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March 3, 2014

Ms. Tara M. Blum, P.E.  
New York State Department of Environmental Conservation, Region 7  
615 Erie Boulevard West  
Syracuse, NY 13204-2400

**Subject: Data Summary/Transmittal Report  
November 2013 Wetland Delineation and Sampling  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, NY**

Dear Ms. Blum:

On behalf of United Technologies Corporation (UTC), AECOM USA, Inc. (AECOM) is pleased to provide this transmittal of data recently collected at the Carrier-Dewitt Landfill site (AOC G) on Thompson Road in Syracuse, New York. The sampling procedures and analytical results described below were commissioned as components of a Remedial Investigation for AOC G. These work elements and procedures were performed in accordance with the Draft Remedial Investigation Work Plan (RIWP) prepared for AOC G in 2013 by EnSafe Inc. and the comments supplied by the New York State Department of Environmental Conservation (NYSDEC) in a memo dated September 12, 2013.

As stated during previous meetings and correspondence between NYSDEC and UTC, it was determined that proceeding with characterization of the wetland area comprising the western portion of the AOC G site (see Site Plan, **Figure 1**) during the latter part of 2013 would be advantageous due to the beneficial vegetative and climatic conditions provided by the fall season. Accordingly, a field team was mobilized to delineate and sample the wetland during the week of November 4, 2013. The remainder of this letter details the wetland delineation and sampling tasks completed during November 2013. It is intended that the upland areas of AOC G (former landfill area) and additional locations within the wetland will be sampled and reported upon during 2014.

## **1.0 November 2013 AOC G Wetland Investigation Scope of Work**

The scope of work completed during November 2013 included:

- Delineation of the AOC G Wetlands,
- Sampling of the AOC G Wetland Area,
- Surveying the AOC G Wetland Area and the Sanders Creek Corridor, and
- Preparing a Data Summary/Transmittal Report.

### **1.1 Delineation of the AOC G Wetlands**

#### **Wetland Delineation and Flagging Methodology**

AECOM delineated the jurisdictional boundary between the upland portions of AOC G and the adjacent wetlands using the three-parameter method documented in the *Corps of Engineers Wetlands Delineation Manual* (Technical Report Y-87-1, 1987) as modified in the *Regional Supplement to the Corps of Engineers Wetland Delineation Manual: Northcentral and Northeast Region, V2* (ERDC/EL TR-12-1, 2012). The delineation was completed during the week ending

November 8, 2013 and was based on the presence of hydrophytic vegetation, hydric soils, and wetland hydrology.

### **Observations and Results**

A map showing the wetland features described below is attached as **Figure 2**.

- The wetland boundary line was started along the eastern edge of a swale ("north swale") that drains northward toward Sanders Creek from a ponded area at the north end of the wetland ("north wetland"). The north swale itself is considered a wetland, but its wetland boundary northward to Sanders Creek (including a feature referred to as the "Hotel Pond") was not delineated.
- The wetland boundary extends south from the north swale along the edge of fill used to construct the UTC-Carrier parking lot past a rip-rap drainage swale from the parking lot and continues along the toe of the landfill slope approximately 1,000 feet to a point where historical earth moving activities (presumably clay mining) has left a complex of upland fill areas separated by low wetland/ponded areas.
- The wetlands-uplands complex was further delineated where it abuts the wetlands in the western part of the UTC property. Identified on the map is a low point where a majority of the wetlands-uplands complex drains into the western wetland area and a low upland peninsula composed of discarded bricks.
- The wetland line continues south along the toe of the landfill slope across the complex of wetlands-uplands to the southwest corner of the landfill, where it meets a drainage swale that drains the truck parking area on the adjacent property south of the landfill.

This was the extent of delineation of the wetland completed during this project phase, as it was expected that this wetland/landfill boundary may be involved in future remedial action permitting.

### **Seasonal/Water Elevation Impacts on Wetland Delineation**

The level of water in the wetlands was higher than that observed during a September 11, 2013 field reconnaissance conducted by AECOM. The lower water elevation aided in delineating wetland hydrology. However, the vegetation was dying back for the winter as a result of previous frosts, so documentation of hydrophytic vegetation was hampered. Additional documentation of selected data points will be completed during the spring field season when hydrophytic vegetation is fully evident.

### **Visual Observation of Water Sources and Outlets (Including Presence or Absence of Outlets for Outlet Flow or Overland Flooding to Sanders Creek and South Branch of Ley Creek)**

- A swale to Sanders Creek (north swale) from the northern wetland was verified.
- The water level in the north wetland at the beginning of the week of November 4, 2013 was 6 to 8 inches below the elevation to initiate flow into the swale to Sanders Creek. At the end of the week the water level in the north wetland was approximately 3 to 4 inches below the threshold into the swale.
- An active outfall (**Figures 1 and 2**, Outfall A) to South Branch of Ley Creek was located on County Sewer District property near the southwest corner of the UTC property. The flow was traced back to the western ponded area on the UTC property. This confirms that water from the landfill area may discharge to Ley Creek and that this drainage pathway is active before discharge through the north swale to Sanders Creek.
- A second trickling outfall to South Branch of Ley Creek (**Figures 1 and 2**, Outfall B), potentially originating on UTC property, was located approximately 800 feet downstream (northwest) from the main outfall.

The elevations of both the northern and southwestern drainage pathways are such that water levels in South Branch of Ley Creek or Sanders Creek are unlikely to rise sufficiently to initiate flow onto the UTC property.

## 1.2 Sampling of the AOC G Wetland Area

Due to the historic usage of the AOC G landfill for disposal of industrial waste material, the RIWP includes activities to investigate potential releases of organic compounds and metals from the landfill into the adjacent wetlands to its north and west. These chemicals may have been present in soils eroded from the landfill and deposited into the wetland sediment by surface water runoff. During the years of operation of the landfill and after its closure, alternating layers of contaminated and clean sediment may have been deposited in the wetlands, so that a layer of contaminated sediment may potentially be buried beneath a layer of uncontaminated sediment. Therefore, during this investigation AECOM collected sediment samples from three depth intervals as described below.

In addition, the infiltration of precipitation into the landfill mass may have dissolved chemicals from the waste and soils and transported these contaminants through groundwater discharge into the wetland. During this wetland sampling event, pore water and surface water samples were not collected because the site was inundated with standing water. A concern that standing water could make it difficult to distinguish between surface water and pore water quality resulted in pore water sampling being deferred until the surface water elevation is lower.

### Choice of Sediment Sampling Locations (Including Sampling of Flow Paths and Transects Outward from Landfill Area to Determine Gradient of Contamination)

- Sediment sample locations were placed on eleven transects (A – K) radiating from the landfill boundary (refer to **Figure 2**), to aid in assessing the presence or absence of a concentration gradient from the landfill into the wetland.
- Nine sediment samples (sample prefix S) were collected along the north swale from the northern wetland to Sanders Creek.
- The Balduck Method was used to estimate the approximate number of sediment sample locations needed to characterize the potential landfill impact on the wetlands. Using an area of approximately 112,500 square yards to estimate the area to be characterized (300 X 375 yards), and an assigned degrees of freedom (Df) of 3:

$$N = Df * 30 * (W * L * (1 / 1.2E06))^{0.33} = 41 \text{ sample locations}$$

where:

N is the number of samples for the area derived by the Balduck Method;  
Df (degrees of freedom) is a theoretical constant and the chosen value (3)  
is conservative;  
W and L are the width and length of the subject wetland area.

- AECOM collected 135 samples from 45 locations, including the nine samples collected in the northern swale to Sanders Creek. Quality control (QC) samples were also collected. The locations where sediments were sampled during November 2013 are shown on **Figure 2**.

### Sample Depth Intervals

AECOM collected sediment samples from three depth intervals (0-6 inches, 6-12 inches, and 18-24 inches) from all sample locations.

### Sample Collection Procedures

- AECOM collected sediment cores using a slide hammer and hand auger. Upon collection, the cores were immediately sealed in plastic zipper bags and transported to the processing team to be logged and sampled.

- After screening with a photoionization detector (PID), volatile organic compound (VOC) samples were collected from each interval prior to homogenization of the interval material for polychlorinated biphenyl (PCB), total organic carbon (TOC), metals, semi-volatile organic compound (SVOC), and polycyclic aromatic hydrocarbon (PAH) sampling.

### **Piezometer Installation**

- Pre-constructed piezometers were installed at ten locations around the toe of the landfill to a depth of 3 feet below ground surface (bgs).
- The piezometers were constructed of 2.5 feet of 1½-inch inside diameter polyvinyl chloride (PVC) 0.010" slotted pipe threaded onto a 3-foot solid riser. The slotted section of pipe was surrounded by filter sand held in place by food-grade plastic mesh. The bottom was capped and the top was fitted with a plug.
- Piezometers were installed at locations A-1, B-2, C-2, D-1, E-1, F-2, H-1, I-1, J-1, and K-1 shown on **Figure 2**. The locations were not underwater at the time of installation.

### **Equipment Decontamination**

The hand augers and slide hammers were decontaminated with an Alconox rinse followed by a de-ionized water rinse between locations. New core-liners were used for each sediment core in the slide hammer.

### **Analytical Methods**

- Each sediment sample was analyzed for VOCs (United States Environmental Protection Agency [US EPA] Method 5035/8260),
- TOC (Lloyd-Kahn method),
- SVOCs (US EPA Method 8270C),
- PCBs (US EPA Method 8082),
- Metals (US EPA method 6010B),
- Percent moisture, and
- T34 PAHs by SIMS (US EPA Method 8270D).

### **1.3 Surveying the AOC G Wetland Area and Sanders Creek Corridor**

A topographic survey of AOC G and adjoining portions of the Sanders Creek Corridor was performed by a New York State Licensed land surveyor.

- Topographic and planimetric features were mapped utilizing photogrammetric mapping techniques based on aerial photography obtained on November 16, 2013.
- Four established benchmarks and 16 survey points were used to provide physical control for the survey. A field edit was performed on January 3, 2014 to verify the accuracy and completeness of the photogrammetric mapping.
- The survey is referenced to the North American datum of 1983, 2011 adjustment (NAD83/2011) and projected on the New York State Plane Coordinate System and vertically to the North American Vertical Datum of 1988 (NAVD88-Geoid 09).

This survey was used as the basis of the maps included with this report, which were prepared at a 1-foot contour interval.

## 2.0 AOC G Wetland Sediment Sampling Data and Discussion

### 2.1 Review of New York State Sediment Screening Methodology

Protection of ecological resources, specifically, fish, wildlife, and habitat thereof within New York State is a responsibility of the NYSDEC Division of Fish, Wildlife and Marine Resources (DFWMR). As part of this responsibility, DFWMR has developed draft guidance for the assessment of sediment quality and the potential for organic and inorganic contamination to adversely affect sediment-dwelling (benthic) organisms.

This document (*Screening and Assessment of Contaminated Sediment*, NYSDEC, DFWMR, Bureau of Habitat, January 24, 2013) is intended to provide information and guidance to DFWMR staff and other interested parties for screening, evaluating and assessing contaminated sediments in New York State; that is, for determining whether or not a given sediment is toxic and whether it potentially poses a risk to aquatic life.

The guidance document describes criteria and procedures for assessing whether or not contaminants present in sediment at a given site are causing toxicity, and if they pose a risk to aquatic life.

Screening sediments is the first, necessary step in the assessment process. In order to screen sediments to assess the potential for risk to aquatic life, a set of numeric criteria is needed for comparing bulk sediment contaminant concentrations. The NYSDEC draft guidance identifies numerical thresholds for various chemical contaminants found in sediments that can be used to identify potential risk to aquatic life. Given no information other than the concentration of a contaminant in sediment, these values allow for a reasonable assessment of the potential for the contaminants to be harmful to aquatic life.

In addition to identifying specific numerical criteria, the draft guidance document explains the technical basis for the derivation of the criteria values selected, and explains how those numeric criteria values can be modified if more information, such as site-specific data, becomes available.

#### Sediment Assessment Procedure

The NYSDEC sediment assessment procedure is a stepwise process:

1. Sediment Guidance Values (SGVs) in Tables 1a (fresh water sediments) and 1b (salt water sediments) are used to make the initial assessment (i.e., Tier 1 screening) of risk to aquatic life from contaminants in sediment.
2. If the concentrations of all of the contaminants present are below the Class A threshold values, the sediment is considered to present a low risk to aquatic life.
3. If the concentration of one or more contaminants exceeds the Class C threshold values, then the sediment is considered highly contaminated, and likely to present a high risk to aquatic life.
4. If the concentration of one or more contaminants lies between the Class A and Class C threshold values and no contaminants exceed the Class C threshold values, then the sediments are considered to be moderately contaminated (Class B). Additional testing and/or evaluation are needed to determine the magnitude and extent of or potential for aquatic life toxicity.
5. For contaminant mixes such as PAHs, the classification of sediment can be based on the mean SGV quotient, assuming none of the individual quotients exceed 1.0. That is, if the mean Class A quotient is less than 1.0, then the sediments may be considered to be Class A; and if the mean Class C threshold is greater than 1.0, then the sediments are considered to be Class C. Sediments with a mean Class A quotient that is greater than 1.0 and a mean Class C quotient that is less than 1.0 are considered to be Class B. Total SGV quotients are used instead of mean SGV quotients if the contaminants are likely to be additive in toxicity.

6. The second tier of sediment assessment is to adjust the SGVs for local conditions, such as TOC. This is applicable only to equilibrium partitioning-based SGVs. Once the site-specific TOC has been measured, the equilibrium-based SGVs can be recalculated and the screening and classification of sediments adjusted.
7. The third tier of sediment assessment is to conduct additional studies to measure and evaluate the potential for toxicity from contaminants in the sediment. The additional studies constitute a weight of evidence approach; that is, multiple lines of evidence are used to evaluate risk. Additional studies should be designed, as a minimum, to be consistent with the sediment quality triad, which includes toxicity testing and benthic macroinvertebrate community analysis. This report does not include a Tier 3 assessment.

## 2.2 Tier 1 Screening Assessment

Bulk sediment concentrations of detected chemicals collected from 45 sample locations at 3 depths (i.e., 0-6 inches, 6-12 inches, and 18-24 inches) were compared to Class A screening values from Table 1a in the DEC sediment guidance document (NYSDEC 2013).

### Inorganics

Twenty-two metals in sediment were analyzed using US EPA method SW6010 and mercury was analyzed using US EPA method SW7471B). Twenty-three metals were detected at least once in site sediments. Calcium, magnesium, iron, potassium, and sodium are considered nutrient minerals and have no screening criteria and were not screened. These minerals are needed by organisms for normal metabolism and internal concentrations can be regulated over a wide range of concentrations. Nine other metals (i.e., aluminum, antimony, barium, beryllium, cobalt, manganese, selenium, thallium, and vanadium) also lack Class A screening values and were not evaluated. **Table 1** presents the results of the Tier 1 screening of metals against Class A criteria. A summary of locations where metals exceedances were most common is presented in **Figure 3**.

The greatest number of different metals exceeding Class A criteria occurred in the 0 to 6 inch interval samples. Nickel exhibited the greatest number of exceedances for all intervals (12 exceedances at 0 to 6 inches, 15 exceedances at 6 to 12 inches, and 27 at 18 to 24 inches). Other exceeding metals included arsenic (3 at 0 to 6 inches, 3 at 6 to 12 inches, and 4 at 18 to 24 inches); cadmium (2 at 0 to 6 inches); chromium (3 at 0 to 6 inches); copper (3 at 0 to 6 inches, 3 at 6 to 12 inches, and 1 at 18 to 24 inches); lead (4 at 0 to 6 inches, and 3 at 6 to 12 inches); mercury (3 at 0 to 6 inches, and 1 at 6 to 12 inches); and zinc (4 at 0 to 6 inches, and 1 at 6 to 12 inches). No metals exceeded NYSDEC Class C criteria at any depth. Therefore, all metals that exceed Class A criteria are considered Class B metals.

Because metals are naturally occurring elements in sediment, concentrations of metals within the natural background range are not expected to cause adverse effects to benthic communities. Natural background concentration ranges for metals in the region of the AOC G landfill are not available for sediment.

Site-specific background concentration ranges for metals may be developed from the existing data set using geochemical enrichment normal probability plots and geochemical correlations between naturally occurring elements, and other statistical tools. Alternatively, sediment samples from locations considered to be similar to the geological environment of the site, but not impacted by metal contamination, could be sampled and analyzed to establish background ranges for site metals.

### Organic Compounds

**Polychlorinated Biphenyls:** PCBs were analyzed using US EPA method SW8082 as seven different Aroclor mixtures. Total PCBs were estimated as the sum of concentrations of detected

Aroclors. This is considered a conservative approach because certain congeners found in more than one mixture are “double counted”.

- Aroclor 1254 and Aroclor 1260 were the only PCBs detected on the site. No other Aroclors were detected. **Table 2** presents the results of the Tier 1 screening of total PCBs against NYSDEC Class A criteria.

Total PCBs were detected at the transect sample locations closest to the toe of the landfill, except at SDC1 in the wetland-upland complex. Exceedance of NYSDEC Class A criteria were found at SDA4 (at 0 to 6 and 6 to 12 inches); SDB1 (at 0 to 6 inches); SDF1 (at 0 to 6, 6 to 12, and 18 to 24 inches); SDH1 (at 0 to 6 inches); SDJ1 (0 to 6 inches); and SDK1 (0 to 6 inches). The distribution of exceedances suggests a landfill source of PCBs. A summary of total PCB concentrations exceeding Class A criteria is presented on **Figure 4**.

- PCBs also exceeded Class A criteria at 2 locations in the Sanders Creek floodplain at S-8 (6 to 12 inches, and 18 to 24 inches) and S-9 (0 to 6 inches, 6 to 12 inches, and 18 to 24 inches).
- Concentrations of total PCBs at SDF1 (0 to 6 inches and 18 to 24 inches) also exceed the NYSDEC Class C criteria, indicating a potential for adverse effects to the benthic community in the area of SDF1. The potential for adverse effects was further evaluated considering the potential bioavailability of PCBs in the sediment.

Using the equilibrium partitioning (EqP) method provided in NYSDEC draft sediment guidance (NYSDEC 2013), carbon-normalized SGVs for Aroclor 1254, Aroclor 1260 and Total PCBs were developed (see **Appendix A** for details). For those locations where Class A and Class C criteria were exceeded, the PCB concentrations were normalized to the TOC content and the result compared to the carbon-normalized SGVs.

- Concentrations of total PCBs at F-1 (0 to 6 inches and 18 to 24 inches) exceed the EqP sediment guidance values indicating a probable adverse effect to the benthic community in the area of F-1. Further risk management action is required for sediments in the area of F-1.
- Concentrations of total PCBs at all other locations where the Class A screening value was exceeded did not exceed the EqP sediment guidance values indicating that risk to the benthic community at those sites is acceptable and no further action is necessary to protect benthic communities from PCB exposure.

**Volatile Organic Compounds:** Samples were analyzed using US EPA method SW8260C for VOCs. The VOCs were screened individually against NYSDEC Class A criteria and the results are presented in **Table 3**. Thirteen VOCs detected in sediment have no NYSDEC Class A criteria. Using the EqP method provided in State guidance (NYSDEC 2013), the missing criteria were developed for 10 of the 13 (**Appendix A**). Chloroform is the only VOC that exceeds a Class A screening criterion. Chloroform was detected at 0.00077 mg/kg at C-2 which slightly exceeds the screening criterion of 0.0007 mg/kg. The relatively small exceedance and low frequency of detection indicate that chloroform exposure is not likely to cause unacceptable adverse effects to benthic communities. Therefore, no VOCs are expected to cause adverse effects to benthic communities and no further action is required regarding VOCs in sediment at the AOC G site.

**Semi-Volatile Organic Compounds:** SVOCs, except PAHs, were analyzed using US EPA method SW8270D. The SVOCs were screened individually against NYSDEC Class A criteria and the results are presented in **Table 4**. Five SVOCs detected in sediment have no NYSDEC Class A criteria. Using the EqP method provided in State guidance (NYSDEC 2013), the 5 missing criteria were developed (**Appendix A**). No SVOCs exceed Class A screening criteria. Therefore, SVOCs are not expected to cause adverse effects to benthic communities and no further action is required regarding SVOCs in sediment at the AOC G site.

**Polycyclic Aromatic Hydrocarbons:** PAHs were analyzed using US EPA method SW8270C SIM for 34 PAHs (18 specific non-alkylated compounds and 16 generic alkylated forms). The PAH

concentrations were screened individually against NYSDEC Class A criteria and the results are presented in **Table 5**. Because PAHs represent a family of compounds with a common mode of toxicity, PAH toxicity is also assessed on a combined basis. NYSDEC proposed the following approach to assess the potential toxic effect expected from exposure to combined PAH concentrations:

- Concentrations of all 34 PAHs and TOC are measured in a sediment sample as micrograms per kilogram ( $\mu\text{g}/\text{kg}$ ) of sediment.
- The concentration of each PAH detected is normalized to the percent of TOC in the sediment to produce a concentration of each PAH detected in units of  $\mu\text{g PAH}/\text{gram (g) TOC}$ .
- The concentration of each individual PAH present is divided by its corresponding SGV (Table 2, Column 5 in NYSDEC 2013). This quotient is described as a Toxic Unit (TU). In this context, a TU is essentially the same as a SGV quotient.
- The resulting TUs for each individual PAH are summed to produce a total TU for the mixture. If the total TU is greater than 1.0, the sediment is considered to be potentially toxic. The PAH concentrations at all locations and depths were screened as TUs and are presented in **Table 5**.
- The total TU for the 34-compound PAH mixture did not exceed 1 at any sample location, indicating that combined PAHs at the AOC G landfill site are not expected to cause adverse effects to benthic communities.
- No further action is required regarding PAHs in sediment at the AOC G landfill site.

### 2.3 Data Usability Assessment

The laboratory data collected during this project was reviewed by a third party chemist to provide an independent verification of its validity and usability. The Data Usability Summary Reports (DUSRs) for this project's Sample Delivery Groups are attached as **Appendix B**. In the opinion of the data validator, the tabulated data attached to this report is acceptable for its intended purpose, as qualified. A few non-detectable analysis results were rejected, as discussed below:

- The values for 2,4-Dinitrophenol (an SVOC) in two samples collected on November 6, 2013 were rejected due to low matrix spike/matrix spike duplicate (MS/MSD) recoveries.
- The non-detect SVOC results for one sample collected on November 7, 2013 were rejected due to exceeded holding times. The PCB results for another sample on this date (equipment blank) were rejected due to low surrogate recoveries.
- Several VOC results for two samples collected on November 8, 2013 were rejected due to low MS/MSD recoveries. The 2,4-Dinitrophenol result for one of these samples was also rejected due to low MS/MSD recoveries.

In AECOM's opinion, these deficiencies do not impact the assessment of sediment quality presented in this report for the following reasons:

- The compound 2,4-Dinitrophenol was not detected in any of the samples collected during the field event, so the rejection of its reported non-detect result for the two samples discussed above does not affect the assessment of this compound.
- Although the non-detect SVOC results for the November 7 sample were rejected, the compounds detected in this sample were at low concentrations, well below the Class A criteria. The rejected non-detect PCB results for the equipment blank on November 7 were not a concern, as PCBs were not detected in this or any of the other equipment blanks analyzed during the sampling event.
- Similarly, the rejected VOC and 2,4-Dinitrophenol non-detect results on November 8 do not affect the assessment of these compounds, as the specific compounds for which the data was rejected were not detected in any other samples during the sampling event, with the exception

of one VOC in a single sample. This detection was at a low concentration, well below the Class A criterion for the compound.

### 3.0 Summary of the AOC G November 2013 Wetland Investigation

The wetland delineation and sampling of the wetland area was completed during the week of November 4, 2013 as reported in this memorandum and summarized below.

- AECOM delineated the jurisdictional boundary between the upland portions of AOC G and the adjacent wetlands using the method documented in the Corps of Engineers Wetlands Delineation Manual (Technical Report Y-87-1, 1987), as modified by Supplement V2 (ERDC/EL TR-12-1, 2012), based on the presence of hydrophytic vegetation, hydric soils, and wetland hydrology.
- Sediment samples were collected from three depth intervals (0-6 inches, 6-12 inches, and 18-24 inches) from forty-three sediment sample locations.
- Each sediment sample was analyzed for VOCs, TOC, SVOCs, PCBs, Metals, percent moisture, and T34 PAHs.

Concentrations of analytes detected in the samples were compared to Class A screening values from Table 1a in the draft NYSDEC sediment guidance document (*Screening and Assessment of Contaminated Sediment*, NYSDEC, DFWMR, Bureau of Habitat, January 24, 2013).

- Metals exceeding the Class A criteria included arsenic, cadmium, chromium, copper, lead, nickel, mercury, and zinc. No metals exceeded NYSDEC Class C criteria at any depth. Therefore, all metals that exceed Class A criteria are considered Class B metals. Site-specific background concentration ranges for metals may be developed from the existing data set or, alternatively, sediment samples from locations considered to be similar to the geological environment of the site, but not impacted by metal contamination, could be sampled and analyzed to establish background ranges for site metals.
- PCBs were detected at the transect sample locations closest to the toe of the landfill, except at the location (SDC1) in the wetland-upland complex. Exceedances of NYSDEC Class A criteria were found at locations SDA4, SDB1, SDF1, SDH1, SDJ1, and SDK1. The distribution of these exceedances suggests a landfill source of PCBs.
- PCBs were also detected in excess of the Class A criteria at 2 locations within the floodplain of Sanders Creek, S-8 and S-9. Because no PCBs were detected in the swale between the AOC G landfill and Sanders Creek, the source of PCBs in the floodplain is assumed to be sediments from Sanders Creek deposited in the floodplain during flooding events that overtopped the creek banks.
- PCBs exceeded the Class C criteria at only one location: SDF1.
- Chloroform, detected at 0.00077 mg/kg at SDC2, was the only VOC that exceeded a Class A screening criterion. Although slightly exceeding the screening criterion, due to its minimal exceedance and low frequency of detection, chloroform exposure is not likely to cause unacceptable adverse effects to benthic communities. Therefore, no VOCs are expected to cause adverse effects to benthic communities and no further action is required regarding VOCs in sediment at the site.
- No SVOCs exceed Class A screening criteria. Therefore, SVOCs are not expected to cause adverse effects to benthic communities and no further action is required regarding SVOCs in sediment at the AOC G site.
- The evaluation of PAHs indicates that PAHs at the site are not expected to cause adverse effects to benthic communities. No further action is required regarding PAHs in sediment at the AOC G landfill site.

## 4.0 Closing

AECOM looks forward to your comments and questions regarding this report. Please feel free to contact Daniel Servetas at (518) 951-2378.

Sincerely,

AECOM USA, Inc.



Daniel Servetas, P.E.  
Project Manager  
[daniel.servetas@aecom.com](mailto:daniel.servetas@aecom.com)



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cc: K. Anders (NYSDOH)  
M. Sergott (NYSDOH)  
J. Wolski (UTC)  
K. McFadden (UTC)  
J. Alberg (AECOM)  
T. Schwendeman (AECOM)

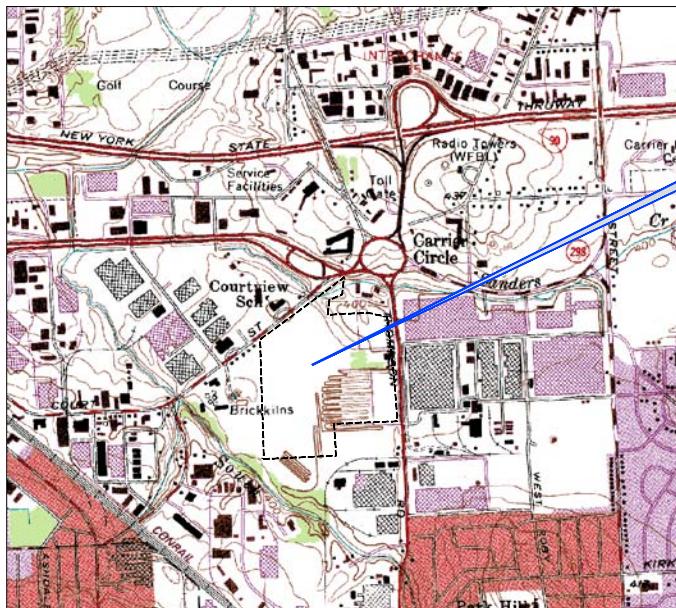
Enc: Figure 1, Site Plan  
Figure 2, Field Activity Summary (D-size)  
Figure 3, Distribution of Metals in Sediments at AOC G  
Figure 4, NYSDEC PCB Sediment Criteria Exceedances at AOC G

Table 1, Screening of Metals in Sediment by Depth  
Table 2, Distribution of PCBs in Sediment by Depth  
Table 3, VOC Screening Table  
Table 4, SVOC Screening Table  
Table 5, PAH Screening Table with Calculated Toxicity Units

Appendix A - Use of Equilibrium Partitioning Theory to Develop Sediment Guidance Values for Nonpolar Organic Compounds

Appendix B - Data Usability Summary Reports

## **Figures**



### SITE LOCATION MAP

NOT TO SCALE



Map References:

1) Base mapping from a field survey and plans titled "Map Showing Existing Conditions, Former Carrier Landfill, AOC G and Sanders Creek, by Thew Associates, dated January 21, 2014, including all notes and references therein.

2) The Topographic and planimetric features shown hereon were mapped utilizing photogrammetric mapping techniques. The aerial photography was obtained on November 16, 2013. A field edit was performed on January 3, 2014 to verify the accuracy and completeness of the photogrammetric mapping.

3) Field observations made by AECOM personnel, various dates.

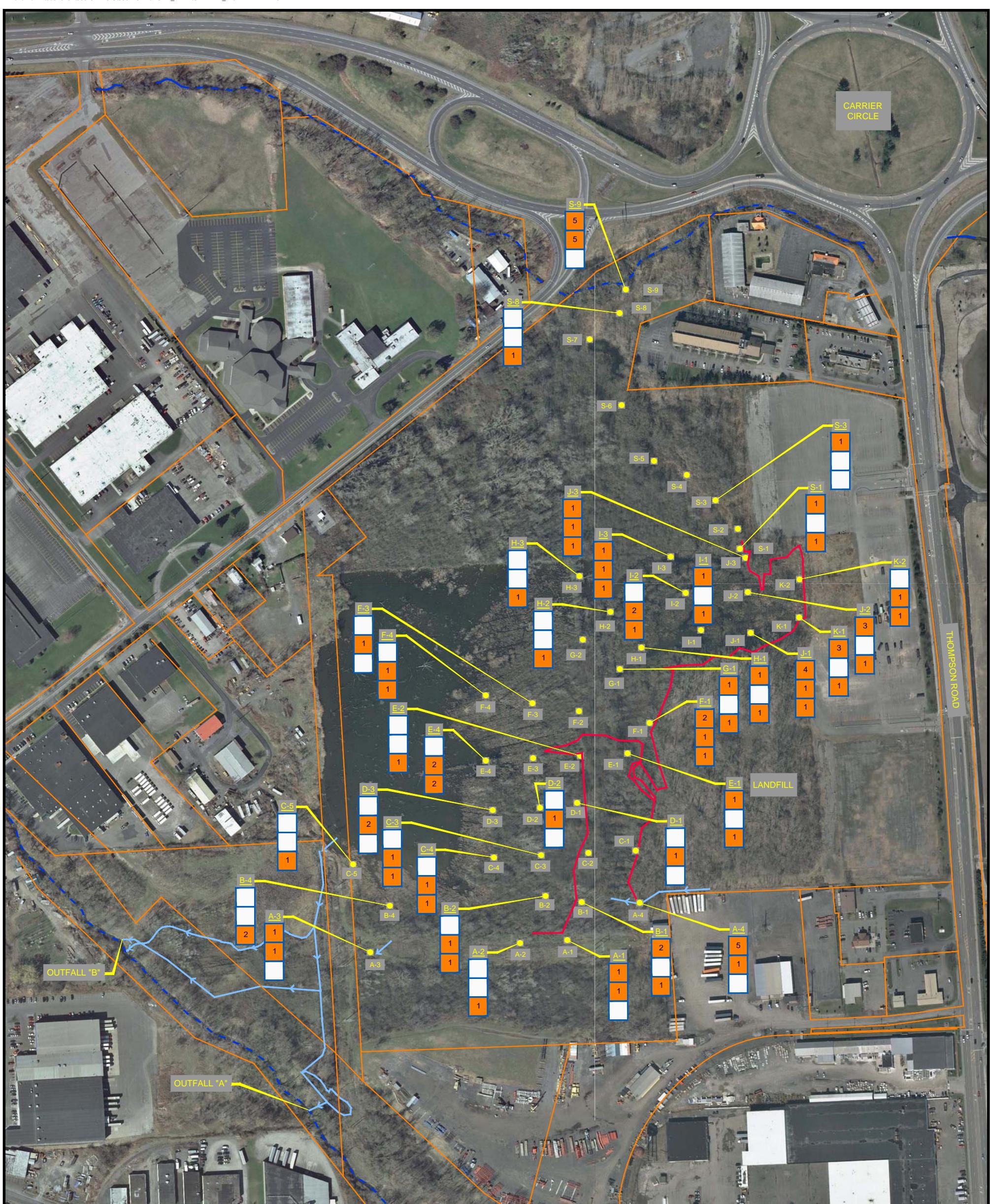
### SITE PLAN

NORTH

0' 75' 150' 300'

Issue Status: Final





Map Reference:  
 Photo from NYS GIS Orthos Online.

#### LEGEND

	APPROXIMATE PROPERTY BOUNDARY (TAX MAP)
	CREEK
	DRAINAGE PATH AND DIRECTION
	WETLAND LINE
	SAMPLE LOCATION (SEE NOTE)

#### SAMPLE KEYS

SAMPLE DEPTH		SAMPLE CONCENTRATION	
	A-4	NO EXCEDANCE AT INTERVAL	
		EXCEEDANCE WITH NUMBER OF METALS	
		EXCEEDING CLASS "A" CRITERIA	2



0° 75° 150° 300°

NOTE:  
 NO DETECTIONS UNLESS IDENTIFIED IN ASSOCIATED CALL-OUT BOX.

Issue Status: Final



Map Reference:  
 Photo from NYS GIS Orthos Online.

#### LEGEND

	APPROXIMATE PROPERTY BOUNDARY (TAX MAP)
	CREEK
	DRAINAGE PATH AND DIRECTION
	WETLAND LINE
	SAMPLE LOCATION (SEE NOTE)

#### SAMPLE KEYS

SAMPLE DEPTH	SAMPLE NAME	SAMPLE CONCENTRATION
0" - 6"	A-4	DETECTION, NO EXCEDENCE
6" - 12"		CLASS "A"
18" - 24"		CLASS "C"

TOTAL PCB CONCENTRATION AT DEPTH INTERVAL



0' 75' 150' 300'

NOTE:  
 NO DETECTIONS UNLESS IDENTIFIED IN ASSOCIATED CALL-OUT BOX.

Issue Status: Final

## **Tables**

Table 1  
Screening of Metals in Sediment by Depth  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria	Transect A				Transect B				Transect C				Transect D			Transect E			Transect F						
		SDA1	SDA2	SDA3	SDA4	SDB1	SDB2 (Dup)	SDB2	SDB4	SDC1	SDC2	SDC3	SDC4	SDC5	SDD1	SDD2	SDD3	SDE1	SDE2	SDE3	SDE4	SDF1	SDF2	SDF3	SDF4	
Sample ID		11/6/2013	11/6/2013	11/7/2013	11/8/2013	11/6/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/6/2013	11/8/2013	11/7/2013	11/6/2013	11/7/2013	11/8/2013	11/6/2013	11/6/2013	11/6/2013	11/8/2013	11/8/2013	11/4/2013	11/7/2013	11/8/2013		
Sample Date																										
Chemical Name	CLASS A	CLASS C																								
<b>Depths: 0-6 inches</b>																										
ALUMINUM	--	--	13200 J	12900 J	12400 J	9020 J	15000 J	12900	13400	8800 J	9890 J	10900 J	11100 J	15100	12500 J	11800 J	12800	12400	12300 J	8180 J	10600	6310	14300 J	12200	11600	11000
ANTIMONY	--	--	< 1.1 UJ	< 0.92 UJ	< 0.94 UJ	< 1.2 UJ	< 1.2 UJ	0.20 J	0.20 J	< 1.1 UJ	< 0.68 UJ	< 0.74 UJ	< 0.90 UJ	< 1.0 UJ	0.23 J	< 0.83 UJ	< 1.0 UJ	< 0.97 UJ	< 0.83 UJ	< 0.81 UJ	< 1.0 UJ	0.26 J	< 1.1 UJ	< 0.29 J	0.25 J	
ARSENIC	<10	<33	4.3	4.2	6.5	8.8	10.3	4.5	4.7	3.1	2.5	3.2	4.4	6.3	4.3	5.0	4.9	4.8	4.8	3.9	6.6	5.2	7.4	5.6	4.3	3.4
BARIUM	--	--	119 J	87.9 J	88.7 J	90.5 J	134 J	85.4 J	86.8 J	70.2 J	71.5 J	76.5 J	71.5 J	103	89.0 J	85.4 J	82.1 J	82.3	82.0 J	71.8 J	73.9	68.7	113 J	95.3 J	99.0	90.6
BERYLLIUM	--	--	0.57 J	0.53 J	0.58 J	0.62 J	0.74 J	0.59 J	0.62 J	0.41 J	0.45 J	0.48 J	0.49 J	0.69	0.54 J	0.53 J	0.53 J	0.52 J	0.40 J	0.54	0.42	0.69 J	0.58 J	0.56	0.53	
CADMUM	<1	<5	< 0.44 U	< 0.37 U	0.066 J	1.5	0.17 J	0.13 J	0.21 J	0.13 J	0.48	0.21 J	0.25 J	0.18 J	0.24 J	0.43	0.094 J	0.17 J	0.041 J	0.33	0.29 J	0.13 J	0.10 J	0.12 J	0.16 J	0.20 J
CALCIUM	Nutrient	86400 J	21500 J	26400 J	163000 J	50400 J	4330	3890	4780 J	3760 J	3290 J	8800 J	4590	3260 J	4840 J	2970	2740	78500 J	4810 J	3310	3890	16900 J	22100	13700	5950	
CHROMIUM, TOTAL	<43	<110	18.9 J	17.9 J	15.8 J	18.8 J	21.3 J	18.3 J	18.6 J	12.1 J	14.9 J	15.7 J	16.2 J	19.0 J	16.6 J	17.7 J	16.6 J	15.7 J	18.5 J	13.3 J	15.2 J	10.4 J	36.3 J	17.4 J	16.3 J	16.1 J
COBALT	--	--	9.8 J	6.6 J	11.3	8.3	13.9 J	6.8 J	6.9 J	5.7 J	5.7 J	6.6 J	5.8 J	6.7	7.5 J	5.8 J	8.4 J	7.8 J	5.6 J	8.8 J	4.5 J	12.1 J	8.3 J	6.9 J		
COPPER	<32	<150	22.7	19.4	25.7	39.6	29.5	20.3	20.4	18.9 J	17.4 J	16.8 J	14.8	22.5	17.1 J	10.7	13.2	22.8 J	20.9 J	22.2	24.3	28.7 J	21.3	18.2	19.4	
IRON	--	--	20000	18000	20700 J	15200 J	24800	17700 J	18400 J	13400 J	12700 J	13500 J	16600 J	29500 J	17400 J	17500 J	18100 J	17400 J	17800 J	12400 J	17900 J	11700 J	24100	19000 J	15100 J	
LEAD	<36	<130	9.1 J	14.3 J	12.0	45.4	21.9 J	27.3 J	28.0 J	22.6	27.0 J	28.9 J	21.1 J	25.2	30.2	28.4 J	20.1 J	20.5	13.3 J	25.4 J	25.6	28.9	25.0 J	23.2 J	20.4	20.3
MAGNESIUM	Nutrient	31300	11400	16700 J	14700 J	23700	4200 J	4340	3830 J	3100 J	3680 J	7030 J	4360	3950 J	4670 J	3760 J	3740	32200 J	3470 J	3500	2120	10500 J	14000 J	9710	5360	
MANGANESE	--	--	417 J	216 J	672 J	562 J	1270 J	288 J	249 J	157 J	105 J	154 J	232 J	541 J	180 J	301 J	156 J	290 J	330 J	188 J	264 J	143 J	434 J	359 J	331 J	201 J
MERCURY	<0.2	<1	0.070	0.060	0.035 J	0.17	0.11	0.10	0.071	0.085	0.12	0.078	0.12	0.086	0.098	0.089	0.069	0.095	0.083	0.076	0.28	0.074	0.066	0.084		
NICKEL	<23	<49	25.6 J	21.3 J	25.4 J	23.2 J	33.4 J	16.1 J	16.9 J	13.4 J	14.5 J	14.7 J	17.1 J	17.4 J	16.1 J	18.2 J	14.4 J	15.5 J	23.0 J	15.4 J	18.1 J	12.2 J	32.8 J	21.0 J	18.9 J	18.4 J
POTASSIUM	Nutrient	1780 J	1180 J	1150 J	1610 J	1680 J	1400 J	1320 J	851 J	1070 J	1100 J	1090 J	1350 J	1130 J	1240 J	1230 J	1180 J	1200 J	1180 J	1200 J	1250 J	1990 J	1370 J	1150 J	1200 J	
SELENIUM	--	--	< 1.1 U	< 0.92 U	< 0.94 U	< 1.2 U	< 1.2 U	< 0.80 J	0.89 J	0.54 J	0.44 J	< 0.90 U	< 1.0 U	0.43 J	0.52 J	0.85 J	0.40 J	< 0.83 U	0.50 J	0.35 J	1.0	< 1.1 U	0.39 J	0.56 J	0.56 J	
SILVER	<1	<2.2	< 0.55 U	< 0.46 U	< 0.47 U	0.21 J	< 0.60 U	< 0.53 U	< 0.50 U	< 0.57 U	< 0.34 U	< 0.37 U	< 0.45 U	< 0.51 U	0.13 J	< 0.41 U	< 0.52 U	< 0.49 U	< 0.41 U	< 0.40 U	< 0.50 U	< 0.43 U	< 0.56 U	< 0.53 U	< 0.53 U	< 0.46 U
SODIUM	Nutrient	< 550 U	< 460 U	< 470 U	< 460 U	< 520 U	< 600 U	< 530 U	< 500 U	< 570 U	< 370 U	< 450 U	< 510 U	< 500 U	< 410 U	< 520 U	< 490 U	< 410 U	< 400 U	< 500 U	< 430 U	68.1 J	< 530 U	< 460 U		
THALLIUM	--	--	0.38 J	0.45 J	< 0.94 U	< 1.2 U	0.55 J	< 1.1 U	< 1.0 U	0.28 J	0.23 J	0.50 J	0.27 J	< 0.99 U	0.16 J	< 1.0 U	0.14 J	< 0.81 U	0.17 J	< 1.0 U	0.28 J	0.37 J	< 1.1 U</td			

Table 1  
Screening of Metals in Sediment by Depth  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria	Transect G		Transect H			Transect I			Transect K		North Swale to Sanders Creek							Sanders Creek				
Sample ID		SDG1	SDG2	SDH1	SDH2	SDH3	SDI1 (Dup)	SDI1	SDI2	SDI3	SDK1	SDK2	SDS1	SDS2	SDS3	SDS4	SDS5	SDS6	SDS7	SDS8 (Dup)	SDS8	SDS9	
Sample Date		11/6/2013	11/7/2013	11/6/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/7/2013	11/5/2013	11/5/2013	11/5/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	
Depths: 0-6 inches																							
ALUMINUM	--	--	14400 J	11600	1870 J	12500	18800	16000 J	12800 J	2420	16300	4480	8840	13700 J	4210 J	14000 J	9680 J	7830 J	11400 J	7090 J	8250 J	7360 J	8420 J
ANTIMONY	--	--	<1.0 UJ	<1.1 UJ	0.49 J	<1.2 UJ	<1.0 UJ	<1.1 UJ	<1.1 UJ	7.7 J	<1.1 UJ	0.32 J	<1.1 UJ	0.44 J	<1.0 UJ	<1.1 UJ	<1.2 UJ	0.27 J	<1.0 UJ	<0.96 UJ	0.15 J	<0.98 UJ	
ARSENIC	<10	<33	8.2	9.4	2.3	4.2	4.8	6.1	5.3	<8.2 U	4.3	2.8	3.5	7.0	3.4	12.0	5.6	6.0	4.4	4.7	9.0	5.6	12.0
BARIUM	--	--	116 J	100 J	34.0 J	80.6 J	125 J	110 J	94.6 J	56.8 J	91.2 J	46.9 J	72.0 J	101 J	44.2 J	92.1 J	91.0 J	61.3 J	105 J	52.3 J	64.0 J	58.2 J	90.8 J
BERYLLIUM	--	--	0.71 J	0.54 J	0.094 J	0.52 J	0.72 J	0.75 J	0.62 J	<3.3 U	0.79	0.28 J	0.42 J	0.67 J	0.21 J	0.56 J	0.46 J	0.38 J	0.54 J	0.33 J	0.38 J	0.39 J	0.49 J
CADMUM	<1	<5	<0.41 U	0.063 J	0.67	0.081 J	0.41	<0.46 U	<0.46 U	0.90 J	<0.44 U	0.41 J	0.21 J	0.56	0.30 J	0.069 J	0.26 J	0.30 J	0.17 J	<0.42 U	<0.38 U	0.23 J	0.80
CALCIUM	Nutrient		30700 J	2630	14000 J	2710	4170	13500 J	5130 J	21900 J	2730 J	44900 J	4370 J	2730 J	3200 J	10600 J	5770 J	5140 J	2990 J	53100 J	34000 J	26400 J	23600 J
CHROMIUM, TOTAL	<43	<110	20.7 J	15.8 J	8.8 J	18.0 J	25.0 J	22.5 J	19.2 J	6.4 J	25.6 J	47.4 J	13.1 J	20.3 J	7.2 J	18.0 J	14.4 J	10.3 J	15.0 J	10.1 J	13.1 J	14.6 J	44.1 J
COBALT	--	--	13.8 J	15.1 J	1.5 J	8.8 J	10.4 J	11.1 J	10.1 J	1.1 J	12.1 J	2.6 J	5.8 J	13.6 J	3.5 J	8.2 J	6.3 J	3.9 J	6.9	5.5	6.2	6.2	9.6
COPPER	<32	<150	26.8	12.1	16.9	14.8	22.4	23.1	19.2	16.6 J	23.2	14.6	17.3	18.1 J	17.8 J	18.1 J	20.8 J	15.9 J	15.9	22.1	18.7	24.0	46.0
IRON	--	--	22800	27500 J	4410	19600 J	20300 J	30600	22200	4880 J	30200 J	7960 J	14000 J	22900	11500	19800	14700	11700	15800 J	12800 J	16000 J	13600 J	17000 J
LEAD	<36	<130	17.5 J	24.5 J	32.8 J	29.8 J	35.7 J	13.0 J	16.1 J	13.1 J	20.2	31.4	17.8	12.9 J	17.8 J	16.6 J	28.6 J	22.4 J	26.5 J	16.9 J	9.5 J	27.4 J	54.7 J
MAGNESIUM	Nutrient		20800	3250 J	1890	3640 J	4790 J	11100	5020	2250 J	5450 J	11000 J	3110 J	4880 J	1920 J	7220 J	3030 J	2080 J	3150 J	32800 J	15600 J	12900 J	13400 J
MANGANESE	--	--	754 J	539 J	160 J	231 J	188 J	1010 J	389 J	240 J	219 J	202 J	298 J	1210 J	281 J	523 J	1420 J	243 J	226 J	475 J	276 J	298 J	841 J
MERCURY	<0.2	<1	0.075	0.10	0.33	0.089	0.16	0.090	0.057	0.072 J	0.069	0.26	0.085	0.045	0.039	0.10	0.12	0.085	0.091	0.054	0.048	0.081	0.086
NICKEL	<23	<49	31.5 J	15.0 J	13.1 J	17.2 J	21.3 J	28.7 J	24.1 J	4.2 J	29.8 J	13.6 J	14.8 J	25.6 J	8.2 J	17.4 J	14.3 J	10.6 J	16.2 J	12.5 J	14.7 J	15.3 J	21.8 J
POTASSIUM	Nutrient		1360 J	1040 J	415 J	1330 J	1700 J	1500 J	1100 J	1910 J	587 J	976 J	1920 J	624 J	1650 J	1050 J	994 J	977 J	1070 J	1090 J	1070 J	981 J	
SELENIUM	--	--	<1.0 U	0.77 J	1.6	0.49 J	0.81 J	<1.1 U	<1.1 U	<8.2 U	<1.1 U	1.6	0.40 J	3.2	1.7	<1.1 U	0.56 J	0.64 J	0.60 J	<1.0 U	<0.96 U	<1.0 U	<0.98 U
SILVER	<1	<2.2	<0.51 U	<0.53 U	<0.58 U	<0.58 U	<0.51 U	<0.57 U	<0.57 U	<4.1 U	<0.55 U	<0.59 U	<0.53 U	<0.51 U	<0.50 U	<0.57 U	<0.59 U	<0.61 U	<0.53 U	<0.52 U	<0.48 U	0.15 J	0.86
SODIUM	Nutrient		<510 U	<530 U	<580 U	<580 U	<510 U	<570 U	<4100 U	<550 U	<530 U	72.8 J	82.5 J	124 J	220 J	<530 U	40.2 J	<480 U	<500 U	<490 U			
THALLIUM	--	--	0.27 J	<1.1 U	<1.2 U	<1.2 U	<1.0 U	0.47 J	0.46 J	<8.2 U	0.38 J	0.28 J	0.33 J	1.0	0.55 J	0.39 J	0.66 J	0.30 J	0.61 J	0.67 J	0.52 J	0.51 J	0.56 J
VANADIUM	--	--	27.1 J	24.3 J	8.3 J	23.4 J	34.8 J	28.2 J	24.7 J	6.9 J	30.8 J	67.3 J	17.5 J	25.3 J	9.9 J	23.2 J	17.9 J	15.3 J	20.9 J	15.0 J	15.8 J	16.7 J	22.1 J
ZINC	<120	<460	57.4 J	53.7 J	34.1 J	62.7 J	97.7 J	55.9 J	53.9 J	89.3 J	73.1 J	153 J	71.5 J	62.6 J	29.9 J	66.6 J	69.3 J	54.2 J	60.8 J	36.9 J	64.1 J	147 J	
Depths: 6-12 inches																							
ALUMINUM	--	--	11000 J	10400	9830 J	11600	11300	10600 J	13200 J	9260	14000	14400	17100	12700 J	6040 J	15600 J	13100 J	13800 J	11200 J	7980 J	8170 J	8440 J	10500
ANTIMONY	--	--	<0.86 UJ	<1.1 UJ	<0.96 UJ	<1.0 UJ	<0.97 UJ	<1.0 UJ	<1.0 UJ	1.3 J	<1.0 UJ	<1.0 UJ	<0.94 UJ	0.41 J	0.15 J	<1.1 UJ	<1.1 UJ	<1.2 UJ	0.26 J	<0.83 UJ	<1.0 UJ	<1.1 UJ	
ARSENIC	<10	<33	4.3	3.8	2.9	2.9	3.4	3.9	5.1	3.3 J	3.9	4.8	7.9	6.3	3.9	5.6	8.8	5.1	4.3	5.6	5.5	6.4	13.9
BARIUM	--	--	93.8 J	71.5 J	64.8 J	83.8 J	77.1 J	87.5 J	96.1 J	81.7 J	87.0 J	128 J	154 J	89.5 J	46.4 J	82.0 J	93.9 J	83.2 J	106 J	63.9 J	62.9 J	67.0 J	77.8 J
BERYLLIUM	--	--	0.50 J	0.42 J	0.45 J	0.46 J	0.46 J	0.58 J	0.61 J	0.63 J	0.63	0.63	0.82	0.64 J	0.29 J	0.60 J	0.65 J	0.63 J	0.53 J	0.38 J	0.43 J	0.44 J	0.52 J
CADMUM	<1	<5	<0.34 U	<0.43 U	<0.39 U	<0.40 U	<0.42 U	0.087 J	<0.41 U	0.99 J	0.073 J	<0.41 U	<0.38 U	0.46	0.14 J	<0.45 U	0.078 J	0.21 J	0.11 J	<0.33 U	0.21 J	0.24 J	0.56
CALCIUM	Nutrient		39500 J	2350	4300 J	2130	2030	6690 J	4160 J	12500 J	2410 J	3710 J	3150 J	4550 J	4070 J	2600 J	4310 J	4190 J	2520 J	25000 J	28500 J	29100 J	
CHROMIUM, TOTAL	<43	<110	16.1 J	13.5 J	21.1 J	15.3 J	15.0 J	15.7 J	18.5 J	21.9 J	19.4 J	18.7 J	25.1 J	19.5 J	8.5 J	20.1 J	16.6 J	14.0 J	11.3 J	16.4 J	16.5 J	37.7 J	
COBALT	--	--	7.8 J	6.6 J	6.8 J	9.9 J	8.4 J	8.0 J	9.3 J	4.5 J	10.6 J	8.3 J	20.8 J	9.8 J	5.6 J	9.2 J	7.1 J	6.1 J	6.7	6.9	7.0	8.0 J	
COPPER	<32	<150	18.8 J	8.0	14.6	9.2	9.8	13.0	32.1	16.7	11.7	21.5	23.2 J	18.9 J	17.1 J	16.2 J	17.5 J	15.1	20.7	28.0	42.9		
IRON	--	--	17400 J	15900 J	14200	15000	14600 J	13900 J	16100	15300	13300	14500 J	13200 J	10300 J	14700 J	8950 J	14400 J	15200 J	9390 J	7050 J	6770 J	7040	
LEAD	<36	<130	9.8 J	13.4 J	21.6 J	14.8 J	12.3 J	11.0 J	68.1 J	12.8 J	10.9 J	13.0	29.5 J	8.4 J	15.9 J	11.8 J	18.7 J	25.7 J	13.0 J	29.9 J	26.8 J	44.0	
MAGNESIUM	Nutrient		24300 J	2900 J	3720	3160 J	3090 J	5580	5030	3210	4140 J	3800 J	6360 J	4450 J	3600 J	4570 J	3730 J	3450 J	2990 J	11200 J	12500 J	14000 J	14900 J
MANGANESE	--	--	344 J	260 J	110 J	238 J	114 J	328 J	433 J	189 J	353 J	242 J	2490 J	1130 J	365 J	363 J	394 J	229 J	170 J	338 J	259 J	544 J	
MERCURY	<0.2	<1	0.030 J	0.029 J	0.030 J	0.033 J	0.029 J	0.046	0.048	0.19	0.044	0.11	0.066	0.094	0.020 J	0.09 J	0.095	0.12	0.088	0.042	0.079	0.32 J	
NICKEL	<23	<49	21.3 J	11.6 J	20.0 J	14.2 J	13.5 J	19.4 J	20.6 J	15.5 J	23.2 J	17.8 J	39.0 J	22.0 J	10.5 J	19.2 J	15.3 J	16.9 J	15.6 J	14.3 J	16.8 J	22.3 J	
POTASSIUM	Nutrient		1020 J	853 J	888 J	1070 J	1030 J	740 J	1100 J	1090 J	1470 J	1040 J	1690 J	1480 J	810 J	1510 J	1410 J	1210 J	964 J	1070 J	1180 J	1240 J	1190 J
SELENIUM	--	--	<0.86 U	0.61 J	0.56 J	0.42 J	0.68 J	<0.97 U	<1.0 U	1.7 J	<1.0 U	<1.0 U	<1.9 U	3.1	1.8	<1.1 U	0.50 J	0.44 J	0.58 J	<0.83 U	<1.0 U	0.39 J	<1.1 U
SILVER	<1	<2.2	<0.43 U	<0.54 U	<0.48 U	<0.50 U	<0.52 U	<0.49 U	<0.51 U	<1.8 U	<0.52 U	<0.51 U	<0.47 U	<0.55 U	<0.51 U	<0.56 U	<0.56 U	<0.61 U	<0.52 U	<0.41 U	0.20 J	0.21 J	0.88
SODIUM	Nutrient		<430 U																				

## Notes

All units are in milligrams per kilogram (mg/kg).

U or < - Metal was analyzed for, but not detected (i.e., concentration less than method detection limit). Presented with Gray text.

U or < - Me  
I Estimate

Black text indicates a detection (i.e., a concentration greater than the method detection limit).

Black text indicates a detection (i.e., a concentration greater than the method detection limit). Exceedances of NYSDEC Class A criteria are highlighted yellow and presented with **Bold** text.

Exceedances of NYSDEC Class A criteria are highlighted yellow and presented with **Bold** text.



Table 2  
 Distribution of PCBs in Sediment by Depth  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Screening Criteria			Transect E												Transect F				Transect G					
				SDE1			SDE2			SDE3			SDE4			SDF1		SDF2	SDF3	SDF4	SDG1		SDG2		
				11/6/2013			11/6/2013			11/8/2013			11/8/2013			11/4/2013		11/7/2013	11/8/2013	11/8/2013	11/6/2013		11/7/2013		
Chemical Name	Sediment Guidance Value ( $\mu\text{g/gOC}$ )	NYSDEC Sediment CLASS A	NYSDEC Sediment CLASS C	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	mg/kg	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	
			<b>Depths: 0-6 inches</b>																						
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	< 0.045 U	--	--	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	< 0.045 U	--	--	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	< 0.045 U	--	--	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	< 0.045 U	--	--	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	< 0.045 U	--	--	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1254 (AROCLOR 1254)	56.5	--	--	< 0.05 U	--	--	< 0.048 U	--	--	< 0.047 U	--	--	< 0.049 U	--	--	<b>4.31</b>	581.87	<b>10.30</b>	< 0.047 U	< 0.045 U	< 0.055 U	< 0.047 U	--	--	< 0.046 U
PCB-1260 (AROCLOR 1260)	86.9	--	--	0.0807	2.5619	0.0295	0.0260 J	0.7879	0.0091	0.0245 J	0.8333	0.0096	0.0359 J	0.5753	0.0066	<b>1.99 J</b>	269.28	<b>3.10</b>	< 0.047 U	< 0.045 U	< 0.055 U	0.0397 J	1.5095	0.0174	< 0.046 U
Total PCBs	70.9	0.1	1	0.0807	--	--	0.0260	--	--	0.0245	--	--	0.0359	--	--	<b>6.3</b>	--	--	< 0.047 U	< 0.045 U	< 0.055 U	0.0397	--	--	< 0.046 U
TOC	--	--	--	31500	--	--	33000	--	--	29400	--	--	62400	--	--	7390	--	--	21300	20800	20600	26300	--	--	19300
Toxicity Units (total)	--	--	--	--	--	0.0295	--	--	0.0091	--	--	0.0096	--	--	0.0066	--	--	<b>13.40</b>	--	--	--	--	--	0.0174	--
			<b>Depths: 6-12 inches</b>																						
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	< 0.04 U	--	--	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	< 0.04 U	--	--	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	< 0.04 U	--	--	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	< 0.04 U	--	--	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	< 0.04 U	--	--	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1254 (AROCLOR 1254)	56.5	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	<b>0.122</b>	7.13	0.13	< 0.046 U	< 0.041 U	< 0.043 U	< 0.042 U	--	--	< 0.046 U
PCB-1260 (AROCLOR 1260)	86.9	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	0.0644 J	3.77	0.04	< 0.046 U	< 0.041 U	< 0.043 U	0.0327 J	1.2433	0.0143	< 0.046 U
Total PCBs	70.9	0.1	1	< 0.042 U	--	--	< 0.043 U	--	--	< 0.041 U	--	--	< 0.055 U	--	--	<b>0.186</b>	--	--	< 0.046 U	< 0.041 U	< 0.043 U	0.0327	--	--	< 0.046 U
TOC	--	--	--	23900	--	--	16400	--	--	19700	--	--	69800	--	--	17100	--	--	15000	2420	9060	13100	--	--	13000
Toxicity Units (total)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	0.17	--	--	--	--	--	0.0143	--
			<b>Depths: 18-24 inches</b>																						
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.038 U	--	--	< 0.041 U	--	--	< 0.041 U	--	--	< 0.062 U	--	--	< 0.043 U	--	--	< 0.041 U	< 0.047 U	< 0.041 U	< 0.039 U	--	--	< 0.041 U
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.038 U	--	--	< 0.041 U	--	--	< 0.041 U	--	--	< 0.062 U	--	--	< 0.043 U	--	--	< 0.041 U	< 0.047 U	< 0.041 U	< 0.039 U	--	--	< 0.041 U
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.038 U	--	--	< 0.041 U	--	--	< 0.041 U	--	--	< 0.062 U	--	--	< 0.043 U	--	--	< 0.041 U	< 0.047 U	< 0.041 U	< 0.			

Table 2  
 Distribution of PCBs in Sediment by Depth  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Screening Criteria			Transect H						Transect I								
				SDH1			SDH2	SDH3			SDI1 (Dup)			SDI1				
				11/6/2013			11/7/2013	11/7/2013			11/6/2013			11/6/2013	11/7/2013	11/5/2013		
Chemical Name	Sediment Guidance Value ( $\mu\text{g/gOC}$ )	NYSDEC Sediment CLASS A	NYSDEC Sediment CLASS C	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	
<b>Depths: 0-6 inches</b>																		
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.12 U			< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.12 U			< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.12 U			< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.12 U			< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.12 U			< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1254 (AROCLOR 1254)	56.5	--	--	<b>0.195 J</b>	1.3732	0.0243	< 0.049 U	< 0.096 U	--	--	< 0.048 U	--	--	< 0.049 U	--	--	< 0.31 U < 0.056 U	
PCB-1260 (AROCLOR 1260)	86.9	--	--	<b>0.343</b>	2.4155	0.0278	< 0.049 U	<b>0.0506 J</b>	0.4217	0.0049	0.0468 J	1.8	0.0207	<b>0.0522</b>	1.8316	0.0211	< 0.31 U < 0.056 U	
Total PCBs	70.9	0.1	1	<b>0.538</b>	--	--	< 0.049 U	<b>0.0506</b>	--	--	0.0468	--	--	<b>0.0522</b>	--	--	< 0.31 U < 0.056 U	
TOC	--	--	--	142000	--	--	37400	120000	--	--	26000	--	--	28500	--	--	413000 32900	
Toxicity Units (total)	--	--	--	--	--	0.0521	--	--	--	0.0049	--	--	0.0207	--	--	0.0211	--	--
<b>Depths: 6-12 inches</b>																		
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1254 (AROCLOR 1254)	56.5	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
PCB-1260 (AROCLOR 1260)	86.9	--	--	< 0.044 U	--	--	< 0.044 U	< 0.046 U	--	--	< 0.043 U	--	--	< 0.046 U	--	--	< 0.14 U < 0.045 U	
Total PCBs	70.9	0.1	1	<b>&lt; 0.044 U</b>	--	--	<b>&lt; 0.044 U</b>	<b>&lt; 0.046 U</b>	--	--	<b>&lt; 0.043 U</b>	--	--	<b>&lt; 0.046 U</b>	--	--	< 0.14 U < 0.045 U	
TOC	--	--	--	11800	--	--	14500	17400	--	--	9710	--	--	17500	--	--	223000 5270	
Toxicity Units (total)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	
<b>Depths: 18-24 inches</b>																		
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1254 (AROCLOR 1254)	56.5	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	--	< 0.062 U < 0.044 U	
PCB-1260 (AROCLOR 1260)	86.9	--	--	< 0.041 U	--	--	< 0.044 U	< 0.041 U	--	--	< 0.042 U	--	--	< 0.043 U	--	0.0649	< 0.044 U	
Total PCBs	70.9	0.1	1	<b>&lt; 0.041 U</b>	--	--	<b>&lt; 0.044 U</b>	<b>&lt; 0.041 U</b>	--	--	<b>&lt; 0.042 U</b>	--	--	<b>&lt; 0.043 U</b>	--	0.0649	< 0.044 U	
TOC	--	--	--	4730	--	--	3370	3550	--	--	7990	--	--	6910	--	--	54400 3830	
Toxicity Units (total)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	

Notes:

U or < - Polychlorinated biphenyl (PCB) was analyzed for, but not detected (i.e., concentration less than method detection limit). Presented with Gray text.

J - Estimated.

Total PCBs values are calculated by summing the detected concentrations. If there are no detected concentrations, total PCBs is reported at the highest non-detect result.

TOC - Total Organic Carbon.

Black text indicates a detection (i.e., a concentration greater than the method detection limit).

Exceedances of NYSDEC Class A criteria are highlighted yellow and presented with **Bold** text.

Exceedances of NYSDEC Class C criteria are highlighted orange and presented with **Bold** text.

mg/kg - milligrams per kilogram.

$\mu\text{g/gOC}$  - microgram per gram of Organic Carbon.

TU - Toxicity Unit.

Table 2  
 Distribution of PCBs in Sediment by Depth  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Screening Criteria			Transect K			North Swale to Sanders Creek							Sanders Creek									
				SDK1		SDK2	SDS1	SDS2	SDS3	SDS4	SDS5	SDS6	SDS7	SDS8 (Dup)			SDS8		SDS9				
	11/5/2013			11/5/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/4/2013	11/5/2013	11/5/2013	11/5/2013			11/5/2013		11/5/2013				
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	NYSDEC Sediment CLASS C	mg/kg	µg/gOC	TU	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	
<b>Depths: 0-6 inches</b>																							
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.08 U	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	< 0.042 U	--	--	< 0.051 U	--	--	< 0.045 U	--	--	
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.08 U	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	< 0.042 U	--	--	< 0.051 U	--	--	< 0.045 U	--	--	
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.08 U	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	< 0.042 U	--	--	< 0.051 U	--	--	< 0.045 U	--	--	
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.08 U	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	< 0.042 U	--	--	< 0.051 U	--	--	< 0.045 U	--	--	
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.08 U	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	< 0.042 U	--	--	< 0.051 U	--	--	< 0.045 U	--	--	
PCB-1254 (AROCLOR 1254)	56.5	--	--	0.0777 J	0.7194	0.0127	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	0.0654	5.23	0.0602	0.0366 J	1	0.0115	0.231	13.05	0.1502	
PCB-1260 (AROCLOR 1260)	86.9	--	--	<b>0.171</b>	1.5833	0.0182	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	0.0654	--	--	0.0366	--	--	<b>0.28</b>	--	--	
Total PCBs	70.9	0.1	1	<b>0.249</b>	--	--	< 0.049 U	< 0.042 U	< 0.048 U	< 0.065 U	< 0.062 U	< 0.051 U	< 0.05 U	0.0654	--	--	0.0366	--	--	17700	--	--	
TOC	--	--	--	108000	--	--	24500	7430	23900	13600	51100	52900	36500	30300	12500	--	--	36600	--	--	17700	--	--
Toxicity Units (total)	--	--	--	--	--	0.0310	--	--	--	--	--	--	--	--	--	0.0602	--	--	0.0115	--	--	0.1502	
<b>Depths: 6-12 inches</b>																							
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	< 0.051 U	--	--	< 0.058 U	--	--
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	< 0.051 U	--	--	< 0.058 U	--	--
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	< 0.051 U	--	--	< 0.058 U	--	--
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	< 0.051 U	--	--	< 0.058 U	--	--
PCB-1248 (AROCLOR 1248)	--	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	< 0.051 U	--	--	< 0.058 U	--	--
PCB-1254 (AROCLOR 1254)	56.5	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	< 0.051 U	--	--	0.0372 J	1.248322	0.0221	<b>0.185 J</b>	5.51	0.0975
PCB-1260 (AROCLOR 1260)	86.9	--	--	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	0.0752	2.07	0.0238	<b>0.113</b>	3.791946	0.0436	<b>0.536</b>	15.95	0.1836
Total PCBs	70.9	0.1	1	< 0.046 U	--	--	< 0.042 U	< 0.052 U	< 0.045 U	< 0.047 U	< 0.046 U	< 0.051 U	< 0.052 U	< 0.04 U	0.0752	--	--	<b>0.15</b>	--	--	<b>0.721</b>	--	--
TOC	--	--	--	22700	--	--	6200	34500	6520	14500	29200	33100	26200	13400	36400	--	--	29800	--	--	33600	--	--
Toxicity Units (total)	--	--	--	--	--	--	--	--	--	--	--	--	--	--	0.0238	--	--	0.0657	--	--	0.2810	--	--
<b>Depths: 18-24 inches</b>																							
PCB-1016 (AROCLOR 1016)	--	--	--	< 0.045 U	--	--	< 0.043 U	< 0.04 U	< 0.052 U	< 0.049 U	< 0.044 U	< 0.046 U	< 0.047 U	< 0.038 U	< 0.041 U	--	--	< 0.041 U	--	--	< 0.042 U	--	--
PCB-1221 (AROCLOR 1221)	--	--	--	< 0.045 U	--	--	< 0.043 U	< 0.04 U	< 0.052 U	< 0.049 U	< 0.044 U	< 0.046 U	< 0.047 U	< 0.038 U	< 0.041 U	--	--	< 0.041 U	--	--	< 0.042 U	--	--
PCB-1232 (AROCLOR 1232)	--	--	--	< 0.045 U	--	--	< 0.043 U	< 0.04 U	< 0.052 U	< 0.049 U	< 0.044 U	< 0.046 U	< 0.047 U	< 0.038 U	< 0.041 U	--	--	< 0.041 U	--	--	< 0.042 U	--	--
PCB-1242 (AROCLOR 1242)	--	--	--	< 0.045 U	--	--	< 0.043 U	< 0.04 U	< 0.052 U	< 0.049 U	< 0.044 U	< 0.046 U	< 0.047 U	< 0.038 U	< 0.041 U</								

Table 3  
VOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect A										Transect B											
		SDA1	SDA1	SDA1	SDA2	SDA2	SDA3	SDA3	SDA4	SDA4	SDB1	SDB1	SDB1	SDB2 (Dup)	SDB2	SDB2	SDB4	SDB4	SDB4	SDB4	SDB4		
Sample ID		11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/8/2013	11/8/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013		
Sample Date		Sample Interval (inches)		Chemical Name		0 - 6		6 - 12		18 - 24		0 - 6		6 - 12		18 - 24		0 - 6		6 - 12		18 - 24	
1,1,1,2-TETRACHLOROETHANE	--	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 U	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,1,1-TRICHLOROETHANE	0.06*	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 U	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 U	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 U	< 0.0027 U	< 0.0020 U
1,1,2,2-TETRACHLOROETHANE	--	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 UJ	< 0.0020 UJ	< 0.0026 UJ	< 0.0030 UJ	< 0.0032 UJ	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 UJ	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 UJ	< 0.0027 U	< 0.0020 UJ
1,1,2-TRICHLOROETHANE	--	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 UJ	< 0.0020 UJ	< 0.0026 UJ	< 0.0030 UJ	< 0.0032 UJ	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 UJ	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 UJ	< 0.0027 U	< 0.0020 UJ
1,1-DICHLOROETHANE	0.0541*	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 U	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 U	< 0.0027 U	< 0.0020 U
1,1-DICHLOROETHENE	0.52	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 U	< 0.0021 UJ	< 0.0035 U	< 0.25 UJ	< 0.0024 UJ	< 0.0020 UJ	< 0.0033 U	< 0.0027 U	< 0.0020 U
1,1-DICHLOROPROPENE	--	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 U	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2,3-TRICHLOROBENZENE	1.72*	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 UJ	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2,3-TRICHLOROPROPANE	--	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 UJ	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2,4-TRICHLOROBENZENE	2.14*	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	R	R	< 0.0063 U	< 0.0067 U	< 0.0053 U	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2,4-TRIMETHYLBENZENE	3.4	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 U	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2-DIBROMO-3-CHLOROPROPANE	--	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 UJ	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U	
1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	--	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 UJ	< 0.0027 U	< 0.0020 U	
1,2-DICHLOROBENZENE	0.28	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 UJ	< 0.0053 UJ	R	R	< 0.0063 U	< 0.0067 U	< 0.0053 UJ	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.0050 U
1,2-DICHLOROETHANE	1.7	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 UJ	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 U	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 U	< 0.0027 U	< 0.0020 U	
1,2-DICHLOROPROPANE	--	< 0.0028 U	< 0.0022 U	< 0.0018 U	< 0.0024 U	< 0.0020 U	< 0.0026 U	< 0.0030 U	< 0.0032 U	< 0.0020 U	< 0.0053 UJ	< 0.0030 UJ	< 0.0025 U	< 0.0027 U	< 0.0021 U	< 0.0021 UJ	< 0.0035 U	< 0.25 U	< 0.0024 U	< 0.0020 U	< 0.0033 U	< 0.0027 U	< 0.0020 U
1,2,4-TRIMETHYLBENZENE (MESITYLENE)	0.832*	< 0.0071 U	< 0.0055 U	< 0.0045 U	< 0.0060 U	< 0.0049 U	< 0.0065 U	< 0.0076 U	< 0.0080 U	< 0.0051 U	< 0.013 UJ	< 0.0074 UJ	< 0.0063 U	< 0.0067 U	< 0.0053 U	< 0.0087 U	< 0.64 U	< 0.0059 U	< 0.0051 U	< 0.0082 U	< 0.0067 U	< 0.005	

**Table 3**  
**VOC Screening Table**  
**November 2013 Event**  
**UTC/Carrier Circle**  
**Carrier-Dewitt Landfill (AOC G)**  
**thompson Road, Syracuse, New York**

## Notes

VOC - Volatile Organic Compound

All units are in milligrams per kilogram (mg/kg).

\* UTC Developed criteria

- STC-Developed criteria.

U or < - Compound was analyzed for, but not detected (i.e., concentration less than 1. Estimated

J - Estimated.  
R - Concentration rejected during validation.

R - Concentration rejected during validation.

AECOM

Table 3  
VOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect E												Transect F												
		SDE1	SDE1	SDE1	SDE2	SDE2	SDE3	SDE3	SDE4	SDE4	SDF1	SDF1	SDF2	SDF2	SDF3	SDF3	SDF4									
		11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/4/2013	11/4/2013	11/7/2013	11/7/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	
Sample ID																										
Sample Date		0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	
Sample Interval (inches)																										
Chemical Name																										
1,1,1,2-TETRACHLOROETHANE	--	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U	
1,1,1-TRICHLOROETHANE	0.06*	0.0019 J	0.00088 J	0.0011 J	<0.0027 U	<0.0024 U	<0.0018 U	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U	
1,1,2,2-TETRACHLOROETHANE	--	<0.0028 UJ	<0.0024 UJ	<0.0018 UJ	<0.0027 UJ	<0.0024 UJ	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U		
1,1,2-TRICHLOROETHANE	--	<0.0028 UJ	<0.0024 UJ	<0.0018 UJ	<0.0027 UJ	<0.0024 UJ	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U		
1,1,2-DICHLOROETHANE	0.0541*	0.0045	0.0027	0.0034	0.0054	0.0126	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U		
1,1-DICHLOROETHENE	0.52	<0.0028 U	<0.0024 U	<0.0018 U	0.0011 J	0.0016 J	0.0013 J	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U	
1,1-DICHLOROPROPENE	--	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U		
1,2,3-TRICHLOROBENZENE	1.72*	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U		
1,2,3-TRICHLOROPROPANE	--	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U	
1,2,4-TRICHLOROBENZENE	2.14*	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U	
1,2,4-TRIMETHYLBENZENE (MESITYLENE)	3.4	0.0051 J	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U	
1,2-DIBROMO-3-CHLOROPROPANE	--	<0.0069 U	<0.0060 U	<0.0046 U	<0.0067 U	<0.0060 U	<0.0045 U	<0.0066 U	<0.0062 U	<0.0048 U	<0.0091 U	<0.0088 U	<0.012 U	<0.0058 U	<0.0046 U	<0.0051 U	<0.0065 U	<0.0062 U	<0.0050 U	<0.0068 U	<0.0062 U	<0.0056 U	<0.0098 U	<0.0054 U	<0.0052 U	
1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	--	<0.0028 U	<0.0024 U	<0.0018 U	<0.0027 UJ	<0.0024 UU	<0.0018 UJ	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U	
1,2-DICHLOROBENZENE	0.28	<0.0028 U	<0.0024 U	<0.0018 U	<0.0027 U	<0.0024 U	<0.0018 U	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U	
1,2-DICHLOROETHANE	1.7	<0.0028 U	<0.0024 U	<0.0018 U	<0.0027 U	<0.0024 U	<0.0018 U	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	<0.0023 U	<0.0018 U	<0.0020 U	<0.0026 U	<0.0025 U	<0.0020 U	<0.0027 U	<0.0025 U	<0.0022 U	<0.0039 U	<0.0022 U	<0.0021 U	
1,2-DICHLOROPROPANE	--	<0.0028 U	<0.0024 U	<0.0018 U	<0.0027 U	<0.0024 U	<0.0018 U	<0.0026 U	<0.0025 U	<0.0019 U	<0.0036 U	<0.0035 U	<0.0049 U	&												

Table 3  
VOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect G						Transect H						Transect I														
		SDG1	SDG1	SDG1	SDG2	SDG2	SDG2	SDH1	SDH1	SDH1	SDH2	SDH2	SDH3	SDH3	SDI1	SDI1 (Dup)	SDI1	SDI1 (Dup)	SDI1	SDI1 (Dup)	SDI2	SDI2	SDI2	SDI3	SDI3	SDI3		
		Sample Date	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013		
Chemical Name		0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	0 - 6	6 - 12	6 - 12	18 - 24	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24			
1,1,1,2-TETRACHLOROETHANE	--	<0.0070 U	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 UJ	<0.0063 U	<0.0045 U	<0.0088 U	<0.0058 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U
1,1,1-TRICHLOROETHANE	0.06*	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 U	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,1,2,2-TETRACHLOROETHANE	--	<0.0028 U	<0.0025 UJ	<0.0036 UJ	<0.0036 U	<0.0027 UJ	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 UJ	<0.0022 UJ	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 UJ	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,1,2-TRICHLOROETHANE	--	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 UJ	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,1-DICHLOROETHANE	0.0541*	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 U	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,1-DICHLOROETHENE	0.52	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 U	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,1-DICHLOROPROPENE	--	<0.0070 U	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U	
1,2,3-TRICHLOROBENZENE	1.72*	<0.0070 U	<0.0064 UJ	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U	
1,2,3-TRICHLOROPROPANE	--	<0.0070 U	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U	
1,2,4-TRICHLOROBENZENE	2.14*	<0.0070 UJ	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U	
1,2,4-TRIMETHYLBENZENE	3.4	<0.0070 U	<b>0.00038 J</b>	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U
1,2-DIBROMO-3-CHLOROPROPANE	--	<0.0070 U	<0.0064 U	<0.0090 U	<0.0089 U	<0.0068 U	<0.0047 U	<0.027 U	<0.0063 U	<0.0045 U	<0.0088 U	<0.0054 U	<0.014 U	<0.0061 U	<0.0062 U	<0.0070 U	<0.0059 U	<0.0058 U	<0.0052 U	<0.0062 U	<0.0056 U	<0.074 U	<0.033 U	<0.0095 U	<0.0073 U	<0.0060 U	<0.0061 U	
1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	--	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 U	<0.0058 U	<0.0025 UJ	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U	
1,2-DICHLOROBENZENE	0.28	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.0036 U	<0.0027 U	<0.0019 U	<0.011 U	<0.0025 U	<0.0018 U	<0.0035 U	<0.0023 U	<0.0022 U	<0.0058 U	<0.0025 U	<0.0025 U	<0.0028 U	<0.0024 U	<0.0023 U	<0.0021 U	<0.0025 U	<0.0022 U	<0.03 U	<0.013 U	<0.0038 U	<0.0029 U	<0.0024 U	<0.0025 U
1,2-DICHLOROETHANE	1.7	<0.0028 U	<0.0025 UJ	<0.0036 U	<0.003																							



Table 3  
VOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Sanders Creek								
		SDS8	SDS8 (Dup)	SDS8	SDS8 (Dup)	SDS8	SDS8 (Dup)	SDS9	SDS9	
		11/5/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	11/5/2013	
Sample Interval (inches)		0 - 6	0 - 6	6 - 12	6 - 12	18 - 24	18 - 24	0 - 6	6 - 12	18 - 24
Chemical Name										
1,1,1,2-TETRACHLOROETHANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,1,1-TRICHLOROETHANE	0.06*	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,1,2,2-TETRACHLOROETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,1,2-TRICHLOROETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,1-DICHLOROETHANE	0.0541*	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,1-DICHLOROETHENE	0.52	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,1-DICHLOROPROPENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,2,3-TRICHLOROBENZENE	1.72*	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,2,3-TRICHLOROPROPANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,2,4-TRICHLOROBENZENE	2.14*	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,2,4-TRIMETHYLBENZENE	3.4	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<b>0.00027 J</b>	<0.0055 U	<b>0.00046 J</b>	<0.0083 U	<0.0072 U
1,2-DIBROMO-3-CHLOROPROPANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,2-DICHLOROBENZENE	0.28	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,2-DICHLOROETHANE	1.7	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,2-DICHLOROPROPANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,3,5-TRIMETHYLBENZENE (MESITYLENE)	0.832*	0.00048 J	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<b>0.00057 J</b>	<0.0083 U	<0.0072 U
1,3-DICHLOROBENZENE	1.8	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
1,3-DICHLOROPROPANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
1,4-DICHLOROBENZENE	0.72	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
2,2-DICHLOROPROPANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
2-CHLOROTOLUENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
4-CHLOROTOLUENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
BENZENE	0.53	<b>0.00070</b>	<b>0.00073</b>	<b>0.0014</b>	<b>0.0011</b>	<b>0.00091</b>	<b>0.00090</b>	<b>0.00092</b>	<b>0.00094</b>	
BROMOBENZENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
BROMOCHLOROMETHANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
BROMODICHLOROMETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
BROMOFORM	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
BROMOMETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
CARBON TETRACHLORIDE	1.07	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
CHLOROBENZENE	0.2	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
CHLOROETHANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
CHLOROFORM	0.000715*	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
CHLOROMETHANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
CIS-1,2-DICHLOROETHYLENE	0.79*	<0.0026 U	<b>0.00086 J</b>	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<b>0.0044</b>	<b>0.0033</b>	<b>0.0011 J</b>
CYMENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
DIBROMOCHLOROMETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
DIBROMOMETHANE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
DICHLORODIFLUOROMETHANE	--	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
ETHYLBENZENE	0.43	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
HEXAACHLOROBUTADIENE	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
ISOPROPYLBENZENE (CUMENE)	--	<0.0066 U	<0.0058 U	<0.0079 U	<0.0066 U	<0.0053 U	<0.0055 U	<0.0070 U	<0.0083 U	<0.0072 U
M,P-XYLENE (SUM OF ISOMERS)	0.05	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U
METHYLENE CHLORIDE	0.041	<0.0026 U	<0.0023 U	<0.0032 U	<0.0026 U	<0.0021 U	<0.0022 U	<0.0028 U	<0.0033 U	<0.0029 U

Table 4  
SVOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect A										Transect B												
		SDA1	SDA1	SDA1	SDA2	SDA2	SDA3	SDA3	SDA4	SDA4	SDB1	SDB1	SDB1	SDB2	SDB2	SDB2	SDB4	SDB4	SDB4	SDB4				
Sample ID	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/8/2013	11/8/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013				
Sample Date																								
Sample Interval (inches)	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24			
Chemical Name																								
1,2,4-TRICHLOROBENZENE	2.14	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U		
1,2-DICHLOROBENZENE	0.28	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.35 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U	
1,3-DICHLOROBENZENE	1.8	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.35 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U	
1,4-DICHLOROBENZENE	0.72	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.35 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U	
2,4,5-TRICHLOROPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2,4,6-TRICHLOROPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2,4-DICHLOROPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2,4-DIMETHYLPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2,4-DINITROPHENOL	--	<1.4 U	<1.2 U	<1.2 U	<1.2 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	R	<1.5 U	<1.5 U	<1.3 U	<1.2 U	<1.6 U	<1.4 U	<1.3 U	<1.6 U	<1.4 U	<1.3 U	<1.6 U	<1.4 U	<1.3 U	
2,4-DINITROTOLUENE	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2,6-DINITROTOLUENE	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2-CHLORONAPHTHALENE	--	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U			
2-CHLOROPHENOL	--	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U			
2-METHYLNAPHTHALENE	0.586*	<0.14 U	<0.12 U	<0.12 U	<0.12 U	<0.13 U	<0.13 U	<0.13 U	<0.12 U	<0.12 U	<0.13 U	<0.12 U	<0.15 U	<0.14 U	<0.12 U	<0.15 U	<0.16 U	<0.14 U	<0.13 U	<0.16 U	<0.14 U	<0.13 U		
2-METHYLPHENOL (O-CRESOL)	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2-NITROANILINE	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
2-NITROPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U			
3,3'-DICHLOROBENZIDINE	--	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U			
3-NITROANILINE	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U		
4,6-DINITRO-2-METHYLPHENOL	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U		
4-BROMOPHENYL PHENYL ETHER	--	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U	<2.1 U	<0.36 U	<0.38 U	<0.32 U	<0.31 U	<0.39 U	<0.35 U	<0.32 U	<0.4 U	<0.35 U	<0.31 U			
4-CHLOROANILINE	--	<0.7 U	<0.61 U	<0.58 U	<0.61 U	<0.62 U	<0.64 U	<0.65 U	<0.63 U	<0.6 U	<4.1 U	<0.73 U	<0.76 U	<0.64 U	<0.62 U	<0.77 U	<0.79 U	<0.7 U	<0.64 U	<0.8 U	<0.71 U	<0.63 U		
4-CHLOROPHENYL PHENYL ETHER	--	<0.35 U	<0.3 U	<0.29 U	<0.3 U	<0.31 U	<0.32 U	<0.33 U	<0.32 U	<0.3 U</td														

Table 4  
SVOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect C												Transect D												
		SDC1 11/6/2013	SDC1 11/6/2013	SDC1 11/6/2013	SDC2 11/6/2013	SDC2 11/6/2013	SDC3 11/6/2013	SDC4 11/8/2013	SDC4 11/8/2013	SDC5 11/7/2013	SDC5 11/7/2013	SDC5 11/7/2013	SDD1 11/6/2013	SDD1 11/6/2013	SDD1 11/6/2013	SDD2 11/7/2013	SDD2 11/7/2013	SDD3 11/8/2013	SDD3 11/8/2013	SDD3 11/8/2013	SDD3 11/8/2013					
Sample ID																										
Sample Date																										
Sample Interval (inches)																										
Chemical Name																										
1,2,4-TRICHLOROBENZENE	2.14	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
1,2-DICHLOROBENZENE	0.28	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
1,3-DICHLOROBENZENE	1.8	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
1,4-DICHLOROBENZENE	0.72	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
2,4,5-TRICHLOROPHENOL	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2,4,6-TRICHLOROPHENOL	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2,4-DIMETHYLPHENOL	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2,4-DINITROPHENOL	--	<1.7 U	<1.4 U	<1.4 U	<1.2 U	<1.3 U	<1.2 U	<1.2 U	<1.3 U	<1.2 U	<1.6 U	<1.3 U	<1.7 U	<1.5 U	<1.2 U	<1.5 U	<1.3 U	<1.2 U	<1.4 U	<1.3 U	<1.2 U	<1.4 U	<1.3 U	<1.2 U	<1.2 U	
2,4-DINITROTOLUENE	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2,6-DINITROTOLUENE	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2-CHLORONAPHTHALENE	--	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
2-CHLOROPHENOL	--	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
2-METHYLNAPHTHALENE	0.586*	<0.17 U	<0.14 U	<0.14 U	<0.12 U	<0.13 U	<0.12 U	<0.12 U	<0.13 U	<0.12 U	<0.16 U	<0.13 U	<0.17 U	<0.15 U	<0.12 U	<0.12 U	<0.13 U	<0.12 U	<0.14 U	<0.13 U	<0.12 U	<0.12 U	<0.12 U	<0.12 U		
2-METHYLPHENOL (O-CRESOL)	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2-NITROANILINE	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
2-NITROPHENOL	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
3,3-DICHLOROBENZIDINE	--	<0.42 U	<0.36 U	<0.34 U	<0.36 U	<0.31 U	<0.3 U	<0.32 U	<0.29 U	<0.29 U	<0.4 U	<0.33 U	<0.33 U	<0.42 U	<0.38 U	<0.3 U	<0.38 U	<0.32 U	<0.3 U	<0.36 U	<0.31 U	<0.34 U	<0.33 U	<0.31 U	<0.31 U	
3-NITROANILINE	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
4,6-DINITRO-2-METHYLPHENOL	--	<0.83 U	<0.72 U	<0.69 U	<0.72 U	<0.61 U	<0.59 U	<0.64 U	<0.59 U	<0.59 U	<0.8 U	<0.66 U	<0.65 U	<0.85 U	<0.61 U	<0.77 U	<0.64 U	<0.6 U	<0.71 U	<0.63 U	<0.68 U	<0.67 U	<0.61 U	<0.61 U	<0.61 U	
4-BROMOPHENYL PHENYL ETHER	--</td																									

Table 4  
SVOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Transect E												Transect F											
		SDE1	SDE1	SDE1	SDE2	SDE2	SDE3	SDE3	SDE4	SDE4	SDF1	SDF1	SDF1	SDF2	SDF2	SDF3	SDF3	SDF4							
		11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/4/2013	11/4/2013	11/4/2013	11/7/2013	11/7/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	11/8/2013	
Sample ID																									
Sample Date																									
Sample Interval (inches)																									
Chemical Name																									
1,2,4-TRICHLOROBENZENE	2.14	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
1,2-DICHLOROBENZENE	0.28	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
1,3-DICHLOROBENZENE	1.8	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
1,4-DICHLOROBENZENE	0.72	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
2,4,5-TRICHLOROPHENOL	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2,4,6-TRICHLOROPHENOL	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2,4-DIMETHYLPHENOL	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2,4-DINITROPHENOL	--	<1.5 U	<1.3 U	<1.2 U	<1.3 U	<1.4 U	<1.2 U	<1.3 U	<1.4 U	<1.2 U	<1.3 U	<1.4 U	<1.5 U	<1.2 U	<1.3 U	<1.4 U	<1.2 U	<1.3 U	<1.4 U	<1.2 U	<1.3 U	<1.4 U	<1.2 U	<1.3 U	<1.2 U
2,4-DINITROTOLUENE	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2,6-DINITROTOLUENE	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2-CHLORONAPHTHALENE	--	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
2-CHLOROPHENOL	--	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
2-METHYLNAPHTHALENE	0.586*	<0.15 U	<0.13 U	<0.12 U	<0.15 U	<0.13 U	<0.12 U	<0.14 U	<0.12 U	0.0827 J	0.0344 J	0.0400 J	<0.67 U	<0.12 U	<0.13 U	<0.15 U	<0.14 U	<0.12 U	<0.14 U	<0.16 U	<0.13 U	<0.12 U	<0.14 U	<0.13 U	<0.12 U
2-METHYLPHENOL (O-CRESOL)	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2-NITROANILINE	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
2-NITROPHENOL	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
3,3-DICHLOROBENZIDINE	--	<0.37 U	<0.32 U	<0.29 U	<0.38 U	<0.33 U	<0.3 U	<0.35 U	<0.32 U	<0.31 U	<0.38 U	<0.43 U	<0.47 U	<1.7 U	<0.3 U	<0.33 U	<0.37 U	<0.35 U	<0.3 U	<0.34 U	<0.31 U	<0.36 U	<0.41 U	<0.32 U	<0.31 U
3-NITROANILINE	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	<0.59 U	<0.66 U	<0.73 U	<0.7 U	<0.59 U	<0.68 U	<0.62 U	<0.71 U	<0.81 U	<0.64 U	<0.62 U
4,6-DINITRO-2-METHYLPHENOL	--	<0.75 U	<0.63 U	<0.58 U	<0.75 U	<0.66 U	<0.6 U	<0.71 U	<0.63 U	<0.61 U	<0.76 U	<0.86 U	<0.94 U	<3.4 U	&lt										

Table 4  
SVOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect		Transect G										Transect H										Transect I									
Sample ID	NYSDEC Screening Criteria - Class A	SDG1	SDG1	SDG1	SDG2	SDG2	SDG2	SDH1	SDH1	SDH1	SDH2	SDH2	SDH3	SDH3	SDH3	SDI1	SDI1 (Dup)	SDI1	SDI1 (Dup)	SDI1	SDI1 (Dup)	SDI2	SDI2	SDI2	SDI3	SDI3	SDI3				
Sample Date	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/7/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/6/2013	11/7/2013	11/7/2013	11/7/2013	11/5/2013	11/5/2013	11/5/2013					
Sample Interval (inches)		0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	0 - 6	6 - 12	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24							
Chemical Name																															
1,2,4-TRICHLOROBENZENE	2.14	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U				
1,2-DICHLOROBENZENE	0.28	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U				
1,3-DICHLOROBENZENE	1.8	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U					
1,4-DICHLOROBENZENE	0.72	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U				
2,4,5-TRICHLOROPHENOL	-	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2,4,6-TRICHLOROPHENOL	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2,4-DICHLOROPHENOL	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2,4-DIMETHYLPHENOL	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2,4-DINITROPHENOL	--	R	<1.2 U	R	<14 U	<14 U	<1.2 U	<3.5 U	<1.3 U	<1.2 U	<1.3 U	<1.3 U	<1.2 U	<1.3 U	<1.3 U	<1.2 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U	<1.3 U			
2,4-DINITROTOLUENE	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.64 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2,6-DINITROTOLUENE	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2-CHLORONAPHTHALENE	--	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U				
2-CHLOROPHENOL	--	<0.36 U	<0.31 U	<0.31 U	<0.36 U	<0.35 U	<0.3 U	<0.87 U	<0.32 U	<0.3 U	<0.38 U	<0.33 U	<0.32 U	<0.72 U	<0.35 U	<0.31 U	<0.37 U	<0.37 U	<0.33 U	<0.32 U	<0.32 U	<2.4 U	<1.1 U	R	<0.43 U	<0.34 U	<0.33 U				
2-METHYLNAPHTHALENE	0.586*	<0.14 U	<0.12 U	<0.14 U	<0.14 U	<0.12 U	<0.35 U	<0.13 U	<0.12 U	<0.15 U	<0.13 U	<0.13 U	<0.12 U	<0.29 U	<0.14 U	<0.13 U	<0.15 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U	<0.13 U				
2-METHYLPHENOL (O-CRESOL)	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2-NITROANILINE	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U	<0.69 U	<0.66 U				
2-NITROPHENOL	--	<0.72 U	<0.62 U	<0.62 U	<0.71 U	<0.7 U	<0.59 U	<1.7 U	<0.64 U	<0.6 U	<0.77 U	<0.66 U	<0.64 U	<1.4 U	<0.69 U	<0.63 U	<0.75 U	<0.74 U	<0.67 U	<0.65 U	<0.63 U	<4.8 U	<2.2 U	R	<0.86 U						

Table 4  
SVOC Screening Table  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
ompson Road, Syracuse, New York

### Note

SVOC - Semi-Volatile Organic Compound

SVOC - Semi-Volatile Organic Compound.  
All units are in milligrams per kilogram (mg/kg).

\* - LTC-Developed criteria

\* - UTC-Developed criteria.

U or < - Compound was a  
↓ Estimated

J - Estimated.  
P - Calculations are based on available data.

Table 4  
 SVOC Screening Table  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria - Class A	Sanders Creek								
		SDS8 11/5/2013	SDS8 (Dup) 11/5/2013	SDS8 11/5/2013	SDS8 (Dup) 11/5/2013	SDS8 11/5/2013	SDS8 (Dup) 11/5/2013	SDS9 11/5/2013	SDS9 11/5/2013	
		0 - 6	0 - 6	6 - 12	6 - 12	18 - 24	18 - 24	0 - 6	6 - 12	18 - 24
1,2,4-TRICHLOROBENZENE	2.14	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
1,2-DICHLOROBENZENE	0.28	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
1,3-DICHLOROBENZENE	1.8	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
1,4-DICHLOROBENZENE	0.72	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
2,4,5-TRICHLOROPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2,4,6-TRICHLOROPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2,4-DICHLOROPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2,4-DIMETHYLPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2,4-DINITROPHENOL	--	<1.5 U	<1.3 U	<1.6 U	<1.5 U	<1.3 U	<1.3 U	<1.3 U	<1.8 U	<1.3 U
2,4-DINITROTOLUENE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2,6-DINITROTOLUENE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2-CHLORONAPHTHALENE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
2-CHLOROPHENOL	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
2-METHYLNAPHTHALENE	0.586*	<0.15 U	<0.13 U	<0.16 U	<0.15 U	<0.13 U	<0.13 U	<0.13 U	0.0240 J	<0.13 U
2-METHYLPHENOL (O-CRESOL)	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2-NITROANILINE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
2-NITROPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
3,3'-DICHLOROBENZIDINE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
3-NITROANILINE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4,6-DINITRO-2-METHYLPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4-BROMOPHENYL PHENYL ETHER	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
4-CHLORO-3-METHYLPHENOL	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4-CHLOROANILINE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4-CHLOROPHENYL PHENYL ETHER	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
4-METHYLPHENOL (P-CRESOL)	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4-NITROANILINE	--	<0.77 U	<0.63 U	<0.81 U	<0.77 U	<0.64 U	<0.63 U	<0.65 U	<0.89 U	<0.63 U
4-NITROPHENOL	--	<1.5 U	<1.3 U	<1.6 U	<1.5 U	<1.3 U	<1.3 U	<1.8 U	<1.3 U	<1.3 U
ACENAPHTHENE	9.82	<0.15 U	<0.13 U	<0.16 U	<0.15 U	<0.13 U	<0.13 U	0.0182 J	0.0724 J	<0.13 U
ACENAPHTHYLENE	9.04	<0.15 U	<0.13 U	<0.16 U	<0.15 U	<0.13 U	0.0243 J	<0.13 U	0.0346 J	<0.13 U
ANTHRACENE	11.88	0.0250 J	<0.13 U	<0.16 U	0.0219 J	<0.13 U	0.0327 J	0.0366 J	0.192	0.0171 J
BENZO(A)ANTHRACENE	16.82	0.12 J	0.0319 J	0.0722 J	0.132 J	0.0239 J	0.102 J	0.164	0.913	0.0778 J
BENZO(A)PYRENE	19.28	0.11 J	0.0273 J	0.0794 J	0.117 J	<0.13 U	0.111 J	0.161	0.943	0.0729 J
BENZO(B)FLUORANTHENE	19.58	0.115 J	0.0323 J	0.0789 J	0.119 J	<0.13 U	0.11 J	0.148	1.04	0.0843 J
BENZO(G,H,I)PERYLENE	21.9	0.0634 J	0.0191 J	0.0486 J	0.0872 J	<0.13 U	0.0683 J	0.0971 J	0.631	0.0525 J
BENZO(K)FLUORANTHENE	19.6	0.101 J	0.0212 J	0.0685 J	0.122 J	<0.13 U	0.0746 J	0.149	0.773	0.0523 J
BENZYL BUTYL PHTHALATE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
BIS(2-CHLOROETHOXY) METHANE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
BIS(2-CHLOROISOPROPYL) ETHER	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
BIS(2-ETHYLHEXYL) PHTHALATE	180*	0.0937 J	0.167 J	0.112 J	0.135 J	0.0359 J	<0.31 U	<0.33 U	0.123 J	0.0423 J
CARBAZOLE	--	<0.15 U	<0.13 U	<0.16 U	0.0201 J	<0.13 U	0.0202 J	0.0310 J	0.178 J	0.0157 J
CHRYSENE	16.86	0.133 J	0.0292 J	0.101 J	0.136 J	0.0208 J	0.11 J	0.176	1	0.0750 J
DIBENZA(A,H)ANTHRACENE	22.44	<0.15 U	<0.13 U	<0.16 U	0.0324 J	<0.13 U	<0.13 U	0.184	<0.13 U	<0.13 U
DIBENZOFURAN	0.831*	<0.15 U	<0.13 U	<0.16 U	<0.15 U	<0.13 U	<0.13 U	0.0469 J	<0.13 U	<0.13 U
DIETHYL PHTHALATE	1.21*	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
DIMETHYL PHTHALATE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
DI-N-BUTYL PHTHALATE	6.47*	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
DI-N-OCTYLPHTHALATE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
FLUORANTHENE	14.16	0.283	0.0568 J	0.191	0.275	0.0425 J	0.22	0.349	1.92	0.162
FLUORENE	10.78	<0.15 U	<0.13 U	<0.16 U	<0.15 U	<0.13 U	<0.13 U	0.0715 J	<0.13 U	<0.13 U
HEXACHLOROBENZENE	--	<0.39 U	<0.32 U	<0.41 U	<0.38 U	<0.32 U	<0.31 U	<0.33 U	<0.44 U	<0.31 U
HEXACHLOROBUTADIENE</td										

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect A																																			
			SDA1			SDA1			SDA1			SDA2			SDA2			SDA3			SDA3			SDA4			SDA4											
			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/7/2013			11/7/2013			11/8/2013			11/8/2013											
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU									
ACENAPHTHENE	491	9.82	0.0685	2.4729	0.0050	0.0031 J	0.2385	0.0005	< 0.0045 U	--	--	< 0.0046 U	--	--	< 0.0051 U	--	--	< 0.0046 U	--	--	0.0029 J	0.5653	0.0012	< 0.0046 U	--	--	0.0529	1.1914	0.0024	0.0182	0.4764	0.0010	< 0.0058 U	--	--			
ACENAPHTHYLENE	452	9.04	0.785	28.3394	0.0627	0.0224	1.7231	0.0038	< 0.0045 U	--	--	0.0103	1.0457	0.0023	0.055	0.3374	0.0007	0.0129	2.0379	0.0045	< 0.0046 U	--	--	0.0031 J	0.6043	0.0013	< 0.0046 U	--	--	0.456	10.2703	0.0227	0.0409	1.0707	0.0024	0.0031 J	0.1131	0.0003
ANTHRACENE	594	11.88	0.753	27.1841	0.0458	0.0174	1.3385	0.0023	< 0.0045 U	--	--	0.0081	0.8223	0.0014	0.045 J	0.2761	0.0005	0.0098	1.5482	0.0026	< 0.0035 J	0.4667	0.0008	0.0065	1.2671	0.0021	< 0.0046 U	--	--	0.44	9.9099	0.0167	0.0643	1.6832	0.0028	< 0.0058 U	--	--
BENZO(A)ANTHRACENE	841	16.82	0.596	21.5162	0.0256	0.0238	1.8308	0.0022	< 0.0045 U	--	--	0.0139	1.4112	0.0017	0.058	0.3558	0.0004	0.0135	2.1327	0.0025	0.0058	0.7733	0.0009	0.0128	2.4951	0.0030	< 0.0046 U	--	--	0.465	10.4730	0.0125	0.151	3.9529	0.0047	0.0043 J	0.1569	0.0002
BENZO(A)PYRENE	964	19.28	0.895	32.3105	0.0335	0.0372	2.8615	0.0030	< 0.0045 U	--	--	0.0200	2.0305	0.0021	0.096	0.5890	0.0006	0.0228	3.6019	0.0037	0.0105	1.4000	0.0015	0.0139	2.7096	0.0028	0.030 J	0.7732	0.0008	0.621	13.9865	0.0145	0.173	4.5288	0.0047	0.0049 J	0.1788	0.0002
BENZO(B)FLUORANTHENE	979	19.58	0.81	29.2419	0.0299	0.0342	2.6308	0.0027	< 0.0045 U	--	--	0.0194	1.9695	0.0020	0.097	0.5951	0.0006	0.0223	3.5229	0.0036	0.0113	1.5067	0.0015	0.0111	2.1637	0.0022	< 0.0046 U	--	--	0.574	12.9279	0.0132	0.144	3.7696	0.0039	0.0040 J	0.1460	0.0001
BENZO(G,H)PERYLENE	1095	21.9	0.838	30.2527	0.0276	0.0332	2.5538	0.0023	< 0.0045 U	--	--	0.0161	1.6345	0.0015	0.087	0.5337	0.0005	0.0199	3.1438	0.0029	0.0064	0.8533	0.0008	0.0078	1.5205	0.0014	< 0.0046 U	--	--	0.52	11.7117	0.0107	0.115	3.0105	0.0027	0.0031 J	0.1131	0.0001
BENZO(K)FLUORANTHENE	980	19.6	0.778	28.0866	0.0287	0.0356	2.7385	0.0028	< 0.0045 U	--	--	0.0189	1.9188	0.0020	0.092	0.5644	0.0006	0.0214	3.3807	0.0034	0.0104	1.3867	0.0014	0.0118	2.3002	0.0023	< 0.0046 U	--	--	0.566	12.7477	0.0130	0.147	3.8482	0.0039	< 0.0058 U	--	--
BENZO(E)PYRENE	967	19.34	0.875	31.5884	0.0327	0.0376	2.8923	0.0030	< 0.0045 U	--	--	0.0185	1.8782	0.0019	0.0101	0.6196	0.0006	0.0228	3.6019	0.0037	0.0130	1.7333	0.0018	< 0.0046 U	--	--	0.54	12.1622	0.0126	0.119	3.1152	0.0032	0.0034 J	0.1241	0.0001			
C1-Benzoanthracenes/Chrysenes	930	18.6	0.517	18.6643	0.0201	0.0213	1.6385	0.0018	< 0.0045 U	--	--	0.0115	1.1675	0.0013	0.049	0.3006	0.0003	0.0122	1.9273	0.0021	0.0145	1.9333	0.0021	0.0067	1.3060	0.0014	< 0.0046 U	--	--	0.358	8.0631	0.0087	0.0693	1.8141	0.0020	0.0037 J	0.1350	0.0001
C1-Fluoranthenes/Pyrenes	769	15.38	0.79	28.5199	0.0371	0.0278	2.1385	0.0028	< 0.0045 U	--	--	0.0162	1.6447	0.0021	0.076	0.4663	0.0006	0.0165	2.6066	0.0034	0.0108	1.4400	0.0019	0.0126	2.4561	0.0032	0.026 J	0.6701	0.0009	0.549	12.3649	0.0161	0.127	3.3246	0.0043	0.0063	0.2299	0.0003
C1-Fluorenes	611	12.22	0.0806	2.9097	0.0048	0.0041 J	0.3154	0.0005	< 0.0045 U	--	--	0.0028 J	0.2843	0.0005	< 0.0046 U	--	--	0.0027 J	0.4265	0.0007	0.0024 J	0.3200	0.0005	< 0.0046 U	--	--	0.0061	1.3761	0.0023	0.0114	0.2984	0.0005	< 0.0058 U	--	--			
C1-Naphthalenes	445	8.9	0.0950	3.4296	0.0077	0.0058	0.4462	0.0010	< 0.0045 U	--	--	0.0039 J	0.3959	0.0009	< 0.0046 U	--	--	0.0033 J	0.5213	0.0012	0.0052	0.6933	0.0016	0.0048	0.9357	0.0021	< 0.0046 U	--	--	0.13	2.9279	0.0066	0.0463	1.2120	0.0027	< 0.0058 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.507	18.3032	0.0273	0.0173	1.3308	0.0020	< 0.0045 U	--	--	0.0117	1.1878	0.0018	0.051	0.3129	0.0005	0.0103	1.6272	0.0024	0.0082	1.0933	0.0016	0.0106	2.0663	0.0031	0.0025 J	0.6443	0.0010	0.503	11.3288	0.0169	0.103	2.6963	0.0040	0.0060	0.2190	0.0003
C2-Benzoanthracenes/Chrysenes	1009	20.18	0.261	9.4224	0.0093	0.0136	1.0462	0.0010	< 0.0045 U	--	--	0.0078	0.7919	0.0008	0.035 J	0.2147	0.0002	0.0076	1.2006	0.0																		

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	NYSDEC Screening Criteria	Transect B																														
		SDB1			SDB1			SDB1			SDB2			SDB2 (Dup)			SDB2			SDB2			SDB4									
		11/6/2013		11/6/2013		11/6/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013		11/7/2013								
		0 - 6		6 - 12		18 - 24		0 - 6		0 - 6		6 - 12		18 - 24		0 - 6		6 - 12		18 - 24		0 - 6		6 - 12		18 - 24						
Chemical Name	Sediment Guidance Value ( $\mu\text{g/gOC}$ )	NYSDEC Sediment CLASS A	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU	mg/kg	$\mu\text{g/gOC}$	TU						
ACENAPHTHENE	491	9.82	0.0050 J	0.0990	0.0002	< 0.0045 U	--	--	< 0.0048 U	--	--	< 0.0059 J	--	--	0.0034 J	0.0912	0.0002	< 0.0052 J	--	--	< 0.0050 U	--	--	< 0.0060 J	--	--	< 0.0045 U	--	--			
ACENAPHTHYLENE	452	9.04	0.0128	0.2535	0.0006	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0098	0.3131	0.0007	0.0125	0.3351	0.0007	0.0028 J	0.1628	0.0004	< 0.0050 U	--	--	0.0093	0.2308	0.0005	< 0.0053 U	--	--	< 0.0045 U	--	--
ANTHRACENE	594	11.88	0.0129	0.2554	0.0004	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0104	0.3326	0.0006	0.0138	0.3700	0.0006	0.0028 J	0.1628	0.0003	< 0.0050 U	--	--	0.0093	0.2308	0.0004	< 0.0053 U	--	--	< 0.0045 U	--	--
BENZO(A)ANTHRACENE	841	16.82	0.0310	0.1639	0.0007	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0253	0.8083	0.0010	0.0333	0.8928	0.0011	0.0042 J	0.2442	0.0003	< 0.0050 U	--	--	0.0333	0.8263	0.0010	0.0067	0.2900	0.0003	0.0030 J	0.6012	0.0007
BENZO(A)PYRENE	964	19.28	0.0378	0.7485	0.0008	0.0031 J	0.1640	0.0002	< 0.0048 U	--	--	0.0295	0.9425	0.0010	0.0373	1.0000	0.0010	0.0048 J	0.2791	0.0003	< 0.0050 U	--	--	0.0404	1.0025	0.0010	0.0075	0.3247	0.0003	0.0033 J	0.6613	0.0007
BENZO(B)FLUORANTHENE	979	19.58	0.0320	0.6337	0.0006	0.0027 J	0.1429	0.0001	< 0.0048 U	--	--	0.0259	0.8275	0.0008	0.0349	0.9357	0.0010	0.0039 J	0.2267	0.0002	< 0.0050 U	--	--	0.0356	0.8834	0.0009	0.0064	0.2771	0.0003	0.0030 J	0.6012	0.0006
BENZO(G,H,I)PERYLENE	1095	21.9	0.0234	0.4634	0.0004	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0196	0.6262	0.0006	0.0248	0.6649	0.0006	0.0030 J	0.1744	0.0002	< 0.0050 U	--	--	0.0250	0.6203	0.0006	0.0042 J	0.1818	0.0002	< 0.0045 U	--	--
BENZO(K)FLUORANTHENE	980	19.6	0.0338	0.6693	0.0007	0.0025 J	0.1323	0.0001	< 0.0048 U	--	--	0.0271	0.8658	0.0009	0.0348	0.9330	0.0010	0.0042 J	0.2442	0.0002	< 0.0050 U	--	--	0.0367	0.9107	0.0009	0.0069	0.2987	0.0003	0.0031 J	0.6212	0.0006
BENZO(E)PYRENE	967	19.34	0.0271	0.5366	0.0006	0.0023 J	0.1217	0.0001	< 0.0048 U	--	--	0.0215	0.6869	0.0007	0.0284	0.7614	0.0008	0.0036 J	0.2093	0.0002	< 0.0050 U	--	--	0.0290	0.7196	0.0007	0.0051 J	0.2208	0.0002	< 0.0045 U	--	--
C1-Benzoanthracenes/Chrysenes	930	18.6	0.0207	0.4099	0.0004	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0160	0.5112	0.0005	0.0202	0.5416	0.0006	0.0034 J	0.1977	0.0002	< 0.0050 U	--	--	0.0184	0.4566	0.0005	0.0047 J	0.2035	0.0002	< 0.0045 U	--	--
C1-Fluoranthenes/Pyrenes	769	15.38	0.0327	0.6475	0.0008	0.0026 J	0.1376	0.0002	< 0.0048 U	--	--	0.0266	0.8496	0.0011	0.0358	0.9598	0.0012	0.0061	0.3837	0.0005	< 0.0050 U	--	--	0.0301	0.7469	0.0010	0.0077	0.3333	0.0004	0.0032 J	0.6413	0.0008
C1-Fluorenes	611	12.22	0.0049 J	0.0970	0.0002	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0057 J	0.1821	0.0003	0.0066	0.1769	0.0003	0.0033 J	0.1919	0.0003	< 0.0050 U	--	--	0.0037 J	0.0918	0.0002	0.0027 J	0.1169	0.0002	< 0.0045 U	--	--
C1-Naphthalenes	445	8.9	0.0093	0.1842	0.0004	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0082	0.2620	0.0006	0.0099	0.2654	0.0006	< 0.0052 J	--	--	< 0.0050 U	--	--	< 0.0060 J	--	--	< 0.0053 J	--	--	< 0.0045 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0256	0.5069	0.0008	0.0026 J	0.1376	0.0002	< 0.0048 U	--	--	0.0265	0.8466	0.0013	0.0338	0.9062	0.0014	0.0067	0.3895	0.0006	< 0.0050 U	--	--	0.0208	0.5161	0.0008	0.0064	0.2771	0.0004	0.0026 J	0.5210	0.0008
C2-Benzoanthracenes/Chrysenes	1009	20.18	0.0125	0.2475	0.0002	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0080	0.2556	0.0003	0.0081	0.2172	0.0002	< 0.0052 J	--	--	< 0.0050 U	--	--	0.0097	0.2407	0.0002	0.0027 J	0.1169	0.0001	< 0.0045 U	--	--
C2-Fluorenes	687	13.74	0.0068	0.1347	0.0002	< 0.0045 U	--	--	< 0.0048 U	--	--	0.0062	0.1981	0.0003	0.0075	0.2011	0.0003	0.0029 J	0.1686	0.0002	< 0.0050 U	--	--	0.0042 J	0.1042	0.0002	< 0.0053 U	--	--	< 0.0045 U	--	--
C2-Naphthalenes	510	10.2	0.0093	0.1842	0.0004	< 0.0045 U	--	--	< 0.0048 U	--	--</td																					

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect C																							
			SDC1			SDC1			SDC1			SDC2			SDC2			SDC2			SDC3					
			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/6/2013					
Sample Date	Sample Interval (inches)	Sediment Guidance Value (µg/gOC)	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24			
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU																					
ACENAPHTHENE	491	9.82	< 0.0066 U	--	--	< 0.0051 U	--	--	< 0.0050 U	--	--	< 0.0056 J	--	--	< 0.0050 U	--	--	< 0.0047 U	--	--	< 0.0049 J	--	--	< 0.0045 U	--	--
ACENAPHTHYLENE	452	9.04	0.0085	0.2457	0.0005	0.0036 J	0.1891	0.0004	< 0.0050 U	--	--	0.0048 J	0.2202	0.0005	< 0.0050 U	--	--	0.0076	0.3858	0.0009	< 0.0045 U	--	--	< 0.0046 U	--	--
ANTHRACENE	594	11.88	0.0088	0.2543	0.0004	0.0040 J	0.1990	0.0003	< 0.0050 U	--	--	0.0043 J	0.1972	0.0003	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0077	0.3909	0.0007	< 0.0045 U	--	--
BENZO(A)ANTHRACENE	841	16.82	0.0264	0.7630	0.0009	0.0120	0.5970	0.0007	< 0.0050 U	--	--	0.0094	0.4312	0.0005	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0178	0.9036	0.0011	< 0.0045 U	--	--
BENZO(A)PYRENE	964	19.28	0.0336	0.9711	0.0010	0.0149	0.7413	0.0008	0.0029 J	0.2266	0.0002	0.0120	0.5505	0.0006	0.0030 J	0.4532	0.0005	< 0.0047 U	--	--	0.0205	1.0406	0.0011	0.0023 J	0.2312	0.0002
BENZO(B)FLUORANTHENE	979	19.58	0.0317	0.9162	0.0009	0.0134	0.6667	0.0007	0.0030 J	0.2344	0.0002	0.0108	0.4954	0.0005	0.0025 J	0.3776	0.0004	< 0.0047 U	--	--	0.0183	0.9289	0.0009	< 0.0045 U	--	--
BENZO(G,H,I)PERYLENE	1095	21.9	0.0224	0.6474	0.0006	0.0094	0.4677	0.0004	< 0.0050 U	--	--	0.0085	0.3899	0.0004	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0131	0.6650	0.0006	< 0.0045 U	--	--
BENZO(K)FLUORANTHENE	980	19.6	0.0314	0.9075	0.0009	0.0136	0.6766	0.0007	0.0028 J	0.2188	0.0002	0.0112	0.5138	0.0005	0.0026 J	0.3927	0.0004	< 0.0047 U	--	--	0.0185	0.9391	0.0010	< 0.0045 U	--	--
BENZO(E)PYRENE	967	19.34	0.0258	0.7457	0.0008	0.0108	0.5373	0.0006	< 0.0050 U	--	--	0.0091	0.4174	0.0004	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0150	0.7614	0.0008	< 0.0045 U	--	--
C1-Benzanthracenes/Chrysenes	930	18.6	0.0148	0.4277	0.0005	0.0078	0.3881	0.0004	< 0.0050 U	--	--	0.0073	0.3349	0.0004	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0116	0.5888	0.0006	< 0.0045 U	--	--
C1-Fluoranthenes/Pyrenes	769	15.38	0.0236	0.6821	0.0009	0.0124	0.6169	0.0008	0.0034 J	0.2656	0.0003	0.0124	0.5688	0.0007	0.0034 J	0.5136	0.0007	< 0.0047 U	--	--	0.0202	1.0254	0.0013	0.0034 J	0.3417	0.0004
C1-Fluorenes	611	12.22	0.0038 J	0.1098	0.0002	0.0026 J	0.1294	0.0002	< 0.0050 U	--	--	0.0031 J	0.1422	0.0002	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0034 J	0.1726	0.0003	< 0.0045 U	--	--
C1-Naphthalenes	445	8.9	0.0050 J	0.1445	0.0003	0.0034 J	0.1692	0.0004	< 0.0050 U	--	--	0.0057	0.2615	0.0006	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0056	0.2843	0.0006	< 0.0045 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0165	0.4769	0.0007	0.0103	0.5124	0.0008	0.0038 J	0.2969	0.0004	0.0126	0.5780	0.0009	0.0034 J	0.5136	0.0008	< 0.0047 U	--	--	0.0183	0.9289	0.0014	0.0032 J	0.3216	0.0005
C2-Benzanthracenes/Chrysenes	1009	20.18	0.0080	0.2312	0.0002	0.0039 J	0.1940	0.0002	< 0.0050 U	--	--	0.0048 J	0.2202	0.0002	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0059	0.2995	0.0003	< 0.0045 U	--	--
C2-Fluorenes	687	13.74	0.0041 J	0.1185	0.0002	0.0026 J	0.1294	0.0002	< 0.0050 U	--	--	0.0038 J	0.1743	0.0003	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0043 J	0.2183	0.0003	< 0.0045 U	--	--
C2-Naphthalenes	510	10.2	0.0056 J	0.1618	0.0003	0.0036 J	0.1791	0.0004	< 0.0050 U	--	--	0.0080	0.3670	0.0007	0.0034 J	0.5136	0.0010	0.0024 J	0.5298	0.0010	0.0089	0.4518	0.0009	0.0032 J	0.3216	0.0006
C2-Phenanthrenes/Anthracenes	745	14.9	0.0082	0.2370	0.0003	0.0055	0.2736	0.0004	< 0.0050 U	--	--	0.0067	0.3073	0.0004	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0112	0.5685	0.0008	< 0.0045 U	--	--
C3-Benzanthracenes/Chrysenes	1213	24.26	< 0.0066 U	--	--	< 0.0051 U	--	--	< 0.0050 U	--	--	< 0.0056 U	--	--	< 0.0050 U	--	--	< 0.0047 U	--	--	< 0.0049 U	--	--	< 0.0045 U	--	--
C3-Fluorenes	768	15.36	< 0.0066 U	--	--	< 0.0051 U	--	--	< 0.0050 U	--	--	< 0.0056 U	--	--	< 0.0050 U	--	--	< 0.0047 U	--	--	< 0.0049 U	--	--	< 0.0045 U	--	--
C3-Naphthalenes	581	11.62	0.0068	0.1965	0.0003	0.0034 J	0.1692	0.0003	< 0.0050 U	--	--	0.0066	0.3028	0.0005	0.0031 J	0.4683	0.0008	< 0.0047 U	--	--	0.0063	0.3198	0.0006	0.0030 J	0.3015	0.0005
C3-Phenanthrenes/Anthracenes	830	16.6	0.0049 J	0.1416	0.0002	0.0026 J	0.1294	0.0002	< 0.0050 U	--	--	0.0037 J	0.1697	0.0002	< 0.0050 U	--	--	< 0.0047 U	--	--	0.0059					

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect C														
			SDC4			SDC4			SDC4			SDC5			SDC5		
			11/8/2013			11/8/2013			11/8/2013			11/7/2013			11/7/2013		
			0 - 6	0 - 6	0 - 6	6 - 12	6 - 12	6 - 12	18 - 24	18 - 24	18 - 24	0 - 6	0 - 6	0 - 6	6 - 12	6 - 12	18 - 24
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU
ACENAPHTHENE	491	9.82	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
ACENAPHTHYLENE	452	9.04	0.0110	0.2670	0.0006	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0087	0.2444	0.0005	0.0033 J	0.1610	0.0004
ANTHRACENE	594	11.88	0.0113	0.2743	0.0005	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0079	0.2219	0.0004	0.0031 J	0.1512	0.0003
BENZO(A)ANTHRACENE	841	16.82	0.0259	0.6286	0.0007	0.0037 J	0.6631	0.0008	< 0.0044 U	--	--	0.0210	0.5899	0.0007	0.0069	0.3366	0.0004
BENZO(A)PYRENE	964	19.28	0.0295	0.7160	0.0007	0.0044 J	0.7885	0.0008	< 0.0044 U	--	--	0.0236	0.6629	0.0007	0.0079	0.3854	0.0004
BENZO(B)FLUORANTHENE	979	19.58	0.0235	0.5704	0.0006	0.0035 J	0.6272	0.0006	< 0.0044 U	--	--	0.0202	0.5674	0.0006	0.0066	0.3220	0.0003
BENZO(G,H,I)PERYLENE	1095	21.9	0.0178	0.4320	0.0004	0.0026 J	0.4659	0.0004	< 0.0044 U	--	--	0.0135	0.3792	0.0005	0.0047 J	0.2293	0.0002
BENZO(K)FLUORANTHENE	980	19.6	0.0268	0.6505	0.0007	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0209	0.5871	0.0006	0.0067	0.3268	0.0003
BENZO(E)PYRENE	967	19.34	0.0196	0.4757	0.0005	0.0029 J	0.5197	0.0005	< 0.0044 U	--	--	0.0161	0.4522	0.0005	0.0053 J	0.2585	0.0003
C1-Benzanthracenes/Chrysenes	930	18.6	0.0168	0.4078	0.0004	0.0024 J	0.4301	0.0005	< 0.0044 U	--	--	0.0128	0.3596	0.0004	0.0051 J	0.2488	0.0003
C1-Fluoranthenes/Pyrenes	769	15.38	0.0287	0.6966	0.0009	0.0043 J	0.7706	0.0010	< 0.0044 U	--	--	0.0222	0.6236	0.0008	0.0083	0.4049	0.0005
C1-Fluorenes	611	12.22	0.0039 J	0.0947	0.0002	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
C1-Naphthalenes	445	8.9	0.0063	0.1529	0.0003	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0256	0.6214	0.0009	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0175	0.4916	0.0007	0.0075	0.3659	0.0005
C2-Benzanthracenes/Chrysenes	1009	20.18	0.0078	0.1893	0.0002	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0069	0.1938	0.0002	< 0.0060 U	--	--
C2-Fluorenes	687	13.74	0.0078	0.1893	0.0003	0.0023 J	0.4122	0.0006	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
C2-Naphthalenes	510	10.2	0.0098	0.2379	0.0005	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0045 J	0.1264	0.0002	0.0034 J	0.1659	0.0003
C2-Phenanthrenes/Anthracenes	745	14.9	0.0188	0.4563	0.0006	0.0032 J	0.5735	0.0008	< 0.0044 U	--	--	0.0097	0.2725	0.0004	0.0040 J	0.1951	0.0003
C3-Benzanthracenes/Chrysenes	1213	24.26	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
C3-Fluorenes	768	15.36	0.0054 J	0.1311	0.0002	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 U	--	--	< 0.0060 U	--	--
C3-Naphthalenes	581	11.62	0.0108	0.2621	0.0005	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0057 J	0.1601	0.0003	0.0035 J	0.1707	0.0003
C3-Phenanthrenes/Anthracenes	830	16.6	0.0096	0.2330	0.0003	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0050 J	0.1404	0.0002	< 0.0060 U	--	--
C4-Benzanthracenes/Chrysenes	1213	24.26	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 U	--	--	< 0.0060 U	--	--
C4-Naphthalenes	657	13.14	0.0125	0.3034	0.0005	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 U	--	--	< 0.0060 U	--	--
C4-Phenanthrenes/Anthracenes	670	13.4	0.0033 J	0.0801	0.0001	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
CHRYSENE	843	16.86	0.0334	0.8107	0.0010	0.0046	0.8244	0.0010	< 0.0044 U	--	--	0.0278	0.7809	0.0009	0.0092	0.4488	0.0005
DIBENZ(A,H)ANTHRACENE	1122	22.44	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0038 J	0.1067	0.0001	< 0.0060 U	--	--
FLUORANTHENE	708	14.16	0.0598	1.4515	0.0021	0.0078	1.3978	0.0020	< 0.0044 U	--	--	0.0514	1.4438	0.0020	0.0173	0.8439	0.0012
FLUORENE	539	10.78	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	< 0.0067 J	--	--	< 0.0060 U	--	--
INDENO(1,2,3-C,D)PYRENE	1115	22.3	0.0183	0.4442	0.0004	0.0026 J	0.4659	0.0004	< 0.0044 U	--	--	0.0143	0.4017	0.0004	0.0046 J	0.2244	0.0002
NAPHTHALENE	385	7.7	< 0.0063 U	--	--	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0051 J	0.1433	0.0004	0.0032 J	0.1561	0.0004
PERYLENE	967	19.34	0.0070	0.1699	0.0002	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0061 J	0.1713	0.0002	< 0.0060 U	--	--
PHENANTHRENE	597	11.94	0.0325	0.7888	0.0013	< 0.0046 U	--	--	< 0.0044 U	--	--	0.0256	0.7191	0.0012	0.0109	0.5317	0.0009
PYRENE	698	13.96	0.0515	1.2500	0.0018	0.0067	1.2007	0.0017	< 0.0044 U	--	--	0.0432	1.2135	0.0017	0.0144	0.7024	0.0010
TOC	--	--	41200	--	--	5580	--	--	3100	--	--	35600	--	--	20500	--	--
Total PAH Toxicity Units	--	--															

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect D																												
			SDD1			SDD1			SDD1			SDD2			SDD2			SDD2			SDD3										
			11/6/2013			11/6/2013			11/6/2013			11/7/2013			11/7/2013			11/7/2013			11/8/2013										
Sample Date	Sample Interval (inches)	Sediment Guidance Value (µg/gOC)	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24								
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU																										
ACENAPHTHENE	491	9.82	< 0.0062 U	--	--	< 0.0050 U	--	--	< 0.0048 U	--	--	< 0.0054 J	--	--	< 0.0050 U	--	--	< 0.0049 U	--	--	< 0.0050 J	--	--	< 0.0043 U	--	--	< 0.0042 J	--	--		
ACENAPHTHYLENE	452	9.04	0.0114	0.2365	0.0005	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0107	0.3702	0.0008	0.0028 J	0.2667	0.0006	< 0.0049 U	--	--	0.0080	0.3540	0.0008	< 0.0043 U	--	--	< 0.0042 J	--	--		
ANTHRACENE	594	11.88	0.0110	0.2282	0.0004	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0108	0.3737	0.0008	< 0.0050 U	--	--	< 0.0049 U	--	--	0.0090	0.3982	0.0007	< 0.0043 U	--	--	< 0.0042 U	--	--		
BENZO(A)ANTHRACENE	841	16.82	0.0337	0.6992	0.0008	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0252	0.8720	0.0010	0.0048 J	0.4571	0.0005	0.0031 J	0.3418	0.0004	0.0244	1.0796	0.0013	0.0024 J	0.8727	0.0010	< 0.0042 J	--	--		
BENZO(A)PYRENE	964	19.28	0.0429	0.8900	0.0009	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0295	1.0208	0.0011	0.0057	0.5429	0.0006	0.0037 J	0.4079	0.0004	0.0267	1.1814	0.0012	0.0030 J	1.0909	0.0011	< 0.0042 U	--	--		
BENZO(B)FLUORANTHENE	979	19.58	0.0387	0.8029	0.0008	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0264	0.9135	0.0009	0.0049 J	0.4667	0.0005	0.0032 J	0.3528	0.0004	0.0212	0.9381	0.0010	0.0026 J	0.9455	0.0010	< 0.0042 J	--	--		
BENZO(G,H,I)PERYLENE	1095	21.9	0.0273	0.5664	0.0005	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0194	0.6713	0.0006	0.0036 J	0.3429	0.0003	< 0.0049 U	--	--	0.0160	0.7080	0.0006	< 0.0043 U	--	--	< 0.0042 J	--	--		
BENZO(K)FLUORANTHENE	980	19.6	0.0395	0.8195	0.0008	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0270	0.9343	0.0010	0.0050	0.4762	0.0005	0.0033 J	0.3638	0.0004	0.0223	0.9867	0.0010	< 0.0043 U	--	--	< 0.0042 J	--	--		
BENZO(E)PYRENE	967	19.34	0.0311	0.6452	0.0007	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0218	0.7543	0.0008	0.0041 J	0.3905	0.0004	0.0027 J	0.2977	0.0003	0.0179	0.7920	0.0008	< 0.0043 U	--	--	< 0.0042 J	--	--		
C1-Benzanthracenes/Chrysenes	930	18.6	0.0223	0.4627	0.0005	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0166	0.5744	0.0006	0.0036 J	0.3429	0.0004	< 0.0049 U	--	--	0.0141	0.6239	0.0007	< 0.0043 U	--	--	< 0.0042 J	--	--		
C1-Fluoranthenes/Pyrenes	769	15.38	0.0350	0.7261	0.0009	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0286	0.9896	0.0013	0.0076	0.7238	0.0009	0.0054	0.5954	0.0008	0.0258	1.1416	0.0015	0.0035 J	1.2727	0.0017	< 0.0042 J	--	--		
C1-Fluorenes	611	12.22	0.0050 J	0.1037	0.0002	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0050 J	0.1730	0.0003	0.0035 J	0.3333	0.0005	0.0029 J	0.3197	0.0005	0.0032 J	0.1416	0.0002	< 0.0043 U	--	--	0.0025 J	0.6394	0.0010		
C1-Naphthalenes	445	8.9	0.0066	0.1369	0.0003	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0082	0.2837	0.0006	< 0.0050 J	--	--	< 0.0049 U	--	--	< 0.0043 U	--	--	< 0.0042 J	--	--					
C1-Phenanthrenes/Anthracenes	670	13.4	0.0279	0.5788	0.0009	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0261	0.9031	0.0013	0.0081	0.7714	0.0012	0.0055	0.6064	0.0009	0.0229	1.0133	0.0015	< 0.0043 U	--	--	0.0065	1.6624	0.0025		
C2-Benzanthracenes/Chrysenes	1009	20.18	0.0109	0.2261	0.0002	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0095	0.3287	0.0003	< 0.0050 U	--	--	< 0.0049 U	--	--	0.0046 J	0.2035	0.0002	< 0.0043 U	--	--	< 0.0042 J	--	--		
C2-Fluorenes	687	13.74	0.0061 J	0.1266	0.0002	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0063	0.2180	0.0003	< 0.0050 U	--	--	0.0025 J	0.2756	0.0004	0.0054	0.2389	0.0003	0.0025 J	0.9091	0.0013	0.0046	1.1765	0.0017		
C2-Naphthalenes	510	10.2	0.0085	0.1763	0.0003	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0025 J	0.1420	0.0003	0.0170	0.5882	0.0012	0.0197	1.8762	0.0037	0.0083	0.9151	0.0018	0.0074	0.3274	0.0006	< 0.0043 U	--	< 0.0042 J	--	--
C2-Phenanthrenes/Anthracenes	745	14.9	0.0155	0.3216	0.0004	< 0.0050 U	--	--	< 0.0048 U	--	--	0.0143	0.4948	0.0007	< 0.0050 U	--	--	0.0030 J	0.3308	0.0004	0.0162	0.7168	0.0010	0.0046	1.6727	0.0022	0.0071	1.8159	0.0024		
C3-Benzanthracenes/Chrysenes	1213	24.26	< 0.0062 U	--	--	< 0.0050 U	--	--	< 0.0048 U	--	--	< 0.0054 U	--	--	< 0.0050 U	--	--	< 0.0049 U	--	--	< 0.0050 U	--	--	< 0.0043 U	--	--	0.0041 J	1.0486	0.0014		
C3-Fluorenes	768	15.36	< 0.0062 U	--	--	< 0.0050 U	--	--	< 0.0048 U	--	--	< 0.0054 U	--	--	< 0.0050 U	--	--	< 0.0049 U	--	--	0.0043 J	0.1903	0.0002	< 0.0043 U	--	--	0.0041 J</				

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect E																																			
			SDE1			SDE1			SDE1			SDE2			SDE2			SDE3			SDE3			SDE4														
			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/6/2013			11/8/2013			11/8/2013			11/8/2013														
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU									
ACENAPHTHENE	491	9.82	0.0058 J	0.1841	0.0004	< 0.0045 U	--	--	< 0.0044 U	--	--	< 0.0060 U	--	--	< 0.0053 U	--	--	< 0.0044 U	--	--	< 0.0043 U	--	--	< 0.0042 U	--	--	< 0.0058 U	--	--	< 0.0070 U	--	--	< 0.0067 U	--				
ACENAPHTHYLENE	452	9.04	0.0216	0.6857	0.0015	0.0060	0.2510	0.0006	< 0.0044 U	--	--	0.0066	0.2000	0.0004	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0093	0.3163	0.0007	0.0042 J	0.2132	0.0005	< 0.0042 U	--	--	0.0133	0.2131	0.0005	0.0143	0.2049	0.0005	0.0140	0.2393	0.0005
ANTHRACENE	594	11.88	0.0203	0.6444	0.0011	0.0054	0.2259	0.0004	< 0.0044 U	--	--	0.0075	0.2273	0.0004	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0107	0.3639	0.0006	0.0038 J	0.1929	0.0003	< 0.0042 U	--	--	0.0151	0.2420	0.0004	0.0146	0.2092	0.0004	0.0149	0.2547	0.0004
BENZO(A)ANTHRACENE	841	16.82	0.0659	2.0921	0.0025	0.0144	0.6025	0.0007	< 0.0044 U	--	--	0.0276	0.8364	0.0010	< 0.0053 U	--	--	0.0031 J	0.8267	0.0010	0.0268	0.9116	0.0011	0.0085	0.4315	0.0005	< 0.0042 U	--	--	0.0386	0.6186	0.0007	0.0300	0.4298	0.0005	0.0320	0.5470	0.0007
BENZO(A)PYRENE	964	19.28	0.0835	2.6508	0.0027	0.0178	0.7448	0.0008	< 0.0044 U	--	--	0.0326	0.9879	0.0010	0.0032 J	0.1951	0.0002	0.0033 J	0.8800	0.0009	1.0408	0.0011	0.0096	0.4873	0.0005	< 0.0042 U	--	--	0.0440	0.7051	0.0007	0.0348	0.4986	0.0005	0.0369	0.6308	0.0007	
BENZO(B)FLUORANTHENE	979	19.58	0.0747	2.3714	0.0024	0.0153	0.6402	0.0007	< 0.0044 U	--	--	0.0316	0.9576	0.0010	0.0030 J	0.1829	0.0002	0.0035 J	0.9333	0.0010	0.0254	0.8639	0.0009	0.0798	< 0.0042 U	--	--	0.0376	0.6026	0.0006	0.0290	0.4155	0.0004	0.0307	0.5248	0.0005		
BENZO(G,H)PERYLENE	1095	21.9	0.0524	1.6635	0.0015	0.0112	0.4686	0.0004	< 0.0044 U	--	--	0.0268	0.8121	0.0007	< 0.0053 U	--	--	0.0033 J	0.8800	0.0008	0.0186	0.6327	0.0006	0.0058	0.2944	0.0003	< 0.0042 U	--	--	0.0284	0.4551	0.0004	0.0233	0.3338	0.0003	0.0239	0.4085	0.0004
BENZO(K)FLUORANTHENE	980	19.6	0.0752	2.3873	0.0024	0.0167	0.6987	0.0007	< 0.0044 U	--	--	0.0325	0.9848	0.0010	0.0029 J	0.1768	0.0002	0.0034 J	0.9067	0.0009	0.2811	0.9558	0.0010	0.0085	0.4315	0.0004	< 0.0042 U	--	--	0.0397	0.6362	0.0006	0.0315	0.4513	0.0005	0.0333	0.5692	0.0006
BENZO(E)PYRENE	967	19.34	0.0594	1.8857	0.0020	0.0126	0.5272	0.0005	< 0.0044 U	--	--	0.0159	0.4818	0.0005	< 0.0053 U	--	--	0.0031 J	0.8267	0.0009	0.0211	0.7177	0.0007	0.0064	0.3249	0.0003	< 0.0042 U	--	--	0.0319	0.5112	0.0005	0.0263	0.3768	0.0004	0.0269	0.4598	0.0005
C1-Benzoanthracenes/Chrysenes	930	18.6	0.0409	1.2984	0.0014	0.0107	0.4477	0.0005	< 0.0044 U	--	--	0.0135	0.4091	0.0004	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0159	0.5408	0.0006	0.0059	0.2995	0.0003	< 0.0042 U	--	--	0.0297	0.4760	0.0005	0.0292	0.4183	0.0004	0.0318	0.5436	0.0006
C1-Fluoranthenes/Pyrenes	769	15.38	0.0646	2.0508	0.0027	0.0162	0.6778	0.0009	< 0.0044 U	--	--	0.0205	0.6122	0.0008	0.0032 J	0.1951	0.0003	0.0025 J	0.6667	0.0009	0.1915	0.0012	0.0099	0.5025	0.0007	< 0.0042 U	--	--	0.0458	0.7324	0.0010	0.0418	0.5989	0.0008	0.0454	0.7761	0.0010	
C1-Fluorenes	611	12.22	0.0662	0.1968	0.0003	0.0025 J	0.1046	0.0002	< 0.0044 U	--	--	0.0039 J	0.1182	0.0002	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0032 J	0.1088	0.0002	< 0.0043 U	--	--	< 0.0042 U	--	--	0.0067	0.1074	0.0002	0.0069 J	0.0989	0.0002	0.0090	0.1538	0.0003
C1-Naphthalenes	445	8.9	0.0991	0.2889	0.0006	0.0034 J	0.1423	0.0003	< 0.0044 U	--	--	0.0112	0.3394	0.0008	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0154	0.5238	0.0012	< 0.0043 U	--	--	< 0.0042 U	--	--	0.0880	1.3141	0.0030	0.0788	1.1289	0.0025	0.0920	1.5726	0.0035
C1-Phenanthrenes/Anthracenes	670	13.4	0.0476	1.5111	0.0023	0.0137	0.5732	0.0009	< 0.0044 U	--	--	0.0193	0.5848	0.0009	0.0038 J	0.2317	0.0003	0.0027 J	0.7200	0.0011	0.0259	0.8810	0.0013	0.0094	0.4772	0.0007	< 0.0042 U	--	--	0.0696	1.1154	0.0017	0.0699	1.0014	0.0015	0.0843	1.4410	0.0022
C2-Benzoanthracenes/Chrysenes	1009	20.18	0.0187	0.5937	0.0006	0.0057	0.2385	0.0002	< 0.0044 U	--	--	0.0073	0.2212	0.0002	< 0.0053 U	--	--	< 0.0044 U	--	--	0.0076	0.2585	0.0003	0.0035 J	0.1777	0.0002	< 0.0042 U	--	--	0.0166	0.2660	0.0003	0.0205	0.2937	0.0005	0.0188	0.3214	0.0003
C2-Fluorenes</td																																						

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect F																																		
			SDF1			SDF1			SDF1			SDF2			SDF2			SDF3			SDF3			SDF4													
			11/4/2013			11/4/2013			11/4/2013			11/7/2013			11/7/2013			11/7/2013			11/8/2013			11/8/2013													
			0 - 6			6 - 12			18 - 24			0 - 6			6 - 12			18 - 24			0 - 6			18 - 24													
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU								
ACENAPHTHENE	491	9.82	0.0028 J	0.3789	0.0008 < 0.0048 U	--	--	< 0.0053 U	--	--	< 0.0058 U	--	--	< 0.0055 U	--	--	< 0.0047 U	--	--	< 0.0053 J	--	--	< 0.0046 U	--	--	< 0.0049 U	--	--	< 0.0065 U	--	--	< 0.0043 U	--	--			
ACENAPHTHYLENE	452	9.04	0.0159	2.1516	0.0048 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0115	0.5399	0.0012	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0057	0.2740	0.0006	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0080	0.3883	0.0009	< 0.0045 U	--	--	< 0.0043 U	--	--
ANTHRACENE	594	11.88	0.0171	2.3139	0.0039 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0105	0.4930	0.0008	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0056	0.2692	0.0005	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0097	0.4709	0.0008	< 0.0045 U	--	--	< 0.0043 U	--	--
BENZO(A)ANTHRACENE	841	16.82	0.0135 J	1.8268	0.0022 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0224	1.0516	0.0013	0.0031 J	0.2067	0.0002	< 0.0047 U	--	--	0.0138	0.6635	0.0008	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0295	1.4320	0.0017	< 0.0045 U	--	--	0.0031 J	1.2351	0.0015
BENZO(A)PYRENE	964	19.28	0.0157 J	2.1245	0.0022 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0270	1.2676	0.0013	0.0036 J	0.2400	0.0002	< 0.0047 U	--	--	0.0165	0.7933	0.0008	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0339	1.6456	0.0017	< 0.0045 U	--	--	0.0037 J	1.4741	0.0015
BENZO(B)FLUORANTHENE	979	19.58	0.0143 J	1.9350	0.0020 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0226	1.0610	0.0011	0.0031 J	0.2067	0.0002	< 0.0047 U	--	--	0.0143	0.6875	0.0007	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0298	1.4466	0.0015	< 0.0045 U	--	--	0.0031 J	1.2351	0.0013
BENZO(G,H)PERYLENE	1095	21.9	0.0147 J	1.9892	0.0018 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0170	0.7981	0.0007	< 0.0048 U	--	--	< 0.0047 U	--	--	0.0099	0.4760	0.0004	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0208	1.0907	0.0009	< 0.0045 U	--	--	0.0022 J	0.8765	0.0008
BENZO(K)FLUORANTHENE	980	19.6	0.0132 J	1.7862	0.0018 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0249	1.1690	0.0012	0.0032 J	0.2133	0.0002	< 0.0047 U	--	--	0.0150	0.7212	0.0007	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0297	1.4417	0.0015	< 0.0045 U	--	--	< 0.0043 U	--	--
BENZO(E)PYRENE	967	19.34	0.0161 J	2.1786	0.0023 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0194	0.9108	0.0009	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0118	0.5673	0.0006	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0235	1.1408	0.0012	< 0.0045 U	--	--	0.0026 J	1.0359	0.0011
C1-Benzenanthrenes/Chrysenes	930	18.6	0.0254	3.4371	0.0037 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0152	0.7136	0.0008	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0079	0.3798	0.0004	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0157	0.7621	0.0008	< 0.0045 U	--	--	0.0022 J	0.8765	0.0009
C1-Fluoranthenes/Pyrenes	769	15.38	0.0253	3.4235	0.0045 0.0054	0.3158	0.0004	0.0052 J	0.2419	0.0003	0.0264	1.2394	0.0016	0.0047 J	0.3133	0.0004	< 0.0047 U	--	--	0.0151	0.7266	0.0009	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0272	1.3204	0.0017	< 0.0045 U	--	--	0.0041 J	1.6335	0.0021
C1-Fluorennes	611	12.22	0.0135	1.8268	0.0030 0.0046 J	0.2690	0.0004	0.0046 J	0.2140	0.0004	0.0045 J	0.2113	0.0003	0.0031 J	0.2067	0.0003	< 0.0047 U	--	--	0.0028 J	0.1346	0.0002	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0034 J	0.1650	0.0003	< 0.0045 U	--	--	< 0.0043 U	--	--
C1-Naphthalenes	445	8.9	0.0249	3.3694	0.0076 0.0058	0.3392	0.0008	0.0063	0.2930	0.0007	0.0065	0.3052	0.0007	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0053 J	0.2903	0.0007	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0045 U	0.1065	0.001	< 0.0043 U	--	--	< 0.0043 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0459	6.2111	0.0093 0.0096	0.5614	0.0008	0.0074	0.3442	0.0005	0.0229	1.0751	0.0016	0.0047 J	0.3133	0.0005	< 0.0047 U	--	--	0.0116	0.5577	0.0008	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0195	0.9466	0.0014	< 0.0045 U	--	--	< 0.0043 U	--	--
C2-Benzenanthrenes/Chrysenes	1009	20.18	0.0254	3.4371	0.0034 < 0.0048 U	--	--	< 0.0053 U	--	--	0.0070	0.3286	0.0003	< 0.0055 U	--	--	< 0.0047 U	--	--	0.0039 J	0.1875	0.0002	< 0.0046 U	--	--	< 0.0049 U	--	--	0.0073	0.3544	0.0004	< 0.0045 U	--	--	< 0.0043 U	--	--
C2-Fluorennes	687	13.74	0.0286																																		

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect Sample ID Sample Date Sample Interval (inches)	NYSDEC Screening Criteria	Transect G																		
		SDG1 11/6/2013 0 - 6				SDG1 11/6/2013 6 - 12				SDG1 11/6/2013 18 - 24				SDG2 11/7/2013 0 - 6						
		mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	
		Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A																	
ACENAPHTHENE	491	9.82	0.0045 J	0.1711	0.0003	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0037 J	0.1917	0.0004	< 0.0055 U	--	--	< 0.0044 U	--	--
ACENAPHTHYLENE	452	9.04	0.0167	0.6350	0.0014	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0116	0.6010	0.0013	< 0.0055 U	--	--	< 0.0044 U	--	--
ANTHRACENE	594	11.88	0.0176	0.6692	0.0011	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0121	0.6269	0.0011	< 0.0055 U	--	--	< 0.0044 U	--	--
BENZO(A)ANTHRACENE	841	16.82	0.0419	1.5932	0.0019	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0230	1.1917	0.0014	0.0045 J	0.3462	0.0004	< 0.0044 U	--	--
BENZO(A)PYRENE	964	19.28	0.0495	1.8821	0.0020	0.0025 J	0.1908	0.0002	< 0.0048 U	--	--	0.0252	1.3057	0.0014	0.0050 J	0.3846	0.0004	< 0.0044 U	--	--
BENZO(B)FLUORANTHENE	979	19.58	0.0420	1.5970	0.0016	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0216	1.1192	0.0011	0.0045 J	0.3462	0.0004	< 0.0044 U	--	--
BENZO(G,H,I)PERYLENE	1095	21.9	0.0336	1.2776	0.0012	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0151	0.7824	0.0007	0.0035 J	0.2692	0.0002	< 0.0044 U	--	--
BENZO(K)FLUORANTHENE	980	19.6	0.0460	1.7490	0.0018	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0235	1.2176	0.0012	0.0048 J	0.3692	0.0004	< 0.0044 U	--	--
BENZO(E)PYRENE	967	19.34	0.0352	1.3384	0.0014	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0180	0.9326	0.0010	0.0038 J	0.2923	0.0003	< 0.0044 U	--	--
C1-Benzanthracenes/Chrysenes	930	18.6	0.0236	0.8973	0.0010	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0159	0.8238	0.0009	0.0032 J	0.2462	0.0003	< 0.0044 U	--	--
C1-Fluoranthenes/Pyrenes	769	15.38	0.0425	1.6160	0.0021	< 0.0047 U	--	--	0.0029 J	0.9295	0.0012	0.0289	1.4974	0.0019	0.0066	0.5077	0.0007	< 0.0044 U	--	--
C1-Fluorenes	611	12.22	0.0046 J	0.1749	0.0003	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0045 J	0.2332	0.0004	0.0031 J	0.2385	0.0004	< 0.0044 U	--	--
C1-Naphthalenes	445	8.9	0.0079	0.3004	0.0007	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0076	0.3938	0.0009	< 0.0055 J	--	--	< 0.0044 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0334	1.2700	0.0019	< 0.0047 U	--	--	0.0034 J	1.0897	0.0016	0.0262	1.3575	0.0020	0.0062	0.4769	0.0007	< 0.0044 U	--	--
C2-Benzanthracenes/Chrysenes	1009	20.18	0.01	0.3802	0.0004	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0083	0.4301	0.0004	< 0.0055 U	--	--	< 0.0044 U	--	--
C2-Fluorenes	687	13.74	0.0067	0.2548	0.0004	< 0.0047 U	--	--	0.0059	1.8910	0.0028	0.0068	0.3523	0.0005	0.0029 J	0.2231	0.0003	< 0.0044 U	--	--
C2-Naphthalenes	510	10.2	0.0088	0.3346	0.0007	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0091	0.4715	0.0009	0.0048 J	0.3692	0.0007	< 0.0044 U	--	--
C2-Phenanthrenes/Anthracenes	745	14.9	0.0185	0.7034	0.0009	< 0.0047 U	--	--	0.0097	3.1090	0.0042	0.0149	0.7720	0.0010	0.0034 J	0.2615	0.0004	< 0.0044 U	--	--
C3-Benzanthracenes/Chrysenes	1213	24.26	0.0060	0.2281	0.0002	< 0.0047 U	--	--	< 0.0048 U	--	--	< 0.0056 U	--	--	< 0.0055 U	--	--	< 0.0044 U	--	--
C3-Fluorenes	768	15.36	0.0061	0.2319	0.0003	< 0.0047 U	--	--	0.0081	2.5962	0.0034	< 0.0056 U	--	--	< 0.0055 U	--	--	< 0.0044 U	--	--
C3-Naphthalenes	581	11.62	0.0092	0.3498	0.0006	< 0.0047 U	--	--	0.0094	3.0128	0.0052	0.0102	1.5285	0.0009	0.0037 J	0.2846	0.0005	< 0.0044 U	--	--
C3-Phenanthrenes/Anthracenes	830	16.6	0.0109	0.4144	0.0005	< 0.0047 U	--	--	0.0096	3.0769	0.0037	0.0079	0.4093	0.0005	< 0.0055 U	--	--	< 0.0044 U	--	--
C4-Benzanthracenes/Chrysenes	1213	24.26	< 0.0055 U	--	< 0.0047 U	--	--	< 0.0048 U	--	--	< 0.0056 U	--	--	< 0.0055 U	--	--	< 0.0044 U	--	--	
C4-Naphthalenes	657	13.14	0.0089	0.3384	0.0005	< 0.0047 U	--	--	0.0231	7.4038	0.0113	0.0085	0.4404	0.0007	< 0.0055 U	--	--	< 0.0044 U	--	--
C4-Phenanthrenes/Anthracenes	670	13.4	0.0065	0.2471	0.0004	< 0.0047 U	--	--	0.0056	1.7949	0.0027	0.0035 J	0.1813	0.0003	< 0.0055 U	--	--	< 0.0044 U	--	--
CHRYSENE	843	16.86	0.0571	2.1711	0.0026	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0305	1.5803	0.0019	0.0064	0.4923	0.0006	< 0.0044 U	--	--
DIBENZ(A,H)ANTHRACENE	1122	22.44	0.0179	0.6806	0.0006	0.0057	0.4351	0.0004	< 0.0048 U	--	--	0.0039 J	0.2021	0.0002	< 0.0055 U	--	--	< 0.0044 U	--	--
FLUORANTHENE	708	14.16	0.104	3.9544	0.0056	0.0023 J	0.1756	0.0002	< 0.0048 U	--	--	0.0612	3.1710	0.0045	0.0106	0.8154	0.0012	< 0.0044 U	--	--
FLUORENE	539	10.78	0.0072	0.2738	0.0005	< 0.0047 U	--	--	< 0.0048 U	--	--	0.0050 J	0.2591	0.0005	< 0.0055 U	--	--	< 0.0044 U	--	--
INDENO(1,2,3-C,D)PYRENE	1115	22.3	0.0344	1.3080	0.0012	0.0034 J	0.2595	0.0002	< 0.0048 U	--	--	0.0148 J	0.7668	0.0007	0.0031 J	0.2385	0.0002	< 0.0044 U	--	--
NAPHTHALENE	385	7.7	0.0117	0.4449	0.0012	< 0.0047 U	--	--	< 0.0048 U	--	--</									

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect H																								
			SDH1			SDH1			SDH1			SDH2			SDH2			SDH3			SDH3						
			11/6/2013			11/6/2013			11/6/2013			11/7/2013			11/7/2013			11/7/2013			11/7/2013						
Sample Date	Sample Interval (inches)	Sediment Guidance Value (µg/gOC)	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24				
Chemical Name	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU		
ACENAPHTHENE	491	9.82	0.0116 J	0.0817	0.0002 < 0.0049 U	--	--	< 0.0046 U	--	--	< 0.0060 J	--	--	< 0.0051 U	--	--	< 0.0048 U	--	--	0.0089 J	0.0742	0.0002 < 0.0055 U	--	--	< 0.0048 U	--	--
ACENAPHTHYLENE	452	9.04	0.0210	0.1479	0.0003 0.0098 0.8305	0.0018 < 0.0046 U	--	--	0.0077 0.2059	0.0005 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0389 0.3242	0.0007 0.0036 J	0.2069	0.0005 < 0.0048 U	--	--	0.0005 < 0.0048 U	--	--			
ANTHRACENE	594	11.88	0.0351	0.2472	0.0004 0.0097 0.8220	0.0014 < 0.0046 U	--	--	0.0068 0.1818	0.0003 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0346 0.2883	0.0005 < 0.0055 U	--	--	< 0.0048 U	--	--	0.0005 < 0.0048 U	--	--		
BENZO(A)ANTHRACENE	841	16.82	0.12	0.8451	0.0010 0.0194 1.6441	0.0020 0.025 J	0.5285	0.0006 0.0183 0.4893	0.0006 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0837 0.6975	0.0008 0.0054 J	0.3103	0.0004 < 0.0048 U	--	--	0.0004 < 0.0048 U	--	--	0.0004 < 0.0048 U	--	--	
BENZO(A)PYRENE	964	19.28	0.157	1.1056	0.0011 0.0228 1.9322	0.0020 0.028 J	0.5920	0.0006 0.0216 0.5775	0.0006 < 0.0051 U	--	--	< 0.0048 U	--	--	0.101 0.8417	0.0009 0.0065 0.3736	0.0004 < 0.0048 U	--	--	0.0004 < 0.0048 U	--	--	0.0004 < 0.0048 U	--	--		
BENZO(B)FLUORANTHENE	979	19.58	0.158	1.1127	0.0011 0.0219 1.8559	0.0019 0.0228 J	0.5920	0.0006 0.0188 0.5027	0.0005 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0851 0.7092	0.0007 0.0055 0.3161	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--		
BENZO(G,H,I)PERYLENE	1095	21.9	0.121	0.8521	0.0008 0.0165 1.3983	0.0013 < 0.0046 U	--	--	0.0130 0.3476	0.0003 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0582 0.4850	0.0004 0.0044 J	0.2529	0.0002 < 0.0048 U	--	--	0.0002 < 0.0048 U	--	--	0.0002 < 0.0048 U	--	--
BENZO(K)FLUORANTHENE	980	19.6	0.146	1.0282	0.0010 0.0216 1.8305	0.0019 0.026 J	0.5497	0.0006 0.0193 0.5160	0.0005 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0922 0.7683	0.0008 0.0057 0.3276	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--		
BENZO(E)PYRENE	967	19.34	0.128	0.9014	0.0009 0.0178 1.5085	0.0016 0.024 J	0.5074	0.0005 0.0149 0.3984	0.0004 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0708 0.5900	0.0006 0.0044 J	0.2529	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	
C1-Benzanthracenes/Chrysenes	930	18.6	0.0631	0.4444	0.0005 0.0164 1.3898	0.0015 < 0.0046 U	--	--	0.0114 0.3048	0.0003 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0541 0.4508	0.0005 0.0049 J	0.2816	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--
C1-Fluoranthenes/Pyrenes	769	15.38	0.0922	0.6493	0.0008 0.0256 2.1695	0.0028 0.046 J	0.9725	0.0013 0.0205 0.5481	0.0007 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0902 0.7517	0.0010 0.0078 0.4483	0.0006 < 0.0048 U	--	--	0.0006 < 0.0048 U	--	--	0.0006 < 0.0048 U	--	--		
C1-Fluorenes	611	12.22	0.0182	0.1282	0.0002 0.0051 0.4322	0.0007 0.025 J	0.5285	0.0009 0.0034 J	0.0909	0.0001 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0086 J 0.0717	0.0001 < 0.0055 U	--	--	< 0.0048 U	--	--	0.0001 < 0.0048 U	--	--		
C1-Naphthalenes	445	8.9	0.0135 J	0.0951	0.0002 0.0061 0.5169	0.0012 < 0.0046 U	--	--	< 0.0060 J	--	--	< 0.0051 U	--	--	< 0.0048 U	--	--	0.0118 0.0983	0.0002 < 0.0055 U	--	--	< 0.0048 U	--	--	0.0002 < 0.0048 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0811	0.5711	0.0009 0.0241 2.0424	0.0030 0.044 J	0.8668	0.0013 0.0159 0.4251	0.0006 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0689 0.5742	0.0009 0.0068 0.3908	0.0006 < 0.0048 U	--	--	0.0006 < 0.0048 U	--	--	0.0006 < 0.0048 U	--	--		
C2-Benzanthracenes/Chrysenes	1009	20.18	0.0340	0.2394	0.0002 0.0091 0.8136	0.0008 < 0.0046 U	--	--	0.0062 0.1658	0.0002 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0247 0.2058	0.0002 0.0032 J	0.1839	0.0002 < 0.0048 U	--	--	0.0002 < 0.0048 U	--	--	0.0002 < 0.0048 U	--	--
C2-Fluorenes	687	13.74	0.0319	0.2246	0.0003 0.0076 0.6441	0.0009 0.0209 J	0.6131	0.0009 0.0035 J	0.0936	0.0001 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0115 0.0958	0.0001 < 0.0055 U	--	--	< 0.0048 U	--	--	0.0001 < 0.0048 U	--	--		
C2-Naphthalenes	510	10.2	0.0399	0.2810	0.0006 0.0101 0.8559	0.0017 0.0052 1.0994	0.0022 0.0043 J	0.1150	0.0002 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0124 0.1033	0.0002 < 0.0055 U	--	--	< 0.0048 U	--	--	0.0002 < 0.0048 U	--	--			
C2-Phenanthrenes/Anthracenes	745	14.9	0.0448	0.3155	0.0004 0.0152 1.2881	0.0017 0.0034 J	0.7188	0.0010 0.0080 0.2139	0.0003 < 0.0051 U	--	--	< 0.0048 U	--	--	0.0384 0.3200	0.0004 0.0037 J	0.2126	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	0.0003 < 0.0048 U	--	--	
C3-Benzanthracenes/Chrysenes	1213	24.26	< 0.014 U	--	--	< 0.0049 U	--	--	< 0.0046 U	--	--	< 0.0060 U	--	--	< 0.0051 U	--	--	< 0.0048 U	--	--	0.0136						

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect I																																			
			SDI1			SDI1 (Dup)			SDI1			SDI1 (Dup)			SDI1			SDI2			SDI2			SDI3														
			11/6/2013		11/6/2013		11/6/2013		11/6/2013		11/6/2013		11/7/2013		11/7/2013		11/7/2013		11/5/2013		11/5/2013		11/5/2013															
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU									
ACENAPHTHENE	491	9.82	< 0.0053 U	--	--	0.0039 J	0.1500	0.0003	< 0.0054 U	--	--	< 0.0048 U	--	--	< 0.0052 U	--	--	0.0189 J	0.0458	0.0001	0.0185	0.0830	0.0002	< 0.0071 U	--	--	< 0.0067 U	--	--	< 0.0046 U	--	--						
ACENAPHTHYLENE	452	9.04	0.0046 J	0.1614	0.0004	0.0151	0.5808	0.0013	< 0.0054 U	--	--	0.0037 J	0.3811	0.0008	< 0.0052 U	--	--	0.0305 J	0.0738	0.0002	0.0400	0.1794	0.0004	0.0167	0.3070	0.0007	0.0060 J	0.1824	0.0004	0.0090	1.7078	0.0038	< 0.0046 U	--	--			
ANTHRACENE	594	11.88	0.0055	0.1930	0.0003	0.0198	0.7615	0.0013	< 0.0054 U	--	--	0.0042 J	0.4325	0.0007	< 0.0052 U	--	--	0.0444	0.1075	0.0002	0.0458	0.2054	0.0003	0.0136	0.2500	0.0004	0.0068	0.2067	0.0028	< 0.0046 U	--	--						
BENZO(A)ANTHRACENE	841	16.82	0.0124	0.4351	0.0005	0.0476	1.8308	0.0022	0.0038 J	0.2171	0.0003	0.0116	1.1946	0.0014	0.0031 J	0.4486	0.0005	< 0.0043 U	--	--	0.155	0.3753	0.0004	0.13	0.5830	0.0007	0.0342	0.6287	0.0007	0.0177	0.5380	0.0006	3.7951	0.0045	0.0055	1.4360	0.0017	
BENZO(A)PYRENE	964	19.28	0.0147	0.5158	0.0005	0.0570	2.1923	0.0023	0.0046 J	0.2629	0.0003	0.0138	0.4949	0.0006	< 0.0043 U	--	--	0.193	0.4673	0.0005	0.165	0.7399	0.0008	0.0406	0.7463	0.0008	0.0687	0.0007	0.0234	4.4402	0.0046	0.0067	1.7493	0.0018				
BENZO(B)FLUORANTHENE	979	19.58	0.0135	0.4737	0.0005	0.0530	2.0385	0.0021	0.0049 J	0.2800	0.0003	0.0140	1.4418	0.0015	0.0045 J	0.6512	0.0007	< 0.0043 U	--	--	0.178	0.4310	0.0004	0.147	0.6592	0.0007	0.0357	0.6563	0.0007	0.0218	0.6626	0.0007	0.0228	4.3264	0.0044	0.0063	1.6449	0.0017
BENZO(G,H)PERYLENE	1095	21.9	0.0105	0.3684	0.0003	0.0410	1.5769	0.0014	0.0041 J	0.2343	0.0002	0.0115	1.1843	0.0011	0.0033 J	0.4776	0.0004	< 0.0043 U	--	--	0.125	0.3027	0.0003	0.103	0.4619	0.0004	0.0237	0.4357	0.0004	0.0170	0.5167	0.0005	0.0161	3.0550	0.0028	0.049	1.2794	0.0012
BENZO(K)FLUORANTHENE	980	19.6	0.0138	0.4842	0.0005	0.0552	2.1231	0.0022	0.0043 J	0.2457	0.0003	0.0131	1.3491	0.0014	0.0035 J	0.5065	0.0005	< 0.0043 U	--	--	0.176	0.4262	0.0004	0.15	0.6726	0.0007	0.0364	0.6691	0.0007	0.0209	0.6553	0.0006	0.0238	4.5161	0.0046	0.0065	1.6971	0.0017
BENZO(E)PYRENE	967	19.34	0.0115	0.4035	0.0004	0.0438	1.6846	0.0017	0.0041 J	0.2343	0.0002	0.0121	1.2461	0.0013	0.0035 J	0.5065	0.0005	< 0.0043 U	--	--	0.142	0.3438	0.0004	0.118	0.5291	0.0005	0.0280	0.5147	0.0005	0.0182	0.5532	0.0006	0.0180	3.4156	0.0035	0.0053	1.3838	0.0014
C1-Benzoanthracenes/Chrysenes	930	18.6	0.0088	0.3088	0.0003	0.0301	1.5777	0.0012	0.0035 J	0.2000	0.0002	0.0102	1.0505	0.0006	< 0.0043 U	--	--	0.0737	0.1785	0.0002	0.0714	0.3202	0.0003	0.0246	0.4522	0.0005	0.0116	0.3526	0.0004	0.0138	2.6186	0.0028	0.0041 J	1.0705	0.0012			
C1-Fluoranthenes/Pyrenes	769	15.38	0.0162	0.5684	0.0007	0.0500	1.9231	0.0025	0.0062	0.3543	0.0005	0.0138	1.4212	0.0018	0.0054 J	0.7815	0.0010	0.0033 J	0.4130	0.0005	0.12	0.2906	0.0004	0.119	0.5336	0.0007	0.0406	0.7463	0.0010	0.0199	0.6049	0.0008	0.0248	4.7059	0.0061	0.0073	1.9067	0.0025
C1-Fluorenes	611	12.22	0.0041 J	0.1439	0.0002	0.0660	0.2308	0.0004	< 0.0054 U	--	--	0.0041 J	0.4222	0.0007	0.0028 J	0.4052	0.0007	< 0.0043 U	--	--	0.0162	0.0726	0.0001	0.0051 J	0.0938	0.0002	0.0041 J	0.1246	0.0002	0.0035 J	0.6641	0.0011	0.0224 J	0.024 J	0.0011	0.2626	0.0010	
C1-Naphthalenes	445	8.9	0.0035 J	0.1228	0.0003	0.0079	0.3038	0.0007	< 0.0054 U	--	--	0.0044 J	0.4531	0.0010	< 0.0052 J	--	--	< 0.0043 U	--	--	0.0187	0.0839	0.0002	0.0072	0.1324	0.0003	< 0.0067 J	--	--	< 0.0053 U	--	--	< 0.0046 J	--	--			
C1-Phenanthrenes/Anthracenes	670	13.4	0.0146	0.5123	0.0008	0.0416	1.6000	0.0024	0.0069	0.3943	0.0006	0.0143	1.4727	0.0022	0.0072	1.0420	0.0016	0.0032 J	0.4005	0.0006	0.0795	0.1925	0.0003	0.0819	0.3673	0.0005	0.0313	0.5754	0.0009	0.0163	0.4954	0.0007	0.0194	3.6812	0.0055	0.0061	1.5927	0.0024
C2-Benzoanthracenes/Chrysenes	1009	20.18	0.0046 J	0.1614	0.0002	0.0133	0.5115	0.0005	0.0031 J	0.1771	0.0002	0.0073	0.7518	0.0007	0.0038 J	0.5499	0.0005	< 0.0043 U	--	--	0.0398	0.0964	0.0001	0.0329	0.1475	0.0001	0.0147	0.2702	0.0003	0.0116	0.3526	0.0004	0.0041 J					

Table 5  
 PAH Screening Table with Calculated Toxicity Units  
 November 2013 Event  
 UTC/Carrier Circle  
 Carrier-Dewitt Landfill (AOC G)  
 Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Transect K														
			SDK1			SDK1			SDK1			SDK2			SDK2		
			11/5/2013			11/5/2013			11/5/2013			11/5/2013			11/5/2013		
			0 - 6			6 - 12			18 - 24			0 - 6			6 - 12		
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU												
ACENAPHTHENE	491	9.82	0.0226	0.2093	0.0004	< 0.0051 J	--	--	< 0.0055 U	--	--	< 0.0059 J	--	--	< 0.0045 U	--	--
ACENAPHTHYLENE	452	9.04	0.122	1.1296	0.0025	0.0048 J	0.2115	0.0005	< 0.0055 U	--	--	0.0033 J	0.1347	0.0003	0.0084	1.3548	0.0030
ANTHRACENE	594	11.88	0.297	2.7500	0.0046	0.0038 J	0.1674	0.0003	< 0.0055 U	--	--	0.0053 J	0.2163	0.0004	0.0101	1.6290	0.0027
BENZO(A)ANTHRACENE	841	16.82	0.256	2.3704	0.0028	0.0072	0.3172	0.0004	0.0030 J	0.2308	0.0003	0.0138	0.5633	0.0007	0.0230	3.7097	0.0044
BENZO(A)PYRENE	964	19.28	0.259	2.3981	0.0025	0.0089	0.3921	0.0004	0.0040 J	0.3077	0.0003	0.0168	0.6857	0.0007	0.0275	4.4355	0.0046
BENZO(B)FLUORANTHENE	979	19.58	0.258	2.3889	0.0024	0.0092	0.4053	0.0004	0.0047 J	0.3615	0.0004	0.0136	0.5551	0.0006	0.0249	4.0161	0.0041
BENZO(G,H,I)PERYLENE	1095	21.9	0.193	1.7870	0.0016	0.0073	0.3216	0.0003	0.0035 J	0.2692	0.0002	0.0109	0.4082	0.0004	0.0202	3.2581	0.0030
BENZO(K)FLUORANTHENE	980	19.6	0.247	2.2870	0.0023	0.0085	0.3744	0.0004	0.0038 J	0.2923	0.0003	0.0144	0.5878	0.0006	0.0252	4.0645	0.0041
BENZO(E)PYRENE	967	19.34	0.212	1.9630	0.0020	0.0075	0.3304	0.0003	0.0038 J	0.2923	0.0003	0.0111	0.4531	0.0005	0.0210	3.3871	0.0035
C1-Benzanthracenes/Chrysenes	930	18.6	0.0163	0.1509	0.0002	0.0064	0.2819	0.0003	0.0031 J	0.2385	0.0003	0.0070	0.2857	0.0003	0.0153	2.4677	0.0027
C1-Fluoranthenes/Pyrenes	769	15.38	0.0683	0.6324	0.0008	0.0096	0.4228	0.0005	0.0052 J	0.4000	0.0005	0.0119	0.4857	0.0006	0.0250	4.0323	0.0052
C1-Fluorenes	611	12.22	< 0.0093 U	--	--	< 0.0029 J	0.1278	0.0002	0.0032 J	0.2462	0.0004	< 0.0059 J	--	--	0.0040 J	0.6452	0.0011
C1-Naphthalenes	445	8.9	0.0390	0.3611	0.0008	< 0.0051 J	--	--	< 0.0055 U	--	--	< 0.0059 J	--	--	< 0.0045 U	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0630	0.5833	0.0009	0.0092	0.4053	0.0006	0.0052 J	0.4000	0.0006	0.0089	0.3633	0.0005	0.0202	3.2581	0.0049
C2-Benzanthracenes/Chrysenes	1009	20.18	0.0163 J	0.1509	0.0001	0.0064 J	0.2818	0.0003	0.0064 J	0.2385	0.0002	< 0.0070 U	0.2857	0.0003	0.0153	2.4677	0.0024
C2-Fluorenes	687	13.74	< 0.0093 U	--	--	< 0.0039 J	0.1718	0.0003	0.0035 J	0.2692	0.0004	< 0.0059 U	--	--	0.0047	0.7581	0.0011
C2-Naphthalenes	510	10.2	0.0217	0.2009	0.0004	0.0059	0.2599	0.0005	0.0064	0.4923	0.0010	0.0038 J	0.1551	0.0003	0.0054	0.8710	0.0017
C2-Phenanthrenes/Anthracenes	745	14.9	0.0228	0.2111	0.0003	0.0061	0.2687	0.0004	0.0037 J	0.2846	0.0004	0.0041 J	0.1673	0.0002	0.0122	1.9677	0.0026
C3-Benzanthracenes/Chrysenes	1213	24.26	< 0.0093 U	--	--	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	0.0044 J	0.7097	0.0006
C3-Fluorenes	768	15.36	< 0.0093 U	--	--	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	< 0.0045 U	--	--
C3-Naphthalenes	581	11.62	0.0164	0.1519	0.0003	0.0063	0.2775	0.0005	0.0050 J	0.3846	0.0007	0.0031 J	0.1265	0.0002	0.0074	1.1935	0.0021
C3-Phenanthrenes/Anthracenes	830	16.6	0.0230	0.2130	0.0003	0.0048 J	0.2115	0.0003	0.0033 J	0.2538	0.0003	< 0.0059 U	--	--	0.0068	1.0968	0.0013
C4-Benzanthracenes/Chrysenes	1213	24.26	< 0.0093 U	--	--	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	< 0.0045 U	--	--
C4-Naphthalenes	657	13.14	< 0.0093 U	--	--	< 0.0065	0.2863	0.0004	0.0067	0.5154	0.0008	0.0040 J	0.1633	0.0002	0.0073	1.1774	0.0018
C4-Phenanthrenes/Anthracenes	670	13.4	0.0050 J	0.0463	0.0001	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	0.0044 J	0.7097	0.0011
CHRYSENE	843	16.86	0.336	3.1111	0.0037	0.0122	0.5374	0.0006	0.0057	0.4385	0.0005	0.0177	0.7224	0.0009	0.0326	5.2581	0.0062
DIBENZ(A,H)ANTHRACENE	1122	22.44	0.0699	0.6472	0.0006	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	0.0073	1.1774	0.0010
FLUORANTHENE	708	14.16	0.507	4.6944	0.0066	0.0155	0.6828	0.0010	0.0069	0.5308	0.0007	0.0329	1.3429	0.0019	0.0517	8.3387	0.0118
FLUORENE	539	10.78	0.0227	0.2102	0.0004	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	0.0029 J	0.4677	0.0009
INDENO(1,2,3-C,D)PYRENE	1115	22.3	0.207	1.9167	0.0017	0.0074	0.3260	0.0003	0.0035 J	0.2692	0.0002	0.0103	0.4204	0.0004	0.0207	3.3387	0.0030
NAPHTHALENE	385	7.7	0.0636	0.5889	0.0015	< 0.0051 U	--	--	< 0.0055 U	--	--	< 0.0059 U	--	--	0.0052	0.8387	0.0022
PERYLENE	967	19.34	0.0627	0.5806	0.0006	0.0093	0.4097	0.0004	0.0077	0.5923	0.0006	0.0061	0.2449	0.0003	0.0325	5.2419	0.0054
PHENANTHRENE	597	11.94	0.299	2.7685	0.0046	0.0094	0.4141	0.0007	0.0050 J	0.3846	0.0006	0.0192	0.7837	0.0013	0.0288	4.6452	0

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	North Swale to Sanders Creek																																			
			SDS1			SDS1			SDS1			SDS2			SDS2			SDS3			SDS3			SDS4														
			11/4/2013			11/4/2013			11/4/2013			11/4/2013			11/4/2013			11/4/2013			11/4/2013			11/4/2013														
			0 - 6			6 - 12			18 - 24			0 - 6			6 - 12			18 - 24			0 - 6			6 - 12			18 - 24											
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU									
ACENAPHTHENE	491	9.82	< 0.0047 U	--	--	< 0.0060 U	--	--	< 0.0042 U	--	--	0.0056	0.2343	0.0005	< 0.0047 J	--	--	0.0057 J	0.2969	0.0006	< 0.0049 J	--	--	< 0.0054 U	--	--	0.0086	0.5181	0.0011	0.0067 J	0.1311	0.0003	< 0.0055 U	--	--	< 0.0047 U	--	--
ACENAPHTHYLENE	452	9.04	0.0025 J	0.3365	0.0007	0.0036 J	0.1043	0.0002	< 0.0042 U	--	--	0.0146	0.6109	0.0014	0.0024 J	0.3681	0.0008	0.0156	0.0011	0.0074	0.5441	0.0012	< 0.0054 U	--	--	0.0134	0.8072	0.0018	0.0319	0.6243	0.0014	< 0.0055 U	--	--	0.0052	0.5149	0.0011	
ANTHRACENE	594	11.88	0.0024 J	0.3230	0.0005	< 0.0060 U	--	--	< 0.0042 U	--	--	0.0213	0.8912	0.0015	0.0029 J	0.4446	0.0007	0.0105	0.0063	0.4632	0.0008	< 0.0054 U	--	--	0.0106	0.6386	0.0011	0.0291	0.5695	0.0010	< 0.0055 U	--	--	0.0034 J	0.3366	0.0006		
BENZO(A)ANTHRACENE	841	16.82	0.0049	0.6595	0.0008	0.0062	0.1797	0.0002	< 0.0042 U	--	--	0.0345	1.4435	0.0017	0.0057	0.8742	0.0010	0.0188	0.0012	0.0161	1.1838	0.0014	< 0.0054 U	--	--	0.0168	1.0120	0.0012	0.0595	1.1644	0.0014	< 0.0055 U	--	--	0.0079	0.7822	0.0009	
BENZO(A)PYRENE	964	19.28	0.0056 J	0.7537	0.0008	0.0075 J	0.2174	0.0002	< 0.0042 U	--	--	0.0362 J	1.5146	0.0016	0.0060 J	0.9202	0.0010	0.0196 J	0.0208	0.0011	0.0180 J	1.3235	0.0014	< 0.0054 U	--	--	0.0180 J	1.0843	0.0011	0.0686 J	1.3425	0.0014	< 0.0055 U	--	--	0.0096 J	0.9505	0.0010
BENZO(B)FLUORANTHENE	979	19.58	0.0053 J	0.7133	0.0007	0.0074 J	0.2145	0.0002	< 0.0042 U	--	--	0.0317 J	1.3264	0.0014	0.0059 J	0.9049	0.0009	0.0179 J	0.9323	0.0010	0.0185 J	1.3603	0.0014	< 0.0054 U	--	--	0.0186 J	1.1205	0.0011	0.0599 J	1.1722	0.0012	< 0.0055 U	--	--	0.0091 J	0.9010	0.0009
BENZO(G,H)PERYLENE	1095	21.9	0.0038 J	0.5114	0.0005	0.0053 J	0.1536	0.0001	< 0.0042 U	--	--	0.0246 J	1.0293	0.0009	0.0047 J	0.7135	0.0007	0.0134 J	0.9853	0.0009	< 0.0054 U	--	--	0.0140 J	0.8434	0.0008	0.0452 J	0.8845	0.0008	< 0.0055 U	--	--	0.0067 J	0.6634	0.0006			
BENZO(K)FLUORANTHENE	980	19.6	0.0052 J	0.6999	0.0007	0.0068 J	0.1971	0.0002	< 0.0042 U	--	--	0.0316 J	1.3222	0.0013	0.0056 J	0.8589	0.0009	0.0181 J	0.9427	0.0010	0.0175 J	1.2868	0.0013	< 0.0054 U	--	--	0.0168 J	1.0120	0.0010	0.0618 J	1.2094	0.0012	< 0.0055 U	--	--	0.0086 J	0.8515	0.0009
BENZO(E)PYRENE	967	19.34	0.0041 J	0.5518	0.0006	0.0057 J	0.1652	0.0002	< 0.0042 U	--	--	0.0274 J	1.1464	0.0012	0.0052 J	0.7975	0.0008	0.0154 J	0.8021	0.0008	0.0154 J	1.1324	0.0012	< 0.0054 U	--	--	0.0162 J	0.9759	0.0010	0.0508 J	0.9941	0.0010	< 0.0055 U	--	--	0.0077 J	0.7624	0.0008
C1-Benzenethenes/Chrysenes	930	18.6	0.0041 J	0.5518	0.0006	0.0062	0.1797	0.0002	< 0.0042 U	--	--	0.0175	0.7322	0.0008	0.0038 J	0.5828	0.0006	0.0142	0.7396	0.0008	< 0.0054 U	--	--	0.0129	0.7771	0.0008	0.0374	0.7319	0.0008	< 0.0055 U	--	--	0.0061	0.6040	0.0006			
C1-Fluoranthenes/Pyrenes	769	15.38	0.0088	1.1844	0.0015	0.0122	0.3536	0.0005	0.0042	0.8268	0.0011	0.0359	1.5021	0.0020	0.0081	1.2423	0.0016	0.0288	1.5000	0.0020	0.0187	1.3750	0.0018	0.0045 J	0.3103	0.0004	0.0250	1.5060	0.0020	0.0711	1.3914	0.0018	0.0038 J	1.3010	0.0002	0.0122	1.2079	0.0016
C1-Fluorenes	611	12.22	0.0066	0.8883	0.0015	0.0064	0.1855	0.0003	0.0039 J	0.7677	0.0013	0.0082	0.4331	0.0006	0.0039 J	0.5982	0.0010	0.0088	0.4583	0.0008	0.0061	0.4485	0.0007	0.0044 J	0.3034	0.0005	0.0084	0.5060	0.0008	0.0085	0.1663	0.0003	0.0033 J	0.1130	0.0002	0.0068	0.6733	0.0011
C1-Naphthalenes	445	8.9	0.0062	0.8345	0.0019	0.0080	0.2319	0.0005	0.0051	1.0039	0.0023	0.0102	0.4268	0.0010	0.0091	1.3957	0.0031	0.0162	0.8438	0.0019	0.0145	1.0662	0.0024	0.0089	0.6138	0.0014	0.0266	1.6024	0.0036	0.0194	0.3796	0.0009	0.0075	0.2568	0.0006	0.0088	0.8713	0.0020
C1-Phenanthrenes/Anthracenes	670	13.4	0.0131	1.7631	0.0026	0.0150	0.4348	0.0006	0.0064	1.2598	0.0019	0.0092	1.4110	0.0021	0.0292	1.2636	0.0019	0.0092	1.4110	0.0022	0.0065	1.4412	0.0022	0.0065	0.4483	0.0007	0.0274	1.6506	0.0025	0.0539	0.1048	0.0016						

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	North Swale to Sanders Creek																										
			SDS5			SDS5			SDS5			SDS6			SDS6			SDS6			SDS7								
			11/4/2013			11/4/2013			11/4/2013			11/5/2013			11/5/2013			11/5/2013			11/5/2013								
Sample Date	Sample Interval (inches)	Sediment Guidance Value (µg/gOC)	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24	0 - 6	6 - 12	18 - 24						
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU						
ACENAPHTHENE	491	9.82	0.0219	0.4140	0.0008	< 0.0060 U	--	--	< 0.0053 U	--	--	< 0.0061 U	--	--	< 0.0057 U	--	--	< 0.0052 U	--	--	< 0.0046 U	--	--	< 0.0043 U	--				
ACENAPHTHYLENE	452	9.04	0.0353	0.6673	0.0015	0.0164	0.4955	0.0011	< 0.0053 U	--	--	0.0119	0.3260	0.0007	0.0135	0.5153	0.0011	0.0046 J	0.1549	0.0003	0.0120	0.3960	0.0009	0.0069	0.5149	0.0011	0.0038 J	0.3930	0.0009
ANTHRACENE	594	11.88	0.0382	0.7221	0.0012	0.0122	0.3686	0.0008	< 0.0053 U	--	--	0.0138	0.3781	0.0006	0.0151	0.5763	0.0010	0.0047 J	0.1582	0.0003	0.0152	0.5017	0.0008	0.0084	0.6269	0.0011	0.0031 J	0.3200	0.0005
BENZO(A)ANTHRACENE	841	16.82	0.0635	1.2004	0.0014	0.0259	0.7825	0.0009	< 0.0053 U	--	--	0.0380	1.0411	0.0012	0.0399	1.5229	0.0018	0.0124	0.4175	0.0005	0.0469	1.5479	0.0018	0.0243	1.8134	0.0022	0.0065	0.6722	0.0008
BENZO(A)PYRENE	964	19.28	0.0685 J	1.2949	0.0013	0.0301 J	0.9094	0.0009	< 0.0053 U	--	--	0.0462	1.2658	0.0013	0.0485	1.8511	0.0019	0.0148	0.4983	0.0005	0.0588	1.9406	0.0020	0.0308	2.2985	0.0024	0.0080	0.8273	0.0009
BENZO(B)FLUORANTHENE	979	19.58	0.0605 J	1.1437	0.0012	0.0247 J	0.7462	0.0008	< 0.0053 U	--	--	0.0382	1.0466	0.0011	0.0414	1.5802	0.0016	0.0123	0.4141	0.0004	0.0539	1.7789	0.0018	0.0278	2.0746	0.0021	0.0066	0.6825	0.0007
BENZO(G,H,I)PERYLENE	1095	21.9	0.0431 J	0.8147	0.0007	0.0185 J	0.5589	0.0008	< 0.0053 U	--	--	0.0288	0.7890	0.0007	0.0299	1.1412	0.0010	0.0086	0.2896	0.0003	0.0400	1.3201	0.0012	0.0206	1.5373	0.0014	0.0053	0.5481	0.0005
BENZO(K)FLUORANTHENE	980	19.6	0.0627 J	1.1853	0.0012	0.0274 J	0.8278	0.0008	< 0.0053 U	--	--	0.0413	1.1315	0.0012	0.0430	1.6412	0.0017	0.0129	0.4343	0.0004	0.0532	1.7558	0.0018	0.0270	2.0149	0.0021	0.0068	0.7032	0.0007
BENZO(E)PYRENE	967	19.34	0.0515 J	0.9735	0.0010	0.0223 J	0.6737	0.0007	< 0.0053 U	--	--	0.0312	0.8548	0.0009	0.0331	1.2634	0.0013	0.0098	0.3300	0.0003	0.0439	1.4488	0.0015	0.0227	1.6940	0.0018	0.0055	0.5688	0.0006
C1-Benzenanthracenes/Chrysenes	930	18.6	0.0387	0.7316	0.0008	0.0187	0.5650	0.0006	< 0.0053 U	--	--	0.0201	0.5507	0.0006	0.0219	0.8359	0.0009	0.0207	0.2357	0.0003	0.0223	0.7360	0.0008	0.0123	0.9179	0.0010	0.0046	0.4757	0.0005
C1-Fluoranthenes/Pyrenes	769	15.38	0.0840	1.5879	0.0021	0.0349	1.0544	0.0014	< 0.0053 U	--	--	0.0361	0.9890	0.0013	0.0382	1.4580	0.0019	0.0122	0.4108	0.0005	0.0389	1.2838	0.0017	0.0212	1.5821	0.0021	0.0078	0.8066	0.0010
C1-Fluorenes	611	12.22	0.0170	0.3214	0.0005	0.0055 J	0.1662	0.0003	< 0.0053 U	--	--	0.0033 J	0.0904	0.0001	< 0.0057 U	--	--	< 0.0052 U	--	--	< 0.0046 U	--	--	< 0.0043 U	--	--	--	--	
C1-Naphthalenes	445	8.9	0.0458	0.8658	0.0019	0.0101	0.3051	0.0007	0.0048 J	0.2254	0.0005	< 0.0061 J	--	--	< 0.0057 J	--	--	< 0.0052 U	--	--	< 0.0046 U	--	--	< 0.0043 J	--	--	--	--	
C1-Phenanthrenes/Anthracenes	670	13.4	0.0852	1.6106	0.0024	0.0320	0.9668	0.0014	0.0038 J	0.1784	0.0003	0.0238	0.6521	0.0010	0.0265	1.0115	0.0015	0.0089	0.2997	0.0004	0.0233	0.7690	0.0011	0.0134	1.0000	0.0015	0.0059	0.6101	0.0009
C2-Benzenanthracenes/Chrysenes	1009	20.18	0.0387	0.7316	0.0007	0.0187	0.5650	0.0006	< 0.0053 U	--	--	0.0201	0.5507	0.0005	0.0219	0.8359	0.0008	0.0070 J	0.2357	0.0002	0.0223	0.7360	0.0007	0.0123 J	0.9179	0.0009	0.0046 U	0.4757	0.0005
C2-Fluorenes	687	13.74	0.0175	0.3308	0.0005	0.0091	0.2749	0.0004	< 0.0053 U	--	--	0.0056 J	0.1534	0.0002	0.0046 J	0.1756	0.0003	< 0.0052 U	--	--	< 0.0047 J	0.1551	0.0002	0.0030 J	0.2239	0.0003	0.0034 J	0.3516	0.0005
C2-Naphthalenes	510	10.2	0.0473	0.8941	0.0018	0.0146	0.4411	0.0009	0.0064	0.3005	0.0006	0.0061	0.1671	0.0003	0.0089	0.3397	0.0007	0.0038 J	0.1279	0.0003	0.0047 J	0.1551	0.0003	0.0036 J	0.2687	0.0005	0.0029 J	0.2999	0.0006
C2-Phenanthrenes/Anthracenes	745	14.9	0.0438	0.8280	0.0011	0.0209	0.6314	0.0008	< 0.0053 U	--	--	0.0117	0.3205	0.0004	0.0139	0.5305	0.0007	0.0052	0.1751	0.0002	0.0115	0.3795	0.0005	0.0060	0.4478	0.0006	0.0032 J	0.3309	0.0004
C3-Benzenanthracenes/Chrysenes	1213	24.26	0.0110	0.2079	0.0002	< 0.0060 U	--	--	< 0.0053 U	--	--	< 0.0061 U	--	--	< 0.0057 U	--	--	< 0.0052 U	--	--	< 0.0046 U	--</td							

Table 5  
PAH Screening Table with Calculated Toxicity Units  
November 2013 Event  
UTC/Carrier Circle  
Carrier-Dewitt Landfill (AOC G)  
Thompson Road, Syracuse, New York

Transect	Sample ID	NYSDEC Screening Criteria	Sanders Creek																										
			SDS8			SDS8 (Dup)			SDS8			SDS8 (Dup)			SDS8			SDS8 (Dup)			SDS9								
			11/5/2013			11/5/2013			11/5/2013			11/5/2013			11/5/2013			11/5/2013			11/5/2013								
Chemical Name	Sediment Guidance Value (µg/gOC)	NYSDEC Sediment CLASS A	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU	mg/kg	µg/gOC	TU						
ACENAPHTHENE	491	9.82	0.0123	0.3361	0.0007	0.0106	0.8480	0.0017	<0.0057 U	--	--	<0.0056 U	--	--	<0.0047 U	--	--	<0.0045 U	--	--	0.0478	2.7006	0.0055	0.0688	2.0476	0.0042	<0.0047 J	--	--
ACENAPHTHYLENE	452	9.04	0.0240	0.6557	0.0015	0.0248	1.9840	0.0044	0.0140	0.4698	0.0010	0.0136	0.3736	0.0008	0.0056	0.4590	0.0010	0.0203	0.8750	0.0019	0.0941	5.3164	0.0118	0.106	3.1548	0.0070	0.0082	1.2893	0.0029
ANTHRACENE	594	11.88	0.0435	1.1885	0.0020	0.0382	3.0560	0.0051	0.0190	0.6376	0.0011	0.0141	0.3874	0.0007	0.0051	0.4180	0.0007	0.0198	0.8534	0.0014	0.183	10.3390	0.0174	0.233	6.9345	0.0117	0.0079	1.2421	0.0021
BENZO(A)ANTHRACENE	841	16.82	0.132	3.6066	0.0043	0.12	9.6000	0.0114	0.0605	2.0302	0.0024	0.0347	0.9533	0.0011	0.0111	0.9098	0.0011	0.0353	1.5216	0.0018	0.542	30.6215	0.0364	0.738	21.9643	0.0261	0.0183	2.8774	0.0034
BENZO(A)PYRENE	964	19.28	0.15	4.0984	0.0043	0.139	11.1200	0.0115	0.0753	2.5268	0.0026	0.0420	1.1538	0.0012	0.0133	1.0902	0.0011	0.0457	1.9698	0.0020	0.658	37.1751	0.0386	0.89	26.4881	0.0275	0.0224	3.5220	0.0037
BENZO(B)FLUORANTHENE	979	19.58	0.128	3.4973	0.0036	0.119	9.5200	0.0097	0.0693	2.3255	0.0024	0.0361	0.9918	0.0010	0.0112	0.9180	0.0009	0.0388	1.6724	0.0017	0.564	31.8644	0.0325	0.776	23.0952	0.0236	0.0189	2.9717	0.0030
BENZO(G,H,I)PERYLENE	1095	21.9	0.0945	2.5820	0.0024	0.0872	6.9760	0.0061	0.0515	1.7282	0.0016	0.0271	0.7445	0.0007	0.0884	0.6885	0.0006	0.0328	1.4138	0.0013	0.438	24.7458	0.0226	0.58	17.2619	0.0158	0.0141	2.2170	0.0020
BENZO(K)FLUORANTHENE	980	19.6	0.134	3.6612	0.0037	0.124	9.9200	0.0101	0.0706	2.3691	0.0024	0.0377	1.0357	0.0011	0.0122	1.0000	0.0010	0.0408	1.7586	0.0018	0.61	34.4633	0.0352	0.8	23.8095	0.0243	0.0199	3.1289	0.0032
BENZO(E)PYRENE	967	19.34	0.105	2.8689	0.0030	0.0964	7.7120	0.0080	0.0572	1.9195	0.0020	0.0299	0.8214	0.0008	0.0094	0.7705	0.0008	0.0355	1.5302	0.0016	0.463	26.1582	0.0271	0.617	18.3631	0.0190	0.0153	2.4057	0.0025
C1-Benzenanthracenes/Chrysenes	930	18.6	0.0569	1.5546	0.0017	0.0534	4.2720	0.0046	0.0291	0.9765	0.0011	0.0195	0.5357	0.0006	0.0073	0.5984	0.0006	0.0255	1.0991	0.0012	0.247	13.9548	0.0150	0.325	9.6726	0.0104	0.109	1.7138	0.0018
C1-Fluoranthenes/Pyrenes	769	15.38	0.103	2.8142	0.0037	0.0972	7.7760	0.0101	0.0510	1.7114	0.0022	0.0338	0.9286	0.0012	0.0132	1.0820	0.0014	0.0432	1.8621	0.0024	0.397	22.4294	0.0292	0.519	15.4464	0.0201	0.0193	3.0346	0.0039
C1-Fluorenes	611	12.22	0.0054 J	0.1475	0.0002	0.0053	0.4240	0.0007	0.0031 J	0.1040	0.0002	0.0029 J	0.0797	0.0001	<0.0047 U	--	--	0.0045	0.1940	0.0003	0.0196	0.0255	0.7589	0.0012	0.0025 J	0.3931	0.0006		
C1-Naphthalenes	445	8.9	0.0664	0.1749	0.0004	0.0059	0.4720	0.0011	<0.0057 U	--	--	<0.0056 U	--	--	<0.0047 U	--	--	0.0047	0.2026	0.0005	0.0223	1.2599	0.0028	0.0243	0.7232	0.0016	<0.0047 J	--	--
C1-Phenanthrenes/Anthracenes	670	13.4	0.0565	1.5437	0.0023	0.0537	4.2960	0.0064	0.0292	0.9799	0.0015	0.0217	0.5962	0.0009	0.0095	0.7787	0.0012	0.0317	1.3664	0.0020	0.232	13.1073	0.0196	0.299	8.8988	0.0133	0.0130	2.0440	0.0031
C2-Benzenanthracenes/Chrysenes	1009	20.18	0.0569	1.5546	0.0015	0.0534	4.2720	0.0042	0.0291	0.9765	0.0010	0.0195	0.5357	0.0005	0.0073	0.5984	0.0006	0.0255	1.0991	0.0011	0.247	13.9548	0.0138	0.325	9.6726	0.0096	0.0109	1.7138	0.0017
C2-Fluorenes	687	13.74	0.0889	0.2432	0.0004	0.0079	0.6320	0.0009	0.0047 J	0.1577	0.0002	0.0044 J	0.1209	0.0002	0.0028 J	0.2295	0.0003	0.0065	0.2802	0.0004	0.0278	1.5706	0.0023	0.0347	1.0327	0.0015	0.0032 J	0.5031	0.0007
C2-Naphthalenes	510	10.2	0.0109	0.2978	0.0006	0.0087	0.6960	0.0014	0.0086	0.2886	0.0006	0.0057	0.1566	0.0003	0.0038 J	0.3115	0.0006	0.0078	0.3362	0.0007	0.0300	1.6949	0.0033	0.0357	1.0625	0.0021	0.0044 J	0.6918	0.0014
C2-Phenanthrenes/Anthracenes	745	14.9	0.0273	0.7459	0.0010	0.0247	1.9760	0.0027	0.0145	0.4866	0.0007	0.0123	0.3379	0.0005	0.0058	0.4754	0.0006	0.0182	0.7845	0.0011	0.115	6.4972	0.0087	0.151	4.4940	0.0060	0.0071	1.1164	0.0015
C3-Benzenanthracenes/Chrysenes	1213	24.26	0.0126	0.3443	0.0003	0.0133	1.0640	0.0009	0.0073	0.2450	0.0002	<0.0056 U	--	--	<0.0047 U	--	--	0.0096	0.4138	0.0003	0.0578	3.2655	0.0027	0.0769	2.2887	0.0019	<0.0047 U	--</	

## **Appendix A**

### **Use of Equilibrium Partitioning Theory to Develop Sediment Guidance Values for Nonpolar Organic Compounds**

**Appendix A**  
**Use of Equilibrium Partitioning Theory to Develop Sediment Guidance Values  
for Nonpolar Organic Compounds**

The equilibrium partitioning methodology is well-documented in the scientific literature (US EPA, 1991; US EPA SAB, 1992; DiToro et al., 1991). It is based on the principle that nonpolar organic contaminants in sediment will partition between sediment pore water and the organic carbon content of sediment in a constant ratio. That ratio is referred to as the organic carbon partitioning coefficient, or Koc.

The Kow, or *n*-octanol water partitioning coefficient, is a measure of the ratio that describes the partitioning of a nonpolar organic compound between water and octanol. The Kow is a useful surrogate of how nonpolar organic compounds will accumulate in lipid in animal tissue (US EPA 1995). Kows are generally readily available for many common organic compounds, and tend to be similar in value to and vary proportionately with a compound's Koc (Kenaga, 1980; Voice, 1983). US EPA (1991) refers to DiToro (1985) to define the relationship between Kow and Koc as:

$$\text{Log10Koc} = 0.00028 + 0.983 \cdot \text{log10Kow} \quad (1)$$

US EPA (2002) reported that the concentration-response curve for biological effects can be correlated with the concentration of the contaminant in the interstitial pore water, *not* with the concentration of the contaminant in sediment; and that the effects concentration for a chemical in pore water is essentially equal to that reported for water-only exposures. In other words, the toxicity of a nonpolar organic contaminant to sediment-dwelling organisms *is proportional to* the concentration of the contaminant that is freely dissolved in the sediment pore water. For nonpolar organic chemicals, the concentration-response curves correlate equally well with the sediment chemical concentration that is normalized on a sediment organic carbon basis. The higher the Kow of a nonpolar organic compound, the stronger the contaminant will adsorb to the organic carbon content in the sediment. When more organic carbon is present in sediment, the concentration of a nonpolar organic contaminant freely dissolved in sediment pore water will be smaller, and therefore, proportionally less toxic to aquatic organisms.

An equilibrium partitioning-based SGV is derived by multiplying the AWQS/GV from 6 NYCRR Part 703.5 or TOGS 1.1.1 for a nonpolar organic compound by its Koc, as derived from equation 1:

$$\text{SGVoc} = \text{AWQS/GV } \mu\text{g/L} * \text{Koc} \quad (2)$$

This results in a SGV in units of microgram ( $\mu\text{g}$ ) of contaminant per gram of organic carbon in the sediment (SGVoc).

UTC-Carrier used this methodology to develop SGVs for organic chemicals detected at the AOC G landfill wetlands that lacked NYSDEC screening values. For chemicals lacking a NYCRR Part 703.5 or TOGS 1.1.1 AWQS for protection of aquatic life, US EPA Region 3 BTAG Kows and fresh water chronic water quality values were used.

**Table A-1: Calculation of NYSDEC Class A Sediment Guidance Values for Nonpolar Organic Compounds**

Chemical	Log Kow	Kow Reference	Koc	Chronic Freshwater Guideline (µg/L)	FW Guideline Reference	Class A FW Sed Criteria (µg/gOC)	NYSDEC TOC %	NYSDEC EqP FW Sed (mg/kg)
1,1,1-Trichloroethane	2.48	EPA, 2006.	274.2331802	11	EPA, 2014a. EPA Region III	3.02E+00	2	6.03E-02
1,1-Dichloroethane	1.79	EPA, 2006.	57.52412217	47	NOAA Squirt Table, 2008	2.70E+00	2	5.41E-02
1,2,3-Trichlorobenzene	4.096	NYS DEC	10632.80869	1.1	NYS DEC, 1998 TOGS 1.1.1	1.17E+01	2	2.34E-01
1,2,4-Trichlorobenzene	3.99	NYS DEC	8364.692881	210	NYS DEC, 1998 TOGS 1.1.1	1.76E+03	2	3.51E+01
1,2-Dichloroethene (cis)	1.86	EPA, 2006.	67.40001606	590	NYS DEC, 1998 TOGS 1.1.1	3.98E+01	2	7.95E-01
1,3,5-Trimethylbenzene	3.78	EPA, 2006.	5200.199438	8	NOAA Squirt Table, 2008	4.16E+01	2	8.32E-01
2-Methylnaphthalene	3.86	ATSDR, 2005.	6232.467177	4.7	EPA, 2014a. EPA Region III	2.93E+01	2	5.86E-01
Aroclor 1254	6.72	EPA, 2006.	4036825.718	0.014	NOAA Squirt Table, 2008	5.65E+01	2	1.13E+00
Aroclor 1260	6.91	EPA, 2006.	6205974.689	0.014	NOAA Squirt Table, 2008	8.69E+01	2	1.74E+00
bis(2-ethylhexyl)phthalate	7.6	NYS DEC	29585574.03	0.6	NYS DEC, 1998 TOGS 1.1.1	1.78E+04	2	3.55E+02
Chloroform	1.92	EPA, 2006.	77.20403519	1.8	NOAA Squirt Table, 2008	1.39E-01	2	2.78E-03
Dibenzofuran	4.12	EPA, 2006.	11226.38676	3.7	EPA, 2014a. EPA Region III	4.15E+01	2	8.31E-01
Diethylphthalate	2.5	EPA, 2006.	286.9326703	210	EPA, 2014a. EPA Region III	6.03E+01	2	1.21E+00
Di-n-butylphthalate	4.61	EPA, 2006.	34033.76533	19	EPA, 2014a. EPA Region III	6.47E+02	2	1.29E+01
m,p-Xylenes	3.2	NYS DEC	1399.200656	17	NYS DEC, 1998 TOGS 1.1.1	2.38E+01	2	4.76E-01
PCBs (NYSDEC)	6.14	EPA, 2006.	1086175.494	0.014	NOAA Squirt Table, 2008	1.52E+01	2	3.04E-01
Styrene	2.94	EPA, 2006.	776.7835144	72	EPA, 2014a. EPA Region III	5.59E+01	2	1.12E+00
Trichloroethene (TCE)	2.53	NYS DEC	307.0930586	290	NYS DEC, 1998 TOGS 1.1.1	8.91E+01	2	1.78E+00
Vinyl chloride	1.5	EPA, 2006.	29.838707	930	EPA, 2014a. EPA Region III	2.77E+01	2	5.55E-01
Xylene (total)	3.2	NYS DEC	1399.200656	17	NYS DEC, 1998 TOGS 1.1.1	2.38E+01	2	4.76E-01

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## **Appendix B**

### **Data Usability Summary Reports**

Due to the size of the Data Usability Summary Reports, they have been provided on the enclosed CD.