 | < 0.70 | 0.70 | U | 0.25 |
| Bromobenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromochloromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromodichloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Bromoform | | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Bromomethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | 0.67 | 5.0 | J | 0.25 |
| Carbon Disulfide | 60 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Carbon tetrachloride | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Chlorobenzene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroethane | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloroform | 7 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| Chloromethane | 5 | 0.3 | 5.0 | J | 0.25 | 2.4 | 5.0 | J | 0.25 | 16 | 5.0 | | 0.25 |
| cis-1,2-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| cis-1,3-Dichloropropene | | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| Dibromochloromethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dibromomethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Dichlorodifluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Ethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 | < 0.50 | 0.50 | U | 0.20 |
| Isopropylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| m&p-Xylenes | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Methyl Ethyl Ketone (2-Butanone) | | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Methyl t-butyl ether (MTBE) | 10 | 14 | 1.0 | | 0.25 | 100 | 10 | D | 2.5 | 5.1 | 1.0 | | 0.25 |
| Methylene chloride | 5 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 | < 3.0 | 3.0 | U | 1.0 |
| Naphthalene | 10 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| n-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| n-Propylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| o-Xylene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| p-Isopropyltoluene | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| sec-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Styrene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| tert-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrachloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Tetrahydrofuran (THF) | | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 | < 5.0 | 5.0 | U | 2.5 |
| Toluene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| trans-1,2-Dichloroethene | 5 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 | < 5.0 | 5.0 | U | 0.25 |
| trans-1,3-Dichloropropene | 0.4 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 | < 0.40 | 0.40 | U | 0.25 |
| trans-1,4-dichloro-2-butene | 5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 | < 2.5 | 2.5 | U | 2.5 |
| Trichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorofluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorotrifluoroethane | | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |
| Vinyl Chloride | 2 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 | < 1.0 | 1.0 | U | 0.25 |

Notes:
RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 8
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Ground Water Analytical Results
Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | Duplicate | | | |
|----------------------------------|--|------------|------|------|------|
| | | 12/1/2015 | | | |
| | | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,1-Trichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 |
| 1,1,2,2-Tetrachloroethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,1,2-Trichloroethane | 1 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloroethane | 5 | < 5.0 | 5.0 | U | 0.25 |
| 1,1-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,1-Dichloropropene | | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 |
| 1,2,3-Trichloropropane | 0.04 | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trichlorobenzene | | < 1.0 | 1.0 | U | 0.25 |
| 1,2,4-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dibromo-3-chloropropane | 0.04 | < 1.0 | 1.0 | U | 0.50 |
| 1,2-Dibromoethane | | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,2-Dichloroethane | 0.6 | < 0.60 | 0.60 | U | 0.25 |
| 1,2-Dichloropropane | 0.94 | < 1.0 | 1.0 | U | 0.25 |
| 1,3,5-Trimethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichlorobenzene | | < 1.0 | 1.0 | U | 0.25 |
| 1,3-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 |
| 1,4-Dichlorobenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 2,2-Dichloropropane | 5 | < 1.0 | 1.0 | U | 0.25 |
| 2-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 2-Hexanone (Methyl Butyl Ketone) | | < 2.5 | 2.5 | U | 2.5 |
| 2-Isopropyltoluene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 4-Chlorotoluene | 5 | < 1.0 | 1.0 | U | 0.25 |
| 4-Methyl-2-Pentanone | | < 2.5 | 2.5 | U | 2.5 |
| Acetone | 50 | < 5.0 | 5.0 | U | 2.5 |
| Acrolein | | < 5.0 | 5.0 | U | 2.5 |
| Acrylonitrile | 5 | < 5.0 | 5.0 | U | 2.5 |
| Benzene | 1 | < 0.70 | 0.70 | U | 0.25 |
| Bromobenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Bromochloromethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| Bromodichloromethane | | < 1.0 | 1.0 | U | 0.25 |
| Bromoform | | < 5.0 | 5.0 | U | 0.25 |
| Bromomethane | 5 | < 5.0 | 5.0 | U | 0.25 |
| Carbon Disulfide | 60 | < 1.0 | 1.0 | U | 0.25 |
| Carbon tetrachloride | 5 | < 1.0 | 1.0 | U | 0.25 |
| Chlorobenzene | 5 | < 5.0 | 5.0 | U | 0.25 |
| Chloroethane | 5 | < 5.0 | 5.0 | U | 0.25 |
| Chloroform | 7 | < 5.0 | 5.0 | U | 0.25 |
| Chloromethane | 5 | 15 | 5.0 | | 0.25 |
| cis-1,2-Dichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 |
| cis-1,3-Dichloropropene | | < 0.40 | 0.40 | U | 0.25 |
| Dibromochloromethane | | < 1.0 | 1.0 | U | 0.25 |
| Dibromomethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| Dichlorodifluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| Ethylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.20 |
| Isopropylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| m&p-Xylenes | 5 | < 1.0 | 1.0 | U | 0.25 |
| Methyl Ethyl Ketone (2-Butanone) | | < 2.5 | 2.5 | U | 2.5 |
| Methyl t-butyl ether (MTBE) | 10 | 5.1 | 1.0 | | 0.25 |
| Methylene chloride | 5 | < 3.0 | 3.0 | U | 1.0 |
| Naphthalene | 10 | < 1.0 | 1.0 | U | 1.0 |
| n-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| n-Propylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| o-Xylene | 5 | < 1.0 | 1.0 | U | 0.25 |
| p-Isopropyltoluene | | < 1.0 | 1.0 | U | 0.25 |
| sec-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Styrene | 5 | < 1.0 | 1.0 | U | 0.25 |
| tert-Butylbenzene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Tetrachloroethene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Tetrahydrofuran (THF) | | < 5.0 | 5.0 | U | 2.5 |
| Toluene | 5 | < 1.0 | 1.0 | U | 0.25 |
| trans-1,2-Dichloroethene | 5 | < 5.0 | 5.0 | U | 0.25 |
| trans-1,3-Dichloropropene | 0.4 | < 0.40 | 0.40 | U | 0.25 |
| trans-1,4-dichloro-2-butene | 5 | < 2.5 | 2.5 | U | 2.5 |
| Trichloroethene | 5 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorofluoromethane | 5 | < 1.0 | 1.0 | U | 0.25 |
| Trichlorotrifluoroethane | | < 1.0 | 1.0 | U | 0.25 |
| Vinyl Chloride | 2 | < 1.0 | 1.0 | U | 0.25 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 9
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Semi-Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW1 | | | | MW2 | | | | MW3 | | | | MW4 | | | |
|-------------------------------|---|-----------|------|------|------|-----------|------|------|------|-----------|------|------|------|-----------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/2/2015 | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,2,4-Trichlorobenzene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 1,2-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,2-Diphenylhydrazine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 1,3-Dichlorobenzene | 3 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,4-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,5-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,6-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dichlorophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dimethylphenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrophenol | 5 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 |
| 2,6-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 2-Chloronaphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| 2-Chlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Methylnaphthalene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 2-Methylphenol (o-cresol) | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 2-Nitrophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3&4-Methylphenol (m&p-cresol) | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3,3'-Dichlorobenzidine | 5 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 |
| 3-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 4,6-Dinitro-2-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Bromophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 4-Chloro-3-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Chloroaniline | 5 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 |
| 4-Chlorophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitroaniline | 5 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitrophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Acenaphthene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Acetophenone | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Aniline | 5 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 |
| Anthracene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Benzidine | 5 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 |
| Benzoic acid | | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 |
| Benzyl butyl phthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Bis(2-chloroethoxy)methane | 5 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Bis(2-chloroethyl)ether | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-chloroisopropyl)ether | | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Carbazole | | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 |
| Dibenzofuran | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Diethyl phthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Dimethylphthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Di-n-butylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Di-n-octylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Fluoranthene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Fluorene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| Hexachlorocyclopentadiene | 5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Isophorone | 50 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Naphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| N-Nitrosodimethylamine | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| N-Nitrosodi-n-propylamine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| N-Nitrosodiphenylamine | 50 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 |
| Phenol | 50 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Pyrene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 1,2,4,5-Tetrachlorobenzene | | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Acenaphthene | 20 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Acenaphthylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benz(a)anthracene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(a)pyrene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(b)fluoranthene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(ghi)perylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(k)fluoranthene | 0.002 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-ethylhexyl)phthalate | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Chrysene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Dibenz(a,h)anthracene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Hexachlorobenzene | 0.04 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Hexachloroethane | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Indeno(1,2,3-cd)pyrene | 0.002 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachloronitrobenzene | | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachlorophenol | 1 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 |
| Phenanthrene | 50 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pyridine | 50 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

The compound was analyzed for but not detected at or above the MDL.

U The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors.

B This compound was also present in the method blank

J The value is estimated. This flag is used
a) on form 1 when the compound is reported above the MDL, but below the PQL, and
b) on the Tentatively Identified Compounds (TIC) form for all compounds identified.

TABLE 9
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Semi-Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) ($\mu\text{g/L}$) | MW5 | | | | MW7 | | | | MW8 | | | | MW9 | | | |
|-------------------------------|--|-------------------|------|------|------|-------------------|------|------|------|-------------------|------|------|------|-------------------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | |
| | | $(\mu\text{g/L})$ | | | | $(\mu\text{g/L})$ | | | | $(\mu\text{g/L})$ | | | | $(\mu\text{g/L})$ | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,2,4-Trichlorobenzene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 1,2-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,2-Diphenylhydrazine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 1,3-Dichlorobenzene | 3 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,4-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,5-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,6-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dichlorophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dimethylphenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrophenol | 5 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 |
| 2,6-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 2-Chloronaphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| 2-Chlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Methylnaphthalene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 2-Methylphenol (o-cresol) | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 2-Nitrophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3&4-Methylphenol (m&p-cresol) | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3,3'-Dichlorobenzidine | 5 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 |
| 3-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 4,6-Dinitro-2-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Bromophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 4-Chloro-3-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Chloroaniline | 5 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 |
| 4-Chlorophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitroaniline | 5 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitrophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Acenaphthene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Acetophenone | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Aniline | 5 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 |
| Anthracene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Benzidine | 5 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 |
| Benzoic acid | | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 |
| Benzyl butyl phthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Bis(2-chloroethoxy)methane | 5 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Bis(2-chloroethyl)ether | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-chloroisopropyl)ether | | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Carbazole | | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 |
| Dibenzofuran | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Diethyl phthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Dimethylphthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Di-n-butylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Di-n-octylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Fluoranthene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Fluorene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| Hexachlorocyclopentadiene | 5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Isophorone | 50 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Naphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| N-Nitrosodimethylamine | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| N-Nitrosodi-n-propylamine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| N-Nitrosodiphenylamine | 50 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 |
| Phenol | 50 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Pyrene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 1,2,4,5-Tetrachlorobenzene | | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Acenaphthene | 20 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Acenaphthylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.05 | 0.02 | U | 0.02 |
| Benz(a)anthracene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.03 | 0.02 | U | 0.02 |
| Benzo(a)pyrene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.03 | 0.02 | U | 0.02 |
| Benzo(b)fluoranthene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.02 | 0.02 | U | 0.02 |
| Benzo(ghi)perylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.04 | 0.02 | U | 0.02 |
| Benzo(k)fluoranthene | 0.002 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-ethylhexyl)phthalate | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | 0.05 | 0.02 | U | 0.02 |
| Chrysene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Dibenz(a,h)anthracene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Hexachlorobenzene | 0.04 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Hexachloroethane | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Indeno(1,2,3-cd)pyrene | 0.002 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachloronitrobenzene | | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachlorophenol | 1 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 |
| Phenanthrene | 50 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pyridine | 50 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL |

TABLE 9
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Semi-Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | MW10 | | | | MW11D | | | | MW12D | | | | MW13D | | | |
|-------------------------------|---|-----------|------|------|------|-----------|------|------|------|-----------|------|------|------|-----------|------|------|------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,2,4-Trichlorobenzene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 1,2-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,2-Diphenylhydrazine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 1,3-Dichlorobenzene | 3 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 1,4-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,5-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,6-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dichlorophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dimethylphenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrophenol | 5 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 | < 5.0 | 5.0 | U | 2.0 |
| 2,6-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| 2-Chloronaphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| 2-Chlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Methylnaphthalene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 2-Methylphenol (o-cresol) | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 2-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 2-Nitrophenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3&4-Methylphenol (m&p-cresol) | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 3,3'-Dichlorobenzidine | 5 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 | < 5.0 | 5.0 | U | 2.4 |
| 3-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 | < 5.0 | 5.0 | U | 5.0 |
| 4,6-Dinitro-2-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Bromophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| 4-Chloro-3-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| 4-Chloroaniline | 5 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 | < 3.5 | 3.5 | U | 2.3 |
| 4-Chlorophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitroaniline | 5 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitrophenol | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Acenaphthene | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Acetophenone | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Aniline | 5 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 | < 3.5 | 3.5 | U | 5.0 |
| Anthracene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Benzidine | 5 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 | < 4.5 | 4.5 | U | 2.9 |
| Benzoic acid | | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 | < 25 | 25 | U | 10 |
| Benzyl butyl phthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Bis(2-chloroethoxy)methane | 5 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Bis(2-chloroethyl)ether | 1 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-chloroisopropyl)ether | | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Carbazole | | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 | < 25 | 25 | U | 3.8 |
| Dibenzofuran | | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Diethyl phthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Dimethylphthalate | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Di-n-butylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Di-n-octylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 | < 5.0 | 5.0 | U | 1.3 |
| Fluoranthene | 50 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| Fluorene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| Hexachlorocyclopentadiene | 5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 | < 5.0 | 5.0 | U | 1.5 |
| Isophorone | 50 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| Naphthalene | 10 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 | < 5.0 | 5.0 | U | 1.4 |
| N-Nitrosodimethylamine | | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| N-Nitrosodi-n-propylamine | | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 | < 5.0 | 5.0 | U | 1.6 |
| N-Nitrosodiphenylamine | 50 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 | < 5.0 | 5.0 | U | 1.9 |
| Phenol | 50 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Pyrene | 50 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 | < 5.0 | 5.0 | U | 1.7 |
| 1,2,4,5-Tetrachlorobenzene | | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Acenaphthene | 20 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Acenaphthylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benz(a)anthracene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(a)pyrene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(b)fluoranthene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(ghi)perylene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(k)fluoranthene | 0.002 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-ethylhexyl)phthalate | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Chrysene | 0.002 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Dibenz(a,h)anthracene | | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Hexachlorobenzene | 0.04 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 | < 0.40 | 0.40 | U | 0.40 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 | < 0.50 | 0.50 | U | 0.50 |
| Hexachloroethane | 5 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 | < 0.02 | 0.02 | U | 0.02 |
| Indeno(1,2,3-cd)pyrene | 0.002 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachloronitrobenzene | | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pentachlorophenol | 1 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 | < 0.80 | 0.80 | U | 0.80 |
| Phenanthrene | 50 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Pyridine | 50 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 | < 10 | 10 | U | 1.2 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 9
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Semi-Volatile Organic Compounds

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (µg/L) | Duplicate | | | |
|-------------------------------|---|-----------|------|------|------|
| | | 12/1/2015 | | | |
| | | (µg/L) | | | |
| | | Results | RL | Qual | MDL |
| 1,2,4-Trichlorobenzene | | < 5.0 | 5.0 | U | 1.5 |
| 1,2-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 |
| 1,2-Diphenylhydrazine | | < 5.0 | 5.0 | U | 1.6 |
| 1,3-Dichlorobenzene | 3 | < 1.0 | 1.0 | U | 1.0 |
| 1,4-Dichlorobenzene | | < 1.0 | 1.0 | U | 1.0 |
| 2,4,5-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 2,4,6-Trichlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dichlorophenol | | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dimethylphenol | | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrophenol | 5 | < 1.0 | 1.0 | U | 1.0 |
| 2,4-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 2.0 |
| 2,6-Dinitrotoluene | 5 | < 5.0 | 5.0 | U | 1.6 |
| 2-Chloronaphthalene | 10 | < 5.0 | 5.0 | U | 1.4 |
| 2-Chlorophenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 2-Methylnaphthalene | | < 5.0 | 5.0 | U | 1.5 |
| 2-Methylphenol (o-cresol) | 1 | < 1.0 | 1.0 | U | 1.0 |
| 2-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 |
| 2-Nitrophenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 3&4-Methylphenol (m&p-cresol) | | < 1.0 | 1.0 | U | 1.0 |
| 3,3'-Dichlorobenzidine | 5 | < 5.0 | 5.0 | U | 2.4 |
| 3-Nitroaniline | 5 | < 5.0 | 5.0 | U | 5.0 |
| 4,6-Dinitro-2-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 4-Bromophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.5 |
| 4-Chloro-3-methylphenol | 1 | < 1.0 | 1.0 | U | 1.0 |
| 4-Chloroaniline | 5 | < 3.5 | 3.5 | U | 2.3 |
| 4-Chlorophenyl phenyl ether | | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitroaniline | 5 | < 5.0 | 5.0 | U | 1.7 |
| 4-Nitrophenol | | < 1.0 | 1.0 | U | 1.0 |
| Acenaphthene | | < 5.0 | 5.0 | U | 1.5 |
| Acetophenone | | < 5.0 | 5.0 | U | 1.6 |
| Aniline | 5 | < 3.5 | 3.5 | U | 5.0 |
| Anthracene | 50 | < 5.0 | 5.0 | U | 1.6 |
| Benzidine | 5 | < 4.5 | 4.5 | U | 2.9 |
| Benzoic acid | | < 25 | 25 | U | 10 |
| Benzyl butyl phthalate | 50 | < 5.0 | 5.0 | U | 1.3 |
| Bis(2-chloroethoxy)methane | 5 | < 5.0 | 5.0 | U | 1.4 |
| Bis(2-chloroethyl)ether | 1 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-chloroisopropyl)ether | | < 5.0 | 5.0 | U | 1.4 |
| Carbazole | | < 25 | 25 | U | 3.8 |
| Dibenzofuran | | < 5.0 | 5.0 | U | 1.5 |
| Diethyl phthalate | 50 | < 5.0 | 5.0 | U | 1.6 |
| Dimethylphthalate | 50 | < 5.0 | 5.0 | U | 1.6 |
| Di-n-butylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 |
| Di-n-octylphthalate | 50 | < 5.0 | 5.0 | U | 1.3 |
| Fluoranthene | 50 | < 5.0 | 5.0 | U | 1.6 |
| Fluorene | 50 | < 5.0 | 5.0 | U | 1.7 |
| Hexachlorocyclopentadiene | 5 | < 5.0 | 5.0 | U | 1.5 |
| Isophorone | 50 | < 5.0 | 5.0 | U | 1.4 |
| Naphthalene | 10 | < 5.0 | 5.0 | U | 1.4 |
| N-Nitrosodimethylamine | | < 1.0 | 1.0 | U | 1.0 |
| N-Nitrosodi-n-propylamine | | < 5.0 | 5.0 | U | 1.6 |
| N-Nitrosodiphenylamine | 50 | < 5.0 | 5.0 | U | 1.9 |
| Phenol | 50 | < 1.0 | 1.0 | U | 1.0 |
| Pyrene | 50 | < 5.0 | 5.0 | U | 1.7 |
| 1,2,4,5-Tetrachlorobenzene | | < 0.50 | 0.50 | U | 0.50 |
| Acenaphthene | 20 | < 0.10 | 0.10 | U | 0.10 |
| Acenaphthylene | | < 0.02 | 0.02 | U | 0.02 |
| Benz(a)anthracene | 0.002 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(a)pyrene | | < 0.02 | 0.02 | U | 0.02 |
| Benzo(b)fluoranthene | 0.002 | < 0.02 | 0.02 | U | 0.02 |
| Benzo(ghi)perylene | | < 0.02 | 0.02 | U | 0.02 |
| Benzo(k)fluoranthene | 0.002 | < 1.0 | 1.0 | U | 1.0 |
| Bis(2-ethylhexyl)phthalate | 5 | < 0.02 | 0.02 | U | 0.02 |
| Chrysene | 0.002 | < 0.02 | 0.02 | U | 0.02 |
| Dibenz(a,h)anthracene | | < 0.02 | 0.02 | U | 0.02 |
| Hexachlorobenzene | 0.04 | < 0.40 | 0.40 | U | 0.40 |
| Hexachlorobutadiene | 0.5 | < 0.50 | 0.50 | U | 0.50 |
| Hexachloroethane | 5 | < 0.02 | 0.02 | U | 0.02 |
| Indeno(1,2,3-cd)pyrene | 0.002 | < 0.10 | 0.10 | U | 0.10 |
| Pentachloronitrobenzene | | < 0.10 | 0.10 | U | 0.10 |
| Pentachlorophenol | 1 | < 0.80 | 0.80 | U | 0.80 |
| Phenanthrene | 50 | < 0.10 | 0.10 | U | 0.10 |
| Pyridine | 50 | < 10 | 10 | U | 1.2 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|---|--|
| | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |

TABLE 10
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Pesticides and PCBs

| Compound | NYSDEC Ambient Water Quality Standards (AQWS) (µg/L) | MW1 | | | | MW2 | | | | MW3 | | | | MW4 | | | | MW5 | | | | MW7 | | | | MW8 | | | | | | | |
|--------------------|---|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|---------|-------|------|-------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| PCB-1016 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1221 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1232 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1242 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1248 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1254 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1260 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1262 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1268 | 0.09 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| 4,4-DDD | 0.3 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| 4,4-DDE | 0.2 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| 4,4-DDT | 0.11 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| a-BHC | 0.94 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| a-Chlordane | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Alachlor | | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 |
| Aldrin | | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| b-BHC | 0.04 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| Chlordane | 0.05 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| d-BHC | 0.04 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| Dieldrin | 0.004 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Endosulfan I | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endosulfan II | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endosulfan Sulfate | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin aldehyde | 5 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin ketone | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| gamma-BHC | 0.05 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| g-Chlordane | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Heptachlor | 0.04 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Heptachlor epoxide | 0.03 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Methoxychlor | 35 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Toxaphene | | < 0.20 | 0.20 | U | 0.20 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.20 | 0.20 | U | 0.20 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.20 | 0.20 | U | 0.20 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

| | |
|-----|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response to the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 10
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Pesticides and PCBs

| Compound | NYSDEC Ambient Water Quality Standards (AQWS) (µg/L) | MW9 | | | | MW10 | | | | MW11D | | | | MW12D | | | | MW13D | | | | Duplicate | | | | | | | |
|--------------------|--|------------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|-----------|-------|------|-------|---------|-------|------|-------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | | | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| PCB-1016 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1221 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1232 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1242 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1248 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1254 | 0.09 | 1.3 | 0.25 | | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1260 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1262 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| PCB-1268 | 0.09 | < 0.25 | 0.25 | U | 0.25 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| 4,4-DDD | 0.3 | < 0.025 | 0.025 | U | 0.025 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| 4,4-DDE | 0.2 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| 4,4-DDT | 0.11 | < 0.15 | 0.15 | U | 0.15 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| a-BHC | 0.94 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| a-Chlordane | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Alachlor | | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 | < 0.075 | 0.075 | U | 0.075 |
| Aldrin | | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| b-BHC | 0.04 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| Chlordane | 0.05 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 | < 0.050 | 0.050 | U | 0.050 |
| d-BHC | 0.04 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| Dieldrin | 0.004 | < 0.050 | 0.050 | U | 0.050 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Endosulfan I | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endosulfan II | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endosulfan Sulfate | | < 0.15 | 0.15 | U | 0.15 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin | | < 0.15 | 0.15 | U | 0.15 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin aldehyde | 5 | < 0.020 | 0.020 | U | 0.020 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Endrin ketone | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| gamma-BHC | 0.05 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 | < 0.005 | 0.005 | U | 0.005 |
| g-Chlordane | | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Heptachlor | 0.04 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Heptachlor epoxide | 0.03 | < 0.015 | 0.015 | U | 0.015 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 | < 0.010 | 0.010 | U | 0.010 |
| Methoxychlor | 35 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 | < 0.10 | 0.10 | U | 0.10 |
| Toxaphene | | < 0.20 | 0.20 | U | 0.20 | < 0.20 | 0.20 | U | 0.20 | < 0.25 | 0.25 | U | 0.25 | < 0.20 | 0.20 | U | 0.20 | < 0.20 | 0.20 | U | 0.20 | < 0.20 | 0.20 | U | 0.20 | < 0.20 | 0.20 | U | 0.20 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater §

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 11
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Dissolved Metals

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (mg/L) | MW1 | | | | MW2 | | | | MW3 | | | | MW4 | | | | MW5 | | | | MW7 | | | | MW8 | | | |
|-----------|---|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| Aluminum | NS | 0.005 | 0.011 | B | 0.005 | < 0.011 | 0.011 | U | 0.005 | 0.008 | 0.011 | B | 0.005 | < 0.011 | 0.011 | U | 0.005 | 0.009 | 0.011 | B | 0.005 | 0.006 | 0.011 | B | 0.005 | 0.007 | 0.011 | B | 0.005 |
| Antimony | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 |
| Arsenic | 0.025 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 |
| Barium | 1 | 0.037 | 0.011 | | 0.001 | 0.047 | 0.011 | | 0.001 | 0.173 | 0.011 | | 0.001 | 0.297 | 0.011 | | 0.001 | 0.139 | 0.011 | | 0.001 | 0.078 | 0.011 | | 0.001 | 0.044 | 0.011 | | 0.001 |
| Beryllium | 0.003 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 |
| Cadmium | 0.005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 |
| Calcium | NS | 251 | 0.11 | | 0.11 | 177 | 0.11 | | 0.11 | 215 | 0.11 | | 0.11 | 51.9 | 0.01 | | 0.01 | 239 | 0.11 | | 0.11 | 223 | 0.11 | | 0.11 | 250 | 0.11 | | 0.11 |
| Chromium | 0.05 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 |
| Cobalt | NS | 0.004 | 0.005 | B | 0.001 | < 0.005 | 0.005 | U | 0.001 | 0.003 | 0.005 | B | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | 0.005 | 0.005 | | 0.001 | 0.004 | 0.005 | B | 0.001 |
| Copper | 0.2 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | 0.023 | 0.011 | | 0.002 | < 0.011 | 0.011 | U | 0.002 |
| Iron | 0.5 | 0.02 | 0.02 | | 0.02 | 0.36 | 0.01 | | 0.01 | < 0.02 | 0.02 | U | 0.02 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 |
| Lead | 0.025 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 |
| Magnesium | 35 | 688 | 0.11 | | 0.11 | 411 | 0.11 | | 0.11 | 293 | 0.11 | | 0.11 | 43.9 | 0.01 | | 0.01 | 288 | 0.11 | | 0.11 | 393 | 0.11 | | 0.11 | 612 | 0.11 | | 0.11 |
| Manganese | 0.3 | 0.777 | 0.005 | | 0.001 | 0.672 | 0.005 | | 0.001 | 1.25 | 0.005 | | 0.001 | 0.643 | 0.005 | | 0.001 | 0.403 | 0.005 | | 0.001 | 6.25 | 0.053 | | 0.011 | 1.04 | 0.005 | | 0.001 |
| Mercury | 0.0007 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 |
| Nickel | 0.1 | 0.006 | 0.004 | | 0.001 | < 0.004 | 0.004 | U | 0.001 | 0.002 | 0.004 | B | 0.001 | < 0.004 | 0.004 | U | 0.001 | 0.002 | 0.004 | B | 0.001 | 0.009 | 0.004 | | 0.001 | 0.003 | 0.004 | B | 0.001 |
| Potassium | NS | 233 | 1.1 | | 0.11 | 144 | 1.1 | | 0.11 | 87.2 | 1.1 | | 0.11 | 51.2 | 1.1 | | 0.11 | 105 | 1.1 | | 0.11 | 129 | 1.1 | | 0.11 | 214 | 1.1 | | 0.11 |
| Selenium | 0.01 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 |
| Silver | 0.05 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 |
| Sodium | 2 | 5290 | 11 | | 11 | 3190 | 11 | | 11 | 1950 | 11 | | 11 | 122 | 1.1 | | 1.1 | 2310 | 11 | | 11 | 2990 | 11 | | 11 | 4450 | 11 | | 11 |
| Thallium | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 |
| Vanadium | NS | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 |
| Zinc | 2 | 0.008 | 0.021 | B | 0.002 | 0.005 | 0.021 | B | 0.002 | 0.005 | 0.021 | B | 0.002 | 0.003 | 0.021 | B | 0.002 | 0.005 | 0.021 | B | 0.002 | 0.021 | 0.021 | B | 0.002 | 0.005 | 0.021 | B | 0.002 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

Bold = Result detected above detection limit

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 11
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Dissolved Metals

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (mg/L) | MW9 | | | | MW10 | | | | MW11D | | | | MW12D | | | | MW13D | | | | Duplicate | | | |
|-----------|---|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| Aluminum | NS | 0.01 | 0.011 | B | 0.005 | 0.007 | 0.011 | B | 0.005 | 0.008 | 0.011 | B | 0.005 | 0.008 | 0.011 | B | 0.005 | < 0.011 | 0.011 | U | 0.005 | < 0.011 | 0.011 | U | 0.005 |
| Antimony | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 | < 0.003 | 0.003 | U | 0.003 |
| Arsenic | 0.025 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | 0.005 | 0.003 | | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 | < 0.003 | 0.003 | U | 0.004 |
| Barium | 1 | 0.054 | 0.011 | | 0.001 | 0.082 | 0.011 | | 0.001 | 0.02 | 0.011 | | 0.001 | 0.075 | 0.011 | | 0.001 | 0.02 | 0.011 | | 0.001 | 0.02 | 0.011 | | 0.001 |
| Beryllium | 0.003 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 |
| Cadmium | 0.005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 | < 0.004 | 0.004 | U | 0.0005 |
| Calcium | NS | 226 | 0.11 | | 0.11 | 240 | 0.11 | | 0.11 | 28.4 | 0.01 | | 0.01 | 71.1 | 0.01 | | 0.01 | 35.1 | 0.01 | | 0.01 | 34.8 | 0.01 | | 0.01 |
| Chromium | 0.05 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 | < 0.001 | 0.001 | U | 0.001 |
| Cobalt | NS | 0.002 | 0.005 | B | 0.001 | 0.004 | 0.005 | B | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 |
| Copper | 0.2 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 | 0.003 | 0.011 | B | 0.002 | < 0.011 | 0.011 | U | 0.002 | < 0.011 | 0.011 | U | 0.002 |
| Iron | 0.5 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 | 0.02 | 0.02 | | 0.02 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 |
| Lead | 0.025 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 |
| Magnesium | 35 | 426 | 0.11 | | 0.11 | 247 | 0.11 | | 0.11 | 37.7 | 0.01 | | 0.01 | 148 | 0.11 | | 0.11 | 55.9 | 0.01 | | 0.01 | 55.3 | 0.01 | | 0.01 |
| Manganese | 0.3 | 1.18 | 0.005 | | 0.001 | 5.46 | 0.053 | | 0.011 | 0.063 | 0.005 | | 0.001 | 0.609 | 0.005 | | 0.001 | 0.099 | 0.005 | | 0.001 | 0.099 | 0.005 | | 0.001 |
| Mercury | 0.0007 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 |
| Nickel | 0.1 | 0.008 | 0.004 | | 0.001 | 0.008 | 0.004 | | 0.001 | 0.002 | 0.004 | B | 0.001 | 0.002 | 0.004 | B | 0.001 | 0.002 | 0.004 | B | 0.001 | 0.002 | 0.004 | B | 0.001 |
| Potassium | NS | 146 | 1.1 | | 0.11 | 82.5 | 1.1 | | 0.11 | 36 | 1.1 | | 0.11 | 60.7 | 1.1 | | 0.11 | 37.1 | 1.1 | | 0.11 | 35.7 | 1.1 | | 0.11 |
| Selenium | 0.01 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 |
| Silver | 0.05 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 | < 0.005 | 0.005 | U | 0.001 |
| Sodium | 2 | 2,900 | 11 | | 11 | 1870 | 11 | | 11 | 719 | 11 | | 11 | 1420 | 11 | | 11 | 832 | 11 | | 11 | 799 | 11 | | 11 |
| Thallium | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 |
| Vanadium | NS | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | 0.002 | 0.011 | B | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 | < 0.011 | 0.011 | U | 0.001 |
| Zinc | 2 | 0.028 | 0.021 | | 0.002 | 0.007 | 0.021 | B | 0.002 | 0.004 | 0.021 | B | 0.002 | 0.004 | 0.021 | B | 0.002 | 0.005 | 0.021 | B | 0.002 | 0.004 | 0.021 | B | 0.002 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater

Bold = Result detected above detection limit

| | |
|-----|--|
| U | The compound was analyzed for but not detected at or above the MDL. |
| U | The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

Table 12
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Total Metals

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (mg/L) | MW1 | | | | MW2 | | | | MW3 | | | | MW4 | | | | MW5 | | | | MW7 | | | | MW8 | | | |
|-----------|--|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| Aluminum | 0.1 | 0.545 | 0.020 | | 0.0048 | 0.41 | 0.020 | | 0.0048 | 2.24 | 0.020 | | 0.0048 | 0.372 | 0.020 | | 0.0048 | 0.911 | 0.020 | | 0.0048 | < 0.020 | 0.020 | U | 0.0048 | < 0.020 | 0.020 | U | 0.0048 |
| Antimony | 0.003 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Arsenic | 0.025 | < 0.008 | 0.008 | U | 0.002 | < 0.008 | 0.008 | U | 0.002 | 0.005 | 0.008 | B | 0.002 | < 0.008 | 0.008 | U | 0.002 | 0.002 | 0.008 | B | 0.002 | 0.002 | 0.008 | B | 0.002 | < 0.008 | 0.008 | U | 0.002 |
| Barium | 1 | 0.042 | 0.020 | | 0.002 | 0.053 | 0.020 | | 0.002 | 0.211 | 0.020 | | 0.002 | 0.367 | 0.020 | | 0.002 | 0.146 | 0.020 | | 0.002 | 0.076 | 0.020 | | 0.002 | 0.05 | 0.020 | | 0.002 |
| Beryllium | 0.003 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Cadmium | 0.005 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | 0.001 | 0.005 | B | 0.0010 |
| Calcium | NS | 258 | 1.0 | | 0.30 | 165 | 1.0 | | 0.30 | 226 | 1.0 | | 0.30 | 51.2 | 1.0 | | 0.30 | 246 | 1.0 | | 0.30 | 236 | 1.0 | | 0.30 | 262 | 1.0 | | 0.30 |
| Chromium | 0.05 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | 0.008 | 0.002 | | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Cobalt | NS | 0.003 | 0.010 | B | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.006 | 0.010 | B | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.005 | 0.010 | B | 0.002 | 0.004 | 0.010 | B | 0.002 |
| Copper | 0.2 | 0.008 | 0.010 | B | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.011 | 0.010 | | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.004 | 0.010 | B | 0.002 | 0.02 | 0.010 | | 0.002 | < 0.010 | 0.010 | U | 0.002 |
| Iron | 0.5 | 9.07 | 0.02 | | 0.02 | 34.3 | 1.0 | | 1.0 | 20.2 | 1.0 | | 1.0 | 8.79 | 0.02 | | 0.02 | 2.19 | 0.02 | | 0.02 | 0.24 | 0.02 | | 0.02 | 48.8 | 1.0 | | 1.0 |
| Lead | 0.025 | 0.014 | 0.004 | | 0.002 | < 0.004 | 0.004 | U | 0.002 | 0.023 | 0.004 | | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 |
| Magnesium | 35 | 662 | 1.0 | | 0.10 | 399 | 1.0 | | 0.10 | 286 | 1.0 | | 0.10 | 41.3 | 1.0 | | 0.10 | 269 | 1.0 | | 0.10 | 422 | 1.0 | | 0.10 | 580 | 1.0 | | 0.10 |
| Manganese | 0.3 | 0.758 | 0.010 | | 0.002 | 0.602 | 0.010 | | 0.002 | 1.25 | 0.010 | | 0.002 | 0.665 | 0.010 | | 0.002 | 0.404 | 0.010 | | 0.002 | 5.61 | 0.50 | | 0.10 | 1.1 | 0.010 | | 0.002 |
| Mercury | 0.0007 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 |
| Nickel | 0.1 | 0.006 | 0.008 | B | 0.002 | < 0.008 | 0.008 | U | 0.002 | 0.006 | 0.008 | B | 0.002 | < 0.008 | 0.008 | U | 0.002 | 0.002 | 0.008 | B | 0.002 | 0.006 | 0.008 | B | 0.002 | 0.003 | 0.008 | B | 0.002 |
| Potassium | NS | 209 | 10 | | 10 | 109 | 10 | | 10 | 84 | 10 | | 10 | 53 | 10 | | 10 | 96 | 10 | | 10 | 123 | 10 | | 10 | 188 | 10 | | 10 |
| Selenium | 0.01 | 0.018 | 0.002 | | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.01 | 0.01 | U | 0.01 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | 0.004 | 0.002 | | 0.001 |
| Silver | 0.05 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.003 | 0.010 | B | 0.002 | < 0.010 | 0.010 | U | 0.002 |
| Sodium | 2 | 5,390 | 10 | | 10 | 2,850 | 10 | | 10 | 1,900 | 10 | | 10 | 132 | 10 | | 10 | 2,390 | 10 | | 10 | 3,220 | 10 | | 10 | 4,830 | 10 | | 10 |
| Thallium | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 |
| Vanadium | NS | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 | < 0.020 | 0.020 | U | 0.002 |
| Zinc | 2 | 0.013 | 0.020 | B | 0.002 | 0.005 | 0.020 | B | 0.002 | 0.022 | 0.020 | | 0.002 | 0.009 | 0.020 | B | 0.002 | 0.013 | 0.020 | B | 0.002 | 0.017 | 0.020 | B | 0.002 | 0.004 | 0.020 | B | 0.002 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit
 ND = Non-detect

Bold/highlighted - Indicated exceedance of the NYSDEC Groundwater Standard

Bold = Result detected above detection limit

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response fo the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

Table 12
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Groundwater Analytical Results
Total Metals

| Compound | NYSDEC Ambient Water Quality Standards (AWQS) (mg/L) | MW9 | | | | MW10 | | | | MW11D | | | | MW12D | | | | MW13D | | | | Duplicate | | | |
|-----------|--|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|--------------|--------|------|---------|
| | | 12/2/2015 | | | | 12/2/2015 | | | | 12/2/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | | 12/1/2015 | | | |
| | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | | (µg/L) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| Aluminum | 0.1 | 19.4 | 1.0 | | 0.24 | 0.016 | 0.020 | B | 0.0048 | 3.06 | 0.020 | | 0.0048 | 37.9 | 1.0 | | 0.24 | 0.997 | 0.020 | | 0.0048 | 1.05 | 0.020 | | 0.0048 |
| Antimony | 0.003 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Arsenic | 0.025 | 0.034 | 0.008 | | 0.002 | 0.021 | 0.008 | | 0.002 | 0.004 | 0.008 | B | 0.002 | < 0.008 | 0.008 | U | 0.002 | < 0.008 | 0.008 | U | 0.002 | < 0.008 | 0.008 | U | 0.002 |
| Barium | 1 | 0.13 | 0.020 | | 0.002 | 0.112 | 0.020 | | 0.002 | 0.056 | 0.020 | | 0.002 | 0.311 | 0.020 | | 0.002 | 0.026 | 0.020 | | 0.002 | 0.026 | 0.020 | | 0.002 |
| Beryllium | 0.003 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 | 0.003 | 0.002 | | 0.002 | < 0.002 | 0.002 | U | 0.002 | < 0.002 | 0.002 | U | 0.002 |
| Cadmium | 0.005 | 0.002 | 0.005 | B | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | 0.002 | 0.005 | B | 0.0010 | < 0.005 | 0.005 | U | 0.0010 | < 0.005 | 0.005 | U | 0.0010 |
| Calcium | NS | 289 | 1.0 | | 0.30 | 249 | 1.0 | | 0.30 | 36 | 0.020 | | 0.006 | 113 | 1.0 | | 0.30 | 35.6 | 1.0 | | 0.30 | 35.7 | 1.0 | | 0.30 |
| Chromium | 0.05 | 0.047 | 0.002 | | 0.002 | < 0.002 | 0.002 | U | 0.002 | 0.018 | 0.002 | | 0.002 | 0.094 | 0.002 | | 0.002 | 0.004 | 0.002 | | 0.002 | 0.004 | 0.002 | | 0.002 |
| Cobalt | NS | 0.019 | 0.010 | | 0.002 | 0.004 | 0.010 | B | 0.002 | 0.003 | 0.010 | B | 0.002 | 0.039 | 0.010 | | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 |
| Copper | 0.2 | 0.096 | 0.010 | | 0.002 | < 0.010 | 0.010 | U | 0.002 | 0.028 | 0.010 | | 0.002 | 0.132 | 0.010 | | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | B | 0.002 |
| Iron | 0.5 | 48.4 | 1.0 | | 1.0 | 18.2 | 1.0 | | 1.0 | 6.67 | 0.02 | | 0.02 | 78.9 | 1.0 | | 1.0 | 2.94 | 0.02 | | 0.02 | 3.08 | 0.02 | | 0.02 |
| Lead | 0.025 | 0.224 | 0.004 | | 0.002 | < 0.004 | 0.004 | U | 0.002 | 0.047 | 0.004 | | 0.002 | 0.053 | 0.004 | | 0.002 | < 0.004 | 0.004 | U | 0.002 | < 0.004 | 0.004 | U | 0.002 |
| Magnesium | 35 | 489 | 1.0 | | 0.10 | 228 | 1.0 | | 0.10 | 40.2 | 1.0 | | 0.10 | 174 | 1.0 | | 0.10 | 60.9 | 1.0 | | 0.10 | 53.2 | 1.0 | | 0.10 |
| Manganese | 0.3 | 2.45 | 0.010 | | 0.002 | 5.35 | 0.50 | | 0.10 | 0.135 | 0.010 | | 0.002 | 1.46 | 0.010 | | 0.002 | 0.114 | 0.010 | | 0.002 | 0.119 | 0.010 | | 0.002 |
| Mercury | 0.0007 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 | < 0.0002 | 0.0002 | U | 0.00015 |
| Nickel | 0.1 | 0.063 | 0.008 | | 0.002 | 0.007 | 0.008 | B | 0.002 | 0.01 | 0.008 | | 0.002 | 0.096 | 0.008 | | 0.002 | < 0.008 | 0.008 | U | 0.002 | 0.006 | 0.008 | B | 0.002 |
| Potassium | NS | 121 | 10 | | 10 | 72 | 10 | | 10 | 34 | 10 | | 10 | 72.3 | 0.2 | | 0.2 | 34 | 10 | | 10 | 36 | 10 | | 10 |
| Selenium | 0.01 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.002 | 0.002 | U | 0.001 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 | < 0.01 | 0.01 | U | 0.01 |
| Silver | 0.05 | < 0.010 | 0.010 | U | 0.002 | 0.002 | 0.010 | B | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 | < 0.010 | 0.010 | U | 0.002 |
| Sodium | 2 | 3,160 | 10 | | 10 | 1,710 | 10 | | 10 | 736 | 10 | | 10 | 1,290 | 10 | | 10 | 780 | 10 | | 10 | 818 | 10 | | 10 |
| Thallium | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 | < 0.0005 | 0.0005 | U | 0.0005 |
| Vanadium | NS | 0.037 | 0.020 | | 0.002 | < 0.020 | 0.020 | U | 0.002 | 0.007 | 0.020 | B | 0.002 | 0.085 | 0.020 | | 0.002 | < 0.020 | 0.020 | U | 0.002 | 0.004 | 0.020 | B | 0.002 |
| Zinc | 2 | 0.306 | 0.020 | | 0.002 | 0.005 | 0.020 | B | 0.002 | 0.038 | 0.020 | | 0.002 | 0.462 | 0.020 | | 0.002 | 0.008 | 0.020 | B | 0.002 | 0.009 | 0.020 | B | 0.002 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

Bold = Result detected above detection limit

| | |
|-----|---|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| B | This compound was also present in the method blank |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response fo the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of a diluted analysis. |
| (*) | See report for comment. |

TABLE 13
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Parameters Detected Above Ambient Groundwater Standards

| COMPOUND | Range in Exceedances | Frequency of Detection | MW1 | MW2 | MW3 | MW4 | MW5 | MW7 | MW8 | MW9 | MW10 | MW11D | MW12D | MW13D | Duplicate |
|-------------------------------|----------------------|------------------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|-------------|
| | | | 12/2/2015 | 12/2/2015 | 12/1/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/2/2015 | 12/1/2015 | 12/1/2015 |
| <i>Sample Results in ug/L</i> | | | | | | | | | | | | | | | |
| Chloromethane | 8.3-39 | 4 | - | - | 39 | - | 8.3 | - | - | - | - | - | - | 16 | 15 |
| <i>Sample Results in ug/L</i> | | | | | | | | | | | | | | | |
| Acenaphthylene | 0.05 | 1 | - | - | - | - | - | - | - | 0.05 | - | - | - | - | - |
| Benz(a)anthracene | 0.03 | 1 | - | - | - | - | - | - | - | 0.03 | - | - | - | - | - |
| Benzo(a)pyrene | 0.03 | 1 | - | - | - | - | - | - | - | 0.03 | - | - | - | - | - |
| Benzo(ghi)perylene | 0.04 | 1 | - | - | - | - | - | - | - | 0.04 | - | - | - | - | - |
| Bis(2-ethylhexyl)phthalate | 0.05 | 1 | - | - | - | - | - | - | - | 0.05 | - | - | - | - | - |
| <i>Sample Results in ug/L</i> | | | | | | | | | | | | | | | |
| PCB-1254 | 1.3 | 1 | - | - | - | - | - | - | - | 1.3 | - | - | - | - | - |
| <i>Sample Results in mg/L</i> | | | | | | | | | | | | | | | |
| Iron (dissolved) | 0.36-21.8 | 2 | - | 0.36 | - | - | - | - | 21.8 | - | - | - | - | - | - |
| Magnesium (dissolved) | 37.7-688 | 13 | 688 | 411 | 293 | 43.9 | 288 | 393 | 612 | 426 | 247 | 37.7 | 148 | 55.9 | 55.3 |
| Manganese (dissolved) | 0.403-5.46 | 11 | 0.777 | 0.672 | 1.25 | 0.643 | 0.403 | 6.25 | 1.04 | 1.18 | 5.46 | - | 0.609 | - | - |
| Sodium (dissolved) | 122-5290 | 13 | 5290 | 3190 | 1950 | 122 | 2310 | 2990 | 4450 | 2,900 | 1870 | 719 | 1420 | 832 | 799 |
| <i>Sample Results in mg/L</i> | | | | | | | | | | | | | | | |
| Aluminum | 0.372-19.4 | 3 | 0.545 | 0.41 | 2.24 | 0.372 | 0.911 | - | - | 19.4 | - | 3.06 | 37.9 | 0.997 | 1.05 |
| Arsenic | 0.034 | 1 | - | - | - | - | - | - | - | 0.034 | - | - | - | - | - |
| Chromium | 0.094 | 1 | - | - | - | - | - | - | - | - | - | - | 0.094 | - | - |
| Iron | 2.19-78.9 | 12 | 9.07 | 34.3 | 20.2 | 8.79 | 2.19 | - | 48.8 | 48.4 | 18.2 | 6.67 | 78.9 | 2.94 | 3.08 |
| Lead | 0.047-0.224 | 3 | - | - | - | - | - | - | - | 0.224 | - | 0.047 | 0.053 | - | - |
| Magnesium | 40.2-662 | 13 | 662 | 399 | 286 | 41.3 | 269 | 422 | 580 | 489 | 228 | 40.2 | 174 | 60.9 | 53.2 |
| Manganese | 0.404-5.61 | 10 | 0.758 | 0.602 | 1.25 | 0.665 | 0.404 | 5.61 | 1.1 | 2.45 | 5.35 | - | 1.46 | - | - |
| Selenium | 0.018 | 1 | 0.018 | - | - | - | - | - | - | - | - | - | - | - | - |
| Sodium | 132-5390 | 13 | 5,390 | 2,850 | 1,900 | 132 | 2,390 | 3,220 | 4,830 | 3,160 | 1,710 | 736 | 1,290 | 780 | 818 |

Notes:

RL - Laboratory Reporting Limit, MDL=Minimum Detection Limit

ND = Non-detect

Bold/highlighted- Indicated exceedance of the NYSDEC Groundwater Standard

TABLE 14
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Gas - Volatile Organic Compounds

| COMPOUNDS | NYSDOH Soil Outdoor Background Levels (µg/m ³) ^(a) | SG1 | | | | SG2 | | | | SG3 | | | | SG4 | | | |
|-------------------------------|---|--------------------------------|----------------|------|------|--------------------------------|------|------|------|--------------------------------|------|------|------|--------------------------------|------|------|------|
| | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,1-Trichloroethane | <2.0 - 2.8 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,2,2-Tetrachloroethane | <1.5 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,2-Trichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1-Dichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1-Dichloroethene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trichlorobenzene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trimethylbenzene | <1.0 | 1.63 | 1.00 | | 1.00 | 1.64 | 1.00 | | 1.00 | 1.08 | 1.00 | | 1.00 | 2.49 | 1.00 | | 1.00 |
| 1,2-Dibromoethane | <1.5 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloropropane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorotetrafluoroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,3,5-Trimethylbenzene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | 1.02 | 1.00 | | 1.00 |
| 1,3-Butadiene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,3-Dichlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,4-Dichlorobenzene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,4-Dioxane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 2-Hexanone | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 4-Ethyltoluene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 4-Isopropyltoluene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 4-Methyl-2-pentanone | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Acetone | NA | 434 | 29.9 | DS | 29.9 | 622 | 29.9 | DS | 29.9 | 572 | 29.9 | DS | 29.9 | 3,250 | 75.0 | DS | 75.0 |
| Acrylonitrile | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Benzene | <1.6 - 4.7 | 2.51 | 1.00 | | 1.00 | 9.93 | 1.00 | | 1.00 | 6.86 | 1.00 | | 1.00 | 2.18 | 1.00 | | 1.00 |
| Benzyl Chloride | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromodichloromethane | <5.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromoform | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromomethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Carbon Disulfide | NA | 13.9 | 1.00 | | 1.00 | 38.3 | 1.00 | | 1.00 | 20 | 1.00 | | 1.00 | 8.81 | 1.00 | | 1.00 |
| Carbon Tetrachloride | <3.1 | 0.28 | 0.25 | | 0.25 | 0.26 | 0.25 | | 0.25 | 0.87 | 0.25 | | 0.25 | 0.8 | 0.25 | | 0.25 |
| Chlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloroethane | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloroform | <2.4 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloromethane | <1.0 - 1.4 | 2.64 | 1.00 | | 1.00 | 3.74 | 1.00 | | 1.00 | 6.19 | 1.00 | | 1.00 | 1.25 | 1.00 | | 1.00 |
| cis-1,2-Dichloroethene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| cis-1,3-Dichloropropene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Cyclohexane | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Dibromochloromethane | <5.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Dichlorodifluoromethane | NA | 1.7 | 1.00 | | 1.00 | 1.63 | 1.00 | | 1.00 | 1.38 | 1.00 | | 1.00 | 1.43 | 1.00 | | 1.00 |
| Ethanol | | 13.3 | 1.00 | | 1.00 | 10.9 | 1.00 | | 1.00 | 16.1 | 1.00 | | 1.00 | 78.7 | 1.00 | E | 1.00 |
| Ethyl Acetate | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Ethylbenzene | <4.3 | 1.19 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Heptane | NA | 561 | 30.0 | D | 30.0 | 2840 | 30.0 | D | 30.0 | 2,240 | 30.0 | D | 30.0 | 111 | 1.00 | | 1.00 |
| Hexachlorobutadiene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Hexane | <1.5 | 1310 | 30.0 | DS | 30.0 | 2520 | 30.0 | DS | 30.0 | 5,880 | 75.0 | DS | 75.0 | 266 | 30.0 | DS | 30.0 |
| Isopropylalcohol | NA | 5.97 | 1.00 | S | 1.00 | 6.29 | 1.00 | S | 1.00 | 10.5 | 1.00 | | 1.00 | 54.3 | 1.00 | | 1.00 |
| Isopropylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Xylene (m&p) | <4.3 | 1.93 | 1.00 | | 1.00 | 3.37 | 1.00 | | 1.00 | 1.68 | 1.00 | | 1.00 | 1.74 | 1.00 | | 1.00 |
| Methyl Ethyl Ketone | | 10.9 | 1.00 | | 1.00 | 14.8 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | 65.1 | 1.00 | | 1.00 |
| MTBE | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Methylene Chloride | <3.4 | 2.51 | 1.00 | S | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| n-Butylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Xylene (o) | <4.3 | 1.03 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Propylene | NA | 61.7 | 29.9 | D | 29.9 | 501 | 29.9 | D | 29.9 | 322 | 29.9 | D | 29.9 | 139 | 29.9 | D | 29.9 |
| sec-Butylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Styrene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Tetrachloroethene | | < 0.25 | 0.25 | U | 0.25 | 0.26 | 0.25 | | 0.25 | < 0.25 | 0.25 | U | 0.25 | 3.52 | 0.25 | | 0.25 |
| Tetrahydrofuran | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Toluene | 1.0 - 6.1 | 2.82 | 1.00 | | 1.00 | 5.57 | 1.00 | | 1.00 | 10.3 | 1.00 | | 1.00 | 2.36 | 1.00 | | 1.00 |
| trans-1,2-Dichloroethene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| trans-1,3-Dichloropropene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Trichloroethene | <1.7 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | 0.42 | 0.25 | | 0.25 | < 0.25 | 0.25 | U | 0.25 |
| Trichlorofluoromethane | NA | 1.01 | 1.00 | | 1.00 | 1.11 | 1.00 | | 1.00 | 1.12 | 1.00 | | 1.00 | 1.11 | 1.00 | | 1.00 |
| Trichlorotrifluoroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Vinyl Chloride | <1.0 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 |
| Total CVOCs | | | 0.28 | | | 0.5 | | | | 1.3 | | | | 4.32 | | | |
| BTEX | | | 9.48 | | | 18.87 | | | | 18.84 | | | | 6.28 | | | |
| Total VOCs | | | 2430.02 | | | 6,580.80 | | | | 9090.50 | | | | 3,990.81 | | | |

Notes:

NA = No guidance value or standard available
(a) = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York, February 2005, Summary of Background Levels for Selected
RL = Laboratory detection limit
Bold = Result detected above detection limit
VOCs = Volatile Organic Compounds
CVOCs = Chlorinated Volatile Organic Compounds
BTEX = Benzene, Toluene, Ethylbenzene, and Xylenes

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of |

TABLE 14
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Gas - Volatile Organic Compounds

| COMPOUNDS | NYSDOH Soil Outdoor Background Levels (µg/m ³) ^(a) | SG5 | | | | SG7 | | | | SG8 | | | | SG9 | | | |
|-------------------------------|---|--------------------------------|-----------------|------|------|--------------------------------|-----------------|------|------|--------------------------------|-----------------|------|------|--------------------------------|---------------|------|------|
| | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,1-Trichloroethane | <2.0 - 2.8 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | 1.02 | 1.00 | | 1.00 |
| 1,1,2,2-Tetrachloroethane | <1.5 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,2-Trichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1-Dichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,1,4-Dichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trichlorobenzene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trimethylbenzene | <1.0 | 2.76 | 1.00 | | 1.00 | 3.35 | 1.00 | | 1.00 | 2.94 | 1.00 | | 1.00 | 2.61 | 1.00 | | 1.00 |
| 1,2-Dibromoethane | <1.5 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloroethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloropropane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorotetrafluoroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,3,5-Trimethylbenzene | <1.0 | 1.01 | 1.00 | | 1.00 | 1.26 | 1.00 | | 1.00 | 1.13 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,3-Butadiene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,3-Dichlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,4-Dichlorobenzene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 1,4-Dioxane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 2-Hexanone | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | 5.61 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 4-Ethyltoluene | NA | 1.69 | 1.00 | | 1.00 | 2.26 | 1.00 | | 1.00 | 1.85 | 1.00 | | 1.00 | 1.6 | 1.00 | | 1.00 |
| 4-Isopropyltoluene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| 4-Methyl-2-pentanone | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Acetone | NA | < 1.00 | 1.00 | U | 1.00 | 582 | 9.99 | DS | 9.99 | 2850 | 75.0 | DS | 75.0 | 118 | 5.01 | DS | 5.01 |
| Acrylonitrile | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Benzene | <1.6 - 4.7 | 10.3 | 1.00 | | 1.00 | 13.6 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | 1.75 | 1.00 | | 1.00 |
| Benzyl Chloride | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromodichloromethane | <5.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromoform | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Bromomethane | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Carbon Disulfide | NA | 26.2 | 1.00 | | 1.00 | 34.5 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Carbon Tetrachloride | <3.1 | 0.35 | 0.25 | | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 |
| Chlorobenzene | <2.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloroethane | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloroform | <2.4 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Chloromethane | <1.0 - 1.4 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| cis-1,2-Dichloroethene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| cis-1,3-Dichloropropene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Cyclohexane | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Dibromochloromethane | <5.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Dichlorodifluoromethane | NA | 2.11 | 1.00 | | 1.00 | 2.08 | 1.00 | | 1.00 | 2.03 | 1.00 | | 1.00 | 27.2 | 1.00 | | 1.00 |
| Ethanol | | 8.93 | 1.00 | S | 1.00 | 14.8 | 1.00 | | 1.00 | 78.7 | 1.00 | E | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Ethyl Acetate | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Ethylbenzene | <4.3 | 2.62 | 1.00 | | 1.00 | 2.87 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Heptane | NA | 3330 | 30.0 | D | 30.0 | 290 | 9.99 | D | 9.99 | 2.51 | 1.00 | | 1.00 | 14.8 | 1.00 | | 1.00 |
| Hexachlorobutadiene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Hexane | <1.5 | 6200 | 75.0 | DS | 75.0 | 824 | 10.0 | DS | 10.0 | < 1.00 | 1.00 | U | 1.00 | 117 | 1.00 | | 1.00 |
| Isopropylalcohol | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Isopropylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Xylene (m&p) | <4.3 | 3.98 | 1.00 | | 1.00 | 5.51 | 1.00 | | 1.00 | 1.51 | 1.00 | | 1.00 | 1.53 | 1.00 | | 1.00 |
| Methyl Ethyl Ketone | | 12.9 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | 40.7 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| MTBE | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Methylene Chloride | <3.4 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| n-Butylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Xylene (o) | <4.3 | < 1.00 | 1.00 | U | 1.00 | 1.59 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Propylene | NA | < 1.00 | 1.00 | U | 1.00 | 270 | 9.99 | D | 9.99 | 7.88 | 1.00 | | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| sec-Butylbenzene | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Styrene | <1.0 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Tetrachloroethene | | 0.39 | 0.25 | | 0.25 | 6.26 | 0.25 | | 0.25 | 2.64 | 0.25 | | 0.25 | 15.8 | 0.25 | | 0.25 |
| Tetrahydrofuran | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Toluene | 1.0 - 6.1 | 44.8 | 1.00 | | 1.00 | 18.8 | 1.00 | | 1.00 | 1.23 | 1.00 | | 1.00 | 1.55 | 1.00 | | 1.00 |
| trans-1,2-Dichloroethene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| trans-1,3-Dichloropropene | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Trichloroethene | <1.7 | < 0.25 | 0.25 | U | 0.25 | 0.43 | 0.25 | | 0.25 | 0.27 | 0.25 | | 0.25 | 9.24 | 0.25 | | 0.25 |
| Trichlorofluoromethane | NA | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | 1.45 | 1.00 | | 1.00 | 11.9 | 1.00 | | 1.00 |
| Trichlorotrifluoroethane | | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 | < 1.00 | 1.00 | U | 1.00 |
| Vinyl Chloride | <1.0 | 0.28 | 0.25 | | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 | < 0.25 | 0.25 | U | 0.25 |
| Total CVOCs | | | 0.74 | | | | 6.69 | | | | 2.91 | | | | 26.06 | | |
| BTEX | | | 61.7 | | | | 42.37 | | | | 2.74 | | | | 4.83 | | |
| Total VOCs | | | 9,648.32 | | | | 2,073.31 | | | | 3,000.45 | | | | 324.00 | | |

Notes:

NA = No guidance value or standard available
(a) = NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York, February 2005, Summary of Background Levels for Selected
RL = Laboratory detection limit
Bold = Result detected above detection limit
VOCs = Volatile Organic Compounds
CVOCs = Chlorinated Volatile Organic Compounds
BTEX = Benzene, Toluene, Ethylbenzene, and Xylenes

| | |
|---|--|
| U | The compound was analyzed for but not detected at or above the MDL. The number immediately preceding the "U" represents the PQL reporting level corrected for percent solids, weight and/or volume calculations, and dilution factors. |
| J | The value is estimated. This flag is used a) on form 1 when the compound is reported above the MDL, but below the PQL, and b) on the Tentatively Identified Compounds (TIC) form for all compounds identified. |
| N | The concentration is based on the response for the nearest internal. This flag is used on the TIC form for all compounds identified. |
| S | This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level. |
| D | The reported concentration is the result of |

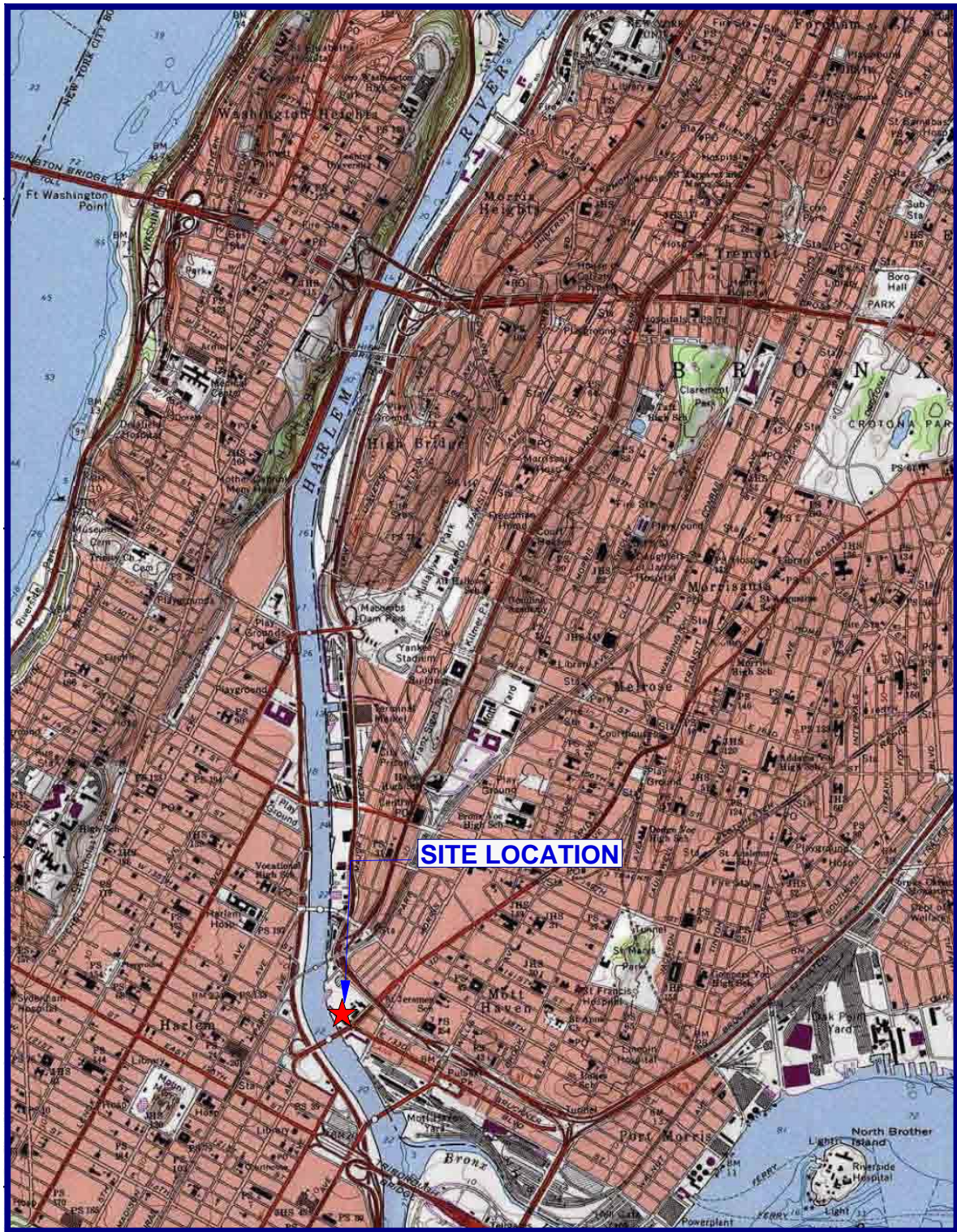
TABLE 14
Former Mugler Shoring
2401 3rd Avenue, Bronx, NY
Soil Gas - Volatile Organic Compounds

| COMPOUNDS | NYSDOH Soil Outdoor Background Levels (µg/m ³) ^(a) | SS1 | | | | SS2 | | | | SS3 | | | | SS4 | | | | SS5 | | | |
|-------------------------------|--|-----------------------------------|------|------|------|-----------------------------------|------|------|------|-----------------------------------|------|------|------|-----------------------------------|------|------|------|-----------------------------------|------|------|------|
| | | 12/2/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/2/2015 (µg/m ³) | | | | 12/1/2015 (µg/m ³) | | | | 12/2/2015 (µg/m ³) | | | |
| | | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL | Results | RL | Qual | MDL |
| 1,1,1,2-Tetrachloroethane | | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,1,1-Trichloroethane | <2.0 - 2.8 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,1,2,2-Tetrachloroethane | <1.5 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,1,2-Trichloroethane | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,1-Dichloroethane | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,1-Dichloroethene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trichlorobenzene | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2,4-Trimethylbenzene | <1.0 | 6.09 | 1.00 | | 1.00 | 3.42 | 1.00 | | 1.00 | 4.06 | 1.00 | | 1.00 | 4.2 | 1.00 | | 1.00 | 3.97 | 1.00 | | 1.00 |
| 1,2-Dibromoethane | <1.5 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorobenzene | <2.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloroethane | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2-Dichloropropane | | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,2-Dichlorotetrafluoroethane | | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,3,5-Trimethylbenzene | <1.0 | 2.56 | 1.00 | | 1.00 | 1.39 | 1.00 | | 1.00 | 1.61 | 1.00 | | 1.00 | 2.19 | 1.00 | | 1.00 | 3.02 | 1.00 | | 1.00 |
| 1,3-Butadiene | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,3-Dichlorobenzene | <2.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,4-Dichlorobenzene | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 1,4-Dioxane | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 2-Hexanone | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 4-Ethyltoluene | NA | 1.51 | 1.00 | | 1.00 | 1.09 | 1.00 | | 1.00 | 1.04 | 1.00 | | 1.00 | 2.57 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 4-Isopropyltoluene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| 4-Methyl-2-pentanone | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.23 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.54 | 1.00 | | 1.00 |
| Acetone | NA | 53.2 | 1.00 | | 1.00 | 23.7 | 1.00 | | 1.00 | 139 | 5.01 | DS | 5.01 | 58.2 | 1.00 | | 1.00 | 102 | 5.01 | DS | 5.01 |
| Acrylonitrile | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Benzene | <1.6 - 4.7 | 1.28 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.1 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.12 | 1.00 | | 1.00 |
| Benzyl Chloride | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Bromodichloromethane | <5.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Bromoform | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Bromomethane | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Carbon Disulfide | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Carbon Tetrachloride | <3.1 | 0.46 | 0.25 | | 0.25 | 0.26 | 0.25 | | 0.25 | 0.43 | 0.25 | | 0.25 | 0.3 | 0.25 | | 0.25 | 0.45 | 0.25 | | 0.25 |
| Chlorobenzene | <2.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Chloroethane | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Chloroform | <2.4 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Chloromethane | <1.0 - 1.4 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.14 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.03 | 1.00 | | 1.00 |
| cis-1,2-Dichloroethene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| cis-1,3-Dichloropropene | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Cyclohexane | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Dibromochloromethane | <5.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Dichlorodifluoromethane | NA | 1.8 | 1.00 | | 1.00 | 2.15 | 1.00 | | 1.00 | 1.86 | 1.00 | | 1.00 | 2.24 | 1.00 | | 1.00 | 1.73 | 1.00 | | 1.00 |
| Ethanol | | 14.2 | 1.00 | | 1.00 | 14.9 | 1.00 | | 1.00 | 98.3 | 5.01 | DS | 5.01 | 4.05 | 1.00 | S | 1.00 | 12.9 | 1.00 | | 1.00 |
| Ethyl Acetate | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.09 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.85 | 1.00 | | 1.00 |
| Ethylbenzene | <4.3 | 1.12 | 1.00 | | 1.00 | 1.51 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.48 | 1.00 | | 1.00 | 2.3 | 1.00 | | 1.00 |
| Heptane | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.22 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.53 | 1.00 | | 1.00 |
| Hexachlorobutadiene | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Hexane | <1.5 | 3.14 | 1.00 | S | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.83 | 1.00 | S | 1.00 | 9.05 | 1.00 | S | 1.00 | 1.76 | 1.00 | S | 1.00 |
| Isopropylalcohol | NA | 1.46 | 1.00 | S | 1.00 | <1.00 | 1.00 | U | 1.00 | 10.1 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 3.02 | 1.00 | S | 1.00 |
| Isopropylbenzene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Xylene (m&p) | <4.3 | 5.55 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 3.66 | 1.00 | | 1.00 | 10.5 | 1.00 | | 1.00 | 9.33 | 1.00 | | 1.00 |
| Methyl Ethyl Ketone | | 5.13 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 8.69 | 1.00 | | 1.00 | 6.57 | 1.00 | | 1.00 | 10.3 | 1.00 | | 1.00 |
| MTBE | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | 6.92 | 1.00 | | 1.00 |
| Methylene Chloride | <3.4 | 1.68 | 1.00 | S | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.92 | 1.00 | S | 1.00 | 1.58 | 1.00 | S | 1.00 | 2.48 | 1.00 | S | 1.00 |
| n-Butylbenzene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Xylene (o) | <4.3 | 2.91 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.92 | 1.00 | | 1.00 | 2.46 | 1.00 | | 1.00 | 5.68 | 1.00 | | 1.00 |
| Propylene | NA | 2.68 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 1.81 | 1.00 | | 1.00 | <1.00 | 1.00 | U | 1.00 | 2.46 | 1.00 | | 1.00 |
| sec-Butylbenzene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Styrene | <1.0 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Tetrachloroethene | | 0.37 | 0.25 | | 0.25 | 4.89 | 0.25 | | 0.25 | 0.26 | 0.25 | | 0.25 | 2.29 | 0.25 | | 0.25 | 0.77 | 0.25 | | 0.25 |
| Tetrahydrofuran | NA | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 | <1.00 | 1.00 | U | 1.00 |
| Toluene | 1. | | | | | | | | | | | | | | | | | | | | |

Table 16
Emergency Contact List

| | | |
|--------------------------|-------------------|------------------|
| General Emergencies | | 911 |
| NYC Police | | 911 |
| NYC Fire Department | | 911 |
| Lincoln Medical Center | | (718) 579-5016 |
| NYSDEC Spills Hotline | | 1-800-457-7362 |
| NYSDEC Project Manager | | (518) 402 - 9656 |
| NYC Department of Health | | (212) 676-2400 |
| National Response Center | | 1-800-424-8802 |
| Poison Control | | 1-800-222-1222 |
| EBC Project Manager | Robert Bennett | (631) 504-6000 |
| EBC BCP Program Manager | Charles Sosik | (631) 504-6000 |
| EBC Site Safety Officer | Kevin Waters | (631) 504-6000 |
| Remedial Engineer | Ariel Czemerinski | (516) 987-1662 |
| Construction Manager | Elliot Lazarus | (516) 808-9085 |

FIGURES

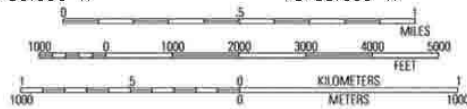


73°57.000' W

73°56.000' W

73°55.000' W

WGS84 73°54.000' W



08/21/13

USGS Central Park Quadrangle 1995, Contour Interval = 10 feet



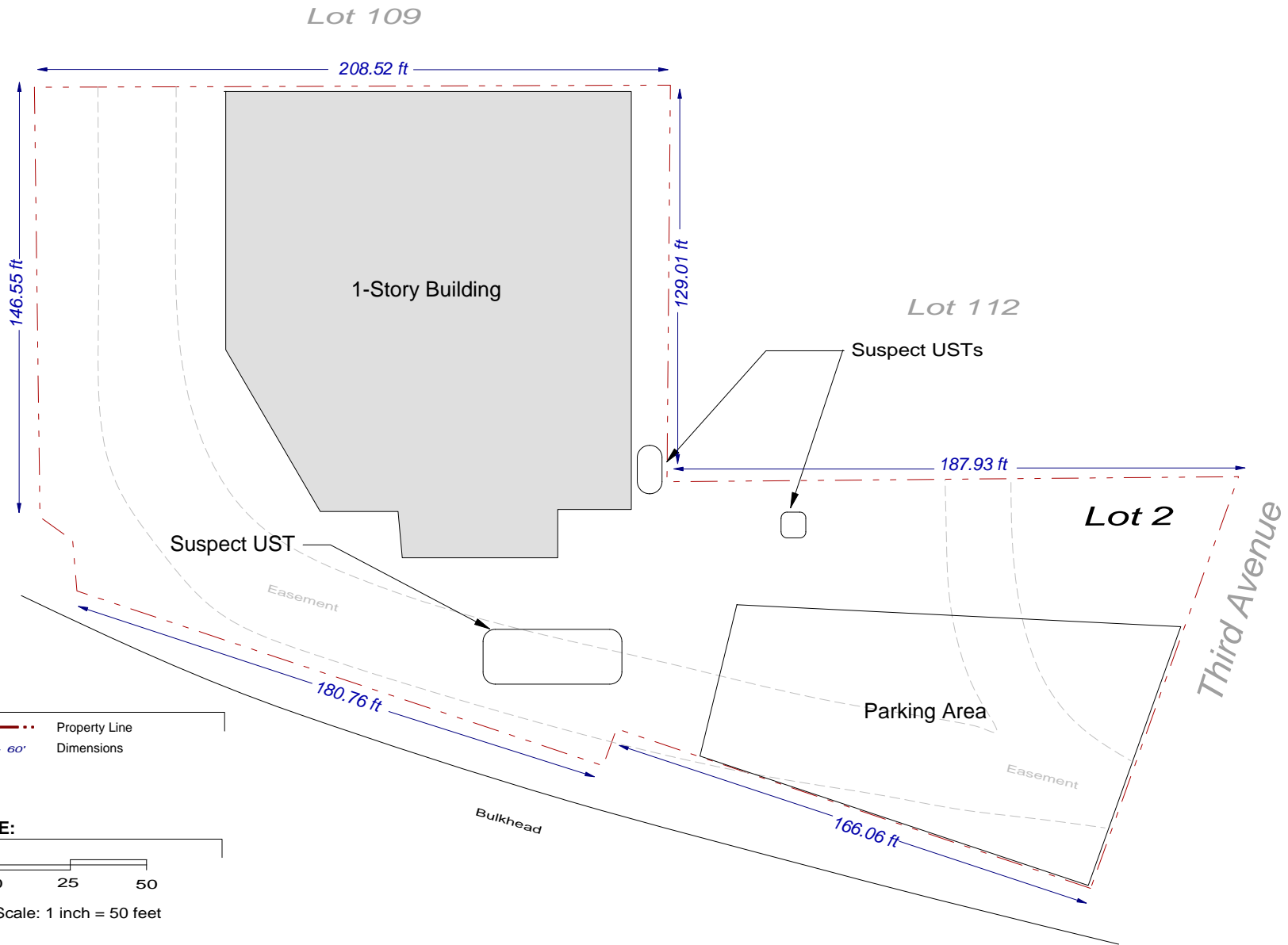
ENVIRONMENTAL BUSINESS CONSULTANTS

Phone 631.504.6000
Fax 631.924.2870

**FORMER MUGLER SHORING INC.
2401 THIRD AVENUE, BRONX NY**

FIGURE 1

SITE LOCATION MAP



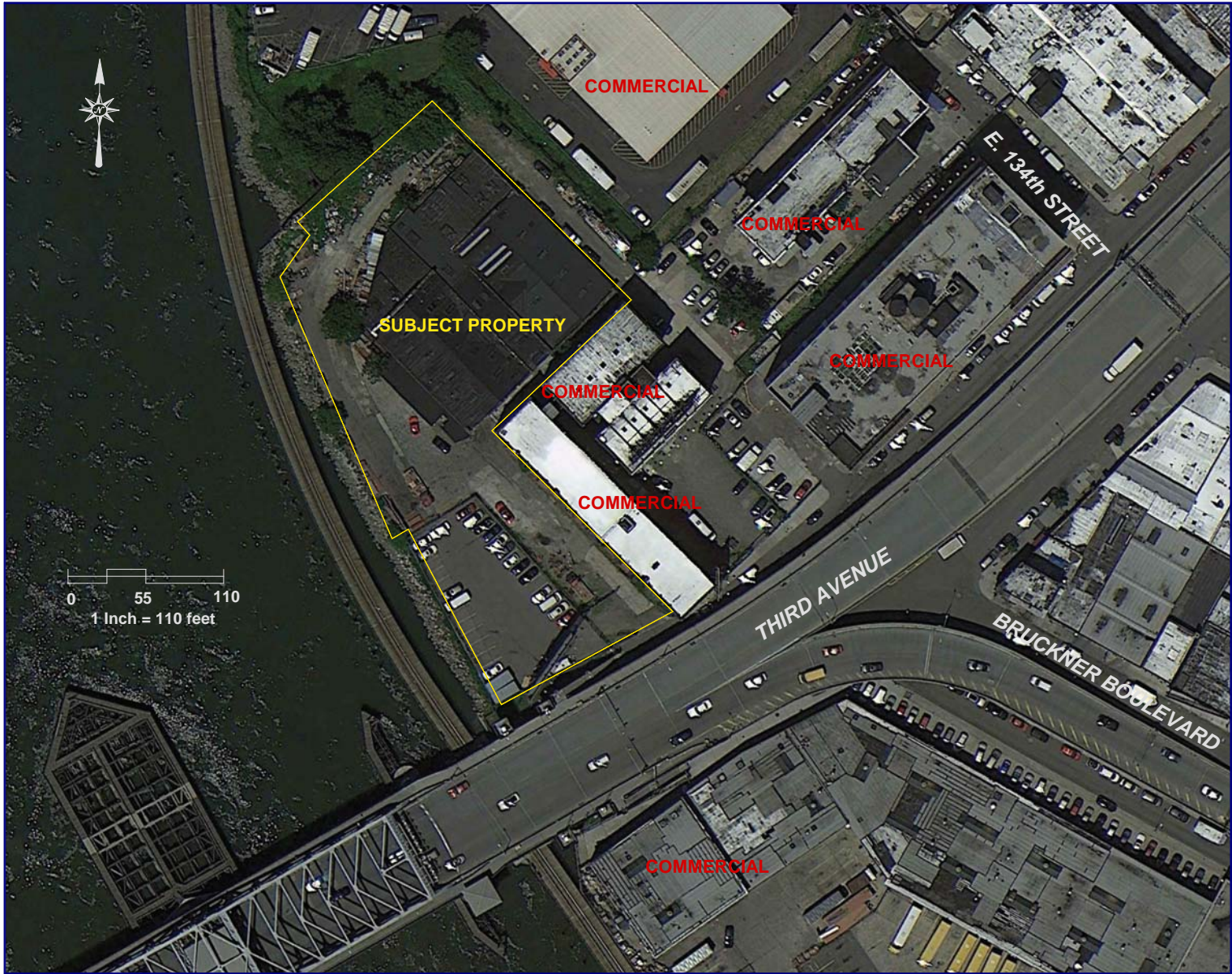
KEY:
- - - Property Line
← 60' Dimensions

SCALE:
0 25 50
Scale: 1 inch = 50 feet

EBC
ENVIRONMENTAL BUSINESS CONSULTANTS
Phone 631.504.6000
Fax 631.924.2870

Figure No.
2

Site Name: **FORMER MUGLER SHORING INC.**
Site Address: **2401 THIRD AVENUE, BRONX, NY**
Drawing Title: **SITE BOUNDARY MAP**



25

0 55 110
1 Inch = 110 feet



ENVIRONMENTAL BUSINESS CONSULTANTS

1808 MIDDLE COUNTRY ROAD, RIDGE, NY 11961

Phone: 631.504.6000

Fax: 631.924.2780

Former Mugler Shoring Inc.

2401 Third Avenue, Bronx NY

FIGURE 3

**PROJECT SITE AND
ADJACENT PROPERTIES**

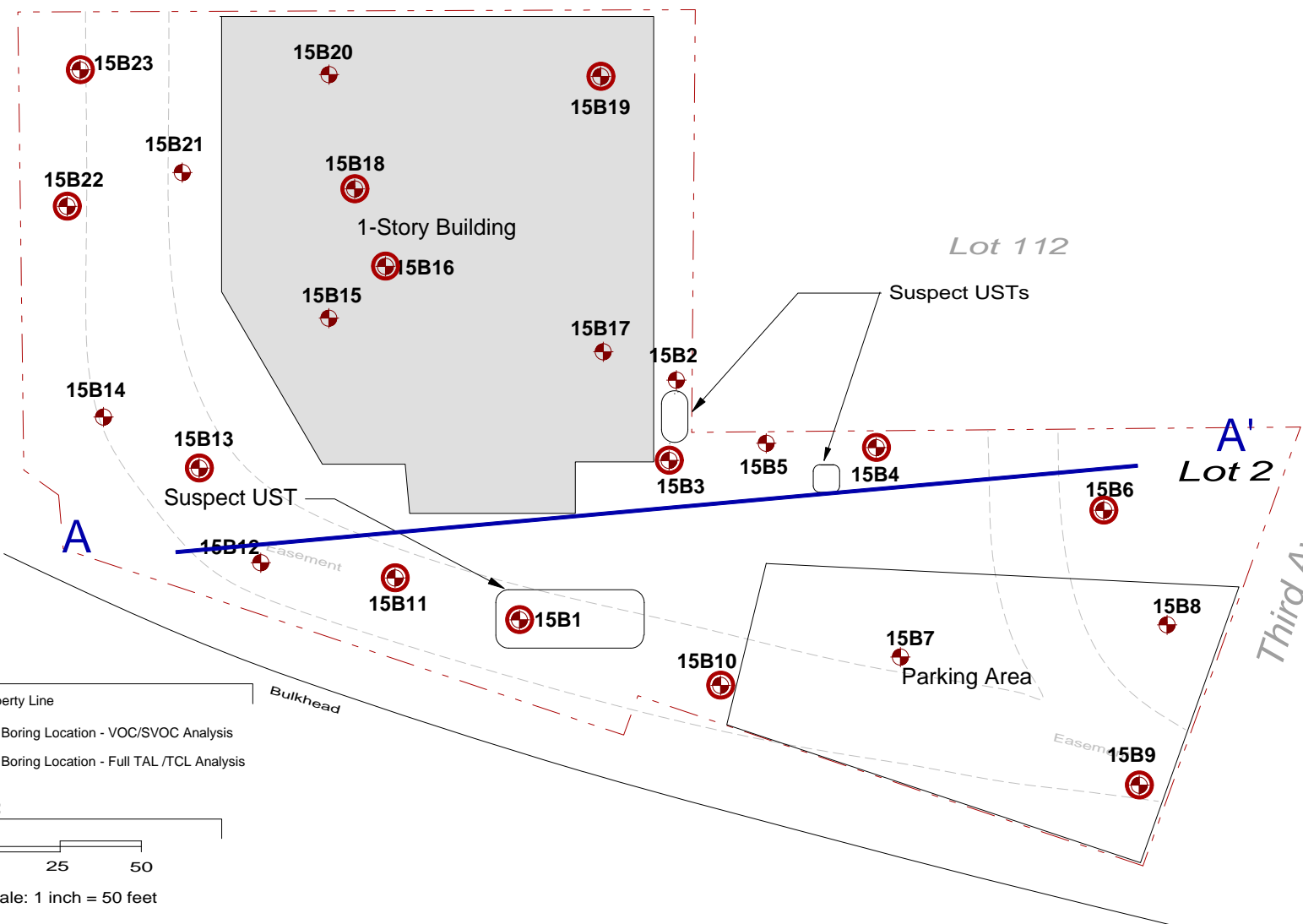


Lot 109

Lot 112

Lot 2

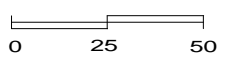
Third Avenue




KEY:

- Property Line
- 15Bx Soil Boring Location - VOC/SVOC Analysis
- 15Bx Soil Boring Location - Full TAL /TCL Analysis

SCALE:

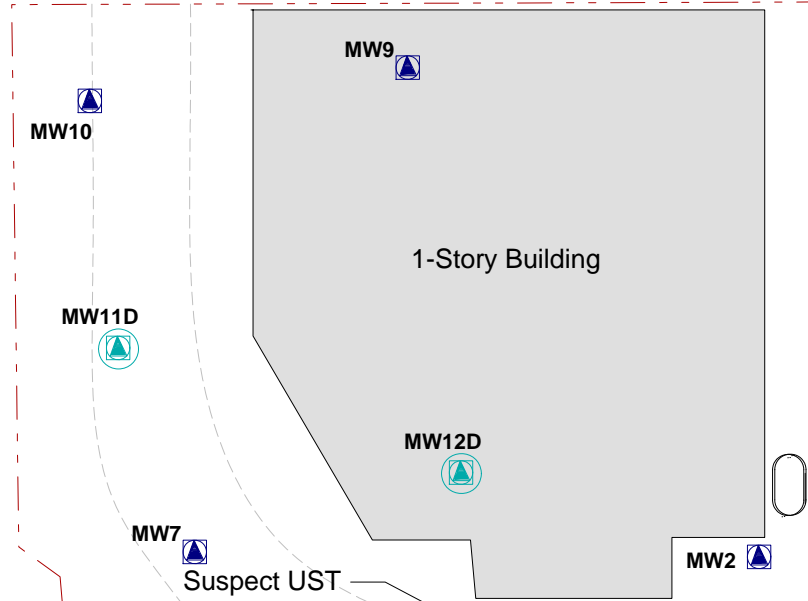


Scale: 1 inch = 50 feet

| | | | | |
|--|--|-------------------|--|---|
|  ENVIRONMENTAL BUSINESS CONSULTANTS | Phone 631.504.6000 Fax 631.924.2870 | Figure No. | | Site Name: FORMER MUGLER SHORING INC. |
| | | 4 | | Site Address: 2401 THIRD AVENUE, BRONX, NY |
| | | | | Drawing Title: SOIL BORING LOCATIONS |



Lot 109






Lot 112

Suspect USTs

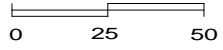
Lot 2

Third Avenue

KEY:

-  Property Line
-  MWx Monitoring Well Location
-  MWxD Deep Monitoring Well Location

SCALE:



Scale: 1 inch = 50 feet

Bulkhead

MW1

MW13D Parking Area

Easement MW5

EBC
ENVIRONMENTAL BUSINESS CONSULTANTS

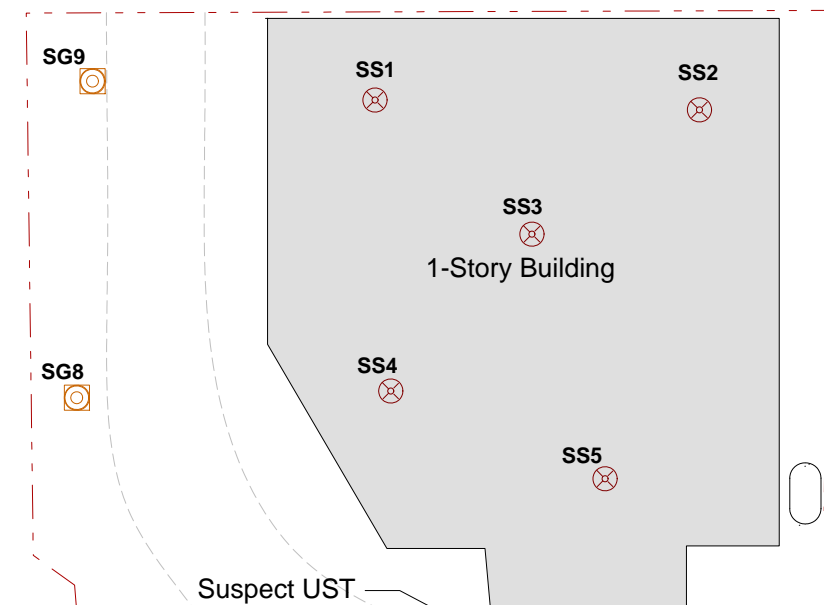
Phone 631.504.6000
Fax 631.924.2870

Figure No.
5

| |
|---|
| Site Name: FORMER MUGLER SHORING INC. |
| Site Address: 2401 THIRD AVENUE, BRONX, NY |
| Drawing Title: MONITORING WELL LOCATIONS |

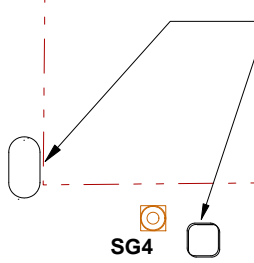


Lot 109

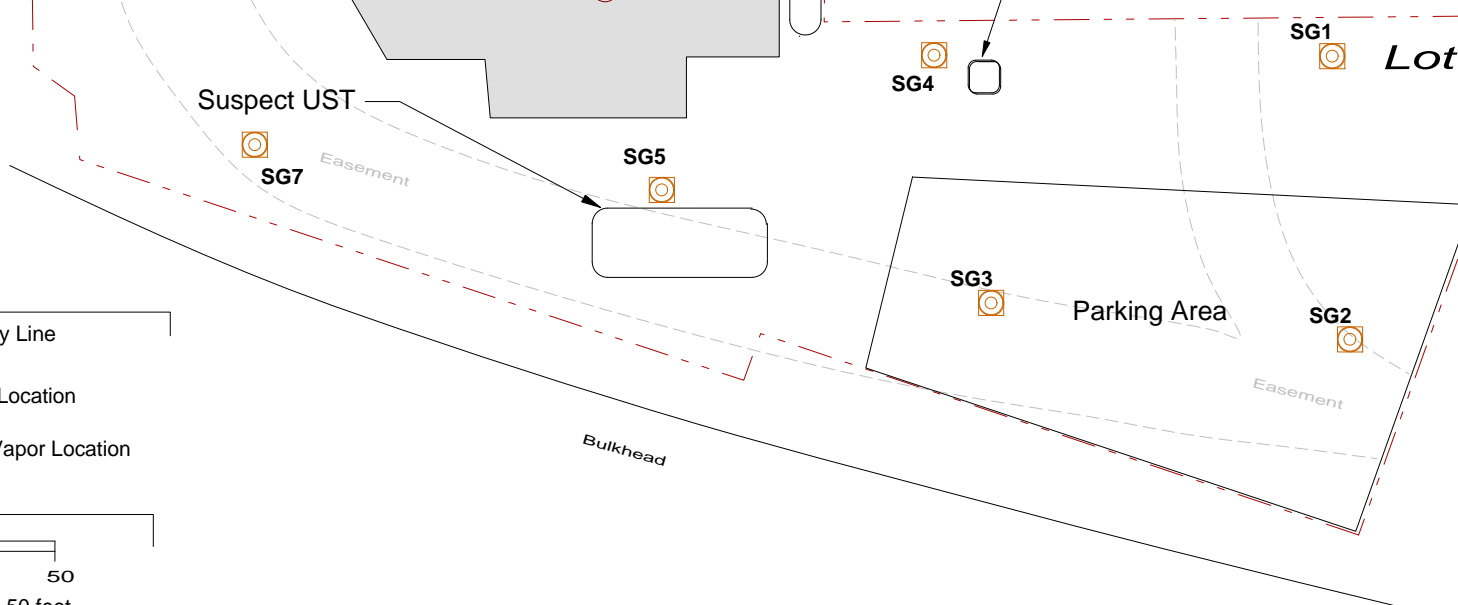


Lot 112

Suspect USTs



Lot 2



Third Avenue

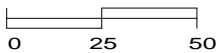
KEY:

--- Property Line

SGx Soil Gas Location

SSx Subslab Vapor Location

SCALE:



Scale: 1 inch = 50 feet

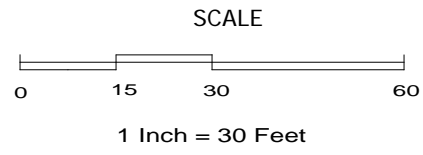
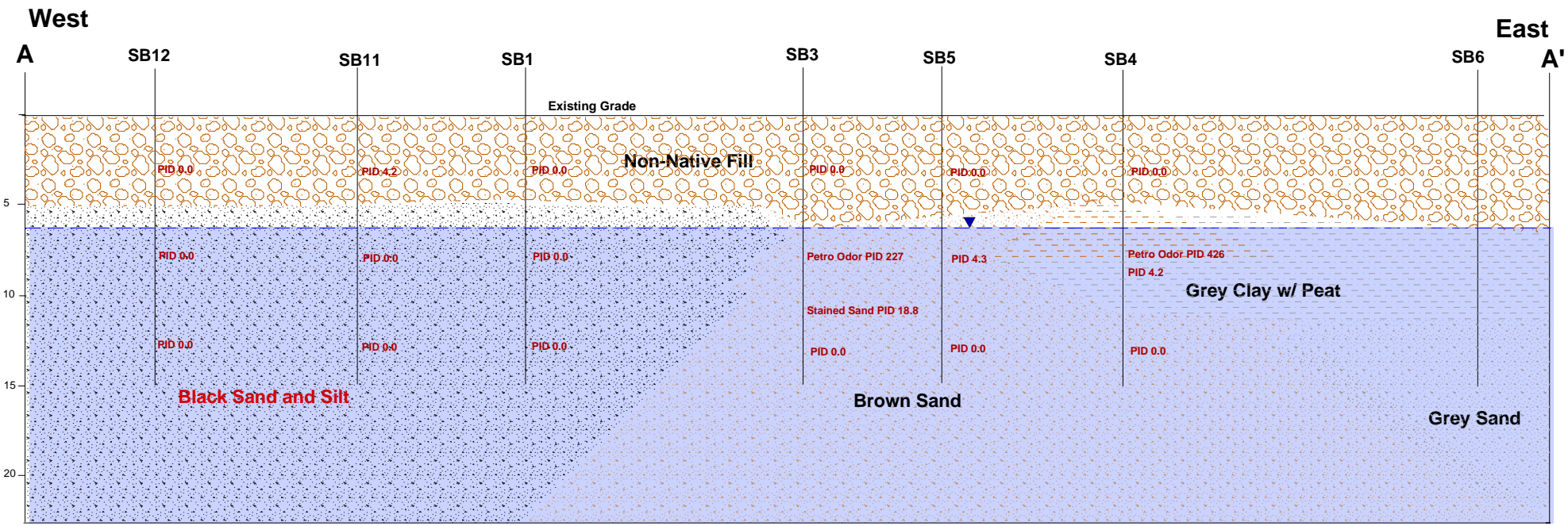
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Figure No.
6

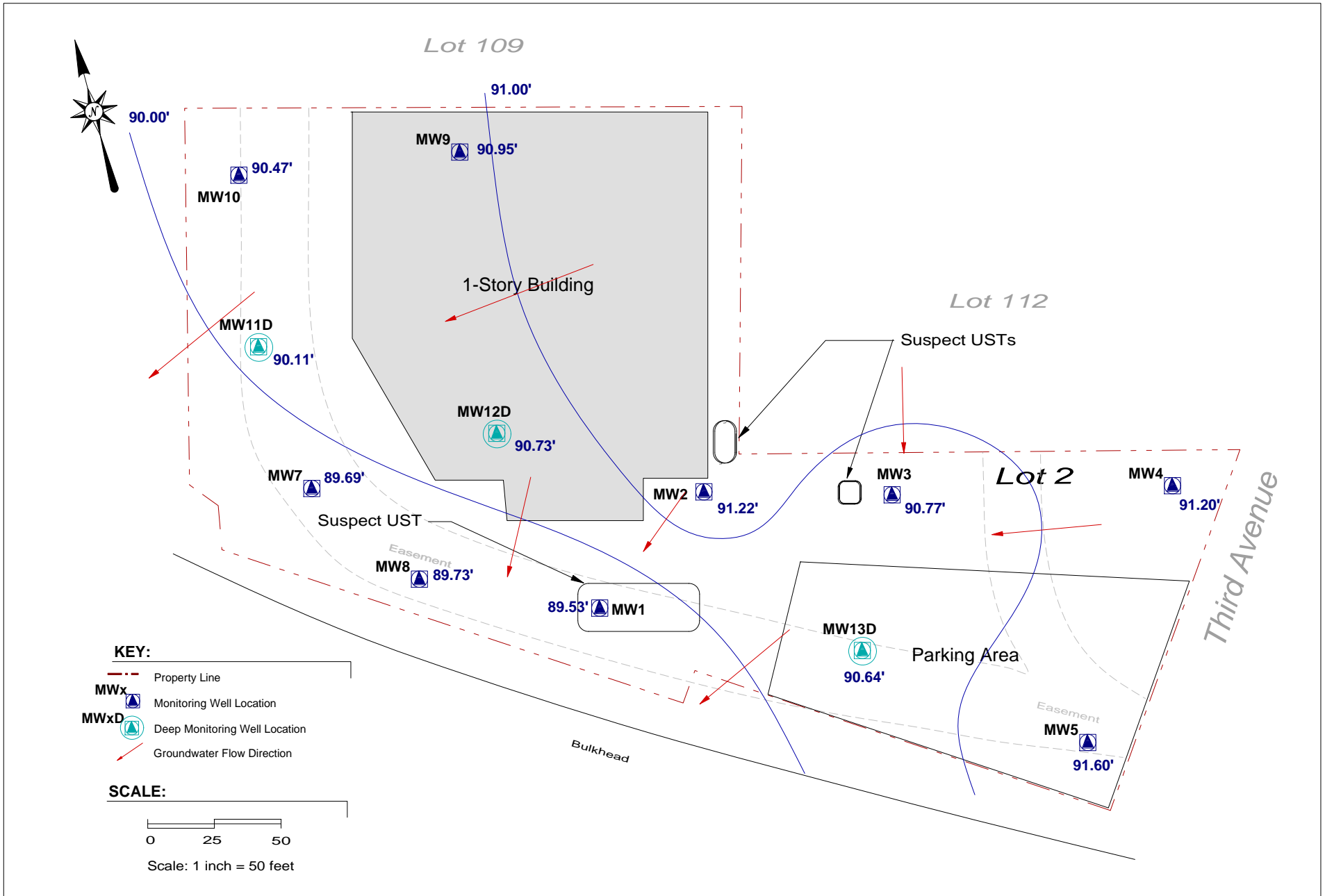
Site Name: **FORMER MUGLER SHORING INC.**


Site Address: **2401 THIRD AVENUE, BRONX, NY**

Drawing Title: **SOIL GAS SAMPLING LOCATIONS**

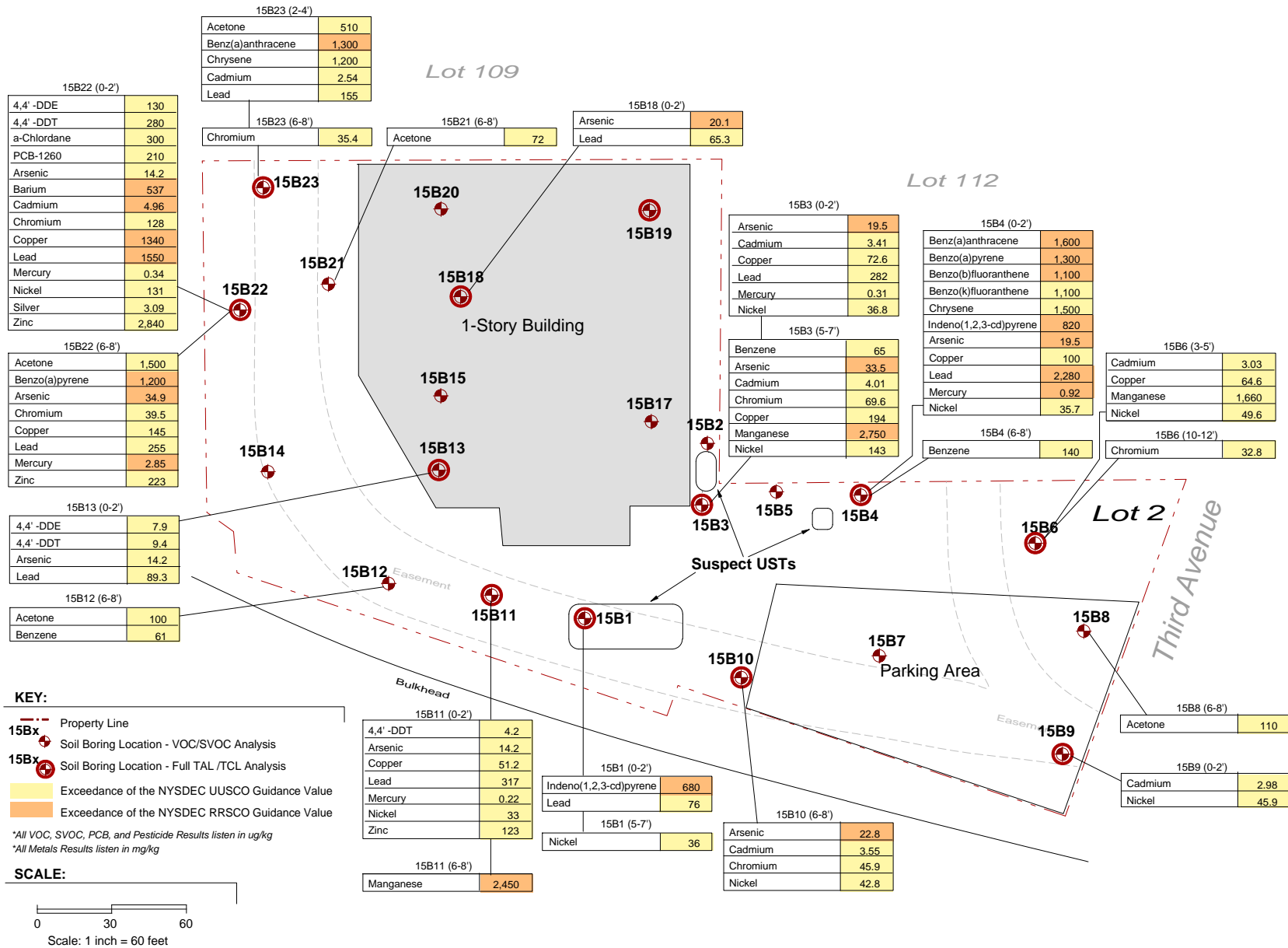


Vertical
Exaggeration
4X



| | | |
|--|-------------------------------|---|
|  | Figure No. 8 | Site Name: FORMER MUGLER SHORING INC. |
| | | Site Address: 2401 THIRD AVENUE, BRONX, NY |
| | | Drawing Title: GROUNDWATER ELEVATION MAP |

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Lot 109

MW10

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Magnesium (dissolved) | 247 |
| Manganese (dissolved) | 5.46 |
| Sodium (dissolved) | 1870 |
| Metals, Total (mg/L) | |
| Iron | 18.2 |
| Magnesium | 228 |
| Manganese | 5.35 |
| Sodium | 1,710 |

MW9

| SVOCs (ug/L) | |
|----------------------------|-------|
| Acenaphthylene | 0.05 |
| Benz(a)anthracene | 0.03 |
| Benzo(a)pyrene | 0.03 |
| Benzo(ghi)perylene | 0.04 |
| Bis(2-ethylhexyl)phthalate | 0.05 |
| Pesticides/PCBs (ug/L) | |
| PCB-1254 | 1.3 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 426 |
| Manganese (dissolved) | 1.18 |
| Sodium (dissolved) | 2,900 |
| Metals, Total (mg/L) | |
| Aluminum | 19.4 |
| Arsenic | 0.034 |
| Iron | 48.4 |
| Lead | 0.224 |
| Magnesium | 489 |
| Manganese | 2.45 |
| Sodium | 3,160 |

MW11D

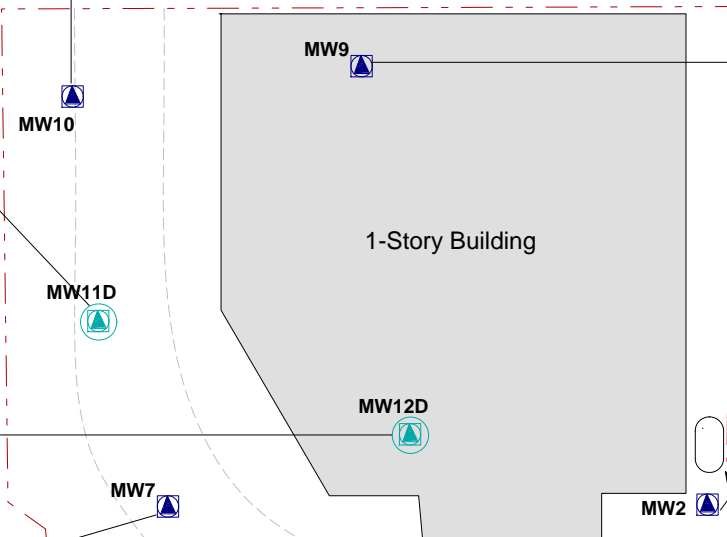
| VOCs (ug/L) | |
|--------------------------|-------|
| Methyl t-butyl ether | 14 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 37.7 |
| Sodium (dissolved) | 719 |
| Metals, Total (mg/L) | |
| Aluminum | 3.06 |
| Iron | 6.67 |
| Lead | 0.047 |
| Magnesium | 40.2 |
| Sodium | 736 |

MW12D

| VOCs (ug/L) | |
|--------------------------|-------|
| Methyl t-butyl ether | 100 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 148 |
| Manganese (dissolved) | 0.609 |
| Sodium (dissolved) | 1420 |
| Metals, Total (mg/L) | |
| Aluminum | 37.9 |
| Chromium | 0.094 |
| Iron | 78.9 |
| Lead | 0.053 |
| Magnesium | 174 |
| Manganese | 1.46 |
| Sodium | 1,290 |

MW7

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Magnesium (dissolved) | 393 |
| Manganese (dissolved) | 6.25 |
| Sodium (dissolved) | 2990 |
| Metals, Total (mg/L) | |
| Magnesium | 422 |
| Manganese | 5.61 |
| Sodium | 3,220 |



MW2

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Iron (dissolved) | 0.36 |
| Magnesium (dissolved) | 411 |
| Manganese (dissolved) | 0.672 |
| Sodium (dissolved) | 3190 |
| Metals, Total (mg/L) | |
| Aluminum | 0.41 |
| Iron | 34.3 |
| Magnesium | 399 |
| Manganese | 0.602 |
| Sodium | 2,850 |

MW3

| VOCs (ug/L) | |
|--------------------------|-------|
| Chloromethane | 39 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 293 |
| Manganese (dissolved) | 1.25 |
| Sodium (dissolved) | 1950 |
| Metals, Total (mg/L) | |
| Aluminum | 2.24 |
| Iron | 20.2 |
| Magnesium | 286 |
| Manganese | 1.25 |
| Sodium | 1,900 |

MW4

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Magnesium (dissolved) | 43.9 |
| Manganese (dissolved) | 0.643 |
| Sodium (dissolved) | 122 |
| Metals, Total (mg/L) | |
| Aluminum | 0.372 |
| Iron | 8.79 |
| Magnesium | 41.3 |
| Manganese | 0.665 |
| Sodium | 132 |

MW8

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Iron (dissolved) | 21.8 |
| Magnesium (dissolved) | 612 |
| Manganese (dissolved) | 1.04 |
| Sodium (dissolved) | 4450 |
| Metals, Total (mg/L) | |
| Iron | 48.8 |
| Magnesium | 580 |
| Manganese | 1.1 |
| Sodium | 4,830 |

MW1

| Metals, Dissolved (mg/L) | |
|--------------------------|-------|
| Magnesium (dissolved) | 688 |
| Manganese (dissolved) | 0.777 |
| Sodium (dissolved) | 5290 |
| Metals, Total (mg/L) | |
| Aluminum | 0.545 |
| Iron | 9.07 |
| Magnesium | 662 |
| Manganese | 0.758 |
| Selenium | 0.018 |
| Sodium | 5,390 |

MW13D

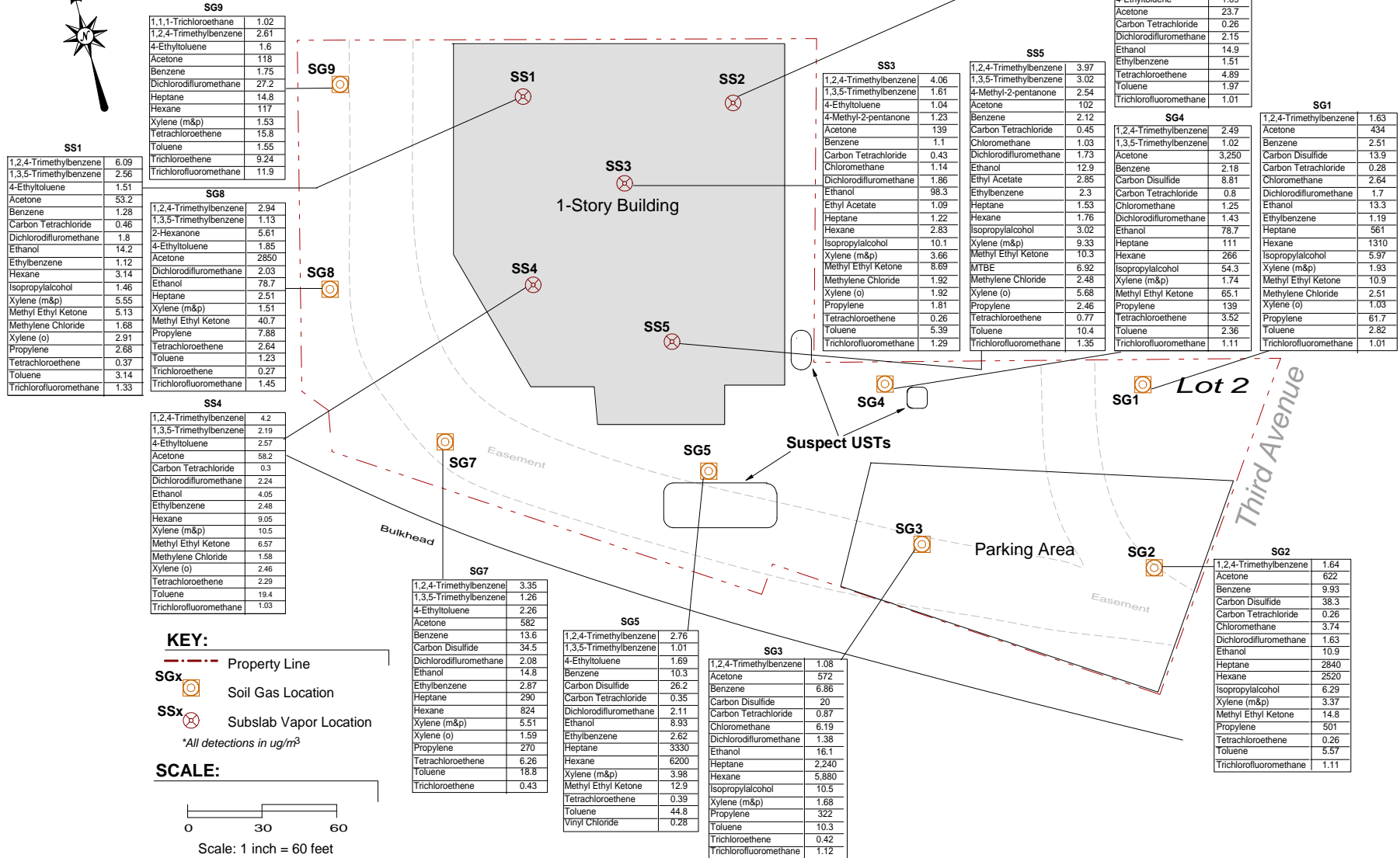
| VOCs (ug/L) | |
|--------------------------|-------|
| Chloromethane | 16 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 55.9 |
| Sodium (dissolved) | 832 |
| Metals, Total (mg/L) | |
| Aluminum | 0.997 |
| Iron | 2.94 |
| Magnesium | 60.9 |
| Sodium | 780 |

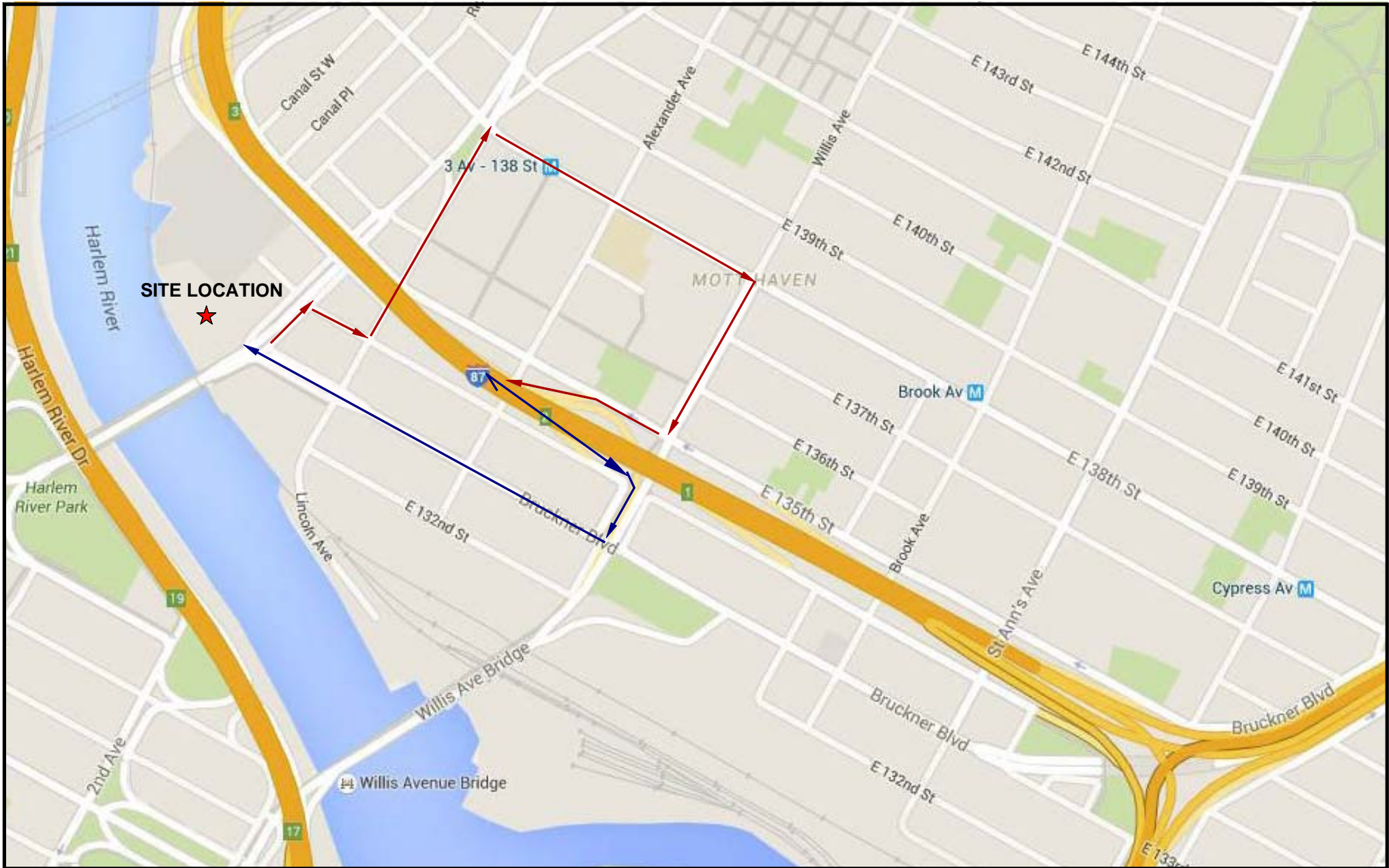
MW5

| VOCs (ug/L) | |
|--------------------------|-------|
| Chloromethane | 8.3 |
| Metals, Dissolved (mg/L) | |
| Magnesium (dissolved) | 288 |
| Manganese (dissolved) | 0.403 |
| Sodium (dissolved) | 2310 |
| Metals, Total (mg/L) | |
| Aluminum | 0.911 |
| Iron | 2.19 |
| Magnesium | 269 |
| Manganese | 0.404 |
| Sodium | 2,390 |



KEY:
 --- Property Line
 MWx Monitoring Well Location
 MWxD Deep Monitoring Well Location

SCALE:
 0 25 55
 Scale: 1 inch = 55 feet





Key

-  Truck Route to the Site
-  Truck Route from the Site



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 1808 MIDDLE COUNTRY ROAD, RIDGE, NY 11961

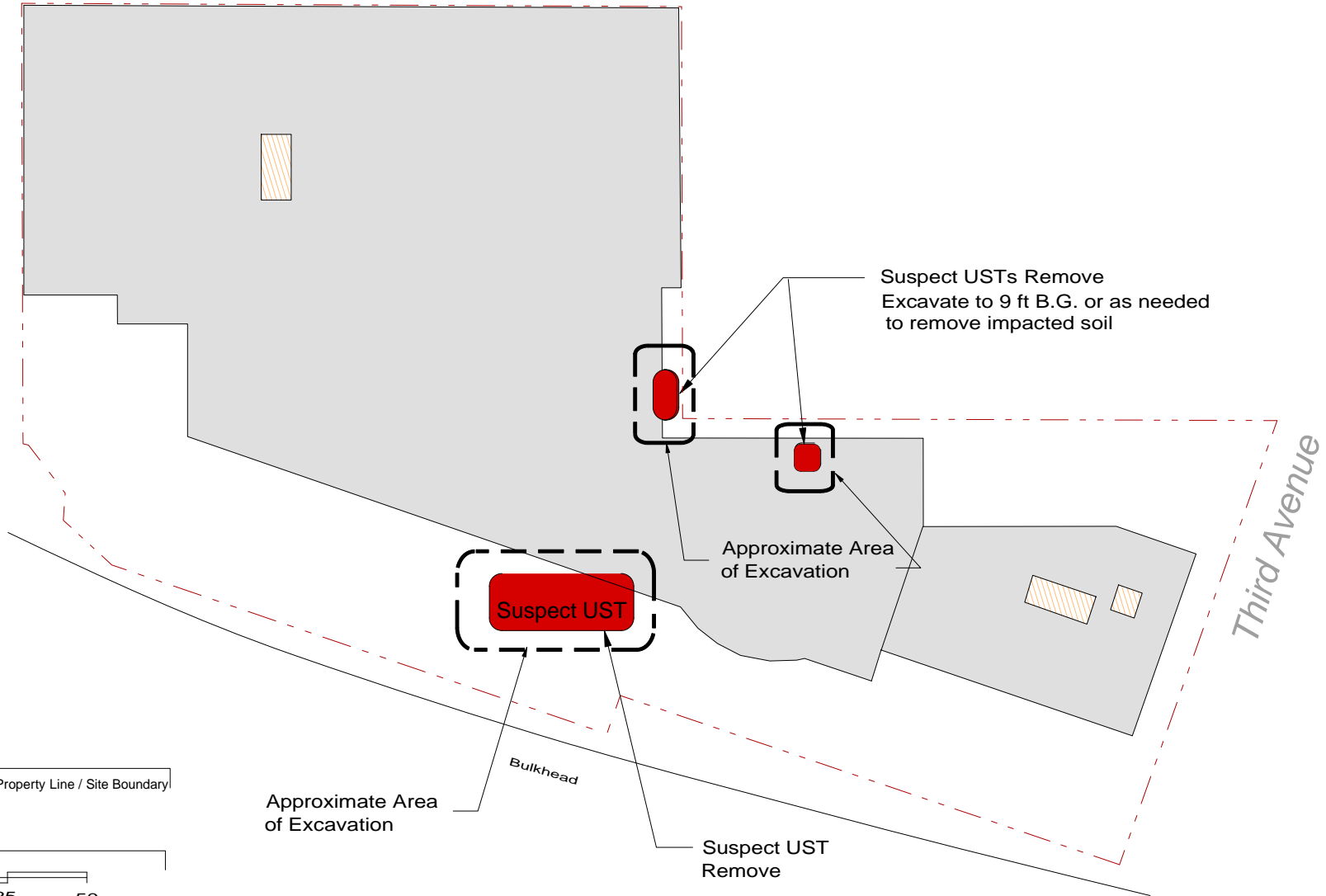
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 Fax 631.924.2780

FORMER MUGLER SHORING
 2401 THIRD AVENUE, BRONX, NY

FIGURE 12 TRUCK ROUTES



Lot 109



KEY:
- - - - - Property Line / Site Boundary

SCALE:
0 25 50
Scale: 1 inch = 50 feet

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Fax 631.924.2870

Figure No.
13

| | |
|----------------|-------------------------------------|
| Site Name: | FORMER MUGLER SHORING INC. |
| Site Address: | 2401 THIRD AVENUE, BRONX, NY |
| Drawing Title: | EXCAVATION PLAN |

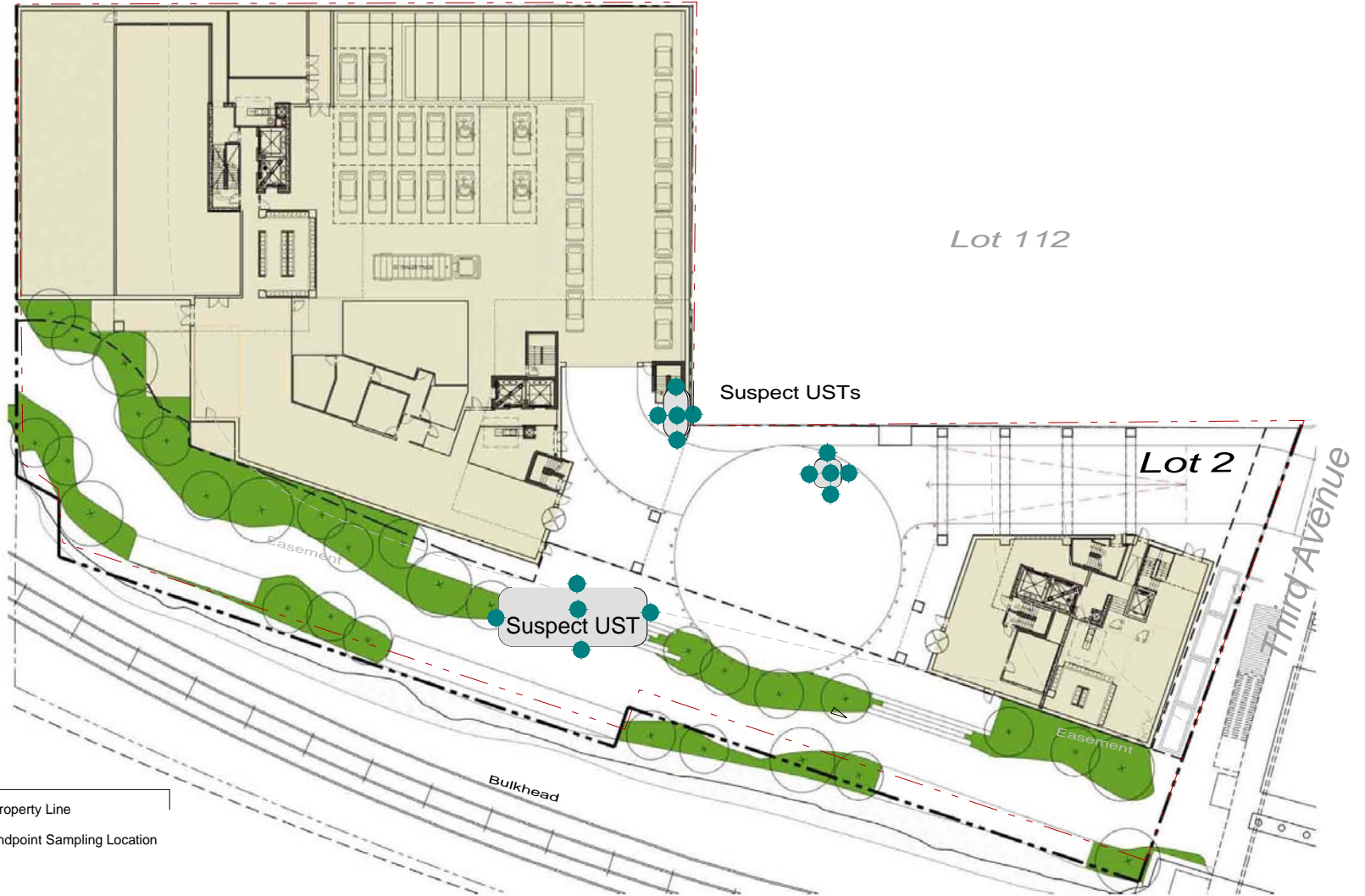


Lot 109



Lot 112

Lot 2

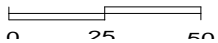
Third Avenue



KEY:

-  Property Line
-  Endpoint Sampling Location

SCALE:



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Figure No.
14

Site Name: **FORMER MUGLER SHORING INC.**
Site Address: **2401 THIRD AVENUE, BRONX, NY**
Drawing Title: **ENDPOINT SAMPLING PLAN**

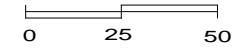


Lot 109

Tower 1
Cap w/ 6" Concrete Slab

Lot 112

SCALE:



Scale: 1 inch = 50 feet





Lot 2
Cap w/ 4-6" Concrete Slab

Tower 2

Third Avenue

Landscape Areas
Top 2 ft to meet RRSCOs

KEY:

-  Property Line
-  Cap with 6" Building Slab
-  Cap with 2' Clean Soil
-  Cap with 4-6" Concrete

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Figure No.
15

Site Name: **FORMER MUGLER SHORING INC.**

Site Address: **2401 THIRD AVENUE, BRONX, NY**

Drawing Title: **CAPPING PLAN**



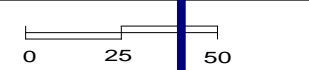
Lot 109

A

B

Lot 112

SCALE:



Scale: 1 inch = 50 feet

C

D

Lot 2

E

Third Avenue

Easement

Easement

KEY:

--- Property Line

EBC
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Figure No.
16

Site Name: **FORMER MUGLER SHORING INC.**

Site Address: **2401 THIRD AVENUE, BRONX, NY**

Drawing Title: **GRID MAP**

ATTACHMENT A
Metes and Bounds Description of Property

EXHIBIT A

Legal Description

ALL THAT CERTAIN PLOT, PIECE OR PARCEL OF LAND, WITH THE BUILDINGS AND IMPROVEMENTS THEREON ERECTED, SITUATE, LYING AND BEING IN THE BOROUGH AND COUNTY OF BRONX, CITY AND STATE OF NEW YORK, BOUNDED AND DESCRIBED AS FOLLOWS:

BEGINNING AT A POINT ON THE WESTERLY SIDE OF THIRD AVENUE, 380.21 FEET SOUTHWESTERLY FROM THE CORNER FORMED BY THE INTERSECTION OF THE WESTERLY SIDE OF THIRD AVENUE AND THE SOUTHERLY SIDE OF EAST 134TH STREET;

RUNNING THENCE NORTHWESTERLY ALONG A COURSE FORMING AN INTERIOR ANGLE OF 70 DEGREES 39 MINUTES 15 SECONDS WITH THE NORTHWESTERLY SIDE OF THIRD AVENUE 119.76 FEET TO A POINT;

THENCE SOUTHWESTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE, 6.00 FEET TO A POINT;

THENCE NORTHWESTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE, 10.00 FEET TO A POINT;

THENCE NORTHEASTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE, 6.00 FEET TO A POINT;

THENCE NORTHWESTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE, 58.17 FEET TO A POINT;

THENCE NORTHEASTERLY ALONG A LINE FORMING AN INTERIOR ANGLE OF 89 DEGREES 44 MINUTES 00 SECONDS ON ITS EASTERLY SIDE WITH THE LAST MENTIONED COURSE 129.01 FEET TO A POINT;

THENCE NORTHWESTERLY ALONG A LINE FORMING AN INTERIOR ANGLE OF 90 DEGREES 02 MINUTES 50 SECONDS ON ITS SOUTHERLY SIDE WITH THE LAST MENTIONED COURSE 107.35 FEET TO A POINT;

THENCE NORTHEASTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE 0.35 FEET TO A POINT;

THENCE NORTHWESTERLY ALONG A LINE FORMING AN INTERIOR ANGLE OF 89 DEGREES 45 MINUTES 50 SECONDS ON ITS SOUTHERLY SIDE WITH THE LAST MENTIONED COURSE 101.17 FEET TO THE SOUTHWESTERLY SIDE OF MOTT HAVEN CANAL;

THENCE SOUTHWESTERLY ALONG THE SOUTHEASTERLY SIDE OF MOTT HAVEN CANAL 146.65' TO A POINT;

THENCE SOUTHEASTERLY FORMING AN INTERIOR ANGLE OF 118 DEGREES 57 MINUTES 57 SECONDS WITH THE LAST MENTIONED COURSE A DISTANCE OF 14.93' TO A POINT;

THENCE SOUTHERLY FORMING AN INTERIOR ANGLE OF 241 DEGREES 4 MINUTES 3 SECONDS WITH THE LAST MENTIONED COURSE A DISTANCE OF 17.72' TO A POINT;

THENCE SOUTHEASTERLY FORMING AN INTERIOR ANGLE OF 108 DEGREES 52 MINUTES 16 SECONDS WITH THE LAST MENTIONED COURSE A DISTANCE OF 180.76 FEET TO A POINT;

THENCE NORTHEASTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE A DISTANCE OF 13.00' TO A POINT;

THENCE SOUTHEASTERLY AT RIGHT ANGLES TO THE LAST MENTIONED COURSE A DISTANCE OF 166.06' TO THE NORTHWESTERLY SIDE OF THIRD AVENUE;

THENCE NORTHEASTERLY ALONG THE NORTHWESTERLY SIDE OF THIRD AVENUE 148.43 FEET TO THE POINT OR PLACE OF BEGINNING.

CONTAINING 1.4 ACRES (61,034.98 SQUARE FEET).

TOGETHER WITH THE BENEFITS OF CERTAIN EASEMENTS OF INGRESS, EGRESS AND REGRESS AS DESCRIBED IN DEED BETWEEN GENHAR REALTY CORP. AND CANAL & 3RD CO. RECORDED 4/6/1961 IN LIBER 2430 AT PAGE 382.

TOGETHER WITH THE BENEFITS OF CERTAIN EASEMENTS OF INGRESS, EGRESS AND REGRESS AS DESCRIBED IN DEED BETWEEN CANAL & 3RD CO. AND MAE-JOE REALTY CORP. INC. RECORDED 4/6/1962 IN LIBER 2479 AT PAGE 383.

TOGETHER WITH THE BENEFITS OF CERTAIN EASEMENTS AS TO GAS, ELECTRICITY AND WATER AS DESCRIBED IN DEED BETWEEN RICHARD C. MUGLER, JR. AND KAI DEVELOPMENT CORP. RECORDED 12/30/1997 IN REEL 1519 PAGE 1925.

BEING THE SAME PREMISES AS DESCRIBED IN DEED FROM RICHARD C. MUGLER, JR. TO 2401 3RD AVENUE, L.L.C. RECORDED IN REEL 1785 AT PAGE 871.

ATTACHMENT B
Health and Safety Plan

FORMER MUGLER SHORING INC. SITE
SITE No. C203052
2401 THIRD AVENUE, BRONX, NY
Block 2319 Lot 2

CONSTRUCTION
HEALTH AND SAFETY PLAN

APRIL 2016

Prepared for:
2401 3rd Ave Associates Property LLC
512 Seventh Avenue 15th Floor
New York, NY 10018


Prepared by:

ENVIRONMENTAL BUSINESS CONSULTANTS
1808 Middle Country Road
Ridge, NY 11961

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Former Mugler Shoring Site
2401 Third Avenue, Bronx, New York

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2401 Third Avenue, Bronx, New York

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Figure 1 Route to Hospital (Appendix D)

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APPENDIX B SITE SAFETY PLAN AMENDMENTS
APPENDIX C CHEMICAL HAZARDS
APPENDIX D HOSPITAL INFORMATION, MAP AND FIELD ACCIDENT REPORT

STATEMENT OF COMMITMENT

This Construction Health and Safety Plan (CHASP) has been prepared to ensure that workers are not exposed to risks from hazardous materials during the Remedial Actions at the Former Mugler Shoring BCP Site, located at 2401 Third Avenue, Bronx, NY.

This CHASP, which applies to persons present at the site actually or potentially exposed to hazardous materials, describes emergency response procedures for actual and potential chemical hazards. This CHASP is also intended to inform and guide personnel entering the work area or exclusion zone. Persons are to acknowledge that they understand the potential hazards and the contents of this Health and Safety policy by signing off on receipt of their individual copy of the document. Contractors and suppliers are retained as independent contractors and are responsible for ensuring the health and safety of their own employees.

1.0 INTRODUCTION AND SITE ENTRY REQUIREMENTS

This document describes the health and safety guidelines developed by Environmental Business Consultants (EBC) for the planned Remedial Action at the Former Mugler Shoring BCP Site, located at 2401 Third Avenue, Bronx, NY, to protect on-site personnel, visitors, and the public from physical harm and exposure to hazardous materials or wastes during remedial activities. In accordance with the Occupational Safety and Health Administration (OSHA) 29 CFR Part 1910.120 Hazardous Waste Operations and Emergency Response Final rule, this CHASP, including the attachments, addresses safety and health hazards related to excavation, loading and other soil disturbance activities and is based on the best information available. The CHASP may be revised by EBC at the request of the owner and/or a regulatory agency upon receipt of new information regarding site conditions. Changes will be documented by written amendments signed by EBC's project manager, site safety officer and/or the EBC health and safety consultant.

Work performed under the remedial action will not involve confined space entry since the excavations will be large and sloped back in accordance with NYCDOB shoring requirements and will not have a limited or restricted means for entry or exit.

1.1 Training Requirements

Personnel entering the exclusion zone or decontamination zone are required to be certified in health and safety practices for hazardous waste site operations as specified in the Federal OSHA Regulations CFR 1910.120e (revised 3/6/90).

Paragraph (e - 3) of the above referenced regulations requires that all on-site management personnel directly responsible for or who supervise employees engaged in hazardous waste operations, must initially receive 8 hours of supervisor training related to managing hazardous waste work.

Paragraph (e - 8) of the above referenced regulations requires that workers and supervisors receive 8 hours of refresher training annually on the items specified in Paragraph (e-1) and/or (e-3).

Additionally all on-site personnel must receive adequate site-specific training in the form of an on-site Health and Safety briefing prior to participating in field work with emphasis on the following:

- Protection of the adjacent community from hazardous vapors and / or dust which may be released during intrusive activities.
- Identification of chemicals known or suspected to be present on-site and the health effects and hazards of those substances.
- The need for vigilance in personnel protection, and the importance of attention to proper use, fit and care of personnel protective equipment.
- Decontamination procedures.
- Site control including work zones, access and security.
- Hazards and protection against heat or cold.
- The proper observance of daily health and safety practices, such as entry and exit of work

- zones and site. Proper hygiene during lunch, break, etc.
- Emergency procedures to be followed in case of fire, explosion and sudden release of hazardous gases.

Health and Safety meetings will be conducted on a daily basis and will cover protective clothing and other equipment to be used that day, potential and chemical and physical hazards, emergency procedures, and conditions and activities from the previous day.

1.2 Medical Monitoring Requirements

Field personnel and visitors entering the exclusion zone or decontamination zone must have completed appropriate medical monitoring required under OSHA 29 CFR 1910.120(f) if respirators or other breathing related PPE is needed. Medical monitoring enables a physician to monitor each employee's health, physical condition, and his fitness to wear respiratory protective equipment and carry out on-site tasks.

1.3 Site Safety Plan Acceptance, Acknowledgment and Amendments

The project superintendent and the site safety officer are responsible for informing personnel (EBC employees and/or owner or owners representatives) entering the work area of the contents of this plan and ensuring that each person signs the safety plan acknowledging the on-site hazards and procedures required to minimize exposure to adverse effects of these hazards. A copy of the Acknowledgement Form is included in **Appendix A**.

Site conditions may warrant an amendment to the CHASP. Amendments to the CHASP are acknowledged by completing forms included in **Appendix B**.

1.4 Key Personnel - Roles and Responsibilities

Personnel responsible for implementing this Health and Safety Plan are:

| Name | Title | Address | Contact Numbers |
|---------------------|-------------------------|---|-----------------|
| Mr. Robert Bennett | EBC – Project Manager | 1808 Middle Country Rd Ridge, NY 11961 | (631) 504-6000 |
| Ms. Chawinie Miller | Health & Safety Manager | 1808 Middle Country Rd Ridge, NY 11961 | (631) 504-6000 |
| Mr. Kevin Waters | Site Safety Officer | 1808 Middle Country Rd Ridge, NY 11961 | (631) 504-6000 |

The project manager is responsible for overall project administration and, with guidance from the site safety officer, for supervising the implementation of this CHASP. The site safety officer will conduct daily (tail gate or tool box) safety meetings at the project site and oversee daily safety issues. Each subcontractor and supplier (defined as an OSHA employer) is also responsible for the health and safety of its employees. If there is any dispute about health and safety or project activities, on-site personnel will attempt to resolve the issue. If the issue cannot be resolved at the site, then the project manager will be consulted.

The site safety officer is also responsible for coordinating health and safety activities related to hazardous material exposure on-site. The site safety officer is responsible for the following:

1. Educating personnel about information in this CHASP and other safety requirements to be observed during site operations, including, but not limited to, decontamination procedures, designation of work zones and levels of protection, air monitoring, fit testing, and emergency procedures dealing with fire and first aid.
2. Coordinating site safety decisions with the project manager.
3. Designating exclusion, decontamination and support zones on a daily basis.
4. Monitoring the condition and status of known on-site hazards and maintaining and implementing the air quality monitoring program specified in this CHASP.
5. Maintaining the work zone entry/exit log and site entry/exit log.
6. Maintaining records of safety problems, corrective measures and documentation of chemical exposures or physical injuries (the site safety officer will document these conditions in a bound notebook and maintain a copy of the notebook on-site).

The person who observes safety concerns and potential hazards that have not been addressed in the daily safety meetings should immediately report their observations/concerns to the site safety officer or appropriate key personnel.

2.0 SITE BACKGROUND AND SCOPE OF WORK

The street address for the Site is 2401 Third Avenue, Bronx, NY. The Site is located in the City of New York in the Port Morris neighborhood of the Borough of the Bronx. The Site is comprised of a single tax parcel covering 61,034.98 (1.4 acres). The subject property is located in the City of New York and Borough of the Bronx (Bronx County). The lot is located on the west side of Third Avenue and is identified as Block 2319 Lot 2 on the NY City tax map. The property has 159 feet of street frontage on Third Avenue and approximately 346 ft of frontage along the north side of the Harlem River.

The property is currently vacant but was must recently used by a shoring company and has historically been used as an iron works, steel fabrication, gear manufacturing, gasket manufacturing and shoring equipment fabrication.

2.1 Previous Investigations

2.1.1 October 2014 – Phase I Environmental Site Assessment (Langan)

A Phase I Environmental Site Assessment Report was prepared by Langan Engineering, Environmental, Surveying and Landscape Architecture, D.P.C. (Langan) on October 28, 2014. Based upon reconnaissance of the subject and surrounding properties, interviews and review of historical records and regulatory agency databases, Langan identified the following recognized environmental conditions:

- **Current and Historical Manufacturing and Industrial Use:**
The Subject Property was historically used for manufacturing purposes since at least 1891. Historical operators include; J.L. Mott Iron Works (1891–1922), Hydraulic Steel Company (1922–1935), General Builders Supply Corporation (1935–1968), Brill Equipment Company (1949–1956), US Gear Manufacturing Company (1965–1971), Ohio Gasket Manufacturing Corporation (1971), and Mugler Inc. (1965–present). Inadvertent releases of solvents, petroleum products, metals, polychlorinated biphenyls (PCB) and/or other chemicals used during manufacturing operations may have adversely impacted soil, groundwater, building components and/or soil vapor. The Subject Property is presently operated by Mugler, Inc. for shoring equipment fabrication, storage, truck loading/unloading, and equipment repairs. As such, there are typical tools, maintenance/repair materials, and miscellaneous equipment used and stored throughout. Discoloration and staining were apparent throughout the interior of the building, suggesting incidental releases of petroleum products during truck maintenance, and the concrete slab was compromised in several areas. Because fractures in the slab provide a conduit for spilled motor oils and/or petroleum products to impact subsurface conditions, current use by Mugler, Inc. constitutes a REC.
- **On-Site Petroleum Bulk Storage:**
The following historical underground storage tanks (USTs) were identified:
 - One 550-gallon gasoline UST, located outside the southeast corner of the building, was reportedly closed-in place circa 1999.

- An area of patched concrete and an apparent abandoned fill port southeast of the building, which is suspected to be associated with another decommissioned UST, was identified during the site reconnaissance.
- One 1,550-gallon gasoline UST was identified on the 1935 through 1946 Sanborn maps.

Based on the known and suspected presence of historic tanks and lack of any documentation of tank closure, the historical USTs are a REC.

- **Current and Historical Use of Surrounding Properties:**
Include a private garage (1946–1951), lubricating oils storage (1935–1983), two chemical corporations (1935–1947), a paint manufacturer (1951), a coal yard (1935–1951), an auto building (1908–1947), an auto house (1935–1944), a private garage (1935–1944), a printing facility (1935–1947), and several auto repair shops (2005–2012). In addition, an active New York State Department of Environmental Conservation (NYSDEC) Brownfield Cleanup Program (BCP) site was identified approximately 660 feet northeast (up-gradient) of the Subject Property. Based on investigations conducted to date, the primary contaminants of concern in soil and groundwater are petroleum related compounds. No information was provided with respect to the off-Site migration of contaminants; however, the Environmental Database Report (EDR) report indicates that the potential exists for off-Site migration of site-related contaminants in soil vapor. Potential petroleum and solvent releases associated with the historical surrounding property uses may have adversely impacted groundwater and/or soil vapor at the Subject Property and is considered a REC.

2.1.2 October 2014 - Phase II Investigation Report (Langan)

A remedial investigation was performed by Langan for the property in July and August of 2014. The Phase II included the following scope of work:

1. Completion of a geophysical investigation;
2. Completion of seven soil borings and collection of 14 soil samples;
3. Installation and sampling of six temporary monitoring wells; and
4. Installation and sampling of six soil vapor points. Installed five soil vapor probes around Site perimeter and collected five samples for chemical analysis.

The findings of the Phase II are summarized as follows:

- The geophysical survey identified two major anomaly areas along the southeast corner of the Subject Property Building. Based on their reflection rates and their proximities, these anomalies are indicative of USTs located approximately 2 to 3 ft bgs. Another anomaly area consistent with disturbed soil was located along the western edge of the Subject Property.
- A layer of historic fill was identified throughout the Subject Property with thicknesses ranging from approximately 5 to 13 feet. This fill layer contained SCO exceedances (SVOC and metal) throughout the majority of the Site.
- The suite of metals and SVOCs and their concentrations detected in soil and groundwater

across the site are attributed to historic fill material.

- One of the seven soil borings (SB-3) was installed next to a suspect 550-gallon gasoline UST that was reportedly decommissioned in 1999. Petroleum-like odors and PID readings (up to 700 ppm) were observed in the soil boring at depths of approximately 8 to 9 ft bgs. Groundwater analytical sampling results from this location confirmed the presence of a petroleum release. Article 12 of the New York State Navigation Law requires that anyone with knowledge of a petroleum release must report the discovery to the New York State Spill Hotline as soon as possible. Therefore, on 13 August, 2014, Langan called in a spill and Spill Case No. 1405230 was opened by NYSDEC.
- Soil vapor concentrations are below NYSDOH AGV values; however, based on the presence of VOCs in the soil vapor, remedial measures should include installation of a soil vapor mitigation system.
- Due to the E-designation of the site, additional investigation including soil, groundwater and soil vapor sampling will be required in order to comply with NYCOER guidelines and obtain permits for construction and occupancy of proposed building structures.
- If site development is planned, a Remedial Action Plan (RAP) and Construction Health and Safety Plan (CHASP) should be prepared and implemented. In accordance with the requirements of the E-Designation program, these documents must be submitted to OER for approval prior to the start of Site work. OER will require additional site investigation to support preparation of the RAP. The RAP and CHASP may need to address the following:
 - Handling, transport and disposal of excess material including fill and debris as required for development with all appropriate requirements.
 - A composite cover system for impacted material that may be left in-place after development. The cover system will prevent direct contact with or inhalation of contaminated material.
 - Odor and soil vapor intrusion should be addressed if odorous or VOC-impacted material is left in-place as part of the development.
 - Registration, removal and closure of known petroleum storage tanks.
 - Permitting and treatment, if required, of dewatering effluent.
 - Health and safety requirements addressing the contaminants of concern including a Community Air Monitoring Plan (CAMP) to monitor odor and dust.
 - A contingency plan for removal of currently unknown underground storage tanks or other subsurface structures

2.2 Redevelopment Plans

The redevelopment project consists of the construction of two new 25-story residential tower buildings identified as the east and the west towers. The two towers flank an entry court which is open to the river and Manhattan views to the south. The East Tower is free-standing, serving as a “portal” on Third Avenue, whereas the West Tower has a 7-story base with common amenities and parking for 200 cars. The residential component consists of 475 rental units in an area of 399,798 gross square feet. Commercial (retail) and community space will be included in the west tower base.

2.3 Description of Remedial Action

Site activities included within the Remedial Action that are included within the scope of this HASP include the following:

1. Excavation of soil/fill exceeding groundwater protection SCOs for those VOC parameters in groundwater above standards to depths as great as 9 feet below grade in the UST source area;
2. Screening for indications of contamination (by visual means, odor, and monitoring with PID) of all excavated soil during any intrusive Site work;
3. Excavation and disposal of historic fill materials as needed for installation of the new buildings basement level foundations.;
4. Excavation and disposal of historic fill materials from planned landscaped (exposed soil) areas;
5. Collection and analysis of end-point samples to evaluate the performance of the remedy with respect to attainment of groundwater protection SCOs for VOCs present in groundwater above standards;
6. Appropriate off-Site disposal of all material removed from the Site in accordance with all Federal, State and local rules and regulations for handling, transport, and disposal;
7. Capping the entire Site with the building foundations, concrete walkways / driveways or 2 ft of soil meeting Restricted Residential SCOs.
8. Import of materials to be used for backfill and cover in compliance with: (1) chemical limits and other specifications, (2) all Federal, State and local rules and regulations for handling and transport of material.
9. Implementation of a Site Management Plan (SMP) for long term maintenance of the Engineering Controls.
10. An Environmental Easement will be filed against the Site to ensure implementation of the SMP.

3.0 HAZARD ASSESSMENT

This section identifies the hazards associated with the proposed scope of work, general physical hazards that can be expected at most sites; and presents a summary of documented or potential chemical hazards at the site. Every effort must be made to reduce or eliminate these hazards. Those that cannot be eliminated must be guarded against using engineering controls and/or personal protective equipment.

3.1 Physical Hazards

3.1.1 Tripping Hazards

An area of risk associated with on-site activities are presented by uneven ground, concrete, curbstones or equipment which may be present at the site thereby creating a potential tripping hazard. During intrusive work, care should be taken to mark or remove any obstacles within the exclusion zone.

3.1.2 Climbing Hazards

During site activities, workers may have to work on excavating equipment by climbing. The

excavating contractor will conform with any applicable NIOSH and OSHA requirements or climbing activities.

3.1.3 Cuts and Lacerations

Field activities that involve excavating activities usually involve contact with various types of machinery. A first aid kit approved by the American Red Cross will be available during all intrusive activities.

3.1.4 Lifting Hazards

Improper lifting by workers is one of the leading causes of industrial injuries. Field workers in the excavation program may be required to lift heavy objects. Therefore, all members of the field crew should be trained in the proper methods of lifting heavy objects. All workers should be cautioned against lifting objects too heavy for one person.

3.1.5 Utility Hazards

Before conducting any excavation, the excavation contractor will be responsible for locating and verifying all existing utilities at each excavation.

3.1.6 Traffic Hazards

All traffic, vehicular and pedestrian, shall be maintained and protected at all times consistent with local, state and federal agency regulations regarding such traffic and in accordance with NYCDOT guidelines. The excavation contractor shall carry on his operations without undue interference or delays to traffic. The excavation contractor shall furnish all labor, materials, guards, barricades, signs, lights, and anything else necessary to maintain traffic and to protect his work and the public, during operations.

3.2 Work in Extreme Temperatures

Work under extremely hot or cold weather conditions requires special protocols to minimize the chance that employees will be affected by heat or cold stress.

3.2.1 Heat Stress

The combination of high ambient temperature, high humidity, physical exertion, and personal protective apparel, which limits the dissipation of body heat and moisture, can cause heat stress.

The following prevention, recognition and treatment strategies will be implemented to protect personnel from heat stress. Personnel will be trained to recognize the symptoms of heat stress and to apply the appropriate treatment.

1. Prevention

- a. Provide plenty of fluids. Available in the support zone will be a 50% solution of fruit punch and water or plain water.
- b. Work in Pairs. Individuals should avoid undertaking any activity alone.
- c. Provide cooling devices. A spray hose and a source of water will be provided to reduce body temperature, cool protective clothing and/or act as a quick-drench shower in case of an exposure incident.
- d. Adjustment of the work schedule. As is practical, the most labor-intensive tasks should be carried out during the coolest part of the day.

2. Recognition and Treatment

a. Heat Rash (or prickly heat):

Cause: Continuous exposure to hot and humid air, aggravated by chafing clothing.

Symptoms: Eruption of red pimples around sweat ducts accompanied by intense itching and tingling.

Treatment: Remove source of irritation and cool skin with water or wet cloths.

b. Heat Cramps (or heat prostration)

Cause: Profuse perspiration accompanied by inadequate replenishment of body water and electrolytes.

Symptoms: Muscular weakness, staggering gait, nausea, dizziness, shallow breathing, pale and clammy skin, approximately normal body temperature.

Treatment: Perform the following while making arrangement for transport to a medical facility. Remove the worker to a contamination reduction zone. Remove protective clothing. Lie worker down on back in a cool place and raise feet 6 to 12 inches. Keep warm, but loosen all clothing. If conscious, provide sips of salt-water solution, using one teaspoon of salt in 12 ounces of water. Transport to a medical facility.

c. Heat Stroke

Cause: Same as heat exhaustion. This is also an extremely serious condition.

Symptoms: Dry hot skin, dry mouth, dizziness, nausea, headache, rapid pulse.

Treatment: Cool worker immediately by immersing or spraying with cool water or sponge bare skin after removing protective clothing. Transport to hospital.

3.2.2 Cold Exposure

Exposure to cold weather, wet conditions and extreme wind-chill factors may result in excessive loss of body heat (hypothermia) and /or frostbite. To guard against cold exposure and to prevent cold injuries, appropriate warm clothing should be worn, warm shelter must be readily available, rest periods should be adjusted as needed, and the physical conditions of on-site field personnel should be closely monitored. Personnel and supervisors working on-site will be made aware of the signs and symptoms of frost bite and hypothermia such as:

- Shivering;
- reduced blood pressure;
- reduced coordination;
- drowsiness;
- impaired judgment;
- fatigue;
- pupils dilated but reactive to light; and,
- numbing of the toes and fingers.

3.3 Chemical Hazards

“Urban fill” materials, present throughout the New York City area typically contain elevated levels of semi-volatile organic compounds and metals. These “contaminants” are not related to a chemical release occurring on the site, but are inherent in the reworked fill material in the area which contains ash and bits of tar and asphalt. Considering the previous sampling results and the past and present use of the site, the following compounds are considered for the site as potential contaminants: volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), pesticides, polychlorinated biphenyl’s (PCBs), and heavy metals such as arsenic, chromium, lead and mercury.

Based on the findings of the Remedial Investigation and the inherent properties of urban fill, the following compounds are considered for the site as potential contaminants: volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), pesticides, and heavy metals.

Volatile organic compounds reported to be present in soil include the following:

| |
|---------|
| benzene |
|---------|

Semi-Volatile organic compounds reported to be present in soil and / or fill materials include the following:

| | | | |
|------------------------|----------------------|----------------|----------|
| Benzo(a)anthracene | Benzo(b)fluoranthene | Benzo(a)pyrene | Chrysene |
| Ideno(1,2,3-cd) pyrene | | | |

PCBs and Pesticides reported to be present in soil and / or fill materials include the following:

| | | | | |
|-----------|-----------|-----------|-------------|------------|
| 4'-4'-DDD | 4'-4'-DDE | 4'-4'-DDT | a-Chlordane | PCB - 1260 |
|-----------|-----------|-----------|-------------|------------|

Metals reported to be present in fill materials include the following:

| | | | | | |
|-----------|---------|---------|----------|--------|------|
| Arsenic | Barium | Cadmium | Chromium | Copper | Lead |
| Manganese | Mercury | Nickel | Silver | Zinc | |

The primary routes of exposure to these contaminants are inhalation, ingestion and absorption. **Appendix C** includes information sheets for suspected chemicals that may be encountered at the site.

3.3.1 Respirable Dust

Dust may be generated from vehicular traffic and/or excavation activities. If visible observation detects elevated levels of dust, a program of wetting will be employed by the site safety officer. If elevated dust levels persist, the site safety office will employ dust monitoring using a particulate monitor (Miniram or equivalent). If monitoring detects concentrations greater than 150 µg/m³ over daily background, the site safety officer will take corrective actions as defined herein, including the use of water for dust suppression and if this is not effective, requiring workers to wear APRs with efficiency particulate air (HEPA) cartridges.

Absorption pathways for dust and direct contact with soils or groundwater will be mitigated with the implementation of latex gloves, hand washing and decontamination exercises when necessary.

3.3.2 Dust Control and Monitoring During Earthwork

Dust generated during excavation activities or other earthwork may contain contaminants identified in soils at the site. Dust will be controlled by wetting the working surface with water. Calcium chloride may be used if the problem cannot be controlled with water. Air monitoring and dust control techniques are specified in a site specific Dust Control Plan (if applicable). Site workers will not be required to wear APR's unless dust concentrations are consistently over 150 $\mu\text{g}/\text{m}^3$ over site-specific background in the breathing zone as measured by a dust monitor unless the site safety officer directs workers to wear APRs. The site safety officer will use visible dust as an indicator to implement the dust control plan.

3.3.3 Organic Vapors

Elevated levels of chlorinated VOCs were detected in soil, soil gas and groundwater samples collected during previous investigations at the site. Therefore, excavation activities may cause the release of organic vapors to the atmosphere. The site safety officer will periodically monitor organic vapors with a Photoionization Detector (PID) during excavation activities to determine whether organic vapor concentrations exceed action levels shown in Section 5 and/or the Community Air Monitoring Plan.

4.0 PERSONAL PROTECTIVE EQUIPMENT

Personal protective equipment (PPE) shall be selected in accordance with the site air monitoring program, OSHA 29 CFR 1910.120(c), (g), and 1910.132. Protective equipment shall be NIOSH approved and respiratory protection shall conform to OSHA 29 CFR Part 1910.133 and 1910.134 specifications; head protection shall conform to 1910.135; eye and face protection shall conform to 1910.133; and foot protection shall conform to 1910.136. The only true difference among the levels of protection from D thru B is the addition of the type of respiratory protection. **It is anticipated that work will be performed in Level D PPE.**

4.1 Level D

Level D PPE shall be donned when the atmosphere contains no known hazards and work functions preclude splashes, immersion, or the potential for inhalation of, or contact with, hazardous concentrations of harmful chemicals. Level D PPE consists of:

- standard work uniform, coveralls, or tyvek, as needed;
- steel toe and steel shank work boots;
- hard hat;
- gloves, as needed;
- safety glasses;
- hearing protection;
- equipment replacements are available as needed.

4.2 Level C

Level C PPE shall be donned when the concentrations of measured total organic vapors in the breathing zone exceed background concentrations (using a portable OVA, or equivalent), but are less than 5 ppm. The specifications on the APR filters used must be appropriate for contaminants identified or expected to be encountered. Level C PPE shall be donned when the identified contaminants have adequate warning properties and criteria for using APR have been met. Level C PPE consists of:

- chemical resistant or coated tyvek coveralls;
- steel-toe and steel-shank workboots;
- chemical resistant overboots or disposable boot covers;
- disposable inner gloves (surgical gloves);
- disposable outer gloves;
- full face APR fitted with organic vapor/dust and mist filters or filters appropriate for the identified or expected contaminants;
- hard hat;
- splash shield, as needed; and,
- ankles/wrists taped with duct tape.

The site safety officer will verify if Level C is appropriate by checking organic vapor concentrations using compound and/or class-specific detector tubes.

- chemical resistant coveralls;
- steel-toe and steel-shank workboots;
- chemical resistant overboots or disposable boot covers;
- disposable inner gloves;
- disposable outer gloves;
- hard hat; and,
- ankles/wrists taped.

The exact PPE ensemble is decided on a site-by-site basis by the Site Safety Officer with the intent to provide the most protective and efficient worker PPE.

4.3 Activity-Specific Levels of Personal Protection

The required level of PPE is activity-specific and is based on air monitoring results (Section 4.0) and properties of identified or expected contaminants. **It is expected that site work will be performed in Level D.** If air monitoring results indicate the necessity to upgrade the level of protection engineering controls (i.e. Facing equipment away from the wind and placing site personnel upwind of drilling locations, active venting, etc.) will be implemented before requiring the use of respiratory protection.

5.0 AIR MONITORING AND ACTION LEVELS

29 CFR 1910.120(h) specifies that monitoring shall be performed where there may be a question of employee exposure to hazardous concentrations of hazardous substances in order to assure proper selection of engineering controls, work practices and personal protective equipment so that employees are not exposed to levels which exceed permissible exposure limits, or published exposure levels if there are no permissible exposure limits, for hazardous substances.

5.1 Air Monitoring Requirements

If excavation work is performed, air will be monitored for VOCs with a portable ION Science 3000EX photoionization detector, or the equivalent. If necessary, Lower Explosive Limit (LEL) and oxygen will be monitored with a Combustible Gas Indicator (CGI). If appropriate, fugitive dust will be monitored using a MiniRam Model PDM-3 aerosol monitor. Air will be monitored when any of the following conditions apply:

- initial site entry;
- during any work where a potential IDLH condition or flammable atmosphere could develop;
- excavation work begins on another portion of the site;
- contaminants, other than those previously identified, have been discovered;
- each time a different task or activity is initiated;
- during trenching and/or excavation work.

The designated site safety officer will record air monitoring data and ensure that air monitoring instruments are calibrated and maintained in accordance with manufacturer's specifications. Instruments will be zeroed daily and checked for accuracy. Monitoring results will be recorded in a field notebook and will be transferred to instrument reading logs.

5.2 Work Stoppage Responses

The following responses will be initiated whenever one or more of the action levels necessitating a work stoppage are exceeded:

- 1 The SSO will be consulted immediately
- 2 All personnel (except as necessary for continued monitoring and contaminant migration, if applicable) will be cleared from the work area (eg from the exclusion zone).
- 3 Monitoring will be continued until intrusive work resumes.

5.3 Action Levels During Excavation Activities

Instrument readings will be taken in the breathing zone above the excavation pit unless otherwise noted. Each action level is independent of all other action levels in determining responses.

| Organic Vapors (PID) | LEL % | Responses |
|--------------------------|-------|---|
| 0-1 ppm above background | 0% | <ul style="list-style-type: none">• Continue excavating• Level D protection• Continue monitoring every 10 minutes |

| | | |
|--|--------|---|
| 1-5 ppm Above Background, Sustained Reading | 1-10% | <ul style="list-style-type: none"> • Continue excavating • Go to Level C protection or employ engineering controls • Continue monitoring every 10 minutes |
| 5-25 ppm Above Background, Sustained Reading | 10-20% | <ul style="list-style-type: none"> • Discontinue excavating, unless PID is only action level exceeded. • Level C protection or employ engineering controls • Continue monitoring for organic vapors 200 ft downwind • Continuous monitoring for LEL at excavation pit |
| >25 ppm Above Background, Sustained Reading | >20% | <ul style="list-style-type: none"> • Discontinue excavating • Withdraw from area, shut off all engine ignition sources. • Allow pit to vent • Continuous monitoring for organic vapors 200 ft downwind. |

Notes: Air monitoring will occur in the breathing zone 30 inches above the excavation pit. Readings may also be taken in the excavation pit but will not be used for action levels.

If action levels for any one of the monitoring parameters are exceeded, the appropriate responses listed in the right hand column should be taken. If instrument readings do not return to acceptable levels after the excavation pit has been vented for a period of greater than one-half hour, a decision will then be made whether or not to seal the pit with suppressant foam.

If, during excavation activities, downwind monitoring PID readings are greater than 5 ppm above background for more than one-half hour, excavation will stop until sustained levels are less than 5 ppm (see Community Air Monitoring Plan).

6.0 SITE CONTROL

6.1 Work Zones

The primary purpose of site controls is to establish the perimeter of a hazardous area, to reduce the migration of contaminants into clean areas, and to prevent access or exposure to hazardous materials by unauthorized persons. When operations are to take place involving hazardous materials, the site safety officer will establish an exclusion zone, a decontamination zone, and a support zone. These zones "float" (move around the site) depending on the tasks being performed on any given day. The site safety officer will outline these locations before work begins and when zones change. The site safety officer records this information in the site log book. **It is expected that the entire fenced in area of the Site will be the exclusion zone, with the decontamination zone the Site entrance.** The support zone will be the office trailer.

Tasks requiring OSHA 40-hour Hazardous Waste Operations and Emergency Response Operations training are carried out in the exclusion zone. The exclusion zone is defined by the site safety officer but will typically be a 50-foot area around work activities. Gross decontamination (as determined by the site Health and Safety Officer) is conducted in the exclusion zone; all other decontamination is performed in the decontamination zone or trailer.

Protective equipment is removed in the decontamination zone. Disposable protective equipment is stored in receptacles staged in the decontamination zone, and non-disposable equipment is decontaminated. All personnel and equipment exit the exclusion zone through the decontamination zone. If a decontamination trailer is provided the first aid equipment, an eye wash unit, and drinking water are kept in the decontamination trailer.

The support zone is used for vehicle parking, daily safety meetings, and supply storage. Eating, drinking, and smoking are permitted only in the support zone. When a decontamination trailer is not provided, the eye wash unit, first aid equipment, and drinking water are kept at a central location designated by the site safety officer.

6.2 General Site Work

An excavation contractor with appropriate experience, personnel and training (40 hr OSHA Hazardous Waste Operations and Emergency Response Operations - HAZWOPER) is required to perform the removal / cleaning of the USTs soil. After this material is removed the contractor will remove historic fill and uncontaminated soil. The excavation contractor's on-site personnel engaged in historic fill and native soil removal will have a minimum of 24 hour HAZWOPER training.

7.0 CONTINGENCY PLAN/EMERGENCY RESPONSE PLAN

Site personnel must be prepared in the event of an emergency. Emergencies can take many forms: illnesses, injuries, chemical exposure, fires, explosions, spills, leaks, releases of harmful contaminants, or sudden changes in the weather.

Emergency telephone numbers and a map to the hospital will be posted in the command post. Site personnel should be familiar with the emergency procedures, and the locations of site safety, first aid, and communication equipment.

7.1 Emergency Equipment On-site

| | |
|---------------------|-------------------------------------|
| Private telephones: | Site personnel. |
| Two-way radios: | Site personnel where necessary. |
| Emergency Alarms: | On-site vehicle horns*. |
| First aid kits: | On-site, in vehicles or office. |
| Fire extinguisher: | On-site, in office or on equipment. |

* Horns: Air horns will be supplied to personnel at the discretion of the project superintendent or site safety officer.

7.2 Emergency Telephone Numbers

| | |
|-------------------------------------|----------------|
| General Emergencies | 911 |
| New York City Police | 911 |
| Lincoln Medical Center | 1-718-579-5016 |
| NYSDEC Spills Division | 1-800-457-7362 |
| NYSDEC Division of Env. Remediation | 1-718-482-4900 |
| NYCDEP | 1-718-699-9811 |
| NYC Department of Health | 1-212-788-4711 |
| NYC Fire Department | 911 |
| National Response Center | 1-800-424-8802 |
| Poison Control | 1-212-340-4494 |
| Site Safety Officer | 1-631-504-6000 |
| Alternate Site Safety Officer | 1-631-504-6000 |

7.3 Personnel Responsibilities During an Emergency

The project manager is primarily responsible for responding to and correcting any emergency situations. However, in the absence of the project manager, the site safety officer shall act as the project manager's on-site designee and perform the following tasks:

- Take appropriate measures to protect personnel including: withdrawal from the exclusion zone, evacuate and secure the site, or upgrade/downgrade the level of protective clothing and respiratory protection;

- Ensure that appropriate federal, state, and local agencies are informed and emergency response plans are coordinated. In the event of fire or explosion, the local fire department should be summoned immediately. If toxic materials are released to the air, the local authorities should be informed in order to assess the need for evacuation;
- Ensure appropriate decontamination, treatment, or testing for exposed or injured personnel;
- Determine the cause of incidents and make recommendations to prevent recurrence; and,
- Ensure that all required reports have been prepared.

The following key personnel are planned for this project:

- Project Manager Mr. Robert Bennett (631) 504-6000
- Construction Superintendent to be named
- Site Safety Officer Mr. Kevin Waters (631) 504-6000

7.4 Medical Emergencies

A person who becomes ill or injured in the exclusion zone will be decontaminated to the maximum extent possible. If the injury or illness is minor, full decontamination will be completed and first aid administered prior to transport. First aid will be administered while waiting for an ambulance or paramedics. A Field Accident Report (**Appendix D**) must be filled out for any injury.

A person transporting an injured/exposed person to a clinic or hospital for treatment will take the directions to the hospital (**Appendix D**) and information on the chemical(s) to which they may have been exposed (**Appendix C**).

7.5 Fire or Explosion

In the event of a fire or explosion, the local fire department will be summoned immediately. The site safety officer or his designated alternate will advise the fire commander of the location, nature and identification of the hazardous materials on-site. If it is safe to do so, site personnel may:

- use fire fighting equipment available on site; or,
- remove or isolate flammable or other hazardous materials that may contribute to the fire.

7.6 Evacuation Routes

Evacuation routes established by work area locations for each site will be reviewed prior to commencing site operations. As the work areas change, the evacuation routes will be altered accordingly, and the new route will be reviewed.

Under extreme emergency conditions, evacuation is to be immediate without regard for equipment. The evacuation signal will be a continuous blast of a vehicle horn, if possible, and/or by verbal/radio communication. When evacuating the site, personnel will follow these instructions:

- Keep upwind of smoke, vapors, or spill location.
- Exit through the decontamination corridor if possible.
- If evacuation through the decontamination corridor is not possible, personnel should remove contaminated clothing once they are in a safe location and leave it near the exclusion zone or in a safe place.
- The site safety officer will conduct a head count to ensure that all personnel have been evacuated safely. The head count will be correlated to the site and/or exclusion zone entry/exit log.
- If emergency site evacuation is necessary, all personnel are to escape the emergency situation and decontaminate to the maximum extent practical.

7.7 Spill Control Procedures

Spills associated with site activities may be attributed to project equipment and include gasoline, diesel and hydraulic oil. In the event of a leak or a release, site personnel will inform their supervisor immediately, locate the source of spillage and stop the flow if it can be done safely. A spill containment kit including absorbent pads, booms and/or granulated speedy dry absorbent material will be available to site personnel to facilitate the immediate recovery of the spilled material. Daily inspections of site equipment components including hydraulic lines, fuel tanks, etc. will be performed by their respective operators as a preventative measure for equipment leaks and to ensure equipment soundness. In the event of a spill, site personnel will immediately notify the NYSDEC (1-800-457-7362), and a spill number will be generated.

7.8 Vapor Release Plan

If work zone organic vapor (excluding methane) exceeds 5 ppm, then a downwind reading will be made either 200 feet from the work zone or at the property line, whichever is closer. If readings at this location exceed 5 ppm over background, the work will be stopped.

If 5 ppm of VOCs are recorded over background on a PID at the property line, then an off-site reading will be taken within 20 feet of the nearest residential or commercial property, whichever is closer. If efforts to mitigate the emission source are unsuccessful for 30 minutes, then the designated site safety officer will:

- contact the local police;
- continue to monitor air every 30 minutes, 20 feet from the closest off-site property. If two successive readings are below 5 ppm (non-methane), off-site air monitoring will be halted.
- All property line and off site air monitoring locations and results associated with vapor releases will be recorded in the site safety log book.

APPENDIX A
SITE SAFETY ACKNOWLEDGEMENT FORM

DAILY BRIEFING SIGN-IN SHEET

Date: _____ Person Conducting Briefing: _____

Project Name and Location: _____

1. AWARENESS (topics discussed, special safety concerns, recent incidents, etc...):

2. OTHER ISSUES (HASP changes, attendee comments, etc...):

3. ATTENDEES (Print Name):

| | |
|-----|-----|
| 1. | 11. |
| 2. | 12. |
| 3. | 13. |
| 4. | 14. |
| 5. | 15. |
| 6. | 16. |
| 7. | 17. |
| 8. | 18. |
| 9. | 19. |
| 10. | 20. |

APPENDIX B
SITE SAFETY PLAN AMENDMENTS

SITE SAFETY PLAN AMENDMENT FORM

Site Safety Plan Amendment #: _____

Site Name: _____

Reason for Amendment: _____

Alternative Procedures: _____

Required Changes in PPE: _____

Project Superintendent (signature)

Date

Health and Safety Consultant (signature)

Date

Site Safety Officer (signature)

Date

APPENDIX C

CHEMICAL HAZARDS

CHEMICAL HAZARDS

The attached International Chemical Safety Cards are provided for contaminants of concern that have been identified in soils and/or groundwater at the site.

International Chemical Safety Cards

BENZENE

ICSC: 0015



Cyclohexatriene
Benzol
C₆H₆
Molecular mass: 78.1

ICSC # 0015
CAS # 71-43-2
RTECS # [CY1400000](#)
UN # 1114
EC # 601-020-00-8
May 06, 2003 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|---|---|---|
| FIRE | Highly flammable. | NO open flames, NO sparks, and NO smoking. | Powder, AFFF, foam, carbon dioxide. |
| EXPLOSION | Vapour/air mixtures are explosive. Risk of fire and explosion: see Chemical Dangers. | Closed system, ventilation, explosion-proof electrical equipment and lighting. Do NOT use compressed air for filling, discharging, or handling. Use non-sparking handtools. Prevent build-up of electrostatic charges (e.g., by grounding). | In case of fire: keep drums, etc., cool by spraying with water. |
| EXPOSURE | | AVOID ALL CONTACT! | |
| •INHALATION | Dizziness. Drowsiness. Headache. Nausea. Shortness of breath. Convulsions. Unconsciousness. | Ventilation, local exhaust, or breathing protection. | Fresh air, rest. Refer for medical attention. |
| •SKIN | MAY BE ABSORBED! Dry skin. Redness. Pain. (Further see Inhalation). | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse skin with plenty of water or shower. Refer for medical attention. |
| •EYES | Redness. Pain. | Face shield, or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | Abdominal pain. Sore throat. Vomiting. (Further see Inhalation). | Do not eat, drink, or smoke during work. | Rinse mouth. Do NOT induce vomiting. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|---|
| Remove all ignition sources. Collect leaking and spilled liquid in sealable containers as far as possible. Absorb remaining liquid in sand or inert absorbent and remove to safe place. Do NOT wash away into sewer. Do NOT let this chemical enter the environment. Personal protection: complete protective clothing including self-contained breathing apparatus. | Fireproof. Separated from food and feedstuffs oxidants halogens | Do not transport with food and feedstuffs. Note: E F symbol T symbol R: 45-46-11-36/38-48/23/24/25-65 S: 53-45 UN Hazard Class: 3 UN Packing Group: II |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0015

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

BENZENE

ICSC: 0015

| | | |
|---|--|--|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: COLOURLESS LIQUID , WITH CHARACTERISTIC ODOUR.</p> <p>PHYSICAL DANGERS: The vapour is heavier than air and may travel along the ground; distant ignition possible. As a result of flow, agitation, etc., electrostatic charges can be generated.</p> <p>CHEMICAL DANGERS: Reacts violently with oxidants, nitric acid, sulfuric acid and halogens causing fire and explosion hazard. Attacks plastic and rubber.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 0.5 ppm as TWA 2.5 ppm as STEL (skin) A1 BEI (ACGIH 2004). MAK: H Carcinogen category: 1 Germ cell mutagen group: 3A (DFG 2004). OSHA PEL: 1910.1028 TWA 1 ppm ST 5 ppm See Appendix F NIOSH REL: Ca TWA 0.1 ppm ST 1 ppm See Appendix A NIOSH IDLH: Ca 500 ppm See: 71432</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation through the skin and by ingestion</p> <p>INHALATION RISK: A harmful contamination of the air can be reached very quickly on evaporation of this substance at 20°C.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: The substance is irritating to the eyes the skin and the respiratory tract Swallowing the liquid may cause aspiration into the lungs with the risk of chemical pneumonitis. The substance may cause effects on the central nervous system , resulting in lowering of consciousness Exposure far above the occupational exposure limit value may result in unconsciousness death</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: The liquid defats the skin. The substance may have effects on the bone marrow immune system , resulting in a decrease of blood cells. This substance is carcinogenic to humans.</p> |
|---|--|--|

| | | |
|-----------------------------------|--|---|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 80°C Melting point: 6°C Relative density (water = 1): 0.88 Solubility in water, g/100 ml at 25°C: 0.18 Vapour pressure, kPa at 20°C: 10 Relative vapour density (air = 1): 2.7</p> | <p>Relative density of the vapour/air-mixture at 20°C (air = 1): 1.2 Flash point: -11°C c.c. Auto-ignition temperature: 498°C Explosive limits, vol% in air: 1.2-8.0 Octanol/water partition coefficient as log Pow: 2.13</p> |
|-----------------------------------|--|---|

| | | |
|----------------------------------|--|---|
| <p>ENVIRONMENTAL DATA</p> | <p>The substance is very toxic to aquatic organisms.</p> |  |
|----------------------------------|--|---|

NOTES

Use of alcoholic beverages enhances the harmful effect. Depending on the degree of exposure, periodic medical examination is indicated. The odour warning when the exposure limit value is exceeded is insufficient.

Transport Emergency Card: TEC (R)-30S1114 / 30GF1-II
NFPA Code: H2; F3; R0

ADDITIONAL INFORMATION

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

ICSC: 0015 **BENZENE**

(C) IPCS, CEC, 1994

| | |
|---------------------------------------|--|
| <p>IMPORTANT LEGAL NOTICE:</p> | <p>Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.</p> |
|---------------------------------------|--|

International Chemical Safety Cards

BENZ(a)ANTHRACENE

ICSC: 0385



1,2-Benzoanthracene
Benzo(a)anthracene
2,3-Benzphenanthrene
Naphthanthracene
 $C_{18}H_{12}$
Molecular mass: 228.3

ICSC # 0385
CAS # 56-55-3
RTECS # [CV9275000](#)
EC # 601-033-00-9
October 23, 1995 Validated



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible. | | Water spray, powder. In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | AVOID ALL CONTACT! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety goggles face shield or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. Wash hands before eating. | Rinse mouth. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|--------------|---|
| Sweep spilled substance into sealable containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Personal protection: complete protective clothing including self-contained breathing apparatus. | Well closed. | T symbol N symbol R: 45-50/53 S: 53-45-60-61 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0385

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

ICSC: 0385

BENZ(a)ANTHRACENE

| | | |
|---|---|---|
| I M P O R T A N T D A T A | PHYSICAL STATE; APPEARANCE: COLOURLESS TO YELLOW BROWN FLUORESCENT FLAKES OR POWDER. | ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation, through the skin and by ingestion. |
| | PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air. | INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly. |
| | CHEMICAL DANGERS: | EFFECTS OF SHORT-TERM EXPOSURE: |
| | OCCUPATIONAL EXPOSURE LIMITS: TLV: A2 (suspected human carcinogen); (ACGIH 2004). MAK: Carcinogen category: 2 (as pyrolysis product of organic materials) (DFG 2005). | EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: This substance is probably carcinogenic to humans. |

| | | |
|----------------------------|---|--|
| PHYSICAL PROPERTIES | Sublimation point: 435°C Melting point: 162°C Relative density (water = 1): 1.274 Solubility in water: none | Vapour pressure, Pa at 20°C: 292 Octanol/water partition coefficient as log Pow: 5.61 |
|----------------------------|---|--|

| | | |
|---------------------------|--|---|
| ENVIRONMENTAL DATA | Bioaccumulation of this chemical may occur in seafood. |  |
|---------------------------|--|---|

NOTES

This substance is one of many polycyclic aromatic hydrocarbons - standards are usually established for them as mixtures, e.g., coal tar pitch volatiles. However, it may be encountered as a laboratory chemical in its pure form. Insufficient data are available on the effect of this substance on human health, therefore utmost care must be taken. Do NOT take working clothes home. Tetraphene is a common name. Card has been partly updated in October 2005 and August 2006: see sections Occupational Exposure Limits, EU classification.

ADDITIONAL INFORMATION

ICSC: 0385

BENZ(a)ANTHRACENE

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

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International Chemical Safety Cards

BENZO(a)PYRENE

ICSC: 0104



Benz(a)pyrene
3,4-Benzopyrene
Benzo(d,e,f)chrysene
 $C_{20}H_{12}$
Molecular mass: 252.3

ICSC # 0104
CAS # 50-32-8
RTECS # [DJ3675000](#)
EC # 601-032-00-3
October 17, 2005 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible. | NO open flames. | Water spray, foam, powder, carbon dioxide. |
| EXPLOSION | | | |
| EXPOSURE | See EFFECTS OF LONG-TERM OR REPEATED EXPOSURE. | AVOID ALL CONTACT! AVOID EXPOSURE OF (PREGNANT) WOMEN! | |
| •INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| •SKIN | MAY BE ABSORBED! | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| •EYES | | Safety goggles or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | | Do not eat, drink, or smoke during work. | Induce vomiting (ONLY IN CONSCIOUS PERSONS!). Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|---------------------------------|---|
| Evacuate danger area! Personal protection: complete protective clothing including self-contained breathing apparatus. Do NOT let this chemical enter the environment. Sweep spilled substance into sealable containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. | Separated from strong oxidants. | T symbol N symbol R: 45-46-60-61-43-50/53 S: 53-45-60-61 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0104

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

BENZO(a)PYRENE

ICSC: 0104

| | | |
|--|---|---|
| <p>I M P O R T A N T A D V I S I O N</p> | <p>PHYSICAL STATE; APPEARANCE: PALE-YELLOW CRYSTALS</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS: Reacts with strong oxidants causing fire and explosion hazard.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: Exposure by all routes should be carefully controlled to levels as low as possible A2 (suspected human carcinogen); (ACGIH 2005). MAK: Carcinogen category: 2; Germ cell mutagen group: 2; (DFG 2005).</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol, through the skin and by ingestion.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE:</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: This substance is carcinogenic to humans. May cause heritable genetic damage to human germ cells. Animal tests show that this substance possibly causes toxicity to human reproduction or development.</p> |
|--|---|---|

| | | |
|-----------------------------------|--|---|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 496°C Melting point: 178.1°C Density: 1.4 g/cm³</p> | <p>Solubility in water: none (<0.1 g/100 ml) Vapour pressure : negligible Octanol/water partition coefficient as log Pow: 6.04</p> |
|-----------------------------------|--|---|

| | | |
|----------------------------------|---|---|
| <p>ENVIRONMENTAL DATA</p> | <p>The substance is very toxic to aquatic organisms. Bioaccumulation of this chemical may occur in fish, in plants and in molluscs. The substance may cause long-term effects in the aquatic environment.</p> |  |
|----------------------------------|---|---|

NOTES

Do NOT take working clothes home. Benzo(a)pyrene is present as a component of polycyclic aromatic hydrocarbons (PAHs) in the environment, usually resulting from the incomplete combustion or pyrolysis of organic matters, especially fossil fuels and tobacco.

ADDITIONAL INFORMATION

| | |
|--|--|
| | |
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| | | |
|-------------------|---------------------|-----------------------|
| ICSC: 0104 | (C) IPCS, CEC, 1994 | BENZO(a)PYRENE |
|-------------------|---------------------|-----------------------|

| | |
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| <p>IMPORTANT LEGAL NOTICE:</p> | <p>Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.</p> |
|---------------------------------------|--|

International Chemical Safety Cards

BENZO(b)FLUORANTHENE

ICSC: 0720



Benz(e)acephenanthrylene
 2,3-Benzofluoranthene
 Benzo(e)fluoranthene
 3,4-Benzofluoranthene
 $C_{20}H_{12}$
 Molecular mass: 252.3

ICSC # 0720
 CAS # 205-99-2
 RTECS # [CU1400000](#)
 EC # 601-034-00-4
 March 25, 1999 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|-------------------------|---|---|
| FIRE | | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | | | |
| EXPOSURE | | AVOID ALL CONTACT! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety spectacles or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|---|
| Sweep spilled substance into covered containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Do NOT let this chemical enter the environment. | Provision to contain effluent from fire extinguishing. Well closed. | T symbol N symbol R: 45-50/53 S: 53-45-60-61 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0720

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

BENZO(b)FLUORANTHENE

ICSC: 0720

| | | |
|----------|---|---|
| I | PHYSICAL STATE; APPEARANCE: COLOURLESS CRYSTALS | ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation |
|----------|---|---|

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PHYSICAL DANGERS:

CHEMICAL DANGERS:

Upon heating, toxic fumes are formed.

OCCUPATIONAL EXPOSURE LIMITS:

TLV: A2 (suspected human carcinogen); (ACGIH 2004).

MAK:

Carcinogen category: 2;
(DFG 2004).

of its aerosol and through the skin.

INHALATION RISK:

Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly.

EFFECTS OF SHORT-TERM EXPOSURE:

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

This substance is possibly carcinogenic to humans. May cause genetic damage in humans.

PHYSICAL PROPERTIES

Boiling point: 481°C
Melting point: 168°C
Solubility in water:
none

Octanol/water partition coefficient as log Pow: 6.12

ENVIRONMENTAL DATA

This substance may be hazardous to the environment; special attention should be given to air quality and water quality.



NOTES

Benzo(b)fluoranthene is present as a component of polycyclic aromatic hydrocarbons (PAH) content in the environment usually resulting from the incomplete combustion or pyrolysis of organic matters, especially fossil fuels and tobacco. ACGIH recommends environment containing benzo(b)fluoranthene should be evaluated in terms of the TLV-TWA for coal tar pitch volatile, as benzene soluble 0.2 mg/m³. Insufficient data are available on the effect of this substance on human health, therefore utmost care must be taken.

ADDITIONAL INFORMATION

ICSC: 0720

BENZO(b)FLUORANTHENE

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

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International Chemical Safety Cards

BENZO(k)FLUORANTHENE

ICSC: 0721



Dibenzo(b,jk)fluorene
8,9-Benzofluoranthene
11,12-Benzofluoranthene
 $C_{20}H_{12}$
Molecular mass: 252.3

ICSC # 0721
CAS # 207-08-9
RTECS # [DF6350000](#)
EC # 601-036-00-5
March 25, 1999 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|-------------------------|---|---|
| FIRE | | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | | | |
| EXPOSURE | | AVOID ALL CONTACT! | |
| •INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| •SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| •EYES | | Safety spectacles or eye protection in combination with breathing protection if powder. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|---|
| Sweep spilled substance into covered containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Do NOT let this chemical enter the environment. | Provision to contain effluent from fire extinguishing. Well closed. | T symbol N symbol R: 45-50/53 S: 53-45-60-61 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0721

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

BENZO(k)FLUORANTHENE

ICSC: 0721

| | | |
|---|---|--|
| I | PHYSICAL STATE; APPEARANCE: YELLOW CRYSTALS | ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol and through the skin. |
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PHYSICAL DANGERS:

CHEMICAL DANGERS:

Upon heating, toxic fumes are formed.

OCCUPATIONAL EXPOSURE LIMITS:

TLV not established.

MAK:

Carcinogen category: 2;
(DFG 2004).

INHALATION RISK:

Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly.

EFFECTS OF SHORT-TERM EXPOSURE:

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

This substance is possibly carcinogenic to humans.

PHYSICAL PROPERTIES

Boiling point: 480°C
Melting point: 217°C
Solubility in water:
none

Octanol/water partition coefficient as log Pow: 6.84

ENVIRONMENTAL DATA

This substance may be hazardous to the environment; special attention should be given to air quality and water quality. Bioaccumulation of this chemical may occur in crustacea and in fish.



NOTES

Benzo(k)fluoranthene is present as a component of polycyclic aromatic hydrocarbons (PAH) content in the environment usually resulting from the incomplete combustion or pyrolysis of organic matters, especially fossil fuels and tobacco. ACGIH recommends environment containing benzo(k)fluoranthene should be evaluated in terms of the TLV-TWA for coal tar pitch volatile, as benzene soluble 0.2 mg/m³. Insufficient data are available on the effect of this substance on human health, therefore utmost care must be taken.

ADDITIONAL INFORMATION

ICSC: 0721

BENZO(k)FLUORANTHENE

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

CHRYSENE

ICSC: 1672



Benzoaphenanthrene
1,2-Benzophenanthrene
1,2,5,6-Dibenzonaphthalene
 $C_{18}H_{12}$
Molecular mass: 228.3

ICSC # 1672
CAS # 218-01-9
RTECS # [GC0700000](#)
UN # 3077
EC # 601-048-00-0
October 12, 2006 Validated



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible. | NO open flames. | Water spray. Dry powder. Foam. Carbon dioxide. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | See EFFECTS OF LONG-TERM OR REPEATED EXPOSURE. | AVOID ALL CONTACT! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety goggles | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|--|--|
| Personal protection: P3 filter respirator for toxic particles. Do NOT let this chemical enter the environment. Sweep spilled substance into sealable containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. | Separated from strong oxidants, Provision to contain effluent from fire extinguishing. Store in an area without drain or sewer access. | T symbol N symbol R: 45-68-50/53 S: 53-45-60-61 UN Hazard Class: 9 UN Packing Group: III Signal: Warning Aqua-Cancer Suspected of causing cancer Very toxic to aquatic life with long lasting effects Very toxic to aquatic life |

SEE IMPORTANT INFORMATION ON BACK


International Chemical Safety Cards

CHRYSENE

ICSC: 1672

| | | |
|--|--|---|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: COLOURLESS TO BEIGE CRYSTALS OR POWDER</p> <p>PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air.</p> <p>CHEMICAL DANGERS: The substance decomposes on burning producing toxic fumes Reacts violently with strong oxidants</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: A3 (confirmed animal carcinogen with unknown relevance to humans); (ACGIH 2006). MAK not established.</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol, through the skin and by ingestion.</p> <p>INHALATION RISK: A harmful concentration of airborne particles can be reached quickly when dispersed</p> <p>EFFECTS OF SHORT-TERM EXPOSURE:</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: This substance is possibly carcinogenic to humans.</p> |
|--|--|---|

| | | |
|-----------------------------------|--|---|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 448°C Melting point: 254 - 256°C Density: 1.3 g/cm³</p> | <p>Solubility in water: very poor Octanol/water partition coefficient as log Pow: 5.9</p> |
|-----------------------------------|--|---|

| | | |
|----------------------------------|--|---|
| <p>ENVIRONMENTAL DATA</p> | <p>The substance is very toxic to aquatic organisms. Bioaccumulation of this chemical may occur in seafood. It is strongly advised that this substance does not enter the environment.</p> |  |
|----------------------------------|--|---|

NOTES

Depending on the degree of exposure, periodic medical examination is suggested. Do NOT take working clothes home. This substance does not usually occur as a pure substance but as a component of polyaromatic hydrocarbon (PAH) mixtures. Human population studies have associated PAH's exposure with cancer and cardiovascular diseases.

Transport Emergency Card: TEC (R)-90GM7-III

ADDITIONAL INFORMATION

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

ICSC: 1672

CHRYSENE

(C) IPCS, CEC, 1994

| | |
|---------------------------------------|--|
| <p>IMPORTANT LEGAL NOTICE:</p> | <p>Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.</p> |
|---------------------------------------|--|

International Chemical Safety Cards

INDENO(1,2,3-cd)PYRENE

ICSC: 0730



o-Phenylenepyrene
2,3-Phenylenepyrene
 $C_{22}H_{12}$
Molecular mass: 276.3

ICSC # 0730
CAS # 193-39-5
RTECS # [NK9300000](#)
March 25, 1999 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|-------------------------|---|---|
| FIRE | | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | | | |
| EXPOSURE | | AVOID ALL CONTACT! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety spectacles or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|-----------------------|
| Sweep spilled substance into covered containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Do NOT let this chemical enter the environment. | Provision to contain effluent from fire extinguishing. Well closed. | R: S: |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0730

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

INDENO(1,2,3-cd)PYRENE

ICSC: 0730

| | | |
|----------|---|--|
| I | PHYSICAL STATE; APPEARANCE: YELLOW CRYSTALS | ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol and through the skin. |
| M | PHYSICAL DANGERS: | INHALATION RISK: |
| P | | |

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CHEMICAL DANGERS:
Upon heating, toxic fumes are formed.

Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly.

OCCUPATIONAL EXPOSURE LIMITS:
TLV not established.
MAK:
Carcinogen category: 2;
(DFG 2004).

EFFECTS OF SHORT-TERM EXPOSURE:

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

This substance is possibly carcinogenic to humans.

PHYSICAL PROPERTIES

Boiling point: 536°C
Melting point: 164°C
Solubility in water:
none

Octanol/water partition coefficient as log Pow: 6.58

ENVIRONMENTAL DATA

This substance may be hazardous to the environment; special attention should be given to air quality and water quality. Bioaccumulation of this chemical may occur in fish.



NOTES

Indeno(1,2,3-cd)pyrene is present as a component of polycyclic aromatic hydrocarbons (PAH) content in the environment usually resulting from the incomplete combustion or pyrolysis of organic matters, especially fossil fuels and tobacco. ACGIH recommends environment containing Indeno(1,2,3-c,d)pyrene should be evaluated in terms of the TLV-TWA for coal tar pitch volatile, as benzene soluble 0.2 mg/m³. Insufficient data are available on the effect of this substance on human health, therefore utmost care must be taken.

ADDITIONAL INFORMATION

ICSC: 0730

INDENO(1,2,3-cd)PYRENE

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

1. PRODUCT AND COMPANY IDENTIFICATION

Product name : 4,4'-DDD PESTANAL,250 MG (2,2-BIS(4-CHL&

Product Number : 35486
Brand : Fluka

Company : Sigma-Aldrich
3050 Spruce Street
SAINT LOUIS MO 63103
USA

Telephone : +1 800-325-5832
Fax : +1 800-325-5052
Emergency Phone # : (314) 776-6555

2. HAZARDS IDENTIFICATION

Emergency Overview

OSHA Hazards

Toxic by ingestion, Harmful by skin absorption., Possible carcinogen.

GHS Label elements, including precautionary statements

Pictogram



Signal word Danger

Hazard statement(s)

H301 Toxic if swallowed.
H312 Harmful in contact with skin.
H351 Suspected of causing cancer.
H400 Very toxic to aquatic life.
H413 May cause long lasting harmful effects to aquatic life.

Precautionary statement(s)

P273 Avoid release to the environment.
P280 Wear protective gloves/protective clothing.
P301 + P310 IF SWALLOWED: Immediately call a POISON CENTER or doctor/physician.

HMIS Classification

Health hazard: 2
Chronic Health Hazard: *
Flammability: 0
Physical hazards: 0

NFPA Rating

Health hazard: 2
Fire: 0
Reactivity Hazard: 0

Potential Health Effects

Inhalation May be harmful if inhaled. May cause respiratory tract irritation.
Skin Harmful if absorbed through skin. May cause skin irritation.
Eyes May cause eye irritation.
Ingestion Toxic if swallowed.

3. COMPOSITION/INFORMATION ON INGREDIENTS

Synonyms : 1,1-Dichloro-2,2-bis(4-chlorophenyl)ethane
4,4'-DDD
TDE

Formula : C₁₄H₁₀Cl₄
Molecular Weight : 320.04 g/mol

| CAS-No. | EC-No. | Index-No. | Concentration |
|--|-----------|-----------|---------------|
| 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | | | |
| 72-54-8 | 200-783-0 | - | - |

4. FIRST AID MEASURES

General advice

Consult a physician. Show this safety data sheet to the doctor in attendance. Move out of dangerous area.

If inhaled

If breathed in, move person into fresh air. If not breathing give artificial respiration. Consult a physician.

In case of skin contact

Wash off with soap and plenty of water. Consult a physician.

In case of eye contact

Rinse thoroughly with plenty of water for at least 15 minutes and consult a physician.

If swallowed

Never give anything by mouth to an unconscious person. Rinse mouth with water. Consult a physician.

5. FIRE-FIGHTING MEASURES

Suitable extinguishing media

Use water spray, alcohol-resistant foam, dry chemical or carbon dioxide.

Special protective equipment for fire-fighters

Wear self contained breathing apparatus for fire fighting if necessary.

6. ACCIDENTAL RELEASE MEASURES

Personal precautions

Use personal protective equipment. Avoid dust formation. Avoid breathing dust. Ensure adequate ventilation. Evacuate personnel to safe areas.

Environmental precautions

Prevent further leakage or spillage if safe to do so. Do not let product enter drains. Discharge into the environment must be avoided.

Methods and materials for containment and cleaning up

Pick up and arrange disposal without creating dust. Keep in suitable, closed containers for disposal.

7. HANDLING AND STORAGE

Precautions for safe handling

Avoid contact with skin and eyes. Avoid formation of dust and aerosols.

Provide appropriate exhaust ventilation at places where dust is formed. Normal measures for preventive fire protection.

Conditions for safe storage

Keep container tightly closed in a dry and well-ventilated place.

8. EXPOSURE CONTROLS/PERSONAL PROTECTION

Contains no substances with occupational exposure limit values.

Personal protective equipment

Respiratory protection

Where risk assessment shows air-purifying respirators are appropriate use a full-face particle respirator type N100 (US) or type P3 (EN 143) respirator cartridges as a backup to engineering controls. If the respirator is the sole means of protection, use a full-face supplied air respirator. Use respirators and components tested and approved under appropriate government standards such as NIOSH (US) or CEN (EU).

Hand protection

Handle with gloves.

Eye protection

Face shield and safety glasses

Skin and body protection

Choose body protection according to the amount and concentration of the dangerous substance at the work place.

Hygiene measures

Avoid contact with skin, eyes and clothing. Wash hands before breaks and immediately after handling the product.

9. PHYSICAL AND CHEMICAL PROPERTIES

Appearance

Form solid

Safety data

| | |
|---|---|
| pH | no data available |
| Melting point | 94.0 - 96.0 °C (201.2 - 204.8 °F) |
| Boiling point | 193.0 °C (379.4 °F) at 1.3 hPa (1.0 mmHg) |
| Flash point | no data available |
| Ignition temperature | no data available |
| Lower explosion limit | no data available |
| Upper explosion limit | no data available |
| Vapour pressure | < 0.00001 hPa (< 0.00001 mmHg) at 25.0 °C (77.0 °F) |
| Density | 1.38 g/cm ³ |
| Water solubility | no data available |
| Partition coefficient: n-octanol/water | log Pow: 6.02 |

10. STABILITY AND REACTIVITY

Chemical stability

Stable under recommended storage conditions.

Conditions to avoid

no data available

Materials to avoid

Strong oxidizing agents

Hazardous decomposition products

Hazardous decomposition products formed under fire conditions. - Carbon oxides, Hydrogen chloride gas

Hazardous decomposition products formed under fire conditions. - Nature of decomposition products not known.

11. TOXICOLOGICAL INFORMATION

Acute toxicity

LD50 Oral - Hamster - > 5,000 mg/kg

TDL0 Oral - Human - 428.5 mg/kg

Remarks: Endocrine:Adrenal cortex hypoplasia.

TDL0 Oral - rat - 6,000 mg/kg

Remarks: Cardiac:Other changes. Gastrointestinal:Other changes. Kidney, Ureter, Bladder:Changes in both tubules and glomeruli.

TDL0 Oral - rat - 14 mg/kg

Remarks: Liver:Changes in liver weight. Endocrine:Estrogenic. Musculoskeletal:Other changes.

TDL0 Oral - rat - 2,100 mg/kg

Remarks: Behavioral:Altered sleep time (including change in righting reflex).

LD50 Dermal - rabbit - 1,200 mg/kg

Remarks: Behavioral:Excitement. Behavioral:Convulsions or effect on seizure threshold. Skin irritation

Skin corrosion/irritation

no data available

Serious eye damage/eye irritation

no data available

Respiratory or skin sensitization

no data available

Germ cell mutagenicity

no data available

Carcinogenicity

This product is or contains a component that has been reported to be possibly carcinogenic based on its IARC, ACGIH, NTP, or EPA classification.

Limited evidence of carcinogenicity in animal studies

IARC: No component of this product present at levels greater than or equal to 0.1% is identified as probable, possible or confirmed human carcinogen by IARC.

ACGIH: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by ACGIH.

NTP: No component of this product present at levels greater than or equal to 0.1% is identified as a known or anticipated carcinogen by NTP.

OSHA: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by OSHA.

Reproductive toxicity

no data available

Specific target organ toxicity - single exposure (GHS)

no data available

Specific target organ toxicity - repeated exposure (GHS)

no data available

Aspiration hazard

no data available

Potential health effects**Inhalation**

May be harmful if inhaled. May cause respiratory tract irritation.

Ingestion

Toxic if swallowed.

Skin

Harmful if absorbed through skin. May cause skin irritation.

Eyes May cause eye irritation.

Signs and Symptoms of Exposure

To the best of our knowledge, the chemical, physical, and toxicological properties have not been thoroughly investigated.

Additional Information

RTECS: KI0700000

12. ECOLOGICAL INFORMATION

Toxicity

Toxicity to fish LC50 - other fish - 1.18 - 9 mg/l - 96.0 h
LC50 - Lepomis macrochirus (Bluegill) - 0.04 - 0.05 mg/l - 96.0 h
LC50 - Oncorhynchus mykiss (rainbow trout) - 0.06 - 0.09 mg/l - 96.0 h
LC50 - Pimephales promelas (fathead minnow) - 3.47 - 5.58 mg/l - 96.0 h

Toxicity to daphnia and other aquatic invertebrates. EC50 - Daphnia pulex (Water flea) - 0.01 mg/l - 48 h

Persistence and degradability

no data available

Bioaccumulative potential

Indication of bioaccumulation.

Mobility in soil

no data available

PBT and vPvB assessment

no data available

Other adverse effects

An environmental hazard cannot be excluded in the event of unprofessional handling or disposal.

Very toxic to aquatic organisms, may cause long-term adverse effects in the aquatic environment.

13. DISPOSAL CONSIDERATIONS

Product

Observe all federal, state, and local environmental regulations. Contact a licensed professional waste disposal service to dispose of this material. Dissolve or mix the material with a combustible solvent and burn in a chemical incinerator equipped with an afterburner and scrubber.

Contaminated packaging

Dispose of as unused product.

14. TRANSPORT INFORMATION

DOT (US)

UN-Number: 2811 Class: 6.1 Packing group: III
Proper shipping name: Toxic solids, organic, n.o.s. (2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane)
Reportable Quantity (RQ): 1 lbs
Marine pollutant: No
Poison Inhalation Hazard: No

IMDG

UN-Number: 2811 Class: 6.1 Packing group: III EMS-No: F-A, S-A
Proper shipping name: TOXIC SOLID, ORGANIC, N.O.S. (2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane)
Marine pollutant: No

IATA

UN-Number: 2811 Class: 6.1 Packing group: III
Proper shipping name: Toxic solid, organic, n.o.s. (2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane)

15. REGULATORY INFORMATION

OSHA Hazards

Toxic by ingestion, Harmful by skin absorption., Possible carcinogen.

DSL Status

This product contains the following components that are not on the Canadian DSL nor NDSL lists.

| | |
|---|--------------------|
| 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | CAS-No. 72-54-8 |
|---|--------------------|

SARA 302 Components

SARA 302: No chemicals in this material are subject to the reporting requirements of SARA Title III, Section 302.

SARA 313 Components

SARA 313: This material does not contain any chemical components with known CAS numbers that exceed the threshold (De Minimis) reporting levels established by SARA Title III, Section 313.

SARA 311/312 Hazards

Acute Health Hazard

Massachusetts Right To Know Components

| | | |
|---|--------------------|---------------|
| 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | CAS-No. 72-54-8 | Revision Date |
|---|--------------------|---------------|

Pennsylvania Right To Know Components

| | | |
|---|--------------------|---------------|
| 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | CAS-No. 72-54-8 | Revision Date |
|---|--------------------|---------------|

New Jersey Right To Know Components

| | | |
|---|--------------------|---------------|
| 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | CAS-No. 72-54-8 | Revision Date |
|---|--------------------|---------------|

California Prop. 65 Components

| | | |
|--|--------------------|---------------|
| WARNING! This product contains a chemical known to the State of California to cause cancer. 2,2-bis(4-Chlorophenyl)-1,1-dichloro-ethane | CAS-No. 72-54-8 | Revision Date |
|--|--------------------|---------------|

16. OTHER INFORMATION

Further information

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The above information is believed to be correct but does not purport to be all inclusive and shall be used only as a guide. The information in this document is based on the present state of our knowledge and is applicable to the product with regard to appropriate safety precautions. It does not represent any guarantee of the properties of the product. Sigma-Aldrich Co., shall not be held liable for any damage resulting from handling or from contact with the above product. See reverse side of invoice or packing slip for additional terms and conditions of sale.

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Search

72-55-9 msds



MSDS 250,000+

MSDS : 2,2-Bis-(4-chlorophenyl)-1,1-dichloroethylene, 99%
 CAS : 72-55-9
 SYNONYMS : p,p'-DDE ; ethylene,1,1-dichloro-2,2-bis-(p-chlorophenyl)- ; DDT dehydrochloride ; DDE; 1-1'-(Dichloroethenylidene)bis(4-chlorobenzene)

[MSDS Safety Sheet](#)

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AdChoices ▶

Catalog of Chemical Suppliers, Buyers, Custom Synthesis Companies And Equipment Manufacturers
 [2,2-Bis-(4-chlorophenyl)-1,1-dichloroethylene, 99% 72-55-9]

Suppliers:

Not Available

Buyers:

Not Available

[Sprayon® LU711 Lubricant](#) Because your environment demands a TRUE Industrial Lubricant Sprayon.com

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AdChoices ▶

**** SECTION 2 - COMPOSITION, INFORMATION ON INGREDIENTS ****

| CAS# | Chemical Name | % | EINECS# |
|---------|---|----|-----------|
| 72-55-9 | 2,2-Bis-(4-chlorophenyl)-1,1-dichloroethy | 99 | 200-784-6 |
| | ethylene | | |

Hazard Symbols: XN

Risk Phrases: 22 33

**** SECTION 3 - HAZARDS IDENTIFICATION ****

EMERGENCY OVERVIEW

Harmful if swallowed. Danger of cumulative effects.Cancer suspect agent.Possible risks of irreversible effects.

Potential Health Effects

Eye:

May cause eye irritation.

Skin:

May cause skin irritation.

Ingestion:

May cause irritation of the digestive tract. May be harmful if swallowed. Ingestion of large amounts may cause liver and/or kidney damage.

Inhalation:

May cause respiratory tract irritation.

Chronic:

May cause cancer according to animal studies. Adverse reproductive effects have been reported in animals. Laboratory experiments have resulted in mutagenic effects.

**** SECTION 4 - FIRST AID MEASURES ****

Eyes:

Flush eyes with plenty of water for at least 15 minutes, occasionally lifting the upper and lower eyelids. Get medical aid.

Skin:

Get medical aid. Flush skin with plenty of water for at least 15 minutes while removing contaminated clothing and shoes. Wash clothing before reuse.

Ingestion:

If victim is conscious and alert, give 2-4 cupfuls of milk or water.

Never give anything by mouth to an unconscious person. Get medical aid immediately.

Inhalation:

Remove from exposure and move to fresh air immediately. If not breathing, give artificial respiration. If breathing is difficult, give oxygen. Get medical aid.

Notes to Physician:

Treat symptomatically and supportively.

**** SECTION 5 - FIRE FIGHTING MEASURES ****

General Information:

As in any fire, wear a self-contained breathing apparatus in pressure-demand, MSHA/NIOSH (approved or equivalent), and full protective gear. Water runoff can cause environmental damage. Dike and collect water used to fight fire. During a fire, irritating and highly toxic gases may be generated by thermal decomposition or combustion. Will burn if involved in a fire.

Extinguishing Media:

For large fires, use water spray, fog or regular foam. For small fires, use dry chemical, carbon dioxide, water spray or regular foam. Cool containers with flooding quantities of water until well after fire is out.

**** SECTION 6 - ACCIDENTAL RELEASE MEASURES ****

General Information: Use proper personal protective equipment as indicated in Section 8.

Spills/Leaks:

Avoid runoff into storm sewers and ditches which lead to waterways. Clean up spills immediately, observing precautions in the Protective Equipment section. Sweep up, then place into a suitable container for disposal. Avoid generating dusty conditions. Provide ventilation.

**** SECTION 7 - HANDLING and STORAGE ****

Handling:

Wash thoroughly after handling. Remove contaminated clothing and wash before reuse. Minimize dust generation and accumulation. Avoid contact with eyes, skin, and clothing. Do not ingest or inhale. Use with adequate ventilation.

Storage:

Keep container closed when not in use. Store in a tightly closed container. Store in a cool, dry, well-ventilated area away from incompatible substances.

**** SECTION 8 - EXPOSURE CONTROLS, PERSONAL PROTECTION ****

Engineering Controls:

Facilities storing or utilizing this material should be equipped with an eyewash facility and a safety shower. Use adequate ventilation to keep airborne concentrations low.

Exposure Limits

CAS# 72-55-9:

Personal Protective Equipment

Eyes:

Wear appropriate protective eyeglasses or chemical safety goggles as described by OSHA's eye and face protection regulations in 29 CFR 1910.133 or European Standard EN166.

Skin:

Wear appropriate protective gloves to prevent skin exposure.

Clothing:

Wear appropriate protective clothing to prevent skin exposure.

Respirators:

A respiratory protection program that meets OSHA's 29 CFR 1910.134 and ANSI Z88.2 requirements or European Standard EN 149 must be followed whenever workplace conditions warrant respirator use.

**** SECTION 9 - PHYSICAL AND CHEMICAL PROPERTIES ****

Physical State: Crystals

Color: white

Odor: None reported.

pH: Not available.

Vapor Pressure: 6.5106 mm Hg @ 20 C

Viscosity: Not available.

Boiling Point: 336 deg C

Freezing/Melting Point: 88.00 - 90.00 deg C

Autoignition Temperature: Not available.

Flash Point: Not available.

Explosion Limits, lower: Not available.

Explosion Limits, upper: Not available.

Decomposition Temperature:

Solubility in water: 0.010 ppm

Specific Gravity/Density:

Molecular Formula: C14H8Cl4

Molecular Weight: 318.02

**** SECTION 10 - STABILITY AND REACTIVITY ****

Chemical Stability:

Stable under normal temperatures and pressures.

Conditions to Avoid:

Incompatible materials, dust generation, strong oxidants.

Incompatibilities with Other Materials:

Strong oxidizing agents - strong bases.

Hazardous Decomposition Products:

Hydrogen chloride, carbon monoxide, carbon dioxide.

Hazardous Polymerization: Has not been reported.

**** SECTION 11 - TOXICOLOGICAL INFORMATION ****

RTECS#:

CAS# 72-55-9: KV9450000

LD50/LC50:

CAS# 72-55-9: Oral, mouse: LD50 = 700 mg/kg; Oral, rat: LD50 = 880 mg/kg.

Carcinogenicity:

2,2-Bis-(4-chlorophenyl)-1,1-dichloroethylene -

California: carcinogen, initial date 1/1/89

Other:

See actual entry in RTECS for complete information.

**** SECTION 12 - ECOLOGICAL INFORMATION ****

Ecotoxicity:

Estimated BCF value = 8,300 based on water solubility. Estimated Koc value = 8,300. There was no movement of DDE reported in soil column mobility experiments.

**** SECTION 13 - DISPOSAL CONSIDERATIONS ****

Dispose of in a manner consistent with federal, state, and local regulations.

**** SECTION 14 - TRANSPORT INFORMATION ****

IATA

Not regulated as a hazardous material.

IMO

Not regulated as a hazardous material.

RID/ADR

Not regulated as a hazardous material.

USA RQ: CAS# 72-55-9: 1 lb final RQ; 0.454 kg final RQ

**** SECTION 15 - REGULATORY INFORMATION ****

European/International Regulations

European Labeling in Accordance with EC Directives

Hazard Symbols: XN

Risk Phrases:

R 22 Harmful if swallowed.

R 33 Danger of cumulative effects.

Safety Phrases:

S 24/25 Avoid contact with skin and eyes.

WGK (Water Danger/Protection)

CAS# 72-55-9: 3

Canada

None of the chemicals in this product are listed on the DSL/NDSL list.

CAS# 72-55-9 is listed on Canada's Ingredient Disclosure List.

US FEDERAL

TSCA

CAS# 72-55-9 is not listed on the TSCA inventory.

It is for research and development use only.

**** SECTION 16 - ADDITIONAL INFORMATION ****

MSDS Creation Date: 9/28/1998 Revision #3 Date: 3/18/2003

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| NAME | CAS |
|---|------------|
| M-Benzoyloxybenzyl Alcohol, 97% | 1700-30-7 |
| Octaphenylcyclotetrasiloxane, 98% | 546-56-5 |
| Cetylpyridinium chloride | 123-03-5 |
| 3,4-Difluorophenol, 99% | 2713-33-9 |
| 1-Benzyl-4-Hydroxypiperidine, 97% | 4727-72-4 |
| 4-tert-Butylbenzoyl chloride | 1710-98-1 |
| Borane-morpholine complex, 97% | 4856-95-5 |
| Benzyl Ether, 99% | 103-50-4 |
| 5-Amino-1-Naphthol (Pract) | 83-55-6 |
| Pyridinium-P-Toluenesulfonate 98% | 24057-28-1 |
| Pyrogallol Red, 98% (Titr.) | 32638-88-3 |
| Amberlite ira 416 | 9002-26-0 |
| 3-Methoxybenzotrile, 98% | 1527-89-5 |
| 1-Adamantanemethanol, 99% | 770-71-8 |
| Inosine, 99% | 58-63-9 |
| Pentafluoropropionic Acid | 422-64-0 |
| Pyruvic Acid | 127-17-3 |
| Potassium hydrogen fluoride, 99+% | 7789-29-9 |
| Aluminum Nitride, 98% Particle Size <10 Micron | 24304-00-5 |
| Nickel(II) hydroxide, c.p., 60-61% Ni | 12054-48-7 |
| 1-Adamantanamine sulfate, 99% | 31377-23-8 |
| S-(Thiobenzoyl)-Thioglycolic Acid, 97% | 942-91-6 |
| N,N-Dimethyl-P-Nitroaniline | 100-23-2 |
| Benzofuroxan | 480-96-6 |
| cis-2-Aminomethyl-1-cyclohexanol hydrochloride, 99% | 24947-68-0 |
| Silver Phosphate, 98% (Titr.) | 7784-09-0 |

| | |
|---|------------|
| 4-Cyano-4-Phenylpiperidine Hydrochloride, 99% (TLC) | 51304-58-6 |
| Methanesulfonamide | 3144-09-0 |
| gamma-Octanoic lactone, 98% | 104-50-7 |
| Cis,cis,cis-1,2,3,4-cyclopentane- tetracarboxylic dianhydride, | 4802-47-5 |
| Tetrachloroethylene Carbonate, 98+% | 22432-68-4 |
| Oxamic Acid, 98% | 471-47-6 |
| 1O,11-Dihydro-5H-Dibenzo(A,D)-Cycloheptene, 98% | 833-48-7 |
| Thallium (I) Sulfate, 99.9+% | 7446-18-6 |
| N-(2,6-Dimethylphenylcarbonyl-Methyl)-Iminodiacetic Acid, 99% | 59160-29-1 |
| P-(Dimethylamino)cinnamic Acid, 99% | 1552-96-1 |
| Biebrich Scarlet, 99% (UV-VIS) | 4196-99-0 |
| 4-Chlorobenzenediazonium hexafluoro- phosphate | 1582-27-0 |
| Ammonium hexachloroiridate(IV), 99.99% | 16940-92-4 |
| Methylamine-d2 deuteriochloride, 98+ atom % D | 593-51-1 |
| 2,2-Bis-(4-chlorophenyl)-1,1-dichloroethylene, 99% | 72-55-9 |
| Nitro red | 56431-61-9 |
| Methyl 2,3-dichlorobenzoate, 98+% | 2905-54-6 |
| Isopropyl Bromoacetate, 98% (GC) | 29921-57-1 |
| 1-Iodo-4-Nitrobenzene, 99% | 636-98-6 |
| 4-Ethylcyclohexanol, 99% cis/trans mixture | 4534-74-1 |
| Fluorescamine | 38183-12-9 |
| Tris(2,2,6,6-Tetramethyl-3,5-Heptanedionato)Dysprosium(III), 99+% | 15522-69-7 |
| 3-Amino-2,2,5,5-Tetramethyl-1-Pyrrolidinyloxy, 99% (Titr.) | 34272-83-8 |
| 3,4-Dihydroxyphenylacetic Acid,98% | 102-32-9 |

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International Chemical Safety Cards

DDT

ICSC: 0034



Dichlorodiphenyltrichloroethane
 1,1,1-Trichloro-2,2-bis(p-chlorophenyl)ethane
 2,2-bis(p-Chlorophenyl)-1,1,1-trichloroethane
 1,1'-(2,2,2-Trichloroethylidene)bis(4-chlorobenzene)
 p,p'-DDT
 $C_{14}H_9Cl_5$
 Molecular mass: 354.5



ICSC # 0034
 CAS # 50-29-3
 RTECS # [KJ3325000](#)
 UN # 2761
 EC # 602-045-00-7
 April 20, 2004 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|---|---|
| FIRE | Combustible. Liquid formulations containing organic solvents may be flammable. Gives off irritating or toxic fumes (or gases) in a fire. | NO open flames. | Powder, water spray, foam, carbon dioxide. |
| EXPLOSION | | | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! STRICT HYGIENE! AVOID EXPOSURE OF (PREGNANT) WOMEN! | |
| •INHALATION | Cough. | Local exhaust or breathing protection. | Fresh air, rest. |
| •SKIN | | Protective gloves. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| •EYES | Redness. | Safety goggles, or eye protection in combination with breathing protection if powder. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | Tremors. Diarrhoea. Dizziness. Headache. Vomiting. Numbness. Paresthesias. Hyperexcitability. Convulsions. | Do not eat, drink, or smoke during work. Wash hands before eating. | Rinse mouth. Give a slurry of activated charcoal in water to drink. Rest. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|---|--|
| Do NOT let this chemical enter the environment. Sweep spilled substance into sealable non-metallic containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Personal protection: P3 filter respirator for toxic particles. | Provision to contain effluent from fire extinguishing. Separated from iron, aluminum and its salts, food and feedstuffs See Chemical Dangers. | Do not transport with food and feedstuffs. Severe marine pollutant. T symbol N symbol R: 25-40-48/25-50/53 S: 1/2-22-36/37-45-60-61 UN Hazard Class: 6.1 UN Packing Group: III |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0034

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

ICSC: 0034

DDT

| | | |
|---|---|---|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: COLOURLESS CRYSTALS WHITE POWDER. TECHNICAL PRODUCT IS WAXY SOLID.</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS: On combustion, forms toxic and corrosive fumes including hydrogen chloride. Reacts with aluminium and iron.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 1 mg/m³ as TWA A3 (ACGIH 2004). MAK: 1 mg/m³ H Peak limitation category: II(8) (DFG 2003). OSHA PEL: TWA 1 mg/m³ skin NIOSH REL: Ca TWA 0.5 mg/m³ See Appendix A NIOSH IDLH: Ca 500 mg/m³ See: 50293</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by ingestion.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly especially if powdered.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: May cause mechanical irritation. The substance may cause effects on the central nervous system, resulting in convulsions and respiratory depression. Exposure at high levels may result in death. Medical observation is indicated.</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: The substance may have effects on the central nervous system and liver. This substance is possibly carcinogenic to humans. Animal tests show that this substance possibly causes toxicity to human reproduction or development.</p> |
|---|---|---|

| | | |
|-----------------------------------|--|---|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 260°C Melting point: 109°C Density: 1.6 g/cm³</p> | <p>Solubility in water: poor Octanol/water partition coefficient as log Pow: 6.36</p> |
|-----------------------------------|--|---|

| | | |
|----------------------------------|---|---|
| <p>ENVIRONMENTAL DATA</p> | <p>The substance is very toxic to aquatic organisms. This substance may be hazardous to the environment; special attention should be given to birds. Bioaccumulation of this chemical may occur along the food chain, for example in milk and aquatic organisms. This substance does enter the environment under normal use. Great care, however, should be given to avoid any additional release, e.g. through inappropriate disposal.</p> |  |
|----------------------------------|---|---|

NOTES

Depending on the degree of exposure, periodic medical examination is indicated. Carrier solvents used in commercial formulations may change physical and toxicological properties. Do NOT take working clothes home. Consult national legislation. Agritan, Azotox, Anofex, Ixodex, Gesapon, Gesarex, Gesarol, Guesapon, Clofenotane, Zeidane, Dicophane, Neocid are trade names.

Transport Emergency Card: TEC (R)-61GT7-III

ADDITIONAL INFORMATION

| | |
|----------------------------|-------------------|
| <p>ICSC: 0034</p> | <p>DDT</p> |
| <p>(C) IPCS, CEC, 1994</p> | |

| | |
|---------------------------------------|--|
| <p>IMPORTANT LEGAL NOTICE:</p> | <p>Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.</p> |
|---------------------------------------|--|

1. PRODUCT AND COMPANY IDENTIFICATION

| | | | |
|--|---|---|---|
| Product name | : | α-Chlordane | |
| Product Number | : | 442449 | |
| Brand | : | Supelco | |
| Product Use | : | For laboratory research purposes. | |
| Supplier | : | Sigma-Aldrich 3050 Spruce Street SAINT LOUIS MO 63103 USA | Manufacturer : Sigma-Aldrich Corporation 3050 Spruce St. St. Louis, Missouri 63103 USA |
| Telephone | : | +1 800-325-5832 | |
| Fax | : | +1 800-325-5052 | |
| Emergency Phone # (For both supplier and manufacturer) | : | (314) 776-6555 | |
| Preparation Information | : | Sigma-Aldrich Corporation Product Safety - Americas Region 1-800-521-8956 | |

2. HAZARDS IDENTIFICATION

Emergency Overview

OSHA Hazards

Toxic by inhalation., Toxic by ingestion, Toxic by skin absorption, Irritant

GHS Classification

Acute toxicity, Inhalation (Category 4)
Acute toxicity, Oral (Category 4)
Acute toxicity, Dermal (Category 3)
Skin irritation (Category 2)
Eye irritation (Category 2A)
Specific target organ toxicity - single exposure (Category 3)
Acute aquatic toxicity (Category 1)

GHS Label elements, including precautionary statements

Pictogram



Signal word

Danger

Hazard statement(s)

| | |
|-------------|-------------------------------------|
| H302 + H332 | Harmful if swallowed or if inhaled. |
| H311 | Toxic in contact with skin. |
| H315 | Causes skin irritation. |
| H319 | Causes serious eye irritation. |
| H335 | May cause respiratory irritation. |
| H400 | Very toxic to aquatic life. |

Precautionary statement(s)

| | |
|--------------------|--|
| P261 | Avoid breathing dust/ fume/ gas/ mist/ vapours/ spray. |
| P273 | Avoid release to the environment. |
| P280 | Wear protective gloves/ protective clothing. |
| P305 + P351 + P338 | IF IN EYES: Rinse cautiously with water for several minutes. Remove contact lenses, if present and easy to do. Continue rinsing. |

P312

Call a POISON CENTER or doctor/ physician if you feel unwell.

HMIS Classification

Health hazard: 2
Flammability: 0
Physical hazards: 0

NFPA Rating

Health hazard: 2
Fire: 0
Reactivity Hazard: 0

Potential Health Effects

Inhalation Toxic if inhaled. Causes respiratory tract irritation.
Skin Toxic if absorbed through skin. Causes skin irritation.
Eyes Causes eye irritation.
Ingestion Toxic if swallowed.

3. COMPOSITION/INFORMATION ON INGREDIENTS

Molecular Weight : 208.29 g/mol

| CAS-No. | EC-No. | Index-No. | Concentration |
|------------------|-----------|-----------|---------------|
| Chlordane | | | |
| 5103-71-9 | 225-825-5 | - | - |

4. FIRST AID MEASURES

General advice

Consult a physician. Show this safety data sheet to the doctor in attendance. Move out of dangerous area.

If inhaled

If breathed in, move person into fresh air. If not breathing, give artificial respiration. Consult a physician.

In case of skin contact

Wash off with soap and plenty of water. Take victim immediately to hospital. Consult a physician.

In case of eye contact

Rinse thoroughly with plenty of water for at least 15 minutes and consult a physician.

If swallowed

Never give anything by mouth to an unconscious person. Rinse mouth with water. Consult a physician.

5. FIRE-FIGHTING MEASURES

Suitable extinguishing media

Use water spray, alcohol-resistant foam, dry chemical or carbon dioxide.

Special protective equipment for fire-fighters

Wear self contained breathing apparatus for fire fighting if necessary.

Hazardous combustion products

Hazardous decomposition products formed under fire conditions. - Carbon oxides, Hydrogen chloride gas

6. ACCIDENTAL RELEASE MEASURES

Personal precautions

Wear respiratory protection. Avoid dust formation. Avoid breathing vapors, mist or gas. Ensure adequate ventilation. Evacuate personnel to safe areas. Avoid breathing dust.

Environmental precautions

Prevent further leakage or spillage if safe to do so. Do not let product enter drains. Discharge into the environment must be avoided.

Methods and materials for containment and cleaning up

Pick up and arrange disposal without creating dust. Sweep up and shovel. Keep in suitable, closed containers for disposal.

7. HANDLING AND STORAGE

Precautions for safe handling

Avoid contact with skin and eyes. Avoid formation of dust and aerosols.

Provide appropriate exhaust ventilation at places where dust is formed. Normal measures for preventive fire protection.

Conditions for safe storage

Keep container tightly closed in a dry and well-ventilated place.

8. EXPOSURE CONTROLS/PERSONAL PROTECTION

Contains no substances with occupational exposure limit values.

Personal protective equipment

Respiratory protection

Where risk assessment shows air-purifying respirators are appropriate use a full-face particle respirator type N99 (US) or type P2 (EN 143) respirator cartridges as a backup to engineering controls. If the respirator is the sole means of protection, use a full-face supplied air respirator. Use respirators and components tested and approved under appropriate government standards such as NIOSH (US) or CEN (EU).

Hand protection

Handle with gloves. Gloves must be inspected prior to use. Use proper glove removal technique (without touching glove's outer surface) to avoid skin contact with this product. Dispose of contaminated gloves after use in accordance with applicable laws and good laboratory practices. Wash and dry hands.

Eye protection

Face shield and safety glasses Use equipment for eye protection tested and approved under appropriate government standards such as NIOSH (US) or EN 166(EU).

Skin and body protection

Complete suit protecting against chemicals, The type of protective equipment must be selected according to the concentration and amount of the dangerous substance at the specific workplace.

Hygiene measures

Avoid contact with skin, eyes and clothing. Wash hands before breaks and immediately after handling the product.

9. PHYSICAL AND CHEMICAL PROPERTIES

Appearance

Form crystalline

Colour colourless

Safety data

pH no data available

Melting/freezing point 93.0 - 94.0 °C (199.4 - 201.2 °F)

Boiling point no data available

Flash point no data available

Ignition temperature no data available

Autoignition temperature no data available

Lower explosion limit no data available

Upper explosion limit no data available

Vapour pressure no data available

| | |
|---|-------------------|
| Density | no data available |
| Water solubility | no data available |
| Partition coefficient: n-octanol/water | no data available |
| Relative vapour density | no data available |
| Odour | no data available |
| Odour Threshold | no data available |
| Evaporation rate | no data available |

10. STABILITY AND REACTIVITY

Chemical stability

Stable under recommended storage conditions.

Possibility of hazardous reactions

no data available

Conditions to avoid

no data available

Materials to avoid

Strong oxidizing agents

Hazardous decomposition products

Hazardous decomposition products formed under fire conditions. - Carbon oxides, Hydrogen chloride gas
Other decomposition products - no data available

11. TOXICOLOGICAL INFORMATION

Acute toxicity

Oral LD50

LD50 Oral - rat - 500.0 mg/kg

Inhalation LC50

Dermal LD50

Other information on acute toxicity

no data available

Skin corrosion/irritation

no data available

Serious eye damage/eye irritation

no data available

Respiratory or skin sensitization

no data available

Germ cell mutagenicity

no data available

Carcinogenicity

IARC: No component of this product present at levels greater than or equal to 0.1% is identified as probable, possible or confirmed human carcinogen by IARC.

ACGIH: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by ACGIH.

NTP: No component of this product present at levels greater than or equal to 0.1% is identified as a known or anticipated carcinogen by NTP.

OSHA: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by OSHA.

Reproductive toxicity

no data available

Teratogenicity

no data available

Specific target organ toxicity - single exposure (Globally Harmonized System)

no data available

Specific target organ toxicity - repeated exposure (Globally Harmonized System)

no data available

Aspiration hazard

no data available

Potential health effects

| | |
|-------------------|---|
| Inhalation | Toxic if inhaled. Causes respiratory tract irritation. |
| Ingestion | Toxic if swallowed. |
| Skin | Toxic if absorbed through skin. Causes skin irritation. |
| Eyes | Causes eye irritation. |

Synergistic effects

no data available

Additional Information

RTECS: Not available

12. ECOLOGICAL INFORMATION

Toxicity

Toxicity to fish LC50 - Lepomis macrochirus (Bluegill) - 0.0074 mg/l - 96 h

Persistence and degradability

no data available

Bioaccumulative potential

Bioaccumulation Lepomis macrochirus (Bluegill) - 24 h
Bioconcentration factor (BCF): 322

Mobility in soil

no data available

PBT and vPvB assessment

no data available

Other adverse effects

An environmental hazard cannot be excluded in the event of unprofessional handling or disposal.

Very toxic to aquatic life.

no data available

An environmental hazard cannot be excluded in the event of unprofessional handling or disposal.

Very toxic to aquatic life.

13. DISPOSAL CONSIDERATIONS

Product

Offer surplus and non-recyclable solutions to a licensed disposal company. Contact a licensed professional waste disposal service to dispose of this material. Dissolve or mix the material with a combustible solvent and burn in a chemical incinerator equipped with an afterburner and scrubber.

Contaminated packaging

Dispose of as unused product.

14. TRANSPORT INFORMATION

DOT (US)

UN-Number: 3077 Class: 9 Packing group: III
Proper shipping name: Environmentally hazardous substances, solid, n.o.s. (Chlordane)
Marine pollutant:
Poison Inhalation Hazard: No

IMDG

UN-Number: 3077 Class: 9 Packing group: III EMS-No: F-A, S-F
Proper shipping name: ENVIRONMENTALLY HAZARDOUS SUBSTANCE, SOLID, N.O.S. (Chlordane)
Marine pollutant: Marine pollutant

IATA

UN-Number: 3077 Class: 9 Packing group: III
Proper shipping name: Environmentally hazardous substance, solid, n.o.s. (Chlordane)

Further information

EHS-Mark required (ADR 2.2.9.1.10, IMDG code 2.10.3) for single packagings and combination packagings containing inner packagings with Dangerous Goods > 5L for liquids or > 5kg for solids.

15. REGULATORY INFORMATION

OSHA Hazards

Toxic by inhalation., Toxic by ingestion, Toxic by skin absorption, Irritant

DSL Status

This product contains the following components that are not on the Canadian DSL nor NDSL lists.

| | |
|-----------|----------------------|
| Chlordane | CAS-No. 5103-71-9 |
|-----------|----------------------|

SARA 302 Components

SARA 302: No chemicals in this material are subject to the reporting requirements of SARA Title III, Section 302.

SARA 313 Components

SARA 313: This material does not contain any chemical components with known CAS numbers that exceed the threshold (De Minimis) reporting levels established by SARA Title III, Section 313.

SARA 311/312 Hazards

Acute Health Hazard

Massachusetts Right To Know Components

No components are subject to the Massachusetts Right to Know Act.

Pennsylvania Right To Know Components

| | | |
|-----------|----------------------|---------------|
| Chlordane | CAS-No. 5103-71-9 | Revision Date |
|-----------|----------------------|---------------|

New Jersey Right To Know Components

| | | |
|-----------|----------------------|---------------|
| Chlordane | CAS-No. 5103-71-9 | Revision Date |
|-----------|----------------------|---------------|

California Prop. 65 Components

This product does not contain any chemicals known to State of California to cause cancer, birth defects, or any other reproductive harm.

16. OTHER INFORMATION

Further information

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1. PRODUCT AND COMPANY IDENTIFICATION

Product name : Aroclor 1262

Product Number : 442463
Brand : Supelco

Supplier : Sigma-Aldrich
3050 Spruce Street
SAINT LOUIS MO 63103
USA

Telephone : +1 800-325-5832
Fax : +1 800-325-5052
Emergency Phone # (For : (314) 776-6555
both supplier and
manufacturer)

Preparation Information : Sigma-Aldrich Corporation
Product Safety - Americas Region
1-800-521-8956

2. HAZARDS IDENTIFICATION**Emergency Overview****OSHA Hazards**

Carcinogen

GHS Classification

Carcinogenicity (Category 1B)

Specific target organ toxicity - repeated exposure (Category 2)

Acute aquatic toxicity (Category 3)

Chronic aquatic toxicity (Category 3)

GHS Label elements, including precautionary statements

Pictogram



Signal word

Danger

Hazard statement(s)

H350

May cause cancer.

H373

May cause damage to organs through prolonged or repeated exposure.

H412

Harmful to aquatic life with long lasting effects.

Precautionary statement(s)

P201

Obtain special instructions before use.

P273

Avoid release to the environment.

P308 + P313

IF exposed or concerned: Get medical advice/ attention.

HMIS Classification

Health hazard: 0

Chronic Health Hazard: *

Flammability: 0

Physical hazards: 0

NFPA Rating

Health hazard: 0

Fire: 0

Reactivity Hazard: 0

Potential Health Effects

Inhalation May be harmful if inhaled. May cause respiratory tract irritation.
Skin May be harmful if absorbed through skin. May cause skin irritation.
Eyes May cause eye irritation.
Ingestion May be harmful if swallowed.

3. COMPOSITION/INFORMATION ON INGREDIENTS

| CAS-No. | EC-No. | Index-No. | Concentration |
|---------------------------|--------|--------------|---------------|
| PCB - Aroclor 1262 | | | |
| 37324-23-5 | - | 602-039-00-4 | - |

4. FIRST AID MEASURES

General advice

Consult a physician. Show this safety data sheet to the doctor in attendance. Move out of dangerous area.

If inhaled

If breathed in, move person into fresh air. If not breathing, give artificial respiration. Consult a physician.

In case of skin contact

Wash off with soap and plenty of water. Consult a physician.

In case of eye contact

Flush eyes with water as a precaution.

If swallowed

Never give anything by mouth to an unconscious person. Rinse mouth with water. Consult a physician.

5. FIRE-FIGHTING MEASURES

Conditions of flammability

Not flammable or combustible.

Suitable extinguishing media

Use water spray, alcohol-resistant foam, dry chemical or carbon dioxide.

Special protective equipment for fire-fighters

Wear self contained breathing apparatus for fire fighting if necessary.

Hazardous combustion products

Hazardous decomposition products formed under fire conditions. - Nature of decomposition products not known.

6. ACCIDENTAL RELEASE MEASURES

Personal precautions

Use personal protective equipment. Avoid breathing vapors, mist or gas. Ensure adequate ventilation. Evacuate personnel to safe areas.

Environmental precautions

Prevent further leakage or spillage if safe to do so. Do not let product enter drains. Discharge into the environment must be avoided.

Methods and materials for containment and cleaning up

Soak up with inert absorbent material and dispose of as hazardous waste. Keep in suitable, closed containers for disposal.

7. HANDLING AND STORAGE

Precautions for safe handling

Avoid contact with skin and eyes. Avoid inhalation of vapour or mist.

Conditions for safe storage

Keep container tightly closed in a dry and well-ventilated place. Containers which are opened must be carefully resealed and kept upright to prevent leakage.

8. EXPOSURE CONTROLS/PERSONAL PROTECTION

Contains no substances with occupational exposure limit values.

Personal protective equipment

Respiratory protection

Where risk assessment shows air-purifying respirators are appropriate use a full-face respirator with multi-purpose combination (US) or type ABEK (EN 14387) respirator cartridges as a backup to engineering controls. If the respirator is the sole means of protection, use a full-face supplied air respirator. Use respirators and components tested and approved under appropriate government standards such as NIOSH (US) or CEN (EU).

Hand protection

Handle with gloves. Gloves must be inspected prior to use. Use proper glove removal technique (without touching glove's outer surface) to avoid skin contact with this product. Dispose of contaminated gloves after use in accordance with applicable laws and good laboratory practices. Wash and dry hands.

Eye protection

Face shield and safety glasses Use equipment for eye protection tested and approved under appropriate government standards such as NIOSH (US) or EN 166(EU).

Skin and body protection

Complete suit protecting against chemicals, The type of protective equipment must be selected according to the concentration and amount of the dangerous substance at the specific workplace.

Hygiene measures

Handle in accordance with good industrial hygiene and safety practice. Wash hands before breaks and at the end of workday.

9. PHYSICAL AND CHEMICAL PROPERTIES

Appearance

| | |
|--------|-------------------|
| Form | liquid |
| Colour | no data available |

Safety data

| | |
|--|-------------------|
| pH | no data available |
| Melting point/freezing point | no data available |
| Boiling point | no data available |
| Flash point | no data available |
| Ignition temperature | no data available |
| Autoignition temperature | no data available |
| Lower explosion limit | no data available |
| Upper explosion limit | no data available |
| Vapour pressure | no data available |
| Density | no data available |
| Water solubility | no data available |
| Partition coefficient: n-octanol/water | no data available |
| Relative vapour density | no data available |

| | |
|------------------|-------------------|
| Odour | no data available |
| Odour Threshold | no data available |
| Evaporation rate | no data available |

10. STABILITY AND REACTIVITY

Chemical stability

Stable under recommended storage conditions.

Possibility of hazardous reactions

no data available

Conditions to avoid

no data available

Materials to avoid

Strong oxidizing agents

Hazardous decomposition products

Hazardous decomposition products formed under fire conditions. - Nature of decomposition products not known.
Other decomposition products - no data available

11. TOXICOLOGICAL INFORMATION

Acute toxicity

Oral LD50

LD50 Oral - rat - 11,300 mg/kg

Inhalation LC50

no data available

Dermal LD50

Other information on acute toxicity

no data available

Skin corrosion/irritation

no data available

Serious eye damage/eye irritation

no data available

Respiratory or skin sensitization

no data available

Germ cell mutagenicity

no data available

Carcinogenicity

Carcinogen

Possible human carcinogen

IARC: No component of this product present at levels greater than or equal to 0.1% is identified as probable, possible or confirmed human carcinogen by IARC.

ACGIH: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by ACGIH.

NTP: No component of this product present at levels greater than or equal to 0.1% is identified as a known or anticipated carcinogen by NTP.

OSHA: No component of this product present at levels greater than or equal to 0.1% is identified as a carcinogen or potential carcinogen by OSHA.

Reproductive toxicity

no data available

Teratogenicity

no data available

Specific target organ toxicity - single exposure (Globally Harmonized System)

no data available

Specific target organ toxicity - repeated exposure (Globally Harmonized System)

May cause damage to organs through prolonged or repeated exposure.

no data available

Aspiration hazard

no data available

Potential health effects

| | |
|-------------------|---|
| Inhalation | May be harmful if inhaled. May cause respiratory tract irritation. |
| Ingestion | May be harmful if swallowed. |
| Skin | May be harmful if absorbed through skin. May cause skin irritation. |
| Eyes | May cause eye irritation. |

Signs and Symptoms of Exposure

To the best of our knowledge, the chemical, physical, and toxicological properties have not been thoroughly investigated.

Synergistic effects

no data available

Additional Information

RTECS: TQ1364000

12. ECOLOGICAL INFORMATION

Toxicity

Toxicity to fish LC50 - *Oncorhynchus clarki* - 50 mg/l - 96 h

Persistence and degradability

Biodegradability Result: - According to the results of tests of biodegradability this product is not readily biodegradable.
Remarks: no data available

Bioaccumulative potential

no data available

Mobility in soil

no data available

PBT and vPvB assessment

no data available

Other adverse effects

An environmental hazard cannot be excluded in the event of unprofessional handling or disposal.

Harmful to aquatic life with long lasting effects.

13. DISPOSAL CONSIDERATIONS

Product

Offer surplus and non-recyclable solutions to a licensed disposal company. Contact a licensed professional waste disposal service to dispose of this material. Dissolve or mix the material with a combustible solvent and burn in a chemical incinerator equipped with an afterburner and scrubber.

Contaminated packaging

Dispose of as unused product.

14. TRANSPORT INFORMATION

DOT (US)

UN number: 2315 Class: 9 Packing group: II
Proper shipping name: Polychlorinated biphenyls, liquid
Reportable Quantity (RQ):
Marine pollutant: No
Poison Inhalation Hazard: No

IMDG

UN number: 2315 Class: 9 Packing group: II EMS-No: F-A, S-A
Proper shipping name: POLYCHLORINATED BIPHENYLS, LIQUID
Marine pollutant: No

IATA

UN number: 2315 Class: 9 Packing group: II
Proper shipping name: Polychlorinated biphenyls, liquid

15. REGULATORY INFORMATION

OSHA Hazards

Carcinogen

SARA 302 Components

SARA 302: No chemicals in this material are subject to the reporting requirements of SARA Title III, Section 302.

SARA 313 Components

SARA 313: This material does not contain any chemical components with known CAS numbers that exceed the threshold (De Minimis) reporting levels established by SARA Title III, Section 313.

SARA 311/312 Hazards

Chronic Health Hazard

Massachusetts Right To Know Components

No components are subject to the Massachusetts Right to Know Act.

Pennsylvania Right To Know Components

| | CAS-No. | Revision Date |
|--------------------|------------|---------------|
| PCB - Aroclor 1262 | 37324-23-5 | 1989-08-11 |

New Jersey Right To Know Components

| | CAS-No. | Revision Date |
|--------------------|------------|---------------|
| PCB - Aroclor 1262 | 37324-23-5 | 1989-08-11 |

California Prop. 65 Components

| | CAS-No. | Revision Date |
|---|------------|---------------|
| WARNING! This product contains a chemical known to the State of California to cause cancer. PCB - Aroclor 1262 | 37324-23-5 | 2008-08-01 |

California Prop. 65 Components

| | CAS-No. | Revision Date |
|---|------------|---------------|
| WARNING! This product contains a chemical known to the State of California to cause birth defects or other reproductive harm. PCB - Aroclor 1262 | 37324-23-5 | 2008-08-01 |

16. OTHER INFORMATION

Further information

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The above information is believed to be correct but does not purport to be all inclusive and shall be used only as a guide. The information in this document is based on the present state of our knowledge and is applicable to the product with regard to appropriate safety precautions. It does not represent any guarantee of the properties of the product. Sigma-Aldrich Co., shall not be held liable for any damage resulting from handling or from contact with the above product. See reverse side of invoice or packing slip for additional terms and conditions of sale.

International Chemical Safety Cards

ARSENIC

ICSC: 0013



Grey arsenic
As
Atomic mass: 74.9

ICSC # 0013
CAS # 7440-38-2
RTECS # [CG0525000](#)
UN # 1558
EC # 033-001-00-X

October 18, 1999 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|---|--|---|
| FIRE | Combustible. Gives off irritating or toxic fumes (or gases) in a fire. | NO open flames. NO contact with strong oxidizers. NO contact with hot surfaces. | Powder, water spray, foam, carbon dioxide. |
| EXPLOSION | Risk of fire and explosion is slight when exposed to hot surfaces or flames in the form of fine powder or dust. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! AVOID ALL CONTACT! AVOID EXPOSURE OF (PREGNANT) WOMEN! | IN ALL CASES CONSULT A DOCTOR! |
| •INHALATION | Cough. Sore throat. Shortness of breath. Weakness. See Ingestion. | Closed system and ventilation. | Fresh air, rest. Artificial respiration may be needed. Refer for medical attention. |
| •SKIN | Redness. | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse skin with plenty of water or shower. |
| •EYES | Redness. | Face shield or eye protection in combination with breathing protection if powder. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | Abdominal pain. Diarrhoea. Nausea. Vomiting. Burning sensation in the throat and chest. Shock or collapse. Unconsciousness. | Do not eat, drink, or smoke during work. Wash hands before eating. | Rinse mouth. Induce vomiting (ONLY IN CONSCIOUS PERSONS!). Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|--|--|
| Evacuate danger area! Sweep spilled substance into sealable containers. Carefully collect remainder, then remove to safe place. Chemical protection suit including self-contained breathing apparatus. Do NOT let this chemical enter the environment. | Separated from strong oxidants, acids, halogens, food and feedstuffs. Well closed. | Do not transport with food and feedstuffs. Marine pollutant. T symbol N symbol R: 23/25-50/53 S: 1/2-20/21-28-45-60-61 UN Hazard Class: 6.1 UN Packing Group: II |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0013

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

ARSENIC

ICSC: 0013

| | | |
|---|--|--|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: ODOURLESS, BRITTLE, GREY, METALLIC-LOOKING CRYSTALS.</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS: Upon heating, toxic fumes are formed. Reacts violently with strong oxidants and halogens, causing fire and explosion hazard. Reacts with acids to produce</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 0.01 mg/m³ as TWA A1 (confirmed human carcinogen); BEI issued (ACGIH 2004). MAK: Carcinogen category: 1; Germ cell mutagen group: 3A; (DFG 2004). OSHA PEL: 1910.1018 TWA 0.010 mg/m³ NIOSH REL: Ca C 0.002 mg/m³ 15-minute See Appendix A NIOSH IDLH: Ca 5 mg/m³ (as As) See: 7440382</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol and by ingestion.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly, when dispersed.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: The substance is irritating to the eyes the skin and the respiratory tract. The substance may cause effects on the gastrointestinal tract cardiovascular system central nervous system kidneys , resulting in severe gastroenteritis, loss of fluid, and electrolytes, cardiac disorders shock convulsions and kidney impairment Exposure above the OEL may result in death. The effects may be delayed. Medical observation is indicated.</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: Repeated or prolonged contact with skin may cause dermatitis. The substance may have effects on the mucous membranes, skin, peripheral nervous system liver bone marrow , resulting in pigmentation disorders, hyperkeratosis, perforation of nasal septum, neuropathy, liver impairment anaemia This substance is carcinogenic to humans. Animal tests show that this substance possibly causes toxicity to human reproduction or development.</p> |
|---|--|--|

| | | |
|-----------------------------------|---|--------------------------------------|
| <p>PHYSICAL PROPERTIES</p> | <p>Sublimation point: 613°C Density: 5.7 g/cm³</p> | <p>Solubility in water: none</p> |
|-----------------------------------|---|--------------------------------------|

| | | |
|----------------------------------|--|---|
| <p>ENVIRONMENTAL DATA</p> | <p>The substance is toxic to aquatic organisms. It is strongly advised that this substance does not enter the environment.</p> |  |
|----------------------------------|--|---|

NOTES

The substance is combustible but no flash point is available in literature. Depending on the degree of exposure, periodic medical examination is suggested. Do NOT take working clothes home. Refer also to cards for specific arsenic compounds, e.g., Arsenic pentoxide (ICSC 0377), Arsenic trichloride (ICSC 0221), Arsenic trioxide (ICSC 0378), Arsine (ICSC 0222).

Transport Emergency Card: TEC (R)-61GT5-II

ADDITIONAL INFORMATION

ICSC: 0013 **ARSENIC**

(C) IPCS, CEC, 1994

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

International Chemical Safety Cards

BARIUM SULFATE

ICSC: 0827



Barium sulphate
Blanc fixe
Artificial barite
BaSO₄
Molecular mass: 233.43

ICSC # 0827

CAS # 7727-43-7

RTECS # [CR0600000](#)

October 20, 1999 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---|--|---|---|
| FIRE | Not combustible. Gives off irritating or toxic fumes (or gases) in a fire. | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | | | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. | Remove contaminated clothes. Rinse skin with plenty of water or shower. |
| • EYES | | Safety spectacles. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. |
| SPILLAGE DISPOSAL | | STORAGE | PACKAGING & LABELLING |
| Sweep spilled substance into containers; if appropriate, moisten first to prevent dusting. Personal protection: P1 filter respirator for inert particles. | | | R: S: |
| SEE IMPORTANT INFORMATION ON BACK | | | |
| ICSC: 0827 | | Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values. | |

International Chemical Safety Cards

BARIUM SULFATE

ICSC: 0827

| | | |
|---|--|--|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: ODOURLESS TASTELESS, WHITE OR YELLOWISH CRYSTALS OR POWDER.</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS: Reacts violently with aluminium powder.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 10 mg/m³ as TWA; (ACGIH 2004). MAK: (Inhalable fraction) 4 mg/m³; (Respirable fraction) 1.5 mg/m³; (DFG 2004). OSHA PEL[†]: TWA 15 mg/m³ (total) TWA 5 mg/m³ (resp) NIOSH REL: TWA 10 mg/m³ (total) TWA 5 mg/m³ (resp) NIOSH IDLH: N.D. See: IDLH INDEX</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a nuisance-causing concentration of airborne particles can, however, be reached quickly.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE:</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: Lungs may be affected by repeated or prolonged exposure to dust particles, resulting in baritosis (a form of benign pneumoconiosis).</p> |
|---|--|--|

| | | |
|-----------------------------------|---|----------------------------------|
| <p>PHYSICAL PROPERTIES</p> | <p>Melting point (decomposes): 1600°C Density: 4.5 g/cm³</p> | <p>Solubility in water: none</p> |
|-----------------------------------|---|----------------------------------|

| | |
|----------------------------------|--|
| <p>ENVIRONMENTAL DATA</p> | |
|----------------------------------|--|

NOTES

Occurs in nature as the mineral barite; also as barytes, heavy spar. Card has been partly updated in October 2005. See section Occupational Exposure Limits.

ADDITIONAL INFORMATION

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|--|--|

| | | |
|--------------------------|----------------------------|------------------------------|
| <p>ICSC: 0827</p> | <p>(C) IPCS, CEC, 1994</p> | <p>BARIUM SULFATE</p> |
|--------------------------|----------------------------|------------------------------|

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

International Chemical Safety Cards

CADMIUM

ICSC: 0020



Cd
Atomic mass: 112.4

ICSC # 0020
CAS # 7440-43-9
RTECS # [EU980000](#)
UN # 2570
EC # 048-002-00-0
April 22, 2005 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Flammable in powder form and spontaneously combustible in pyrophoric form. Gives off irritating or toxic fumes (or gases) in a fire. | NO open flames, NO sparks, and NO smoking. NO contact with heat or acid(s). | Dry sand. Special powder. NO other agents. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! AVOID ALL CONTACT! | IN ALL CASES CONSULT A DOCTOR! |
| • INHALATION | Cough. Sore throat. | Local exhaust or breathing protection. | Fresh air, rest. Refer for medical attention. |
| • SKIN | | Protective gloves. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | Redness. Pain. | Safety goggles or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | Abdominal pain. Diarrhoea. Headache. Nausea. Vomiting. | Do not eat, drink, or smoke during work. | Rest. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|--|--|
| Evacuate danger area! Personal protection: chemical protection suit including self-contained breathing apparatus. Remove all ignition sources. Sweep spilled substance into containers. Carefully collect remainder, then remove to safe place. | Fireproof. Dry. Keep under inert gas. Separated from ignition sources, oxidants acids, food and feedstuffs | Airtight. Unbreakable packaging; put breakable packaging into closed unbreakable container. Do not transport with food and feedstuffs. Note: E T+ symbol N symbol R: 45-26-48/23/25-62-63-68-50/53 S: 53-45-60-61 UN Hazard Class: 6.1 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0020

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

CADMIUM

ICSC: 0020

| | | |
|---|---|---|
| <p>I M P O R T A N T A D V I S I O N</p> | <p>PHYSICAL STATE; APPEARANCE: SOFT BLUE-WHITE METAL LUMPS OR GREY POWDER. MALLEABLE. TURNS BRITTLE ON EXPOSURE TO 80°C AND TARNISHES ON EXPOSURE TO MOIST AIR.</p> <p>PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air.</p> <p>CHEMICAL DANGERS: Reacts with acids forming flammable/explosive gas (hydrogen - see ICSC0001.) Dust reacts with oxidants, hydrogen azide, zinc, selenium or tellurium , causing fire and explosion hazard.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: (Total dust) 0.01 mg/m³ (Respirable fraction) 0.002 mg/m³ as TWA A2 (suspected human carcinogen); BEI issued (ACGIH 2005). MAK: skin absorption (H); Carcinogen category: 1; Germ cell mutagen group: 3A; (DFG 2004). OSHA PEL*: 1910.1027 TWA 0.005 mg/m³ *Note: The PEL applies to all Cadmium compounds (as Cd). NIOSH REL*: Ca See Appendix A *Note: The REL applies to all Cadmium compounds (as Cd). NIOSH IDLH: Ca 9 mg/m³ (as Cd) See: IDLH INDEX</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol and by ingestion.</p> <p>INHALATION RISK: A harmful concentration of airborne particles can be reached quickly when dispersed, especially if powdered.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: The fume is irritating to the respiratory tract Inhalation of fume may cause lung oedema (see Notes). Inhalation of fumes may cause metal fume fever. The effects may be delayed. Medical observation is indicated.</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: Lungs may be affected by repeated or prolonged exposure to dust particles. The substance may have effects on the kidneys , resulting in kidney impairment This substance is carcinogenic to humans.</p> |
|---|---|---|

| | | |
|-----------------------------------|--|--|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 765°C Melting point: 321°C Density: 8.6 g/cm³</p> | <p>Solubility in water: none Auto-ignition temperature: (cadmium metal dust) 250°C</p> |
|-----------------------------------|--|--|

| | |
|----------------------------------|--|
| <p>ENVIRONMENTAL DATA</p> | |
|----------------------------------|--|

NOTES

Reacts violently with fire extinguishing agents such as water, foam, carbon dioxide and halons. Depending on the degree of exposure, periodic medical examination is indicated. The symptoms of lung oedema often do not become manifest until a few hours have passed and they are aggravated by physical effort. Rest and medical observation are therefore essential. Do NOT take working clothes home. Cadmium also exists in a pyrophoric form (EC No. 048-011-00-X), which bears the additional EU labelling symbol F, R phrase 17, and S phrases 7/8 and 43. UN numbers and packing group will vary according to the physical form of the substance.

ADDITIONAL INFORMATION

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

| | |
|----------------------------|-----------------------|
| <p>ICSC: 0020</p> | <p>CADMIUM</p> |
| <p>(C) IPCS, CEC, 1994</p> | |

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

International Chemical Safety Cards

CHROMIUM

ICSC: 0029



Chrome
Cr
Atomic mass: 52.0
(powder)

ICSC # 0029
CAS # 7440-47-3
RTECS # [GB4200000](#)
October 27, 2004 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible under specific conditions. | No open flames if in powder form. | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! | |
| • INHALATION | Cough. | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. | Remove contaminated clothes. Rinse skin with plenty of water or shower. |
| • EYES | Redness. | Safety goggles. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|---------|-----------------------|
| Sweep spilled substance into containers; if appropriate, moisten first to prevent dusting. Personal protection: P2 filter respirator for harmful particles. | | R: S: |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0029

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

CHROMIUM

ICSC: 0029

| | | |
|----------|--|---|
| I | PHYSICAL STATE; APPEARANCE: GREY POWDER | ROUTES OF EXPOSURE: |
| M | PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air. | INHALATION RISK: A harmful concentration of airborne particles can be reached quickly when dispersed. |
| P | | |

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CHEMICAL DANGERS:

Chromium is a catalytic substance and may cause reaction in contact with many organic and inorganic substances , causing fire and explosion hazard.

EFFECTS OF SHORT-TERM EXPOSURE:

May cause mechanical irritation to the eyes and the respiratory tract.

OCCUPATIONAL EXPOSURE LIMITS:

TLV: (as Cr metal, Cr(III) compounds) 0.5 mg/m³ as TWA A4 (ACGIH 2004).
MAK not established.

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

OSHA PEL*: TWA 1 mg/m³ [See Appendix C](#) *Note: The PEL also applies to insoluble chromium salts.

NIOSH REL: TWA 0.5 mg/m³ [See Appendix C](#)

NIOSH IDLH: 250 mg/m³ (as Cr) See: [7440473](#)

PHYSICAL PROPERTIES

Boiling point: 2642°C
Melting point: 1900°C
Density: 7.15 g/cm³

Solubility in water:
none

ENVIRONMENTAL DATA

NOTES

The surface of the chromium particles is oxidized to chromium(III)oxide in air. See ICSC 1531 Chromium(III) oxide.

ADDITIONAL INFORMATION

ICSC: 0029

CHROMIUM

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

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International Chemical Safety Cards

COPPER

ICSC: 0240



Cu
(powder)

ICSC # 0240

CAS # 7440-50-8

RTECS # [GL5325000](#)

September 24, 1993 Validated

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible. | NO open flames. | Special powder, dry sand, NO other agents. |
| EXPLOSION | | | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! | |
| • INHALATION | Cough. Headache. Shortness of breath. Sore throat. | Local exhaust or breathing protection. | Fresh air, rest. Refer for medical attention. |
| • SKIN | Redness. | Protective gloves. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | Redness. Pain. | Safety goggles. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | Abdominal pain. Nausea. Vomiting. | Do not eat, drink, or smoke during work. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|--|-----------------------|
| Sweep spilled substance into containers. Carefully collect remainder. Then remove to safe place. (Extra personal protection: P2 filter respirator for harmful particles). | Separated from - See Chemical Dangers. | R: S: |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0240

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

COPPER

ICSC: 0240

| | | |
|---|---|--|
| <p>I</p> <p>M</p> <p>P</p> | <p>PHYSICAL STATE; APPEARANCE: RED POWDER, TURNS GREEN ON EXPOSURE TO MOIST AIR.</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS:</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation and by ingestion.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed.</p> |
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Shock-sensitive compounds are formed with acetylenic compounds, ethylene oxides and azides. Reacts with strong oxidants like chlorates, bromates and iodates, causing explosion hazard.

EFFECTS OF SHORT-TERM EXPOSURE:
Inhalation of fumes may cause metal fume fever. See Notes.

OCCUPATIONAL EXPOSURE LIMITS:
TLV: 0.2 mg/m³ fume (ACGIH 1992-1993).
TLV (as Cu, dusts & mists): 1 mg/m³ (ACGIH 1992-1993).
Intended change 0.1 mg/m³
Inhal.,
A4 (not classifiable as a human carcinogen);
MAK: 0.1 mg/m³ (Inhalable fraction)
Peak limitation category: II(2) Pregnancy risk group: D (DFG 2005).
OSHA PEL*: TWA 1 mg/m³ *Note: The PEL also applies to other copper compounds (as Cu) except copper fume.
NIOSH REL*: TWA 1 mg/m³ *Note: The REL also applies to other copper compounds (as Cu) except Copper fume.
NIOSH IDLH: 100 mg/m³ (as Cu) See: [7440508](#)

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:
Repeated or prolonged contact may cause skin sensitization.

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| PHYSICAL PROPERTIES | Boiling point: 2595°C Melting point: 1083°C Relative density (water = 1): 8.9 | Solubility in water: none |
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| ENVIRONMENTAL DATA | |
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NOTES

The symptoms of metal fume fever do not become manifest until several hours.

ADDITIONAL INFORMATION

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| ICSC: 0240 | (C) IPCS, CEC, 1994 | COPPER |
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International Chemical Safety Cards

LEAD

ICSC: 0052



Lead metal
Plumbum
Pb
Atomic mass: 207.2
(powder)


ICSC # 0052
CAS # 7439-92-1
RTECS # [OF7525000](#)
October 08, 2002 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---|--|---|---|
| FIRE | Not combustible. Gives off irritating or toxic fumes (or gases) in a fire. | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | See EFFECTS OF LONG-TERM OR REPEATED EXPOSURE. | PREVENT DISPERSION OF DUST! AVOID EXPOSURE OF (PREGNANT) WOMEN! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety spectacles. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | Abdominal pain. Nausea. Vomiting. | Do not eat, drink, or smoke during work. Wash hands before eating. | Rinse mouth. Give plenty of water to drink. Refer for medical attention. |
| SPILLAGE DISPOSAL | | STORAGE | PACKAGING & LABELLING |
| Sweep spilled substance into containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Do NOT let this chemical enter the environment. Personal protection: P3 filter respirator for toxic particles. | | Separated from food and feedstuffs incompatible materials See Chemical Dangers. | R: S: |
| SEE IMPORTANT INFORMATION ON BACK | | | |
| ICSC: 0052 | | Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values. | |

International Chemical Safety Cards

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| <p>I M P O R T A N T T A D A</p> | <p>PHYSICAL STATE; APPEARANCE: BLUISH-WHITE OR SILVERY-GREY SOLID IN VARIOUS FORMS. TURNS TARNISHED ON EXPOSURE TO AIR.</p> <p>PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air.</p> <p>CHEMICAL DANGERS: On heating, toxic fumes are formed. Reacts with oxidants. Reacts with hot concentrated nitric acid, boiling concentrated hydrochloric acid and sulfuric acid. Attacked by pure water and by weak organic acids in the presence of oxygen.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 0.05 mg/m³ A3 (confirmed animal carcinogen with unknown relevance to humans); BEI issued (ACGIH 2004). MAK: Carcinogen category: 3B; Germ cell mutagen group: 3A; (DFG 2004). EU OEL: as TWA 0.15 mg/m³ (EU 2002). OSHA PEL*: 1910.1025 TWA 0.050 mg/m³ See Appendix C *Note: The PEL also applies to other lead compounds (as Pb) -- see Appendix C. NIOSH REL*: TWA 0.050 mg/m³ See Appendix C *Note: The REL also applies to other lead compounds (as Pb) -- see Appendix C. NIOSH IDLH: 100 mg/m³ (as Pb) See: 7439921</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation and by ingestion.</p> <p>INHALATION RISK: A harmful concentration of airborne particles can be reached quickly when dispersed, especially if powdered.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE:</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: The substance may have effects on the blood bone marrow central nervous system peripheral nervous system kidneys , resulting in anaemia, encephalopathy (e.g., convulsions), peripheral nerve disease, abdominal cramps and kidney impairment. Causes toxicity to human reproduction or development.</p> |
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| PHYSICAL PROPERTIES | Boiling point: 1740°C Melting point: 327.5°C | Density: 11.34 g/cm ³ Solubility in water: none |
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| ENVIRONMENTAL DATA | Bioaccumulation of this chemical may occur in plants and in mammals. It is strongly advised that this substance does not enter the environment. |  |
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NOTES

Depending on the degree of exposure, periodic medical examination is suggested. Do NOT take working clothes home.
 Transport Emergency Card: TEC (R)-51S1872

ADDITIONAL INFORMATION

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| ICSC: 0052 | LEAD |
| (C) IPCS, CEC, 1994 | |

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| IMPORTANT LEGAL NOTICE: | Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values. |
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International Chemical Safety Cards

MANGANESE

ICSC: 0174

Mn
Atomic mass: 54.9
(powder)



ICSC # 0174
CAS # 7439-96-5
RTECS # [OO9275000](#)
November 27, 2003 Validated

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Combustible. | NO open flames. | Dry sand, special powder. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! AVOID EXPOSURE OF (PREGNANT) WOMEN! | |
| •INHALATION | Cough. | Local exhaust or breathing protection. | Fresh air, rest. Refer for medical attention. |
| •SKIN | | Protective gloves. | Rinse and then wash skin with water and soap. |
| •EYES | | Safety goggles, or eye protection in combination with breathing protection if powder. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | Abdominal pain. Nausea. | Do not eat, drink, or smoke during work. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|----------------------------|-----------------------|
| Sweep spilled substance into containers. Carefully collect remainder, then remove to safe place. (Extra personal protection: P2 filter respirator for harmful particles.) | Separated from acids. Dry. | |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0174

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards


MANGANESE

ICSC: 0174

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| I | <p>PHYSICAL STATE; APPEARANCE: GREY - WHITE POWDER</p> <p>PHYSICAL DANGERS:</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its aerosol and by ingestion.</p> |
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| <p>M P O R T A N T D A T A</p> | <p>Dust explosion possible if in powder or granular form, mixed with air.</p> <p>CHEMICAL DANGERS: Reacts slowly with water more rapidly with steam and acids forming flammable/explosive gas (hydrogen - see ICSC0001) causing fire and explosion hazard.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 0.2 mg/m³ (as TWA); (ACGIH 2003). MAK: (Inhalable fraction) 0.5 mg/m³; Pregnancy risk group: C; (DFG 2007). OSHA PEL*: C 5 mg/m³ *Note: Also see specific listings for Manganese cyclopentadienyl tricarbonyl and Methyl cyclopentadienyl manganese tricarbonyl. NIOSH REL*: TWA 1 mg/m³ ST 3 mg/m³ *Note: Also see specific listings for Manganese cyclopentadienyl tricarbonyl, Methyl cyclopentadienyl manganese tricarbonyl, and Manganese tetroxide. NIOSH IDLH: 500 mg/m³ (as Mn) See: 7439965</p> | <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: The aerosol is irritating to the respiratory tract .</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: The substance may have effects on the lungs and central nervous system , resulting in increased susceptibility to bronchitis, pneumonitis and neurologic, neuropsychiatric disorders (manganism). Animal tests show that this substance possibly causes toxicity to human reproduction or development.</p> |
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| PHYSICAL PROPERTIES | Boiling point: 1962°C Melting point: 1244°C Density: 7.47 g/cm ³ | Solubility in water: none |
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| ENVIRONMENTAL DATA | This substance may be hazardous in the environment; special attention should be given to aquatic organisms. |  |
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NOTES

Depending on the degree of exposure, periodic medical examination is suggested. The recommendations on this Card also apply to ferro manganese.

ADDITIONAL INFORMATION

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| ICSC: 0174 | MANGANESE |
| (C) IPCS, CEC, 1994 | |

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International Chemical Safety Cards

MERCURY

ICSC: 0056



Quicksilver
Liquid silver
Hg
Atomic mass: 200.6

ICSC # 0056
CAS # 7439-97-6
RTECS # [OV4550000](#)
UN # 2809
EC # 080-001-00-0
April 22, 2004 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|---|---|
| FIRE | Not combustible. Gives off irritating or toxic fumes (or gases) in a fire. | | In case of fire in the surroundings: use appropriate extinguishing media. |
| EXPLOSION | Risk of fire and explosion. | | In case of fire: keep drums, etc., cool by spraying with water. |
| EXPOSURE | | STRICT HYGIENE! AVOID EXPOSURE OF (PREGNANT) WOMEN! AVOID EXPOSURE OF ADOLESCENTS AND CHILDREN! | IN ALL CASES CONSULT A DOCTOR! |
| •INHALATION | Abdominal pain. Cough. Diarrhoea. Shortness of breath. Vomiting. Fever or elevated body temperature. | Local exhaust or breathing protection. | Fresh air, rest. Artificial respiration if indicated. Refer for medical attention. |
| •SKIN | MAY BE ABSORBED! Redness. | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. Refer for medical attention. |
| •EYES | | Face shield, or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| •INGESTION | | Do not eat, drink, or smoke during work. Wash hands before eating. | Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|--|--|
| Evacuate danger area in case of a large spill! Consult an expert! Ventilation. Collect leaking and spilled liquid in sealable non-metallic containers as far as possible. Do NOT wash away into sewer. Do NOT let this chemical enter the environment. Chemical protection suit including self-contained breathing apparatus. | Provision to contain effluent from fire extinguishing. Separated from food and feedstuffs Well closed. | Special material. Do not transport with food and feedstuffs. T symbol N symbol R: 23-33-50/53 S: 1/2-7-45-60-61 UN Hazard Class: 8 UN Packing Group: III |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0056

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.


International Chemical Safety Cards

MERCURY

ICSC: 0056

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| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: ODOURLESS, HEAVY AND MOBILE SILVERY LIQUID METAL.</p> <p>PHYSICAL DANGERS:</p> <p>CHEMICAL DANGERS: Upon heating, toxic fumes are formed. Reacts violently with ammonia and halogens causing fire and explosion hazard. Attacks aluminium and many other metals forming amalgams.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV: 0.025 mg/m³ as TWA (skin) A4 BEI issued (ACGIH 2004). MAK: 0.1 mg/m³ Sh Peak limitation category: II(8) Carcinogen category: 3B (DFG 2003). OSHA PEL[±]: C 0.1 mg/m³ NIOSH REL: Hg Vapor: TWA 0.05 mg/m³ skin Other: C 0.1 mg/m³ skin NIOSH IDLH: 10 mg/m³ (as Hg) See: 7439976</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of its vapour and through the skin, also as a vapour!</p> <p>INHALATION RISK: A harmful contamination of the air can be reached very quickly on evaporation of this substance at 20°C.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: The substance is irritating to the skin. Inhalation of the vapours may cause pneumonitis. The substance may cause effects on the central nervous system and kidneys. The effects may be delayed. Medical observation is indicated.</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: The substance may have effects on the central nervous system kidneys, resulting in irritability, emotional instability, tremor, mental and memory disturbances, speech disorders. Danger of cumulative effects. Animal tests show that this substance possibly causes toxic effects upon human reproduction.</p> |
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| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 357°C Melting point: -39°C Relative density (water = 1): 13.5 Solubility in water: none</p> | <p>Vapour pressure, Pa at 20°C: 0.26 Relative vapour density (air = 1): 6.93 Relative density of the vapour/air-mixture at 20°C (air = 1): 1.009</p> |
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| <p>ENVIRONMENTAL DATA</p> | <p>The substance is very toxic to aquatic organisms. In the food chain important to humans, bioaccumulation takes place, specifically in fish.</p> |  |
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NOTES

Depending on the degree of exposure, periodic medical examination is indicated. No odour warning if toxic concentrations are present. Do NOT take working clothes home.

Transport Emergency Card: TEC (R)-80GC9-II+III

ADDITIONAL INFORMATION

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| ICSC: 0056 | (C) IPCS, CEC, 1994 | MERCURY |
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| <p>IMPORTANT LEGAL NOTICE:</p> | <p>Neither NIOSH, the CEC or the IPCS nor any person acting on behalf of NIOSH, the CEC or the IPCS is responsible for the use which might be made of this information. This card contains the collective views of the IPCS Peer Review Committee and may not reflect in all cases all the detailed requirements included in national legislation on the subject. The user should verify compliance of the cards with the relevant legislation in the country of use. The only modifications made to produce the U.S. version is inclusion of the OSHA PELs, NIOSH RELs and NIOSH IDLH values.</p> |
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International Chemical Safety Cards

NICKEL

ICSC: 0062



Ni
Atomic mass: 58.7
(powder)

ICSC # 0062
CAS # 7440-02-0
RTECS # [QR5950000](#)
EC # 028-002-00-7
October 17, 2001 Peer reviewed

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|--|--|---|
| FIRE | Flammable as dust. Toxic fumes may be released in a fire. | | Dry sand. NO carbon dioxide. NO water. |
| EXPLOSION | Finely dispersed particles form explosive mixtures in air. | Prevent deposition of dust; closed system, dust explosion-proof electrical equipment and lighting. | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! AVOID ALL CONTACT! | |
| • INHALATION | Cough. Shortness of breath. | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. Protective clothing. | Remove contaminated clothes. Rinse and then wash skin with water and soap. |
| • EYES | | Safety spectacles, or eye protection in combination with breathing protection. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | Rinse mouth. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|---|------------------------------|-------------------------------------|
| Vacuum spilled material. Carefully collect remainder, then remove to safe place. Personal protection: P2 filter respirator for harmful particles. | Separated from strong acids. | Xn symbol R: 40-43 S: 2-22-36 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0062

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

NICKEL

ICSC: 0062

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| I | <p>PHYSICAL STATE; APPEARANCE: SILVERY METALLIC SOLID IN VARIOUS FORMS.</p> <p>PHYSICAL DANGERS:</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation of the dust.</p> |
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Dust explosion possible if in powder or granular form, mixed with air.

CHEMICAL DANGERS:

Reacts violently, in powder form, with titanium powder and potassium perchlorate, and oxidants such as ammonium nitrate, causing fire and explosion hazard. Reacts slowly with non-oxidizing acids and more rapidly with oxidizing acids. Toxic gases and vapours (such as nickel carbonyl) may be released in a fire involving nickel.

OCCUPATIONAL EXPOSURE LIMITS:

TLV:
(Inhalable fraction)
1.5 mg/m³ as TWA A5 (not suspected as a human carcinogen); (ACGIH 2004).
MAK: (Inhalable fraction) sensitization of respiratory tract and skin (Sah);
Carcinogen category: 1;
(DFG 2004).
OSHA PEL*†: TWA 1 mg/m³ *Note: The PEL does not apply to Nickel carbonyl.
NIOSH REL*: Ca TWA 0.015 mg/m³ [See Appendix A](#)
*Note: The REL does not apply to Nickel carbonyl.
NIOSH IDLH: Ca 10 mg/m³ (as Ni) See: [7440020](#)

INHALATION RISK:

Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed.

EFFECTS OF SHORT-TERM EXPOSURE:

May cause mechanical irritation. Inhalation of fumes may cause pneumonitis.

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

Repeated or prolonged contact may cause skin sensitization. Repeated or prolonged inhalation exposure may cause asthma. Lungs may be affected by repeated or prolonged exposure. This substance is possibly carcinogenic to humans.

PHYSICAL PROPERTIES

Boiling point: 2730°C
Melting point: 1455°C
Density: 8.9 g/cm³

Solubility in water:
none

ENVIRONMENTAL DATA

NOTES

At high temperatures, nickel oxide fumes will be formed. Depending on the degree of exposure, periodic medical examination is suggested. The symptoms of asthma often do not become manifest until a few hours have passed and they are aggravated by physical effort. Rest and medical observation are therefore essential. Anyone who has shown symptoms of asthma due to this substance should avoid all further contact with this substance.

ADDITIONAL INFORMATION

ICSC: 0062

NICKEL

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

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International Chemical Safety Cards

SILVER

ICSC: 0810



Argentum
C.I. 77820
Ag

ICSC # 0810
CAS # 7440-22-4
RTECS # [VW3500000](#)
September 10, 1997 Validated

| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|------------------------------------|--|---|
| FIRE | Not combustible, except as powder. | | |
| EXPLOSION | | | |
| EXPOSURE | | PREVENT DISPERSION OF DUST! | |
| • INHALATION | | Local exhaust or breathing protection. | Fresh air, rest. |
| • SKIN | | Protective gloves. | Rinse skin with plenty of water or shower. |
| • EYES | | Safety spectacles, or eye protection in combination with breathing protection if powder. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | | Do not eat, drink, or smoke during work. | |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|-----------------------|
| Sweep spilled substance into containers; if appropriate, moisten first to prevent dusting. Carefully collect remainder, then remove to safe place. Do NOT let this chemical enter the environment. | Separated from ammonia, strong hydrogen peroxide solutions, strong acids. | |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 0810

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

SILVER

ICSC: 0810

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| I | PHYSICAL STATE; APPEARANCE: WHITE METAL, TURNS DARK ON EXPOSURE TO OZONE, HYDROGEN SULFIDE OR SULFUR. | ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation and by ingestion. |
| M | | |
| P | PHYSICAL DANGERS: | INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed. |
| O | CHEMICAL DANGERS: Shock-sensitive compounds are formed with acetylene. | |
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Reacts with acids causing fire hazard. Contact with strong hydrogen peroxide solution will cause violent decomposition to oxygen gas. Contact with ammonia may cause formation of compounds that are explosive when dry.

OCCUPATIONAL EXPOSURE LIMITS:

TLV (metal): 0.1 mg/m³ (ACGIH 1997).
 EU OEL: 0.1 mg/m³ as TWA (EU 2000).
 OSHA PEL: TWA 0.01 mg/m³
 NIOSH REL: TWA 0.01 mg/m³
 NIOSH IDLH: 10 mg/m³ (as Ag) See: [IDLH INDEX](#)

EFFECTS OF SHORT-TERM EXPOSURE:

Inhalation of high amounts of metallic silver vapours may cause lung damage with pulmonary oedema.

EFFECTS OF LONG-TERM OR REPEATED EXPOSURE:

The substance may cause a grey-blue discoloration of the eyes, nose, throat and skin (argyria/argyrosis).

PHYSICAL PROPERTIES

Boiling point: 2212°C
 Melting point: 962°C

Relative density (water = 1): 10.5
 Solubility in water: none

ENVIRONMENTAL DATA

This substance may be hazardous to the environment; special attention should be given to aquatic organisms.



NOTES

Card has been partially updated in March 2008: see Occupational Exposure Limits.

ADDITIONAL INFORMATION

ICSC: 0810

SILVER

(C) IPCS, CEC, 1994

IMPORTANT LEGAL NOTICE:

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International Chemical Safety Cards

ZINC POWDER

ICSC: 1205



Blue powder
Merrillite
Zn
Atomic mass: 65.4
(powder)

ICSC # 1205
CAS # 7440-66-6
RTECS # [ZG8600000](#)
UN # 1436 (zinc powder or dust)
EC # 030-001-00-1
October 24, 1994 Peer reviewed



| TYPES OF HAZARD/ EXPOSURE | ACUTE HAZARDS/ SYMPTOMS | PREVENTION | FIRST AID/ FIRE FIGHTING |
|---------------------------|---|--|---|
| FIRE | Highly flammable. Many reactions may cause fire or explosion. Gives off irritating or toxic fumes (or gases) in a fire. | NO open flames, NO sparks, and NO smoking. NO contact with acid(s), base (s) and incompatible substances (see Chemical Dangers). | Special powder, dry sand, NO other agents. NO water. |
| EXPLOSION | Risk of fire and explosion on contact with acid(s), base(s), water and incompatible substances. | Closed system, ventilation, explosion-proof electrical equipment and lighting. Prevent build-up of electrostatic charges (e.g., by grounding). Prevent deposition of dust. | In case of fire: cool drums, etc., by spraying with water but avoid contact of the substance with water. |
| EXPOSURE | | PREVENT DISPERSION OF DUST! STRICT HYGIENE! | |
| • INHALATION | Metallic taste and metal fume fever. Symptoms may be delayed (see Notes). | Local exhaust. | Fresh air, rest. Refer for medical attention. |
| • SKIN | Dry skin. | Protective gloves. | Rinse and then wash skin with water and soap. |
| • EYES | | Safety spectacles. | First rinse with plenty of water for several minutes (remove contact lenses if easily possible), then take to a doctor. |
| • INGESTION | Abdominal pain. Nausea. Vomiting. | Do not eat, drink, or smoke during work. Wash hands before eating. | Rinse mouth. Refer for medical attention. |

| SPILLAGE DISPOSAL | STORAGE | PACKAGING & LABELLING |
|--|---|---|
| Extinguish or remove all ignition sources. Do NOT wash away into sewer. Sweep spilled substance into containers. then remove to safe place. Personal protection: self-contained breathing apparatus. | Fireproof. Separated from acids, bases oxidants Dry. | Airtight. F symbol N symbol R: 15-17-50/53 S: 2-7/8-43-46-60-61 UN Hazard Class: 4.3 UN Subsidiary Risks: 4.2 |

SEE IMPORTANT INFORMATION ON BACK

ICSC: 1205

Prepared in the context of cooperation between the International Programme on Chemical Safety & the Commission of the European Communities (C) IPCS CEC 1994. No modifications to the International version have been made except to add the OSHA PELs, NIOSH RELs and NIOSH IDLH values.

International Chemical Safety Cards

ZINC POWDER

ICSC: 1205

| | | |
|---|---|---|
| <p>I M P O R T A N T D A T A</p> | <p>PHYSICAL STATE; APPEARANCE: ODOURLESS GREY TO BLUE POWDER.</p> <p>PHYSICAL DANGERS: Dust explosion possible if in powder or granular form, mixed with air. If dry, it can be charged electrostatically by swirling, pneumatic transport, pouring, etc.</p> <p>CHEMICAL DANGERS: Upon heating, toxic fumes are formed. The substance is a strong reducing agent and reacts violently with oxidants. Reacts with water and reacts violently with acids and bases forming flammable/explosive gas (hydrogen - see ICSC0001) Reacts violently with sulfur, halogenated hydrocarbons and many other substances causing fire and explosion hazard.</p> <p>OCCUPATIONAL EXPOSURE LIMITS: TLV not established.</p> | <p>ROUTES OF EXPOSURE: The substance can be absorbed into the body by inhalation and by ingestion.</p> <p>INHALATION RISK: Evaporation at 20°C is negligible; a harmful concentration of airborne particles can, however, be reached quickly when dispersed.</p> <p>EFFECTS OF SHORT-TERM EXPOSURE: Inhalation of fumes may cause metal fume fever. The effects may be delayed.</p> <p>EFFECTS OF LONG-TERM OR REPEATED EXPOSURE: Repeated or prolonged contact with skin may cause dermatitis.</p> |
|---|---|---|

| | | |
|-----------------------------------|---|--|
| <p>PHYSICAL PROPERTIES</p> | <p>Boiling point: 907°C Melting point: 419°C Relative density (water = 1): 7.14</p> | <p>Solubility in water: reaction Vapour pressure, kPa at 487°C: 0.1 Auto-ignition temperature: 460°C</p> |
|-----------------------------------|---|--|

| | |
|----------------------------------|--|
| <p>ENVIRONMENTAL DATA</p> | |
|----------------------------------|--|

NOTES

Zinc may contain trace amounts of arsenic, when forming hydrogen, may also form toxic gas arsine (see ICSC 0001 and ICSC 0222). Reacts violently with fire extinguishing agents such as water, halons, foam and carbon dioxide. The symptoms of metal fume fever do not become manifest until several hours later. Rinse contaminated clothes (fire hazard) with plenty of water.

Transport Emergency Card: TEC (R)-43GWS-II+III
NFPA Code: H0; F1; R1;

ADDITIONAL INFORMATION

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

ICSC: 1205

ZINC POWDER

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

APPENDIX D
HOSPITAL INFORMATION AND MAP
FIELD ACCIDENT REPORT

FIELD ACCIDENT REPORT

This report is to be filled out by the designated Site Safety Officer after EVERY accident.

PROJECT NAME _____ PROJECT. NO. _____

Date of Accident _____ Time _____ Report By _____

Type of Accident (Check One):

Vehicular Personal Property

Name of Injured _____ DOB or Age _____

How Long Employed _____

Names of Witnesses _____

Description of Accident _____

Action Taken _____

Did the Injured Lose Any Time? _____ How Much (Days/Hrs.)? _____

Was Safety Equipment in Use at the Time of the Accident (Hard Hat, Safety Glasses, Gloves, Safety Shoes, etc.)? _____

(If not, it is the EMPLOYEE'S sole responsibility to process his/her claim through his/her Health and Welfare Fund.)

INDICATE STREET NAMES, DESCRIPTION OF VEHICLES, AND NORTH ARROW

HOSPITAL INFORMATION AND MAP

The hospital nearest the site is:

Lincoln Medical and Mental Health Center

234 E 149th St
 Bronx, NY 10451
 (718) 579-5016

0.9 Miles – About 6 Minutes

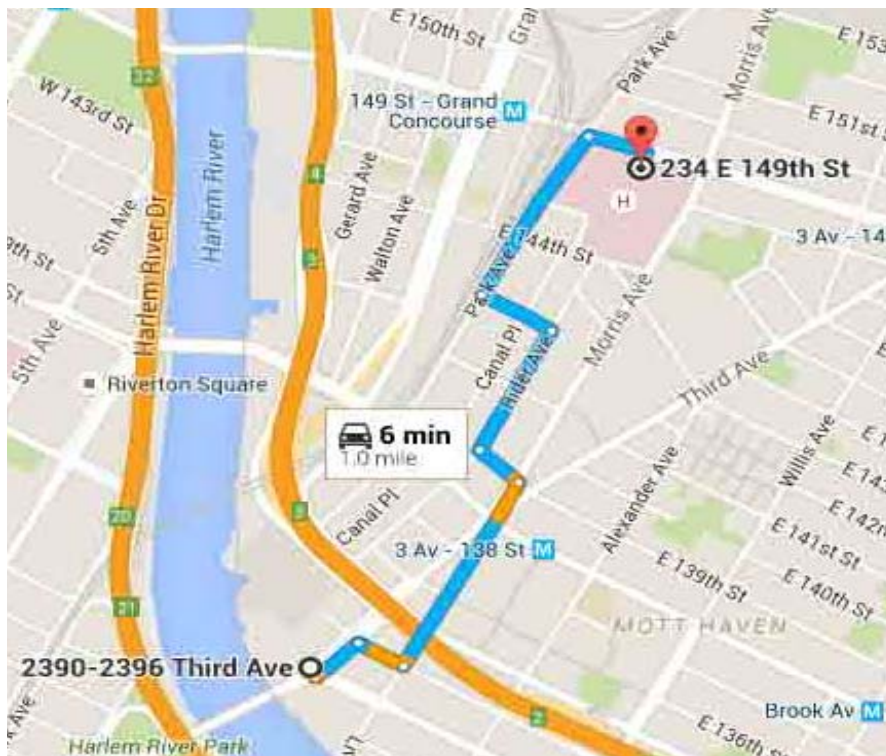
Take Lincoln Ave, Rider Ave and Park Ave to E 149th St

6 min (0.9 mi)

- ↑ 1. Head south on Third Ave toward Bruckner Blvd 82 ft
- ↶ 2. Turn left to stay on Third Ave 335 ft
- ↷ 3. Turn right onto E 134th St 299 ft
- ↶ 4. Turn left at the 1st cross street onto Lincoln Ave 0.3 mi
- ↶ 5. Turn left onto E 139th St 312 ft
- ↷ 6. Turn right onto Rider Ave 0.2 mi
- ↶ 7. Turn left onto E 141st St 459 ft
- ↷ 8. Turn right onto Park Ave 0.2 mi

- ↷ Turn right onto E 149th St 46 s (348 ft)

i Destination will be on the right



ATTACHMENT C
Quality Assurance Project Plan

**QUALITY ASSURANCE PROJECT PLAN
FORMER MUGLER SHORING SITE
2401 Third Avenue, Bronx, NY**

Prepared on behalf of:

2401 3rd Ave Associates Property LLC
512 Seventh Avenue 15th Floor
New York, NY 10018

Prepared by:

EBC
ENVIRONMENTAL BUSINESS CONSULTANTS
RIDGE, NY 11961

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QUALITY ASSURANCE PROJECT PLAN FORMER MUGLER SHORINGSITE

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1.0 INTRODUCTION

This Quality Assurance Project Plan (QAPP) has been prepared in accordance with DER-10 to detail procedures to be followed during the course of the sampling and analytical portion of the project, as required by the approved work plan.

To ensure the successful completion of the project each individual responsible for a given component of the project must be aware of the quality assurance objectives of his / her particular work and of the overall project. The EBC Project Director, Charles Sosik will be directly responsible to the client for the overall project conduct and quality assurance/quality control (QA/QC) for the project. The Project Director will be responsible for overseeing all technical and administrative aspects of the project and for directing QA/QC activities. Mr. Patrick Recio will serve as the Quality Assurance Officer (QAO) and in this role may conduct:

- conduct periodic field and sampling audits;
- interface with the analytical laboratory to resolve problems; and
- interface with the data validator and/or the preparer of the DUSR to resolve problems.

Robert Bennett will serve as the Project Manager and will be responsible for implementation of the IRM and coordination with field sampling crews and subcontractors. Reporting directly to the Project Manager will be the Field Operations Officer, Kevin Waters; who will serve as the on-Site qualified environmental professional who will record observations, direct the field crew and be responsible for the collection and handling of all samples.

1.1 Organization

Project QA will be maintained under the direction of the Project Manager, in accordance with this QAPP. QC for specific tasks will be the responsibility of the individuals and organizations listed below, under the direction and coordination of the Project Manager

| GENERAL RESPONSIBILITY | SCOPE OF WORK | RESPONSIBILITY OF QUALITY CONTROL |
|------------------------|---|-----------------------------------|
| Field Operations | Supervision of Field Crew, sample collection and handling | K. Waters, EBC |
| Project Manager | Implementation of the RI according to the RIWP. | R. Bennett, EBC |
| Laboratory Analysis | Analysis of soil samples by NYSDEC ASP methods Laboratory | NYSDOH-Certified Laboratory |
| Data review | Review for completeness and compliance | 3 rd party validation |

2.0 QUALITY ASSURANCE PROJECT PLAN OBJECTIVES

2.1 Overview

Overall project goals are defined through the development of Data Quality Objectives (DQOs), which are qualitative and quantitative Statements that specify the quality of the data required to support decisions; DQOs, as described in this section, are based on the end uses of the data as described in the work plan.

In this plan, Quality Assurance and Quality Control are defined as follows:

- Quality Assurance - The overall integrated program for assuring reliability of monitoring and measurement data.
- Quality Control - The routine application of procedures for obtaining prescribed standards of performance in the monitoring and measurement process.

2.2 QA / QC Requirements for Analytical Laboratory

Samples will be analyzed by a New York State Department of Health (NYSDOH) certified laboratory that is certified in the appropriate categories. Data generated from the laboratory will be used to evaluate contaminants such as chlorinated and other volatile organic compounds (VOCs) in soil, soil gas and groundwater. The QA requirements for all subcontracted analytical laboratory work performed on this project are described below. QA elements to be evaluated include accuracy, precision, sensitivity, representativeness, and completeness. The data generated by the analytical laboratory for this project are required to be sensitive enough to achieve required quantification limits as specified in NYSDEC Analytical Services Protocol (NYSDEC ASP, 07/2005) and useful for comparison with clean-up objectives. The analytical results meeting the required quantification limits will provide data sensitive enough to meet the data quality objectives of this remedial program as described in the work plan. Reporting of the data must be clear, concise, and comprehensive. The QC elements that are important to this project are completeness of field data, sample custody, sample holding times, sample preservation, sample storage, instrument calibration and blank contamination.

2.2.1 Instrument Calibration

Calibration curves will be developed for each of the compounds to be analyzed. Standard concentrations and a blank will be used to produce the initial curves. The development of calibration curves and initial calibration response factors must be consistent with method requirements presented in the most recent version of NYSDEC ASP 07/2005).

2.2.2 Continuing Instrument Calibration

The initial calibration curve will be verified every 12 hrs by analyzing one calibration standard. The standard concentration will be the midpoint concentration of the initial calibration curve. The calibration check compound must come within 25% relative percent difference (RPD) of the average response factor obtained during initial calibration. If the RPD is greater than 25%, then corrective action must be taken as provided in the specific methodology.

2.2.3 Method Blanks

Method blank or preparation blank is prepared from an analyte-free matrix which includes the same reagents, internal standards and surrogate standards as the related samples. It is carried through the

entire sample preparation and analytical procedure. A method blank analysis will be performed once for each 12 hr period during the analysis of samples for volatiles. An acceptable method blank will contain less than two (2) times the CRQL of methylene chloride, acetone and 2-butanone. For all other target compounds, the method blank must contain less than or equal to the CRQL of any single target compound. For non-target peaks in the method blank, the peak area must be less than 10 percent of the nearest internal standard. The method blank will be used to demonstrate the level of laboratory background and reagent contamination that might result from the analytical process itself.

2.2.4 Trip Blanks.

Trip blanks consist of a single set of sample containers filled at the laboratory with deionized, laboratory-grade water. The water used will be from the same source as that used for the laboratory method blank. The containers will be carried into the field and handled and transported in the same way as the samples collected that day. Analysis of the trip blank for VOCs is used to identify contamination from the air, shipping containers, or from other items coming in contact with the sample bottles. (The bottles holding the trip blanks will be not opened during this procedure.) A complete set of trip blanks will be provided with each shipment of samples to the certified laboratory.

2.2.5 Surrogate Spike Analysis

For organic analyses, all samples and blanks will be spiked with surrogate compounds before purging or extraction in order to monitor preparation and analyses of samples. Surrogate spike recoveries shall fall within the advisory limits in accordance with the NY5DEC ASP protocols for samples falling within the quantification limits without dilution.

2.2.6 Matrix Spike / Matrix Spike Duplicate / Matrix Spike Blank (MS/MSDIMSB) Analysis

MS, MSD and MSB analyses will be performed to evaluate the matrix effect of the sample upon the analytical methodology along with the precision of the instrument by measuring recoveries. The MS / MSD / MSB samples will be analyzed for each group of samples of a similar matrix at a rate of 5% (one for every 20 field samples). The RPD will be calculated from the difference between the MS and MSD. Matrix spike blank analysis will be performed to indicate the appropriateness of the spiking solution(s) used for the MS/MSD. 10% of the samples of each matrix should be sampled and analyzed as Duplicates.

2.3 Accuracy

Accuracy is defined as the nearness of a real or the mean (x) of a set of results to the true value. Accuracy is assessed by means of reference samples and percent recoveries. Accuracy includes both precision and recovery and is expressed as percent recovery (% REC). The MS sample is used to determine the percent recovery. The matrix spike percent recovery (% REC) is calculated by the following equation:

$$\%REC = \frac{SSR - SR}{SA} \times 100$$

Where:

SSR = spike sample results

SR = sample results

SA = spike added from spiking mix

2.4 Precision

Precision is defined as the measurement of agreement of a set of replicate results among themselves without a Precision is defined as the measurement of agreement of a set of replicate results among themselves without assumption of any prior information as to the true result. Precision is assessed by means of duplicate/replicate sample analyses.

Analytical precision is expressed in terms of RPD. The RPD is calculated using the following formula:

$$RPD = \frac{D^1 - D^2}{(D^1 + D^2)/2} \times 100$$

Where:

RPD = relative percent difference

D¹ = first sample value

D² = second sample value (duplicate)

2.5 Sensitivity

The sensitivity objectives for this plan require that data generated by the analytical laboratory achieve quantification levels low enough to meet the required detection limits specified by NYSDEC ASP and to meet all site-specific standards, criteria and guidance values (SGCs) established for this project.

2.6 Representativeness

Representativeness is a measure of the relationship of an individual sample taken from a particular site to the remainder of that site and the relationship of a small aliquot of the sample (i.e., the one used in the actual analysis) to the sample remaining on site. The representativeness of samples is assured by adherence to sampling procedures described in the Remedial Investigation Work Plan.

2.7 Completeness

Completeness is a measure of the quantity of data obtained from a measurement system as compared to the amount of data expected from the measurement system. Completeness is defined as the percentage of all results that are not affected by failing QC qualifiers, and should be between 70 and 100% of all analyses performed. The objective of completeness in laboratory reporting is to provide a thorough data support package. The laboratory data package provides documentation of sample analysis and results in the form of summaries, QC data, and raw analytical data. The laboratory will be required to submit data packages that follow NYSDEC ASP Category B reporting format which, at a minimum, will include the following components:

1. All sample chain-of-custody forms.
2. The case narrative(s) presenting a discussion of any problems and/or procedural changes required during analyses. Also presented in the case narrative are sample summary forms.
3. Documentation demonstrating the laboratory's ability to attain the contract specified detection limits for all target analytes in all required matrices.
4. Tabulated target compound results and tentatively identified compounds.
5. Surrogate spike analysis results (organics).
6. Matrix spike/matrix spike duplicate/matrix spike blank results.
7. QC check sample and standard recovery results
8. Blank results (field, trip, and method).
9. Internal standard area and RT summary.

2.8 Laboratory Custody Procedures

The following elements are important for maintaining the field custody of samples:

- Sample identification
- Sample labels
- Custody records
- Shipping records
- Packaging procedures

Sample labels will be attached to all sampling bottles before field activities begin; each label will contain an identifying number. Each number will have a suffix that identifies the site and where the sample was taken. Approximate sampling locations will be marked on a map with a description of the sample location. The number, type of sample, and sample identification will be entered into the field logbook. A chain-of-custody form, initiated at the analytical laboratory will accompany the sample bottles from the laboratory into the field. Upon receipt of the bottles and cooler, the sampler will sign and date the first received blank space. After each sample is collected and appropriately identified, entries will be made on the chain-of-custody form that will include:

- Site name and address
- Samplers' names and signatures

2.9 Sample Handling and Decontamination Procedures

Collected samples will be appropriately packaged, placed in coolers and shipped via overnight courier or delivered directly to the analytical laboratory by field personnel. Samples will be containerized in appropriate laboratory provided glassware and shipped in plastic coolers. Samples will be preserved through the use of ice or cold-pak(s) to maintain a temperature of 4°C.

Dedicated disposable sampling materials will be used for both soil and groundwater samples (if collected), eliminating the need to prepare field equipment (rinsate) blanks. However, if non-disposable equipment is used, (stainless steel scoop, etc.) field rinsate blanks will be prepared at the rate of 1 for every eight samples collected. No field filtering will be conducted; any required filtration will be completed by the laboratory.

Decontamination of non-dedicated sampling equipment will consist of the following:

- Gently tap or scrape to remove adhered soil;
- Rinse with tap water;
- Wash withalconox® detergent solution and scrub ;
- Rinse with tap water;
- Rinse with distilled or deionized water.

Prepare field blanks by pouring distilled or deionized water over decontaminated equipment and collecting the water in laboratory provided containers. Trip blanks will accompany samples each time they are transported to the laboratory. Matrix spike and matrix spike duplicates (MS/MSD) will be collected at the rate of one per 20 samples submitted to the laboratory and duplicate samples will be collected at a rate of one per ten samples submitted to the laboratory.

3.0 ANALYTICAL PROCEDURES

3.1 Laboratory Analysis

Samples will be analyzed by the NYSDOH ELAP laboratory for one or more of the following parameters: VOCs + TICs in soil / groundwater by USEPA Method 8260C, SVOCs + TICs in soil / groundwater by USEPA Method 8270D, Target Analyte List (TAL) Metals 6010 in soil and groundwater, pesticides / PCBs by USEPA Method 8081B/8082A and VOCs in air by USEPA Method TO15 (Table 2). If any modifications or additions to the standard procedures are anticipated and if any nonstandard sample preparation or analytical protocol is to be used, the modifications and the nonstandard protocol will be explicitly defined and documented. Prior approval by EBC's PM will be necessary for any nonstandard analytical or sample preparation protocol used by the laboratory, i.e., dilution of samples or extracts by greater than a factor of five (5).

4.0 DATA REDUCTION, REVIEW, AND REPORTING

4.1 Overview

The process of data reduction, review, and reporting ensures the assessments or a conclusion based on the final data accurately reflects actual site conditions. This plan presents the specific procedures, methods, and format that will be employed for data reduction, review and reporting of each measurement parameter determined in the laboratory and field. Also described in this section is the process by which all data, reports, and work plans are proofed and checked for technical and numerical errors prior to final submission.

4.2 Data Reduction

Standard methods and references will be used as guidelines for data handling, reduction, validation, and reporting. All data for the project will be compiled and summarized with an independent verification at each step in the process to prevent transcription/typographical errors. Any computerized entry of data will also undergo verification review.

Sample analysis will be provided by a New York State certified environmental laboratory. Laboratory reports will include ASP category B deliverables for use in the preparation of a data usability summary report (DUSR). All results will be provided in accordance with the NYSDEC Environmental Information Management System (EIMS) electronic data deliverable (EDD) format. Analytical results shall be presented on standard NYSDEC ASP-B forms or equivalents, and include the dates the samples were received and analyzed, and the actual methodology used. Note that if waste characterization samples are analyzed they will be in results only format and will not be evaluated in the DUSR.

Laboratory QA/QC information required by the method protocols will be compiled, including the application of data QA/QC qualifiers as appropriate. In addition, laboratory worksheets, laboratory notebooks, chains-of-custody, instrument logs, standards records, calibration records, and maintenance records, as applicable, will be provided in the laboratory data packages to determine the validity of data. Specifics on internal laboratory data reduction protocols are identified in the laboratory's SOPs.

Following receipt of the laboratory analytical results by EBC, the data results will be compiled and presented in an appropriate tabular form. Where appropriate, the impacts of QA/QC qualifiers resulting from laboratory or external validation reviews will be assessed in terms of data usability.

4.3 Laboratory Data Reporting

All sample data packages submitted by the analytical laboratory will be required to be reported in conformance to the NYSDEC ASP (7/2005), Category B data deliverable requirements as applicable to the method utilized. All results will be provided in accordance with the NYSDEC Environmental Information Management System (EIMS) electronic data deliverable (EDD) format. Note that waste characterization samples, if analyzed, will be in results only format and will not be evaluated in the DUSR.

5.0 CORRECTIVE ACTION

Review and implementation of systems and procedures may result in recommendations for corrective action. Any deviations from the specified procedures within approved project plans due to unexpected site-specific conditions shall warrant corrective action. All errors, deficiencies, or other problems shall be brought to the immediate attention of the EBC PM, who in turn shall contact the Quality Assurance/Data Quality Manager or his designee (if applicable).

Procedures have been established to ensure that conditions adverse to data quality are promptly investigated, evaluated and corrected. These procedures for review and implementation of a change are as follows:

- Define the problem.
- Investigate the cause of the problem.
- Develop a corrective action to eliminate the problem, in consultation with the personnel who defined the problem and who will implement the change.
- Complete the required form describing the change and its rationale (see below for form requirements).
- Obtain all required written approvals.
- Implement the corrective action.
- Verify that the change has eliminated the problem.

During the field investigation, all changes to the sampling program will be documented in field logs/sheets and the EBC PM advised.

If any problems occur with the laboratory or analyses, the laboratory must immediately notify the PM, who will consult with other project staff. All approved corrective actions shall be controlled and documented.

All corrective action documentation shall include an explanation of the problem and a proposed solution which will be maintained in the project file or associated logs. Each report must be approved by the necessary personnel (e.g., the PM) before implementation of the change occurs. The PM shall be responsible for controlling, tracking, implementing and distributing identified changes.

**TABLE 1
SUMMARY OF
SAMPLING PROGRAM RATIONALE AND ANALYSIS**

| Matrix | Location | Approximate Number of Samples | Frequency | Rationale for Sampling | Laboratory Analysis | Duplicates | Matrix Spikes | Spike Duplicates | Trip Blanks |
|--------|--------------------------------------|-------------------------------|---|---|--|------------|------------------|------------------|-------------|
| Soil | UST Areas | 15 | 1 per 900 square feet of excavation base | Endpoint Verification of footing excavations | VOCs EPA Method 8260B, SVOCs EPA Method 8270, | 1 per day | 1 per 20 samples | 1 per 20 samples | 1 per trip |
| | Excavated Petroleum Impacted Soil | 1 | 1 per 800 cy | Waste Characterization for disposal if not stockpiled on site | VOCs EPA Method 8260B, PAHs EPA Method 8270, RCRA metals, pesticides and PCBs by EPA 8081/8082, other as per disposal facility | 0 | 0 | 0 | 0 |
| Soil | Excavated Historic Fill Material | 19 | 1 per 800 cy | Waste Characterization for disposal if not stockpiled on site | VOCs EPA Method 8260B, PAHs EPA Method 8270, RCRA metals, pesticides and PCBs by EPA 8081/8082, other as per disposal facility | 0 | 0 | 0 | 0 |
| | Excavated Uncontaminated Native Soil | 9 | 7 Grabs for 1st 1,000 cy, 2 for each additional 1,000 cy As per CP51 | Clean Verification for disposal if not stockpiled. | VOCs EPA Method 8260B | 0 | 0 | 0 | 0 |
| Soil | Excavated Uncontaminated Native Soil | 3 | 2 Composites for 1st 1,000 cy, 1 for each additional 1,000 cy As per CP51 | Clean Verification for disposal if not stockpiled. | SVOCs, pesticides/and PCBs by EPA 8081/8082, and RCRA metals. | 0 | 0 | 0 | 0 |

**TABLE 2
SAMPLE COLLECTION AND ANALYSIS PROTOCOLS**

| Sample Type | Matrix | Sampling Device | Parameter | Sample Container | Sample Preservation | Analytical Method# | CRQL / MDLH | Holding Time |
|--------------------|---------------|------------------------|---------------------|--------------------------|----------------------------|--------------------------------------|--------------------------------|---------------------|
| Soil | Soil | Scoop Direct into Jar | VOCs | (1) 2 oz Jar | Cool to 4° C | EPA Method 8260C (test method 5035A) | Compound specific (1-5 ug/kg) | 14 days |
| Soil | Soil | Scoop Direct into Jar | SVOCs | (1) 8 oz jar | Cool to 4° C | EPA Method 8270D | Compound specific (1-5 ug/kg) | 14 day ext/40 days |
| Soil | Soil | Scoop Direct into Jar | Pest/PCBs | from 8oz jar above | Cool to 4° C | EPA Method 8081B/8082A | Compound specific (1-5 ug/kg) | 14 day ext/40 days |
| Soil | Soil | Scoop Direct into Jar | Metals | from 8oz jar above | Cool to 4° C | TAL Metals 6010 | Compound specific (01-1 mg/kg) | 6 months |
| Groundwater | Water | Pump tubing | VOCs | (3) 40 ml vials | Cool to 4° C 1:1 HCL | EPA Method 8260C | Compound specific (1-5 ug/L) | 14 days |
| Groundwater | Water | Pump tubing | SVOCs | (1) 1 Liter Amber Bottle | Cool to 4° C | EPA Method 8270D | Compound specific (1-5 ug/L) | 14 days |
| Groundwater | Water | Pump tubing | Pesticides and PCBs | (2) 1 Liter Amber Bottle | Cool to 4° C | EPA Method 8081B / 8082A | Compound specific (1-5 ug/L) | 14 days |
| Groundwater | water | Pump tubing | Total Metals | (1) 100 ml | HNO3 | TAL Metals 6010 | Compound specific (1-5 mg/L) | 6 months |
| Groundwater | water | Pump tubing | Dissolved Metals | (1) 100 ml | None | TAL Metals 6010 | Compound specific (1-5 mg/L) | 6 months |

Notes:

All holding times listed are from Verified Time of Sample Receipt (VTSR) unless noted otherwise. * Holding time listed is from time of sample collection.

The number in parentheses in the "Sample Container" column denotes the number of containers needed.

Triple volume required when collected MS/MSD samples

The number of trip blanks are estimated.

CRQL / MDL = Contract Required Quantitation Limit / Method Detection Limit

NA = Not available or not applicable.

ATTACHMENT D
Community Air Monitoring Plan

COMMUNITY AIR MONITORING PLAN

FORMER MUGLER SHORING SITE
2401 THIRD AVENUE
BRONX, NY

APRIL - 2016

**COMMUNITY AIR MONITORING PLAN
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APPENDICES

Appendix A Action Limit Report

1.0 INTRODUCTION

This Community Air Monitoring Plan (CAMP) has been prepared for the excavation and construction activities to be performed under a Remedial Action Work Plan (RAWP) at 2401 Third Avenue, in Bronx, NY. The CAMP provides measures for protection for the downwind community (i.e., off-site receptors including residences, businesses, and on-site workers not directly involved in the investigation activities) from potential airborne contaminant releases resulting from excavation activities at the site.

Compliance with this CAMP is required during all activities associated with soil disturbance activities that have the potential to generate airborne particulate matter and volatile organic compounds (VOCs). These activities include excavation and loading of affected soil. This CAMP has been prepared to ensure that remedial activities do not adversely affect passersby, residents, or workers in the area immediately surrounding the Site and to preclude or minimize airborne migration of site-related contaminants to off-site areas.

1.1 Regulatory Requirements

This CAMP was established in accordance with the following requirements:

- New York State Department of Health's (NYSDOH) Generic Community Air Monitoring Plan as presented in DER-10 Technical Guidance for Site Investigation and Remediation (NYSDEC May 3, 2010). This guidance specifies that a community air-monitoring program shall be implemented to protect the surrounding community and to confirm that the work does not spread contamination off-site through the air.

2.0 AIR MONITORING

Petroleum volatile organic compounds (VOCs), chlorinated VOCs, semi-volatile organic compounds (SVOCs), metals and pesticides are the constituents of concern at the Site. The appropriate method to monitor air for these constituents during remediation activities is through real-time VOC and air particulate (dust) monitoring.

2.1 Meteorological Data

At a minimum, wind direction will be evaluated at the start of each workday, noon of each workday, and the end of each workday. These readings will be utilized to position the monitoring equipment in appropriate upwind and downwind locations.

2.2 Community Air Monitoring Requirements

To establish ambient air background concentrations, air will be monitored at several locations around the site perimeter before activities begin. These points will be monitored periodically in series during the site work. When the excavation area is within 20 feet of potentially exposed populations or occupied structures, the perimeter monitoring points will be located to represent the nearest potentially exposed individuals at the downwind location and will take into account the locations of ventilation system intakes of nearby structures.

Fugitive respirable dust will be monitored using a MiniRam Model PDM-3 aerosol monitor (or equivalent). Air will be monitored for VOCs with a portable Ionscience 3000 photoionization detector (PID), or equivalent. All air monitoring data will be documented in a site log book by the designated site safety officer. The site safety officer or delegate must ensure that air monitoring instruments are calibrated and maintained in accordance with manufacturer's specifications. All instruments will be zeroed daily and checked for accuracy. A daily log will be kept. If additional monitoring is required, the protocols will be developed and appended to this plan

3.0 VOC MONITORING, RESPONSE LEVELS, AND ACTIONS

Volatile organic compounds (VOCs) will be monitored at the downwind perimeter of the immediate work area (i.e., the exclusion zone) on a continuous basis or as otherwise specified. Upwind concentrations should be measured at the start of each workday and periodically thereafter to establish background conditions. The monitoring work should be performed using equipment appropriate to measure the types of contaminants known or suspected to be present.

The equipment should be calibrated at least daily for the contaminant(s) of concern or for an appropriate surrogate. The equipment should be capable of calculating 15-minute running average concentrations, which will be compared to the levels specified below.

- If the ambient air concentration of total organic vapors at the downwind perimeter of the work area or exclusion zone exceeds 5 parts per million (ppm) above background for the 15-minute average, work activities must be temporarily halted and monitoring continued. If the total organic vapor level readily decreases (per instantaneous readings) below 5 ppm over background, work activities can resume with continued monitoring.
- If total organic vapor levels at the downwind perimeter of the work area or exclusion zone persist at levels in excess of 5 ppm over background but less than 25 ppm, work activities must be halted, the source of vapors identified, corrective actions taken to abate emissions, and monitoring continued. After these steps, work activities can resume provided that the total organic vapor level 200 feet downwind of the exclusion zone or half the distance to the nearest potential receptor or residential/commercial structure, whichever is less - but in no case less than 20 feet, is below 5 ppm over background for the 15-minute average.
- If the organic vapor level is above 25 ppm at the perimeter of the work area, activities must be shutdown. All 15-minute readings must be recorded and be available for State (DEC and DOH) personnel to review. Instantaneous readings, if any, used for decision purposes should also be recorded.

All readings will be recorded and made available for NYSDEC and NYSDOH personnel to review. If an exceedance of the Action Limits occurs, an Action Limit Report, as shown in Appendix A, will be completed.

3.1 Potential Corrective Measures and VOC Suppression Techniques

If the 15-minute integrated VOC level at the downwind location persists at a concentration that exceeds the upwind level by more than 5 ppm but less than 25 ppm during remediation activities, then vapor suppression techniques will be employed. The following techniques, or others, may be employed to mitigate the generation and migration of fugitive organic vapors:

- limiting the excavation size;
- limiting the drop-height when loading soil into trucks;
- spraying chemical odorants onto the soil;
- covering soil stockpiles with 6-mil plastic sheeting or tarps;
- hauling waste materials in properly tarped containers; and/or
- applying vapor suppressant foam.

4.0 PARTICULATE MONITORING

Air monitoring for particulates (i.e., dust) will be performed continuously during excavation and loading activities using both air monitoring equipment and visual observation at upwind and downwind locations. Monitoring equipment capable of measuring particulate matter smaller than 10 microns (PM₁₀) and capable of integrating (averaging) over periods of 15 minutes or less will be set up at upwind (i.e., background) and downwind locations, at heights approximately four to five feet above land surface (i.e., the breathing zone). Monitoring equipment will be MIE Data Ram monitors, or equivalent. The audible alarm on the particulate monitoring device will be set at 90 micrograms per cubic meter ($\mu\text{g}/\text{m}^3$). This setting will allow proactive evaluation of worksite conditions prior to reaching the action level of 100 $\mu\text{g}/\text{m}^3$ above background. The monitors will be calibrated at least once per day prior to work activities and recalibrated as needed thereafter. In addition, fugitive dust migration will be visually assessed during all intrusive work activities.

The following summarizes particulate action levels and the appropriate responses:

- If the downwind PM-10 particulate level is 100 $\mu\text{g}/\text{m}^3$ greater than background (upwind perimeter) for the 15-minute period, or if airborne dust is observed leaving the work area, then dust suppression techniques must be employed. Work may continue with dust suppression techniques provided that downwind PM-10 particulate levels do not exceed 150 $\mu\text{g}/\text{m}^3$ above the upwind level and provided that no visible dust is migrating from the work area.
- If, after implementation of dust suppression techniques, downwind PM-10 particulate levels are greater than 150 $\mu\text{g}/\text{m}^3$ above the upwind level, work must be stopped and an evaluation of activities initiated. Work can resume provided that dust suppression measures (as described in Section 2.3.1 below) and other controls are successful in reducing the downwind PM-10 particulate concentration to within 150 $\mu\text{g}/\text{m}^3$ of the upwind level and in preventing visible dust migration.

All readings will be recorded and be available for NYSDEC and NYSDOH personnel to review. If an exceedance of the Action Limits occurs, an Action Limit Report as shown in **Appendix A** will be completed.

4.1 Potential Particulate Suppression Techniques

If the integrated particulate level at the downwind location exceeds the upwind level by more than 100 $\mu\text{g}/\text{m}^3$ at any time during remediation activities, then dust suppression techniques will be employed. The following techniques, or others, may be employed to mitigate the generation and migration of fugitive dusts:

- limiting the excavation size;
- spraying water onto the excavation faces and equipment;
- covering soil stockpiles with plastic sheeting or tarps;
- use of gravel paths / roadways;
- hauling waste materials in properly tarped containers; and/or
- limiting vehicle speeds onsite.

Work may continue with dust suppression techniques provided that downwind PM₁₀ levels are not more than 150 µg/m³ greater than the upwind levels.

There may also be situations where the dust is generated by remediation activities and migrates to downwind locations, but is not detected by the monitoring equipment at or above the action level. Therefore, if dust is observed leaving the working area, dust suppression techniques such as those listed above will be employed.

If dust suppression techniques do not lower particulates to below 150 µg/m³, or visible dust persists, work will be suspended until appropriate corrective measures are identified and implemented to remedy the situation.

All air monitoring readings will be recorded in the field logbook and will be available for the NYSDEC and NYSDOH personnel to review.

5.0 DATA QUALITY ASSURANCE

5.1 Calibration

Instrument calibration shall be documented on instrument calibration and maintenance sheets or in the designated field logbook. All instruments shall be calibrated as required by the manufacturer. Calibration checks may be used during the day to confirm instrument accuracy. Duplicate readings may be taken to confirm individual instrument response.

5.2 Operations

All instruments shall be operated in accordance with the manufacturer's specifications. Manufacturers' literature, including an operations manual for each piece of monitoring equipment will be maintained on-site by the SSO for reference.

5.3 Data Review

The SSO will interpret all monitoring data based the established criteria and his/her professional judgment. The SSO shall review the data with the PM to evaluate the potential for worker exposure, upgrades/downgrades in level of protection, comparison to direct reading instrumentation and changes in the integrated monitoring strategy.

Monitoring and sampling data, along with all sample documentation will be periodically reviewed by the PM.

6.0 RECORDS AND REPORTING

All air readings must be recorded on daily air monitoring log sheets and made available for review by personnel from NYSDEC and NYSDOH.

APPENDIX A
ACTION LIMIT REPORT

ATTACHMENT E
Citizen Participation Plan



New York State Department of Environmental Conservation

Brownfield Cleanup Program

Citizen Participation Plan for **FORMER MUGLER SHORING SITE**

**2401 Third Avenue
Bronx, NY 10451**

September 2015

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Note: The information presented in this Citizen Participation Plan was current as of the date of its approval by the New York State Department of Environmental Conservation. Portions of this Citizen Participation Plan may be revised during the site’s investigation and cleanup process.

Applicant: **2401 3rd Ave Associates Property LLC**
Site Name: **Former Mugler Shoring (“Site”)**
Site Address: **2401 Third Avenue**
Site County: **Bronx**
Site Number: **C203052**

1. What is New York’s Brownfield Cleanup Program?

New York’s Brownfield Cleanup Program (BCP) works with private developers to encourage the voluntary cleanup of contaminated properties known as “brownfields” so that they can be reused and developed. These uses include recreation, housing, and business.

A *brownfield* is any real property that is difficult to reuse or redevelop because of the presence or potential presence of contamination. A brownfield typically is a former industrial or commercial property where operations may have resulted in environmental contamination. A brownfield can pose environmental, legal, and financial burdens on a community. If a brownfield is not addressed, it can reduce property values in the area and affect economic development of nearby properties.

The BCP is administered by the New York State Department of Environmental Conservation (NYSDEC) which oversees Applicants that conduct brownfield site investigation and cleanup activities. An Applicant is a person who has requested to participate in the BCP and has been accepted by NYSDEC. The BCP contains investigation and cleanup requirements, ensuring that cleanups protect public health and the environment. When NYSDEC certifies that these requirements have been met, the property can be reused or redeveloped for the intended use.

For more information about the BCP, go online at: <http://www.dec.ny.gov/chemical/8450.html>.

2. Citizen Participation Activities

Why NYSDEC Involves the Public and Why It Is Important

NYSDEC involves the public to improve the process of investigating and cleaning up contaminated sites, and to enable citizens to participate more fully in decisions that affect their health, environment, and social well-being. NYSDEC provides opportunities for citizen involvement and encourages early two-way communication with citizens before decision-makers form or adopt final positions.

Involving citizens affected and interest in site investigation and cleanup programs is important for many reasons. These include:

- Promoting the development of timely, effective site investigation and cleanup programs that protect public health and the environment;
- Improving public access to, and understanding of, issues and information related to a particular site and that Site’s investigation and cleanup process;

- Providing citizens with early and continuing opportunities to participate in NYSDEC's site investigation and cleanup process;
- Ensuring that NYSDEC makes site investigation and cleanup decisions that benefit from input that reflects the interests and perspectives found within the affected community; and
- Encouraging dialogue to promote the exchange of information among the affected/interested public, State agencies, and other interested parties that strengthens trust among the parties, increases understanding of site and community issues and concerns, and improves decision-making.

This Citizen Participation (CP) Plan provides information about how NYSDEC will inform and involve the public during the investigation and cleanup of the Site identified above. The public information and involvement program will be carried out with assistance, as appropriate, from the Applicant.

Project Contacts

Appendix A identifies NYSDEC project contact(s) to whom the public should address questions or request information about the site's investigation and cleanup program. The public's suggestions about this CP Plan and the CP program for the Site are always welcome. Interested people are encouraged to share their ideas and suggestions with the project contacts at any time.

Locations of Reports and Information

The locations of the reports and information related to the Site's investigation and cleanup program also are identified in Appendix A. These locations provide convenient access to important project documents for public review and comment. Some documents may be placed on the NYSDEC website. If this occurs, NYSDEC will inform the public in fact sheets distributed about the Site and by other means, as appropriate.

Site Contact List

Appendix B contains the site contact list. This list has been developed to keep the community informed about, and involved in, the site's investigation and cleanup process. The site contact list will be used periodically to distribute fact sheets that provide updates about the status of the project. These will include notifications of upcoming activities at the Site (such as fieldwork), as well as availability of project documents and announcements about public comment periods.

The site contact list includes, at a minimum:

- Chief executive officer and planning board chairperson of each county, city, town and village in which the Site is located;
- Residents, owners, and occupants of the Site and properties adjacent to the Site;
- The public water supplier which services the area in which the Site is located;
- Any person who has requested to be placed on the site contact list;

- The administrator of any school or day care facility located on or near the Site for purposes of posting and/or dissemination of information at the facility; and
- Location(s) of reports and information.

The site contact list will be reviewed periodically and updated as appropriate. Individuals and organizations will be added to the site contact list upon request. Such requests should be submitted to the NYSDEC project contact(s) identified in Appendix A. Other additions to the site contact list may be made at the discretion of the NYSDEC project manager, in consultation with other NYSDEC staff as appropriate.

CP Activities

The table at the end of this section identifies the CP activities, at a minimum, that have been and will be conducted during the Site’s investigation and cleanup program. The flowchart in Appendix D shows how these CP activities integrate with the site investigation and cleanup process. The public is informed about these CP activities through fact sheets and notices distributed at significant points during the program. Elements of the investigation and cleanup process that match up with the CP activities are explained briefly in Section 5.

- **Notices and fact sheets** help the interested and affected public to understand contamination issues related to a site, and the nature and progress of efforts to investigate and clean up a site.
- **Public forums, comment periods and contact with project managers** provide opportunities for the public to contribute information, opinions and perspectives that have potential to influence decisions about a site’s investigation and cleanup.
- **Document repositories** allow the public to access and review project documents including investigation and cleanup work plans and final reports.

The public is encouraged to contact project staff at any time during the Site’s investigation and cleanup process with questions, comments, or requests for information. This CP Plan may be revised due to changes in major issues of public concern identified in Section 3 or in the nature and scope of investigation and cleanup activities. Modifications may include additions to the site contact list and changes in planned citizen participation activities.

Technical Assistance Grant

NYSDEC must determine if the Site poses a significant threat to public health or the environment. This determination generally is made using information developed during the investigation of the Site, as described in Section 5.

If the Site is determined to be a significant threat, a qualifying community group may apply for a Technical Assistance Grant (TAG). The purpose of a TAG is to provide funds to the qualifying group to obtain independent technical assistance. This assistance helps the TAG recipient to

interpret and understand existing environmental information about the nature and extent of contamination related to the Site and the development/implementation of a remedy.

An eligible community group must certify that its membership represents the interests of the community affected by the Site, and that its members' health, economic well-being or enjoyment of the environment may be affected by a release or threatened release of contamination at the Site.

For more information about TAGs, go online at <http://www.dec.ny.gov/regulations/2590.html>.

Note: The table identifying the citizen participation activities related to the Site's investigation and cleanup program follows on the next page:

| Citizen Participation Requirements (Activities) | Timing of CP Activity(ies) |
|--|---|
| Application Process: | |
| <ul style="list-style-type: none"> • Prepare site contact list • Establish document repositories | At time of preparation of application to participate in the BCP. |
| <ul style="list-style-type: none"> • Publish notice in Environmental Notice Bulletin (ENB) announcing receipt of application and 30-day public comment period • Publish above ENB content in local newspaper • Mail above ENB content to site contact list • Conduct 30-day public comment period | When NYSDEC determines that BCP application is complete. The 30-day public comment period begins on date of publication of notice in ENB. End date of public comment period is as stated in ENB notice. Therefore, ENB notice, newspaper notice, and notice to the site contact list should be provided to the public at the same time. |
| After Execution of Brownfield Site Cleanup Agreement: | |
| <ul style="list-style-type: none"> • Prepare Citizen Participation (CP) Plan | Before start of Remedial Investigation |
| Before NYSDEC Approves Remedial Investigation (RI) Work Plan: | |
| <ul style="list-style-type: none"> • Distribute fact sheet to site contact list about proposed RI activities and announcing 30-day public comment period about draft RI Work Plan • Conduct 30-day public comment period | Before NYSDEC approves RI Work Plan. If RI Work Plan is submitted with application, public comment periods will be combined and public notice will include fact sheet. Thirty-day public comment period begins/ends as per dates identified in fact sheet. |
| After Applicant Completes Remedial Investigation: | |
| <ul style="list-style-type: none"> • Distribute fact sheet to site contact list that describes RI results | Before NYSDEC approves RI Report |
| Before NYSDEC Approves Remedial Work Plan (RWP): | |
| <ul style="list-style-type: none"> • Distribute fact sheet to site contact list about proposed RWP and announcing 45-day public comment period • Public meeting by NYSDEC about proposed RWP (if requested by affected community or at discretion of NYSDEC project manager) • Conduct 45-day public comment period | Before NYSDEC approves RWP. Forty-five day public comment period begins/ends as per dates identified in fact sheet. Public meeting would be held within the 45-day public comment period. |
| Before Applicant Starts Cleanup Action: | |
| <ul style="list-style-type: none"> • Distribute fact sheet to site contact list that describes upcoming cleanup action | Before the start of cleanup action. |
| After Applicant Completes Cleanup Action: | |
| <ul style="list-style-type: none"> • Distribute fact sheet to site contact list that announces that cleanup action has been completed and that summarizes the Final Engineering Report • Distribute fact sheet to site contact list announcing issuance of Certificate of Completion (COC) | At the time NYSDEC approves Final Engineering Report. These two fact sheets are combined if possible if there is not a delay in issuing the COC. |

3. Major Issues of Public Concern

This section of the CP Plan identifies major issues of public concern that relate to the Site. Additional major issues of public concern may be identified during the course of the Site's investigation and cleanup process.

The Site is located in an Environmental Justice Area. Environmental justice is defined as the fair treatment and meaningful involvement of all people regardless of race, color, national origin, or income with respect to the development, implementation, and enforcement of environmental laws, regulations, and policies.

The Department has received comments regarding the need to perform a Fish & Wildlife Impact Analysis (FWIA) due to the site's close proximity to the Harlem River, and that the Department is in agreement with the issues raised and a FWIA will be performed. The Site is located in an area with a large Hispanic-American population nearby. Therefore, all future fact sheets will be translated into Spanish.

The major issues of concern to the public will be potential impacts of nuisance odors and dust during the removal of affected soil at the Site. Another example of a major issue of public concern would be the impact of increased truck traffic on the surrounding neighborhood. Construction safety issues will also be addressed.

This work will be performed in accordance with procedures which will be specified under a detailed Remedial Program which considers and takes preventive measures for exposures to future residents of the property and those on adjacent properties during construction. Detailed plans to monitor the potential for exposure including a Health and Safety Plan (HASP) and a Community Air Monitoring Plan (CAMP) are required components of the remedial program. Implementation of these plans will be under the direct oversight of the NYSDEC and the New York State Department of Health (NYSDOH).

These plans will specify the following worker and community health and safety activities during remedial activity at the Site:

- On-site air monitoring for worker protection;
- Perimeter air monitoring for community protection;
- The use of odor, vapor, and dust controls, such as water or foam sprays, as needed;
- Monitoring and control of soil, sediments, and water generated during remediation; and
- Truck routes which avoid residential streets.

The HASP and the CAMP will be prepared as part of the Remedial Action Work Plan (RAWP) and will be available for public review at the document repository as identified in Appendix A (page 11).

Furthermore, the Applicant has prepared a Scoping Sheet for Major Issues of Public Concern which will assist them in identifying any concerns. Experience from similar projects, 311 complaints and other construction projects in the area will help in identifying such issues.

4. Site Information

Appendix C contains a map identifying the location of the Site.

Site Description

The Site to be remediated and redeveloped is located in the South Bronx (Bronx County) and is comprised of a single tax parcel covering 61,034.98 (1.4 acres). The subject property is located in the City of New York and Borough of the Bronx (Bronx County). The lot is located on the west side of Third Avenue and is identified as Block 2319 Lot 2 on the NY City tax map. The property has 159 feet of street frontage on Third Avenue and approximately 346 ft of frontage along the north side of the Harlem River. The lot is developed with a one-story 19,450 sf commercial building which according to the NYC Department of Buildings was constructed in 1931.

The land use in the immediate vicinity of the Site includes underutilized or vacant, commercial properties to the north and east, the Third Avenue Bridge to the southeast and the Harlem River to the southwest.

The area surrounding the property is highly urbanized and predominantly consists of heavy commercial / industrial / warehouse properties to the north along a corridor adjacent to the Harlem River. Multi-use residential / commercial (retail) properties are present to the east along Bruckner Boulevard and a large housing project is located to the northeast.

The Lot is currently zoned M1-3/R8 residential / commercial. The M1-3/R8 is a Special Mixed Use District established to encourage investment in, and enhance the vitality of, existing neighborhoods with mixed residential and industrial uses in close proximity and create expanded opportunities for new mixed use communities. New residential and non-residential uses (commercial, community facility and light industrial) can be developed as-of-right and be located side-by-side or within the same building. Pairing an M1 district with an R3 through R10 district ensures a balanced variety of uses.

History of Site Use, Investigation, and Cleanup

The Site is currently owned by the 2401 3rd Ave Associates Property LLC. The property consists of a 1-story commercial / industrial building which is currently vacant. The Requestor purchased the property in January 2015. The property was most recently occupied by a company.

The Subject Property was historically used for manufacturing purposes since at least 1891. Historical operators include; J.L. Mott Iron Works (1891–1922), Hydraulic Steel Company (1922–1935), General Builders Supply Corporation (1935–1968), Brill Equipment Company (1949–1956), US Gear Manufacturing Company (1965–1971), Ohio Gasket Manufacturing Corporation (1971), and Mugler Inc. (1965–2015).

A Phase II investigation performed at the Site in October 2014 historic fill materials and petroleum contamination. The primary contaminants identified included petroleum volatile

organic compounds (VOCs) in soil at a single location, toluene in the vicinity of a suspect 550 gallon underground storage tank. As a result, a spill was reported to the New York State Department of Environmental Conservation (NYSDEC). The NYSDEC assigned spill number 1405230 to the case.

The depth to groundwater at the site is approximately 13 feet below grade. Soil at the site is described as historic fill materials to a depth which ranges from approximately 5 feet to 13 feet below the surface followed by native brown sand with cobble size rocks.

5. Investigation and Cleanup Process

Application

The Applicant has applied for and been accepted into New York's Brownfield Cleanup Program (BCP) as a Volunteer. This means that the Applicant was not responsible for the disposal or discharge of the contaminants or whose ownership or operation of the Site took place after the discharge or disposal of contaminants. The Volunteer must fully characterize the nature and extent of contamination on-site, and must conduct a qualitative exposure assessment, a process that characterizes the actual or potential exposures of people, fish and wildlife to contaminants on the Site and to contamination that has migrated from the Site.

The Applicant proposes that the Site will be used for the construction of two new 25-story residential tower buildings identified as the east and the west towers. The two towers flank an entry court which is open to the river and Manhattan views to the south. The East Tower is free-standing, serving as a "portal" on Third Avenue, whereas the West Tower has a 7-story base with common amenities and parking for 200 cars. The residential component consists of 475 rental units in an area of 399,798 gross square feet. Commercial (retail) and community space will be included in the west tower base. The proposed use is consistent with existing M1-3/R8 zoning of the property.

To achieve this goal, the Applicant will conduct investigation and cleanup activities at the Site with oversight provided by NYSDEC. The Brownfield Cleanup Agreement (BCA) executed by NYSDEC and the Applicant sets forth the responsibilities of each party in conducting these activities at the Site.

Investigation

The Applicant completed a partial site investigation before it entered into the BCP. For the partial investigation, NYSDEC will determine if the data are useable. The Applicant will now conduct an investigation of the Site officially called a "remedial investigation" (RI). This investigation will be performed with NYSDEC oversight. Upon receipt of the RI, the NYSDEC will determine if the investigation goals and requirements of the BCP have been met or if additional work is needed before a remedy can be selected.

The site investigation has several goals:

- 1) Define the nature and extent of contamination in soil, surface water, groundwater and any

- other parts of the environment that may be affected;
- 2) Identify the source(s) of the contamination;
 - 3) Assess the impact of the contamination on public health and the environment; and
 - 4) Provide information to support the development of a proposed remedy to address the contamination or the determination that cleanup is not necessary.

When the investigation is complete, the Applicant will prepare and submit a report that summarizes the results. This report also will recommend whether cleanup action is needed to address site-related contamination. The investigation report is subject to review and approval by NYSDEC.

NYSDEC will use the information in the investigation report to determine if the Site poses a significant threat to public health or the environment. If the Site is a significant threat, it must be cleaned up using a remedy selected by NYSDEC from an analysis of alternatives prepared by the Applicant and approved by NYSDEC. If the Site does not pose a significant threat, the Applicant may select the remedy from the approved analysis of alternatives.

Remedy Selection

When the investigation of the Site has been determined to be complete, the project likely would proceed in one of two directions:

1. The Applicant may recommend in its investigation report that no action is necessary at the Site. In this case, NYSDEC would make the investigation report available for public comment for 45 days. NYSDEC then would complete its review, make any necessary revisions, and, if appropriate, approve the investigation report. NYSDEC would then issue a Certificate of Completion (COC) (described below) to the Applicant.

or

2. The Applicant may recommend in its investigation report that action needs to be taken to address site contamination. After NYSDEC approves the investigation report, the Applicant may then develop a cleanup plan, officially called a Remedial Work Plan. The Remedial Work Plan describes the Applicant's proposed remedy for addressing contamination related to the Site.

When the Applicant submits a proposed Remedial Work Plan for approval, NYSDEC would announce the availability of the proposed plan for public review during a 45-day public comment period.

Cleanup Action

NYSDEC will consider public comments, and revise the draft cleanup plan if necessary, before approving the proposed remedy. The New York State Department of Health (NYSDOH) must concur with the proposed remedy. After approval, the proposed remedy becomes the selected remedy.

The Applicant may then design and perform the cleanup action to address the site contamination. NYSDEC and NYSDOH oversee the activities. When the Applicant completes cleanup activities, it will prepare a Final Engineering Report (FER) that certifies that cleanup requirements have been achieved or will be achieved within a specific time frame. NYSDEC will review the report to be certain that the cleanup is protective of public health and the environment for the intended use of the Site.

Certificate of Completion

When NYSDEC is satisfied that cleanup requirements have been achieved or will be achieved for the Site, it will approve the FER. NYSDEC then will issue a COC to the Applicant. The COC states that cleanup goals have been achieved, and relieves the Applicant from future liability for site-related contamination, subject to certain conditions. The Applicant would be eligible to redevelop the Site after it receives a COC.

Site Management

Site management is the last phase of the site cleanup program. This phase begins when the COC is issued. Site management may be conducted by the Applicant under NYSDEC oversight, if contamination will remain in place. Site management incorporates any institutional and engineering controls required to ensure that the remedy implemented for the Site remains protective of public health and the environment. All significant activities are detailed in a Site Management Plan (SMP).

An institutional control is a non-physical restriction on use of the Site, such as a deed restriction that would prevent or restrict certain uses of the property. An institutional control may be used when the cleanup action leaves some contamination that makes the Site suitable for some, but not all uses.

An engineering control is a physical barrier or method to manage contamination. Examples include: caps, covers, barriers, fences, and treatment of water supplies.

Site management also may include the operation and maintenance of a component of the remedy, such as a system that is pumping and treating groundwater. Site management continues until NYSDEC determines that it is no longer needed.

Appendix A

Project Contacts and Locations of Reports and Information

Project Contacts

For information about the site's investigation and cleanup program, the public may contact any of the following project staff:

New York State Department of Environmental Conservation (NYSDEC):

Sadique Ahmed P.E.
New York State Department of
Environmental Conservation
625 Broadway
Albany, New York 12233-7016
Tel: (518) 402-9656
Email: saduque.ahmed@dec.ny.gov

Thomas Panzone
Regional Citizen Participation Specialist
NYSDEC Region 2
Office of Communications Services
47-40 21st Street
Long Island City, NY 11101-5407
Tel: (718) 482-4953
Email: thomas.panzone@dec.ny.gov

New York State Department of Health (NYSDOH):

Steven Karpinski
New York State Department of Health
Bureau of Environmental Exposure Investigation
Empire State Plaza – Corning Tower Room 1787
Albany, New York 12237
Tel: (518) 402-7860
Email: beei@health.ny.gov

Locations of Reports and Information

The facilities identified below are being used to provide the public with convenient access to important project documents:

New York Public Library - Mott Haven Branch

321 East 140th Street
Bronx, NY 10454
(718) 665-4878

Hours:

Sunday: Closed

Monday: 10am- 6pm

Tuesday: 10am- 7pm

Wednesday & Thursday: 10am- 7pm

Friday: 10am- 5pm

Saturday: 10am – 5pm

Appendix B - Site Contact List

Local Government Contacts:

City of New York

Hon. Bill de Blasio
Mayor of New York City
City Hall
New York, NY 10007

Hon. Ruben Diaz, Jr.
Bronx Borough President
851 Grand Concourse, Suite 301
Bronx, New York 10451
Telephone (718) 590-3500

George Rodriguez
Chair, Bronx Community Board 1
3024 Third Avenue
Bronx, NY 10455
TEL: (718) 585-7117
FAX: (718) 292-0558

Cedric Loftin
District Manager, Bronx Community Board 1
3024 Third Avenue
Bronx, NY 10455
TEL: (718) 585-7117
FAX: (718) 292-0558

Hon. Melissa Mark - Viverito
New York City Council Speaker – District 8
105 East 116th Street
New York, NY 10029
Telephone (212) 828-9800
Fax (212) 722-6378

Carl Weisbrod
Chair of City Planning (Zoning)
22 Reade St.
Third Floor
New York, NY 10007

Carol Samol
Director, NYC Planning Commission – Bronx Office
1 Fordham Plz.
Bronx, New York 10458
Telephone (718) 220-8500

Constance Moran
New York City Department of Transportation
Bronx Borough Commissioner
55 Water Street, 9th Floor
New York, NY 10041
212-748-6680

Bronx County Clerk's Office
Luis M. Diaz, County Clerk
851 Grand Concourse, Room 118
Bronx, New York 10451
Telephone (866) 797-7214

Hon. Letitia James
Public Advocate
1 Centre Street, 15th Floor
New York, NY 10007

Hon. Scott M. Stringer
Office of the Comptroller
1 Centre Street
New York, NY 10007

Julie Stein
Office of Environmental Planning & Assessment
NYC Dept. of Environmental Protection
96-05 Horace Harding Expressway
Flushing, NY 11373

Daniel Walsh
NYC Office of Environmental Remediation
100 Gold Street – 2nd Floor
New York, NY 10038

Nilda Mesa, Director
NYC Office of Environmental Sustainability
100 Gold Street – 2nd Floor
New York, NY 10038

New York State

Senator José M. Serrano
1916 Park Avenue Suite 202,
New York, NY 10037
Telephone (212) 828-5829
Fax (212) 828-2420

Hon.. Carmen E. Arroyo
NYS Assemblymember
384 East 149th Street, Suite 301
Bronx, New York 10455
Telephone (718) 292-2901

Federal

Hon. Charles Schumer
U.S. Senator
780 Third Avenue, Suite 2301
New York, NY 10017

Hon. Kirsten Gillibrand
U.S. Senator
780 Third Avenue, Suite 2601
New York, NY 10017

Rep. José E. Serrano
Congressional District: 15
1231 Lafayette Avenue, 4th Floor
Bronx, New York 10474
Telephone (718) 620-0084
Fax (718) 620-0658

Contact information for the identified owners, as listed in the New York City ACRIS Database,
are as follows:

North

1. Owner / Tenant
CUBESMART EAST 135TH, LLC
PTA
P.O. BOX 320099
ALEXANDRIA, VA 22320-4099

2. Owner
GLS REAL ESTATE CO.,
2413 3RD AVE.
BRONX, NY 10451-6330

Occupant / Tenant
2413 3RD AVE.
BRONX, NY 10451-6330

3. Owner
KAI DEVELOPMENT CORP.
211 W. 58TH ST. APT. 1
NEW YORK, NY 10019-1418

Occupant / Tenant
2403 3RD AVE
BRONX, NY 10451-6330

4. Owner / Tenant
FIVE BORO STORAGE INC.
220 E. 134TH ST. APT. FRNT A.
BRONX, NY 10451-6410

5. Owner
MADHATTERS REALTY INC.
2417 3RD AVE.
BRONX, NY 10451-6331

Occupant / Tenant
2417 3RD AVE
BRONX, NY 10451-6331

East

6. Owner
101 LINCOLN ASSOCIATES PROPERTY LLC
C/O THE CHETRIT GROUP LLC
512 SEVENTH AVENUE APT. 15TH FLOOR
NEW YORK, NY 10018

Occupant / Tenant
101 LINCOLN AVENUE
BRONX, NY 10454

Local News Media

Bronx Times
900 E. 132nd Street
Bronx, NY 10454
(718) 597-1116

New York Daily News
4 New York Plaza
New York, NY 10004

New York Post
1211 Avenue of the Americas
New York, NY 10036-8790

NY 1 News
75 Ninth Avenue
New York, NY 10011

Hoy Nueva York
1 MetroTech Center, 18th Floor
Brooklyn, NY 11201

El Diario La Prensa
1 MetroTech Center, 18th Floor
Brooklyn, NY 11201

Impacto New York
225 West 35th Street, Suite 305
New York, NY 10001

La Voz Hispana NY
159 East 116th Street
New York, NY 10029

Public Water Supplier

Hon. Emily Lloyd, Commissioner
New York City Department of Environmental Protection
59-17 Junction Boulevard
Flushing, NY 11373

Requested Contacts

No requests have been made at this time.

Schools and Daycare Facilities

There are no Schools or Daycare facilities within a quarter mile of the project Site. Schools nearest the Site include:

- 1 City of New York: PS 154 Jonathan D Hyatt
333 East 135th Street, Bronx, NY 10454
Principal: Dr. Alison Coviello
(718) 292-4742

Community, Civic, Religious and other Educational Institutions

Bronx Terminal Market
Attn: Executive Director
610 Exterior Street
Bronx, NY 10451

Mitchell Houses
Management Development Office
NYCHA
303 East 135th Street
Bronx, NY 10454

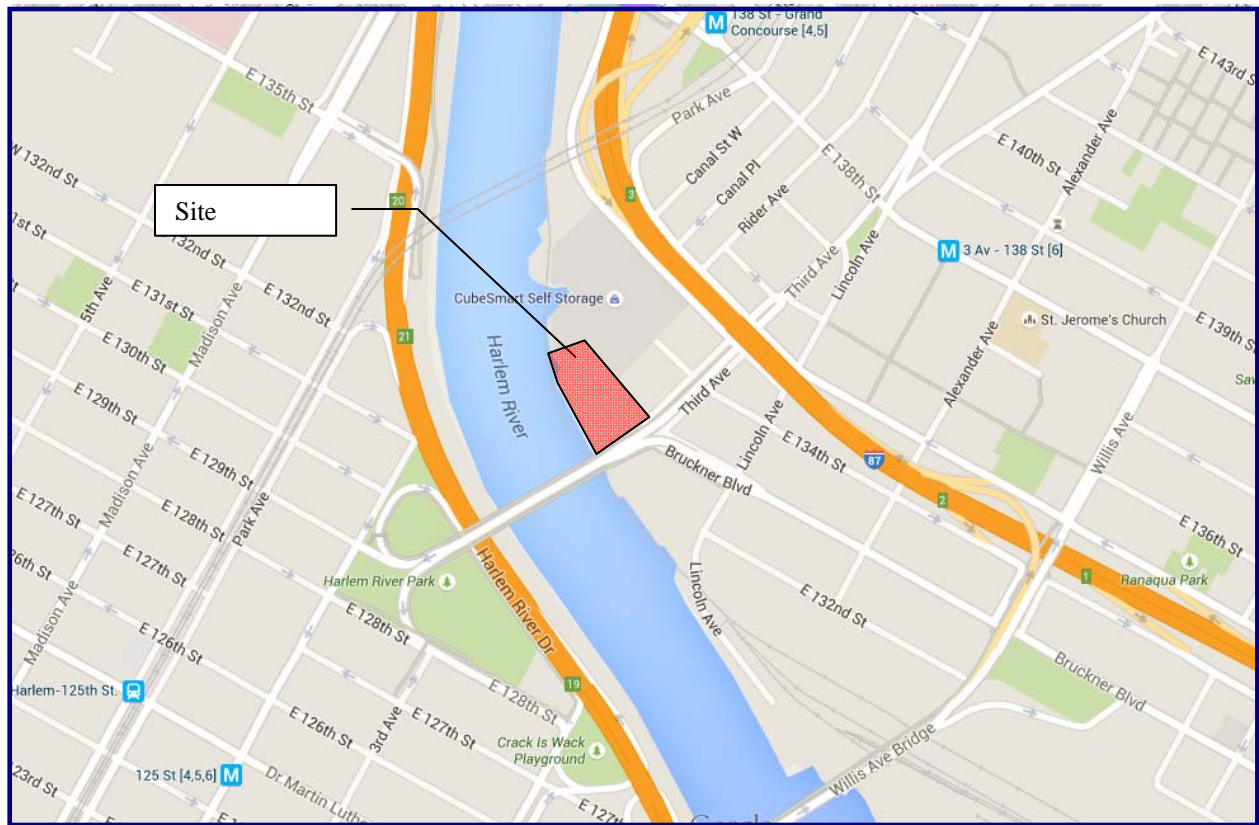
Mitchell Houses
Attn: President, Resident Association
NYCHA
303 East 135th Street
Bronx, NY 10454

Office of Media Relations
NYCHA
250 Broadway
New York, NY 10007

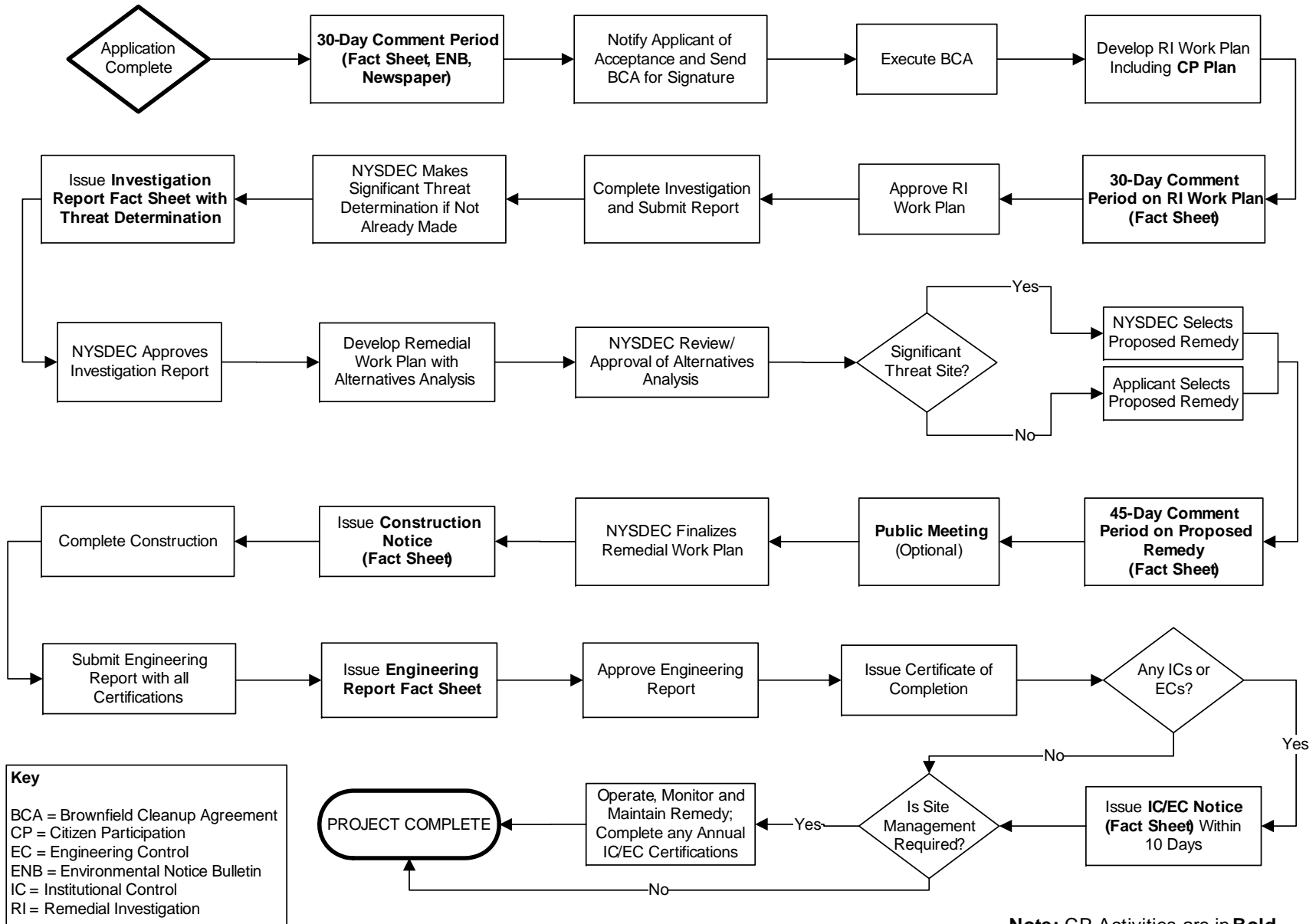
Mitchell Community Center
Attn: Executive Director
210 Alexander Avenue
Bronx, NY 10454

St. Jerome's RC Church
Attn: Pastor
230 Alexander Avenue
Bronx, NY 10454

Appendix C - Site Location Map



Appendix D– Brownfield Cleanup Program Process



ATTACHMENT F
Resumes



ENVIRONMENTAL BUSINESS CONSULTANTS

Charles B. Sosik, PG, PHG, Principal

Professional Experience

24 years

Education

MS, Hydrogeology, Adelphi University, NY
BS, Geology, Northern Arizona University, AZ

Areas of Expertise

- Brownfields Redevelopment
- Hazardous Waste Site Investigations
- Pre-purchase Site Evaluations and Support
- Regulatory Negotiations
- Remedial Planning and "Cost to Cure" Analysis
- Strategic Planning
- Real Estate Transactions
- NYC "E" Designations

Professional Certification

- Professional Geologist, NH
- Professional Geologist, Hydrogeologist, WA
- OSHA 40-hr HAZMAT
- OSHA 8-hr. Supervisor

Professional Affiliation / Committees

- NYS Council of Professional Geologists (NYSCPG)
 - Association of Groundwater Scientists & Engineers (AGSE)
 - NYS RBCA Advisory Committee
 - Massachusetts LSP Association
 - New Hampshire Association of Professional Geologists
 - Interstate Technology Regulatory Council/MTBE Team
 - Environmental Business Association, Brownfields Task Force
 - Part 375 Working Group
-

PROFILE

Mr. Sosik has 24 years of experience in environmental consulting. He specializes in advising clients on managing environmental compliance with federal, state, and municipal agencies and has successfully directed numerous investigation and remediation projects involving petroleum, pesticides, chlorinated solvents, heavy metals and radiologically activated media. His work included extensive three-dimensional investigations on MTBE, which have been used effectively to help shape public policy. He also has experience in applying models to groundwater related problems and has completed several large-scale projects to determine fate and transport of contaminants, establish spill scenarios, and closure criteria. His experience and expertise in the area of contaminant hydrogeology has resulted in requests from environmental attorneys, property owners and New York State to serve as an expert witness and technical advisor on a variety of legal disputes.

For the past 10 years Mr. Sosik has been primarily engaged in providing environmental consulting to developers responding to the extensive rezoning of former industrial and commercial properties, which is currently taking place throughout New York City. These services include everything from pre-purchase evaluations and contract negotiations to gaining acceptance in and moving projects through the NYS Brownfields Program. Mr. Sosik has taken a pro-active role in the continued development of the NYS Brownfields Program and related policy, by attending numerous working seminars, active participation in work groups and task forces and by providing commentary to draft versions of new guidance documents. Throughout his professional career, Mr. Sosik has remained committed to developing innovative cost- efficient solutions to environmental issues, specifically tailored to the needs of his clients.

SELECTED PROJECTS

Scavenger Waste Treatment Facility (SWTF), Suffolk County, NY

Water Treatment Plant EIS - Focused EIS - In response to requests from the Suffolk County Council on Environmental Quality and the Brookhaven Conservation Advisory Council, Mr. Sosik prepared a focused EIS to evaluate the potential impacts to an important surface water resource from the proposed facility including cumulative and synergistic effects with established contaminant plumes in the area.

Advanced Residential Communities, Rockville Centre, NY

Brownfield Project – As the senior project manager on this large scale, high profile redevelopment project, Mr. Sosik was asked to develop a plan to accelerate the regulatory process in the face of general community opposition. Through numerous discussions with the BCP management team, He was able to condense the schedule and review period, through the submission of supporting documents (Investigation Report, Remedial Work Plan) with the BCP application package. Community opposition, which focused on the environmental condition of the site as a means to block the project, was used to advantage in expediting approval of the aggressive interim remedial

plan. This will allow the developer to begin remedial work approximately 5 months ahead of schedule.

Former Temco Uniform site, West Haverstraw, NY

Brownfield Project – Mr. Sosik took over management of this project from another consultant following transition of this VCP site to the BCP. Mr. Sosik used the opportunity to renegotiate and revise the scope of work to allow a more cost effective and focused investigation plan without re-writing or resubmitting the RIWP. During the NYSDEC's review of the transition package, he met with and coordinated changes with the NYSDEC Project Manager to gain approval. The result saved the client a significant amount of money, but perhaps more importantly in this case, did so without loss of time.

Grovick Properties, Jackson Heights, NY

Brownfield Project – This Brownfield property is somewhat unique in that it had been investigated and partially remediated by the NYSDEC through the petroleum spill fund. The client was interested in purchasing the property and redeveloping it as office and retail space. Mr. Sosik reviewed the NYSDEC investigation and developed a



Charles B. Sosik, PG, PHG, Principal

supplemental plan to meet the requirements of an RI under the BCP program. By performing this limited amount of field work "up-front" he was able to complete an RI Report and Remedial Plan and submit both with the BCP application package. The NYSDEC and NYSDOH approved the RI Report and the Remedial Plan with minor changes. This cut 120 days from the review process and allowed the client to arrange financing and move his project forward knowing what the clean-up costs would be at the outset.

Metro Management, Bronx, NY

Brownfield Project – The site of a former gas station, the developer had planned to construct a 12-story affordable housing apartment complex with first floor retail space. Since the site was located in an Environmental zone, potential tax credits of 22% for site development, remediation and tangible property could be realized under the BCP. In a pre-application meeting with the NYSDEC, Mr. Sosik realized that the department did not believe the site was eligible for the BCP, since it had been previously investigated and closed under the spills program.

Mr. Sosik assisted the developer in securing financing, and due to the demands of an aggressive construction schedule developed an Interim Remedial Measure (IRM), based on chemical oxidation treatment. Working closely with the clients environmental counsel, Mr. Sosik was able to get the IRM approved without a public comment period. Implementation of the IRM is currently underway.

The project was awarded the 2009 NYC Brownfield Award for Innovation.

Brandt Airflex, NY

Technical Consulting Services - Mr. Sosik provided senior level technical advice and strategic planning in developing an off-site RI/FS for the site, in negotiating a tax reduction for the property due to the environmental condition and in preparing a cost to cure estimate for settlement between business partners. After achieving a favorable tax consideration and settlement agreement for his client

Allied Aviation Services, Dallas, Fort Worth, Airport, Dallas, TX

Jet Fuel Investigation - Mr. Sosik developed and managed an investigative plan to quickly identify the extent and source of jet fuel which was discharging from the Airport's storm drain system to a creek a mile away. Through the use of a refined conceptual model, accelerated investigative techniques and a flexible work plan, he was able to identify the source of the fuel and the migration route within a single week. He then identified remedial options and successfully negotiated a risk based plan with the Texas regulatory agency that had issued a notice of enforcement action against the facility.

KeySpan – Former LILCO Facilities, Various NY Locations

Pesticide Impact Evaluation - Mr. Sosik developed, negotiated and implemented a site screening procedure to evaluate impact to public health and the environment as the result of past herbicide use at 211 utility sites. Using an unsaturated zone leaching model (PRZM) on a small subset of the sites, he was able to establish mass loading schedules for the remaining sites. This was combined with public well data in a GIS environment to perform queries with respect to mass

loading, time transport and proximity to vulnerable public supply wells. Using this approach Mr. Sosik was able to show that there were no concerns for future impact. This effort satisfied the public health and resource concerns of the state environmental agency and county health department in a reasonable amount of time and at a fraction of the cost of a full scale investigation.

Former Computer Circuits (Superfund) Site, Hauppauge, NY

CERCLA RI/FS - As Senior Project Manager for the site, he played a major role in regaining control of the investigation activities for the PRP. This action prevented the USEPA from initiating an extensive investigation at the site using a RAC II contractor allowing the client to perform a more efficient investigation. He was involved in all negotiations with EPA and was the project lead in developing a revised site characterization plan (work plan, field sampling plan, quality assurance plan, etc.). By carefully managing all phases of the investigation and continued interaction with each of the three regulatory agencies involved, Mr. Sosik was able to keep the project focused and incrementally reinforce the clients position. The estimated cost of the revised investigation is expected to save the client 1.5 to 2 million dollars.

Sun Oil, Seaford, NY

Remediation Consulting Services & Project Management - Under an atmosphere of regulatory distrust, political pressure and mounting public hostility toward the client, Mr. Sosik conducted an off-site 3-D investigation to define the extent of contamination and the potential impact on public health. By designing and implementing an aggressive source area remediation program and personal interaction with the public and regulatory agencies, he was able to successfully negotiate a limited off-site remediation favorable to the client. Source area remediation was completed within 6 months and the project successfully closed without damage to the client's public image or working relationship with the regulatory agencies.

Con Edison, Various Locations, NY

Hydrogeologic Consulting Services - Under a general consulting contract, Mr. Sosik conducted detailed subsurface hydrogeologic investigations at five locations to assist in the development of groundwater contingency planning. He also developed and implemented work plans to investigate and remediate existing petroleum, cable fluid, and PCB releases at many of the generating facilities and substations. An important aspect of his role was in assisting the client in strategic planning and negotiations with the regulatory agency.

Keyspan - Tuthill Substation, Aqueboque, NY

Accelerated Site Characterization - Using accelerated site characterization techniques, Mr. Sosik presented the project as a case study in establishing the transport of an herbicide and its metabolites applied at utility sites in the 1980's. The results were then used to establish a screening method for evaluating 211 similar sites controlled by the client in a reasonable and efficient manner.

NYSDEC Spill, East Moriches, NY

Spill Release Analysis - With recognized expertise in the area of gasoline plume development on Long Island, Mr. Sosik was asked by



Charles B. Sosik, PG, PHG, Principal

the State to establish the release date (and principal responsible party) of an extensive petroleum spill, which impacted a residential neighborhood. He used multiple lines of evidence, and a new EPA model (HSSM), which he has helped to refine, to reconstruct the release scenario and spill date, in support of the State Attorney General's cost recovery effort from the PRP.

Minmilt Realty, Farmingdale, NY

Fate & Transport Modeling - He completed an RI/FS at this location for a PCE plume that had been in transit for over 30 years. Mr. Sosik applied a conservative model to evaluate time/concentration impacts under a variety of transport scenarios to a municipal wellfield located 13,000 feet away. Through the use of the model and careful interpretation of an extensive data set compiled from several sources, Mr. Sosik was able to propose a plan which was both acceptable to the regulator and favorable to the client.

Sebonack Golf Course Project, Town of Southampton, NY

IPM Pesticide Study - Provided professional hydrogeologic services in support of the EIS prepared for the development of the site. The proposed development included an 18-hole golf course, clubhouse, dormitory facility, cottages, associated structures, and a 6,000 square foot research station for Southampton College. Mr. Sosik performed an extensive evaluation (using a pesticide-leaching model) on the effects of pesticide and nitrogen loading to groundwater as part of the projects commitment to an Integrated Pest Management (IPM) approach.

NYSDEC, Spills Division, Regions 1 - 4

Petroleum Spills Investigation & Remediation - As a prime contractor/consultant for the NYSDEC in Regions 1-4, Mr. Sosik has managed the investigation and remediation of numerous petroleum spills throughout the State. Many of these projects required the development of innovative investigation and remediation techniques to achieve project goals. He was also involved in many pilot projects and research studies to evaluate innovative investigation techniques such as accelerated site characterization, and alternative approaches to remediation such as monitored natural attenuation and risk based corrective action.

Sun Oil, E. Meadow, NY

Exposure Assessment - Performed to seek closure of the spill file, despite the presence of contaminants above standards, Mr. Sosik determined after the extended assessment that the level of remaining contamination would not pose a future threat to human health or the environment. He used multiple lines of evidence, and a fate and

transport model to show that degradation processes would achieve standards within a reasonable time.

Sand & Gravel Mine, NY

Property Development - As part of the development of a sand and gravel mine, Mr. Sosik provided environmental consulting services to assist in obtaining a mining permit, which would result in the construction of a 150-acre lake. Specifically, Mr. Sosik investigated if the proposed lake would reduce groundwater quantity to domestic and public well fields, and/or accelerate the migration of potential surface contaminants to the lower part of the aquifer. After assuming the lead role in negotiations with the regulatory agency, Mr. Sosik was able to obtain a permit for the client by adequately addressing water quality and quantity issues, and by preparing a monitoring plan and spill response plan, acceptable to all parties.

NYSDEC, Mamaroneck, NY

Site Characterization / Source Identification - In a complex hydrogeologic setting consisting of contaminant transport through fractured metamorphic bedrock and variable overburden materials, Mr. Sosik was able to develop and implement a sub-surface investigation to differentiate and separate the impact associated with each of two sources. The results of this investigation were successful in encouraging the spiller to accept responsibility for the release.

Riverhead Municipal Water District, NY

Site Characterization / Remedial Planning - Using accelerated characterization techniques, he implemented a 3-D site investigation to identify two service stations 4,000 ft. away as the source of contamination impacting a municipal wellfield. In accordance with the strict time table imposed by the need to return the wellfield to production by early spring, he designed and implemented a multi-point (9 RW, 6 IW) recovery and injection well system using a 3-d numerical flow model, and completed the project on time. Using a contaminant transport model, Mr. Sosik developed clean-up goals which were achieved in 9 months of operation, well below the projected 3 to 5 year project duration.

Montauk Fire Department, NY

Site Assessment - Mr. Sosik performed a limited investigation and used a 2-D flow model to demonstrate that the property could not have been the source of contamination which had impacted an adjacent wellfield as per the results of a previous investigation. This small focused effort successfully reversed a \$500,000, and rising, claim against the department by the water district and the NYSDEC.

PREVIOUS EXPERIENCE

P.W. Grosser Consulting, Bohemia, NY

Senior Project Manager, 1999-2006

Environmental Assessment & Remediation, Patchogue, NY

Senior Project Manager, 1994-1999

Miller Environmental Group, Calverton, NY

Project Manager, 1989-1994

DuPont Biosystems, Aston, PA

Hydrogeologist, 1988-1989



Charles B. Sosik, PG, PHG, Principal

EXPERT WITNESS TESTIMONY AND DEPOSITIONS

Fact Witness -Testimony on relative age of petroleum spill based on nature and extent of residual and dissolved components at the Delta Service Station in Uniondale, NY Fall/1999

Expert Witness / Expert Report for defendant in cost recovery case by NYS Attorney General regarding a Class II Inactive Hazardous Waste (State Superfund) project by the NYSDEC (October 2004 – present, Report: March 2005, Deposition: April 2005)

Expert Witness / Fact Witness for plaintiff seeking compensation for partial expenses incurred during the investigation and remediation of a USEPA CERCLA site due to the release and migration of contaminants from an "upgradient" industrial property. (Deposition May 2005, case settled April 2007).

Expert Witness / Fact Witness for NYS Attorney General with respect to cost recovery for a NYSDEC petroleum spill site in Holtzville, NY (Deposition April 2005 - case settled).

Expert Witness – Statement of opinion and expert testimony at trial for plaintiff seeking damages from a major oil corporation for contamination under a prior leasing agreement in Rego Park, NY. Case decided in favor of plaintiff. Trial July 2007, in favor of Plaintiff. Qualified as Expert Witness.

Expert Witness / Fact Witness for NYS Attorney General with respect to cost recovery for a NYSDEC petroleum spill site in Lindenhurst, NY (Trial date December 2009, in favor of plaintiff. Qualified as Expert Witness.

Expert Witness / Fact Witness for defendant with respect to cost recovery and third party responsibility for a NYSDEC petroleum spill site. (Expert Statement of Fact – October 2005).

Expert Witness for plaintiff seeking damages related to a petroleum spill from the previous owner/operator of a gas station in College Point, NY. Case settled 2009.

Expert Witness for plaintiff (municipal water supply purveyor) seeking damages from major oil companies and manufacturer of MTBE at various locations in Suffolk County, NY. Expert reports July 2007, August 2007 and October 2007, Case settled August, 2008.

Expert Witness - Deposition for NYS Attorney General regarding NYSDEC cost recovery for a petroleum spill site at Sag Harbor, NY. August 2002

Expert Witness - for NYS Attorney General regarding NYSDEC cost recovery for a petroleum spill site at Riverhead, NY. Case settled July 2008.

Expert Witness for defendant responding to a claim from adjacent commercial property owner on the origin of chlorinated solvents on plaintiff's property located in Cedarhurst, NY. Expert opinion submitted to lead counsel on March 6, 2009, case settled April 2009.

Expert Report - for Attorney General on modeling performed to determine the spill release scenario at a NYSDEC petroleum spill site in East Moriches, NY. June 2000.

MODELING EXPERIENCE (PARTIAL LISTING)

Table with 3 columns: PROJECT, MODEL, APPLICATION. Rows include Riverhead Water District, NYSDEC - Region 1, AMOCO, Keyspan Energy, Saboneck Golf Club, Suffolk County Department of Public Works, SCDPW SUNY Waste Water Treatment Plant, and Water Authority of Great Neck North.

PUBLICATIONS / PROFESSIONAL PAPERS

- Smart Pump & Treat Strategy for MTBE Impacting a Public Water Supply (14th Annual Conference on Contaminated Soils Proceedings, 1998)
Transport & Transformation of BTEX & MTBE in a Sand Aquifer (Groundwater Monitoring & Remediation 05/1998)
Characteristics of Gasoline Releases in the Water Table Aquifer of Long Island (Petroleum Hydrocarbons Conference Proceedings, 1999)
Field Applications of the Hydrocarbon Spill Screening Model (HSSM) (USEPA Interactive Modeling Web Course www.epa.gov/athens/software/training/webcourse Authored module on model application and applied use of calculators, 02/2000)
Comparative Evaluation of MTBE Sites on Long Island, US EPA Workshop on MTBE Bioremediation (Cincinnati, 02/2000)
Comparison of Four MTBE Plumes in the Upper Glacial Aquifer of Long Island (American Geophysical Union, San Francisco, 12/1996)
Analysis and Simulation of the Gasoline Spill at East Patchogue, New York (American Geophysical Union, San Francisco, 12/1998)



AMC Engineering
99 Jericho Turnpike, Suite 300J
Jericho, NY 11590
Phone: (516) 417-8588

ARIEL CZEMERINSKI, P.E.

Mr. Czemerinski is a New York State Professional Engineer and CEO of AMC Engineering PLLC an EBC affiliate. Mr. Czemerinski has with 20 years of experience in the chemical and environmental areas. Areas of expertise include environmental compliance, permitting, remedial system design, process and plant safety, and management of a production facility. Mr. Czemerinski is a Registered Professional Engineer in NY, IN, IL, and MI.

Professional Experience

AMC: 14

Prior: 6 years

Education

Master of Science in Chemical Engineering, Columbia University, New York, NY, Feb. 1990.
Bachelor of Science in Chemical Engineering, University Of Buenos Aires, Buenos Aires, Argentina, May 1987

Areas of Expertise

- Vapor Intrusion - Barrier and Sub Slab Venting System Design
- Environmental Assessment Statements and Environmental Impact Assessments under CEQR, ULURP
- Remedial Program Design and Management
- Environmental Compliance, Clean Water Act, Clean Air Act, Hazardous Materials
- Dewatering & Treatment System Design
- NYCDEP Sewer Discharge Permitting
- Transfer Station Permitting and Compliance
- Chemical Process Design and Optimization
- Wastewater Treatment Systems and Permitting, SPEDES, Air
- Zoning Regulations and Permitting
- Safety and Environmental Training
- Waste Management Plans

Professional Certifications

- OSHA 40-hr HAZWOPER
- OSHA 8-hr HAZWOPER Supervisor



ENVIRONMENTAL BUSINESS CONSULTANTS

Robert Bennett, Project Manager

Professional Experience

EBC: February 2015 - Present

Prior: 7 years

Education

Bachelor of Science, Environmental Science, State University of New York College at Oneonta, Oneonta, NY

Associates in Applied Sciences, Field Biology, State University of New York College at Delhi, Delhi, NY

Areas of Expertise

- Phase I / Phase II Property Assessments
- Waste Characterization / Soil Management
- Brownfield Closure and Planning Board
- Remedial Investigations
- Groundwater, Soil and Soil Vapor Remediation
- Indoor Air Quality (IAQ) Investigations
- Lead-Based Pain Risk Assessor
- Asbestos-Containing Materials Investigator/Inspector
- Landfill Closure and Monitoring
- Dredging Monitoring and Management
- Hazardous Materials Assessments
- Title V & NY Air Permitting and Registrations
- NYS / Nassau & Suffolk County Sanitary Code Compliance

Professional Certification

- OSHA 40-hr HAZWOPER
- OSHA 10-hr Construction Safety
- NYSDOH Asbestos Inspector
- NYCDEP Asbestos Investigator
- EPA Lead-Based Paint Inspector & Risk Assessor

PROFILE

Mr. Bennett has 8 years experience as an environmental consultant and is responsible for assessment and investigative services for a wide variety of projects, including industrial and commercial properties, mass transit facilities, parking structures, and sanitary and wastewater treatment facilities. Mr. Bennett has conducted Phase I, II and III Environmental Site Assessments for commercial, industrial, and residential properties in New York, New Jersey, and Massachusetts.



ENVIRONMENTAL BUSINESS CONSULTANTS

Robert Bennett, Project Manager

Mr. Bennett conducts research and provides support for various projects on a daily basis and coordinates with clients, regulatory agencies, attorneys and sub-contractors to provide cost-effective business solutions for a plethora of environmental concerns. Mr. Bennett's field experience includes tank removal and installations, dredging oversight and monitoring, asbestos and lead inspections, compliance audits, spill management and closure, soil and groundwater sampling, and both the oversight and operation of soil boring and well installation equipment. In addition, Mr. Bennett has performed project research, data reduction and evaluation, and has prepared reports for both regulatory and client use.

PREVIOUS EXPERIENCE

Dvirka & Bartilucci Engineers and Architects, P.C., Woodbury, NY
Environmental Scientist II, 2014-2015

Gannett Fleming Engineers and Architects, P.C., Woodbury, NY
Environmental Scientist, 2012-2014

Apex Companies L.L.C., Bohemia, NY
Environmental Scientist / Project Manager, 2008-2012

SELECT PROJECT EXPERIENCE

Project: Fulton Street Redevelopment Project - 1134 Fulton Street, Brooklyn NY
Description: NYC E-Designation. Soil contaminated with chlorinated solvents and heavy metals requiring excavation, soil management and disposal as well as a Soil Vapor Extraction System under a Remedial Action Work Plan, Soil / Materials Management Plan, Construction Health and Safety Plan and Community Air Monitoring Plan
Client: Porter Avenue Holdings
Authority: NYSDEC, NYSDOH & NYCOER
Role: Mr. Bennett served as the Project Manager for the project.

Project: Redevelopment Project - 391 Meeker Avenue, Brooklyn NY
Description: NYC E-Designation. Historic Fill Material requiring excavation, soil management and disposal under a Remedial Action Work Plan, Soil / Materials Management Plan, Construction Health and Safety Plan and Community Air Monitoring Plan
Client: Draftex Architectural Drafting & As Built Services
Authority: NYCOER
Role: Mr. Bennett serves as the Project Manager for the project.

Project: Redevelopment Project - 1555-1557 Fulton Street, Brooklyn NY



ENVIRONMENTAL BUSINESS CONSULTANTS

Robert Bennett, Project Manager

Description: NYC E-Designation. Historic Fill Material requiring excavation, soil management and disposal under a Remedial Action Work Plan, Soil / Materials Management Plan, Construction Health and Safety Plan and Community Air Monitoring Plan
Client: Waterfront Property Management, LLC.
Authority: NYCOER
Role: Mr. Bennett serves as the Project Manager for the project.

Project: Governor's Office of Storm Recovery (GOSR) New York Rising Buyout and Acquisition Program / Superstorm Sandy Relief Program
Location: Long Island and New York City
Type: Phase I Environmental Site Assessments (ESAs) and Property Evaluation
Contamination: Asbestos, Lead, Mold and PCBs
Role: Environmental Scientist II responsible for the creation and review of a high volume of Phase I ESAs

Project: WMATA Metrorail System Assessment Program
Location: Washington D.C. Area
Type: Hazardous materials inspection and evaluation for planning and engineering design purposes.
Contamination: Asbestos, Lead and PCBs
Role: Environmental Scientist and Inspection Team Leader

Project: Armonk Square Redevelopment Plan
Location: Armonk Square, Armonk, NY
Type: Monitoring well and recovery well installation. Sub-slab depressurization system (SSDS) installation and operational modifications.
Contamination: Chlorinated Solvents
Role: Environmental Scientist responsible for the planning and oversight of monitoring well and recovery well installation. Planning, oversight, and modifications to SSDS.

Project: Newtown Creek Dredging Project for NYCDEP
Location: NYCDEP Newtown Creek Wastewater Treatment Facility, Brooklyn, NY
Type: Navigational waterway dredging
Contamination: Hazardous and biological pollutants in bottom sediment.
Role: Environmental Scientist responsible for the implementation and operation of engineering controls and turbidity monitoring.

Project: Boring / Coring Program, Northeast U.S. Region
Location: New Bedford Harbor, New Bedford, MA. Long Island and Massachusetts.



ENVIRONMENTAL BUSINESS CONSULTANTS

Robert Bennett, Project Manager

- Type: Bathymetric surveys. Borings and Corings advanced through deep sediment and bedrock to determine the proper allocation dredge areas and confined aquatic disposal zones. Additionally, Vibracore drilling was conducted in shallow and easily accessible areas.
- Contamination: PCBs
- Role: Environmental Scientist / Project Manager serving as an on-site geologist to interpret and record geological investigations.
- Project: New York State Air Permit Facilities
- Location: Westchester, Orange and Rockland County, NY
- Type: Title V Air Permits, state registration and permitting for multiple industrial laundering facilities.
- Contamination: Hazardous Air Pollutants
- Role: Environmental Scientist / Project Manager responsible for all air permitting work for a NY-branch office.
- Project: Dredging Oversight and Water Quality Monitoring
- Location: New Bedford Harbor, New Bedford, MA
- Type: Bathymetric surveys. Supervised maintenance dredging and confined aquatic disposal zone excavation operations. Turbidity and sediment flocculation monitoring.
- Contamination: PCBs
- Role: Environmental Scientist providing project oversight, coordinating daily with Mass DEP and sub-contractors. Documenting geological data.
- Project: Stormwater Abatement System Inspections, Repairs and Reporting
- Location: Multiple retailer locations throughout New York State
- Type: Stormwater drainage system and stormwater control structure inspections and repairs
- Contamination: PCBs
- Role: Environmental Scientist / Project Manager assigned to coordinate and perform routine inspections of drainage systems and stormwater control structures. Made repairs to stormwater appurtenances where necessary.
- Project: ConEdison Truck-flush facility, effluent discharge monitoring.
- Location: Multiple ConEdison truck-flush facilities located throughout New York City, NY.
- Type: Compliance sampling and evaluation with regard to New York City Sewer Effluent Limitations.
- Contamination: Oil & Grease, Metals, Pesticides/PCBs, VOCs, SVOCs
- Role: Effluent sampling. Coordinating with client and laboratory to conduct quarterly sampling events.
- Project: RCRA Closure Support



ENVIRONMENTAL BUSINESS CONSULTANTS

Robert Bennett, Project Manager

Location: Pall Corporation Former Headquarters, East Hills, NY
Type: Environmental closure of a medical equipment manufacturing facility
Contamination: Formic Acid, Dimethylacetamide (DMAC)
Role: Environmental Scientist / Project Manager responsible for the supervision of the removal of all process tanks, piping and associated appurtenances. Accomplished final decommissioning activities. RCRA Closure Report.

Project: Brownfield Closure Support
Location: Multiple locations throughout New York City
Type: Remedial investigations. Interim remedial measures. Soil vapor intrusion studies. RCRA Closure.
Contamination: VOCs, SVOCs, Oil & Grease, Pesticides/PCBs, Metals
Role: Environmental Scientist / Project Manager responsible for preparing and conducting remedial investigations, interim remedial measures, soil vapor intrusion studies and RCRA closure.

Project: Mirant Bowline Power Plant Asbestos Survey
Location: West Haverstraw, NY
Type: Asbestos inspection. Personal exposure monitoring. Asbestos labeling Program. Reporting.
Contamination: Asbestos
Role: Environmental Scientist / Project Manager serving as a team leader to conduct large scale asbestos inspection, labeling program and reporting.

Project: Estee Lauder SPCC Facilities
Location: Multiple manufacturing facilities throughout Long Island
Type: Spill Prevention Control & Countermeasures (SPCC) inspections, evaluation and reporting.
Contamination: N/A
Role: Environmental Scientist / Project Manager responsible for conducting inspections, facility engineering review, and reporting.

Project: Nassau and Suffolk County Sanitary Code Facility Compliance Audits
Location: Multiple medical equipment manufacturing facilities throughout Long Island.
Type: Article XI and XII Sanitary Code Compliance Audits and multiple medical equipment manufacturing facilities.
Contamination: N/A
Role: Environmental Scientist / Project Manager responsible for conducting inspections, facility engineering review, and reporting.

PUBLICATIONS

Dredging and Beach Nourishment Public Notices
(Cape Cod Times, 2008-2010)



ENVIRONMENTAL BUSINESS CONSULTANTS

Kevin Waters, Field Manager

Professional Experience

EBC: October 2010

Prior: 5 years

Education

Bachelor of Science, Geology, State University of New York, Stony Brook

Areas of Expertise

- Field Operations
- Phase II and RI Implementation, Site Characterization Studies
- Health & Safety Monitoring and Oversight
- Waste Characterization / Soil Management
- Site Logistics

Professional Certification

- OSHA 40-hr HAZWOPER
- OSHA 8-hr HAZWOPER Supervisor

PROFILE

Mr. Waters has 10 years experience as an environmental consultant and has worked on a wide range of environmental projects. Mr. Waters has conducted Phase II and III Environmental Site Assessments for commercial, industrial, and residential properties in New York.

Mr. Waters' field experience includes soil, air and groundwater sampling, operations and maintenance of groundwater remediation systems, tank removals, spill management and closure, and oversight of monitoring well installations. In addition, Mr. Waters has prepared reports for both regulatory and client use.

PREVIOUS EXPERIENCE

P.W. Grosser Consulting, Bohemia, NY

Field Hydrogeologist, 2003-2008

SELECT PROJECT EXPERIENCE

| | |
|----------------|--|
| Project: | Former Gas Station / car wash to mixed use affordable housing / commercial |
| Location: | Bronx, NY, Southern Boulevard |
| Type: | NYS BCP, NYC E-Site Hazmat, Former gas station / gar wash |
| Contamination: | Petroleum - Gasoline |
| Role: | Field Operations Manager, Health and Safety Officer |



ENVIRONMENTAL BUSINESS CONSULTANTS

Kevin Waters, Field Manager

SELECT PROJECT EXPERIENCE

Project: Former Uniforms for Industry Site – Richmond Hill Senior Living Residences / Richmond Place
Location: Jamaica Ave, Richmond Hill Queens, NY
Type: NYS BCP, NYC E-Site Hazmat, Noise, Former industrial Laundry
Contamination: Chlorinated Solvents, Historic Fill, Petroleum - Fuel oil/Mop oil
Role: Field Operations Manager, Health and Safety Monitoring and Field Oversight

Project: Rikers Island – West Intake Facility
Location: NYC Department of Corrections, Rikers Island, NY
Type: Municipal Construction Project
Contamination: Hazardous levels of lead, heavy metals in Historic fill
Role: Field Operations Manager, Health and Safety Monitoring and Field Oversight

Project: Residential Redevelopment Project
Location: Williamsburg Section of Brooklyn, Wallabout Street
Type: NYC E-Designation Site
Contamination: Hazardous levels of lead, heavy metals, SVOCs in Historic fill
Role: Implement RI Work Plan, Supervise sample collection in all media



ENVIRONMENTAL BUSINESS CONSULTANTS

Kevin R. Brussee, Senior Project Manager

Professional Experience

EBC: January 2008

Prior: 6 years

Education

Bachelor of Science, Environmental Science, Plattsburgh State University, NY

Master of Science, Environmental Studies, University of Massachusetts, Lowell

Areas of Expertise

- Management of Site Investigations / Remedial Oversight NYC "E" Designation Sites
- Management of RI Investigations / RAWP Implementation NYS BCP Sites
- NYSDEC Spill Site Investigations
- Phase I / Phase II Property Assessments
- Waste Characterization / Soil Management

Professional Certification

- OSHA 40-hr HAZWOPER
- OSHA 8-hr HAZWOPER Supervisor

PROFILE

Mr. Brussee has 10 years experience as an environmental consultant/contractor and has worked on and managed a wide range of environmental projects. Mr. Brussee has conducted Phase I, II and III Environmental Site Assessments for commercial, industrial, and residential properties in New York, New Jersey, Maryland and Delaware.

Mr. Brussee's field experience includes tank removal and installations, spill management and closure, soil and groundwater sampling, and both the oversight and operation of soil boring and well installation equipment. In addition, Mr. Brussee has performed project research, data reduction and evaluation, and has prepared reports for both regulatory and client use.

PREVIOUS EXPERIENCE

Eastern Environmental Solutions, Inc., Manorville, NY

Project Manager, 2006-2008

EA Engineering, Science & Technology

Hydrogeologist, 2005-2006

P.W. Grosser Consulting, Bohemia, NY

Field Hydrogeologist, 2002-2003



ENVIRONMENTAL BUSINESS CONSULTANTS

Kevin R. Brussee, Senior Project Manager

SELECT PROJECT EXPERIENCE

- Project:** Former Dico G, Auto and Truck Repair Site - Bronx Park Apartments, redevelopment from commercial to mixed use
Location: Bronx, NY, White Plains Road
Type: NYS BCP Site, Former gas station, repair shop & junk yard
Contamination: Petroleum - Gasoline
Role: Project Manager, during Site Management Phase
- Project:** Former Uniforms for Industry Site – Richmond Hill Senior Living Residences / Richmond Place
Location: Jamaica Ave, Richmond Hill Queens, NY
Type: NYS BCP, NYC E-Site Hazmat, Noise, Former industrial Laundry
Contamination: Chlorinated Solvents, Historic Fill, Petroleum - Fuel oil/Mop oil
Role: Project Manager, RAWP implementation
- Project:** Former Gas Station / car wash to mixed use affordable housing / commercial
Location: Bronx, NY, Southern Boulevard
Type: NYS BCP, NYC E-Site Hazmat, Former gas station / gar wash
Contamination: Petroleum - Gasoline
Role: Project Manager, RAWP implementation
- Project:** Redevelopment of former industrial property to residential
Location: Williamsburg section of Brooklyn, NY, Bedford Ave
Type: NYC E-Designation Site, Former dye manufacturing plant
Contamination: Hazardous levels of heavy metals, fuel oil tanks
Role: Project Manager, RAWP implementation
- Project:** Former Domsey Fiber Corp Site
Location: Williamsburg section of Brooklyn, NY, Kent Ave
Type: NYC E-Designation Site, Former commercial property
Contamination: Chlorinated solvents, fuel oil and Historic fill
Role: Project Manager, RIWP Development and Implementation, RAWP development and implementation, waste characterization and soil management

PUBLICATIONS

Chemical Stress Induced by Copper, Examination of a Biofilm System;
(Water Science Technology, 2006; 54(9): 191-199.)



ENVIRONMENTAL BUSINESS CONSULTANTS

Chawinie Reilly, Project Manager / Industrial Hygienist

Professional Experience

EBC: March 2013

Prior: 8 years

Education

Bachelor of Science, Environmental Health and Safety, Stony Brook University, NY

Areas of Expertise

- Phase I / Property Condition Assessments
- Occupational Health and Safety Sampling
- Indoor Air Quality (IAQ) Investigations
- Mold Investigations and Remediation
- Soil and Ground Water Investigations
- Noise Studies
- Lead Paint and Asbestos Surveys
- Hazardous Materials Assessments

Professional Certification

- OSHA 40-hr HAZWOPER
- NYS Asbestos Inspector
- NYC Asbestos Investigator
- USEPA Lead Inspector
- USEPA Lead Risk Assessor
- OSHA 10-hr Construction Health and Safety
- Hazard Analysis and Critical Control Point (HACCP) Certified

PROFILE

Mrs. Reilly has 9 year's experience as an environmental consultant/contractor and has worked on and managed a wide range of environmental projects. Ms. Miller has conducted Phase Is and Property Condition Assessments for commercial, industrial, and residential properties in New York, New Jersey and Connecticut. In addition, Ms. Miller has conducted various IAQ, asbestos, mold and occupational health and safety sampling investigations for a variety of city, state, federal and private clients.

PREVIOUS EXPERIENCE

The Louis Berger Group, New York, New York
Industrial Hygienist, 2008-2013

AEI Consultants, Jersey City, New Jersey
Environmental Scientist, 2005-2008

ATTACHMENT G
BCP Signage Specifications



Department of
Environmental
Conservation

New York State Brownfield Cleanup Program

FORMER MUGLER SHORING INC. SITE

BCP Site No. C203052

2401 3rd Ave Associates Property LLC

Governor Andrew M. Cuomo

NYSDEC Acting Commissioner Basil Seggos

Mayor Bill de Blasio

Transform the Past... Build for the Future.

Sign Requirements

Size: Horizontal format – 96” wide by 48” high

Construction

Materials: Aluminum or wood blank sign boards with vinyl sheeting.

Inserts: “New York State and DEC logo”, “Program Name”, “Site Name”, “Site No.”, “Name of Party Performing Remedial Activities or New York State Department of Environmental Conservation”, “Governor”, “DEC Commissioner”, “Municipal Executive”, “Transform the Past...Build for the Future”.

Color Scheme: All body font should be black or green Pantone 350 C or CMYK 80/43/83/42. If blue is desired, use following values: Pantone 288 C or CMYK 100/87/27/19.

New York State and DEC logo: use eps file [here](#) (it is high resolution and scalable. If vendor needs a different format, use jpg file [here](#) . Both utilize the correct color.

Text:

Program Name (choose one):

| | | |
|---------------------------------------|---|--|
| State Superfund Program | } | Green text (PANTONE 350C or CMYK 100/43/83/42 |
| Brownfield Cleanup Program | | |
| 1996 Clean Water/Clean Air Bond Act – | | |
| Environmental Restoration Program | | |
| Voluntary Cleanup Program | | |
| Petroleum Remediation Program | | |

Site Name: Blue text (PANTONE 288C or CMYK 100/87/27/19)

Site Number: Blue text (PANTONE 288C or CMYK 100/87/27/19)

Name of Party Performing Remedial Activities or New York State Department of Environmental Conservation: Green text (PANTONE 350C or CMYK 100/43/ 83/42

Governor: Black text

DEC Commissioner: Black text

Municipal Executive: Black text

Transform the Past....Build for the Future: Blue text (PANTONE 288C or CMYK 100/87/27/19)

Type

Specifications:

All type is Ariel.

Format is: Center each line of copy with initial caps and small Letters.

Production

Notes:


96" wide x 48" high aluminum blanks will be covered with vinyl sheeting to achieve background color. Copy and logo will be silk screened on this surface.

See Attached

Format:

Next page.

8'

| | | | |
|----|--|---|----------------------|
| 2" |  <p>NEW YORK STATE OF OPPORTUNITY.</p> | <p>Department of Environmental Conservation</p> <p>Logo (Use eps or .jpg file) Green Text (See Key) White Background</p> | Green Text (See Key) |
| 6" | Program Name | | Blue Text (See Key) |
| 2" | Site Name | | Blue Text (See Key) |
| 4" | Site No. | | Blue Text (See Key) |
| 2" | Remedial Party or New York State Department of Environmental Conservation | | Green Text (See Key) |
| 2" | Governor | | Black Text |
| 2" | Commissioner | | Black Text |
| 2" | Municipal Executive | | Black Text |
| 4" | Transform the Past.... Build for the Future | | Blue Text (See Key) |
| 2" | | | |

4'

Color Key for Text

Green Text = Pantone 350C or CMYK 80/43/83/42

Blue Text = Pantone 288C or CMYK 100/87/27/19

Project Sign Format

ATTACHMENT H
Estimated Remedial Costs

FORMER MUGLAR SHORING
2401 Third Avenue
Bronx, NY

Summary of Project Costs

NYS Brownfields Cleanup Program
Costs by Task

| TASK - ENVIRONMENTAL REMEDIATION | Alternative 1 - Track 1 | Alternative 2 - Track 4 |
|---|--------------------------------|--------------------------------|
| BCP Entry Documents | Completed | Completed |
| Supplemental Investigation And RI Report | Completed | Completed |
| Remedial Work Plan, Remedy Scoping & Coordination | \$ 18,450.00 | \$ 18,450.00 |
| Remedial Program Implementation | \$ 3,266,179.00 | \$ 727,655.00 |
| Final Engineering Report, Site Management Plan & IC/ECs | \$ 18,200.00 | \$ 50,450.00 |
| <i>Subtotal</i> | \$ 3,302,829.00 | \$ 796,555.00 |
| <i>15% Contingency</i> | \$ 495,424.35 | \$ 119,483.25 |
| <i>Total</i> | \$ 3,798,253.35 | \$ 916,038.25 |

**FORMER MUGLER SHORING
REMEDIAL PROGRAM COST ANALYSIS
REMEDIAL ACTION
ALTERNATIVE 1**

| Description | Category | Quant. | Unit | Unit Cost | Capital Cost |
|---|--|--------|---------------|---------------|------------------------|
| Excavation and Disposal of Impacted Soil | | | | | \$ 3,257,179.00 |
| Includes costs associated with disposal of excavated soil from petroleum hotspot (and entire site to 8 ft. Includes backfill of the portions of the site as needed to return to grade as per the development plans.l | Supervision, oversight & air monitoring | 90 | day | \$ 850.00 | \$ 76,500.00 |
| | Project Management & Coordination PM | 200 | hr | \$ 135.00 | \$ 27,000.00 |
| | Project Management & Coordination Princ. | 40 | hr | \$ 195.00 | \$ 7,800.00 |
| | Project Management & Coordination Field Manager | 100 | hr | \$ 105.00 | \$ 10,500.00 |
| | Petroleum Impacted Soil Non Haz disposal (150 cy) | 225 | tons | \$ 55.00 | \$ 12,375.00 |
| | Soil Non Haz disposal (18,084 cy) | 27126 | tons | \$ 55.00 | \$ 1,491,930.00 |
| | Excvate top 8 feet of Site (18,084 cy) | 90 | days | \$ 3,500.00 | \$ 315,000.00 |
| | Backfill Materials (fill excavation and restore grade at Site) | 15861 | cy | \$ 29.00 | \$ 459,969.00 |
| | End point sample analysis | 67 | ea | \$ 550.00 | \$ 36,850.00 |
| | Waste Characterization Analysis | 22 | sample | \$ 1,500.00 | \$ 33,000.00 |
| | Waste Characterization Sample Collection | 5 | day | \$ 850.00 | \$ 4,250.00 |
| | Deliverables and EDDs | 2 | LS | \$ 2,500.00 | \$ 5,000.00 |
| | Data Validation | 67 | ea | \$ 115.00 | \$ 7,705.00 |
| | DUSRs | 2 | ea | \$ 2,500.00 | \$ 5,000.00 |
| | Testing of backfill materials | 26 | ea | \$ 550.00 | \$ 14,300.00 |
| | Dewatering west third of site | 1 | LS | \$ 250,000.00 | \$ 250,000.00 |
| Shoring | 1 | LS | \$ 500,000.00 | \$ 500,000.00 | |
| Reporting | | | | | \$ 9,000.00 |

**FORMER MUGLER SHORING
REMEDIAL PROGRAM COST ANALYSIS
REMEDIAL ACTION
ALTERNATIVE 2**

| Description | Category | Quant. | Unit | Unit Cost | Capital Cost |
|--|--|--------|--------|-------------|---------------|
| Excavation and Disposal of Impacted Soil | | | | | |
| Includes costs associated with disposal of excavated soil from petroleum hotspot (150 cy) . Includes excavation and backfill of hotspot excavation and historic fill from building areas (34,391 sf). The soil to go for non-hazardous disposal of all excavated soil either as urban fill or petroleum impacted soil. Non-hazardous soil to be excavated as part of construction. | Supervision, oversight & air monitoring | 50 | day | \$ 850.00 | \$ 42,500.00 |
| | Project Management & Coordination PM | 60 | hr | \$ 135.00 | \$ 8,100.00 |
| | Project Management & Coordination Princ. | 20 | hr | \$ 195.00 | \$ 3,900.00 |
| | Project Management & Coordination Field Manager | 60 | hr | \$ 105.00 | \$ 6,300.00 |
| | Petroleum Impacted Soil Non Haz disposal (150 cy) | 225 | tons | \$ 55.00 | \$ 12,375.00 |
| | Soil disposal from building areas (5,094 cy) | 7641 | tons | \$ 55.00 | \$ 420,255.00 |
| | Backfill Materials | 1000 | cy | \$ 26.00 | \$ 26,000.00 |
| | Excavate hotspot, building basement areas and exposed soil areas | 50 | days | \$ 3,500.00 | \$ 175,000.00 |
| | End point sample analysis | 15 | ea | \$ 200.00 | \$ 3,000.00 |
| | Waste Characterization Analysis | 10 | sample | \$ 1,500.00 | \$ 15,000.00 |
| | Waste Characterization Sample Collection | 1 | day | \$ 850.00 | \$ 850.00 |
| | Testing of backfill materials | 3 | ea | \$ 550.00 | \$ 1,650.00 |
| | Deliverables and EDDs | 1 | LS | \$ 2,500.00 | \$ 2,500.00 |
| | Data Validation | 15 | ea | \$ 115.00 | \$ 1,725.00 |
| | DUSRs | 1 | ea | \$ 2,500.00 | \$ 2,500.00 |
| Reporting | | | | | |
| | | | | | \$ 6,000.00 |