



**Environmental Resources Group**

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## PHASE II ENVIRONMENTAL SITE ASSESSMENT

GREEN BAXTER COURT

1701-1747 GREEN ROAD

ANN ARBOR, MICHIGAN 48105

ERG PROJECT 1127.001

PREPARED FOR:

NORSTAR DEVELOPMENT USA, L.P.  
733 BROADWAY  
ALBANY, NEW YORK 12207

AUGUST 14, 2013



**Environmental Resources Group**

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August 14, 2013

Mr. Richard Higgins  
Norstar Development USA, L.P.  
733 Broadway  
Albany, New York 12207

**Re: Phase II Environmental Site Assessment  
Green Baxter Court  
1701-1747 Green Road, Ann Arbor, Michigan  
Environmental Resources Group, LLC Project 1127.001**

Dear Mr. Higgins:

Environmental Resources Group, LLC (ERG) has completed the Phase II Environmental Site Assessment (ESA) of the Green Baxter Court Housing Development in Ann Arbor, Washtenaw County, Michigan. The results of the Phase II ESA are presented in the attached Report.

The Report was prepared for the exclusive use of NORSTAR DEVELOPMENT USA, L.P., ANN ARBOR HOUSING COMMISSION, and the MICHIGAN STATE HOUSING DEVELOPMENT AUTHORITY, each of whom may rely on the Report's contents.

We are pleased to provide this service and hope that we can be of service in the future. Should you have any questions or require further information, please do not hesitate to call Mr. Foerg at (248) 773-7986.

Sincerely,  
**Environmental Resources Group, LLC**

A handwritten signature in black ink that reads "Andrew J. Foerg". The signature is fluid and cursive, with "Andrew" and "J." being more stylized and "Foerg" being more clearly legible.

Andrew J. Foerg, CPG  
Senior Project Manager

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## **1.0 INTRODUCTION**

Norstar Development USA, L.P. (Norstar), retained Environmental Resources Group, LLC (ERG) to conduct a Phase II Environmental Site Assessment (Phase II ESA) of a property located at 1701-1747 Green Road, Ann Arbor, Michigan (subject property). This Phase II ESA was conducted in accordance with Michigan State Housing Development Authority (MSHDA) Environmental Review Requirements for 2013 and American Society for Testing and Materials (ASTM) Designation E 1903 "Standard Guide for Environmental Site Assessments: Phase II Environmental Site Assessment Process."

This Phase II ESA scope of work is intended to evaluate the recognized environmental conditions (RECs) identified in ERG's July 2013 Phase I ESA, which are presented in Section 2.5.

The Report was prepared for the exclusive use of NORSTAR DEVELOPMENT USA, L.P., ANN ARBOR HOUSING COMMISSION, and the MICHIGAN STATE HOUSING DEVELOPMENT AUTHORITY, each of whom may rely on the Report's contents.

## **2.0 BACKGROUND**

### **2.1 SITE DESCRIPTION AND PHYSICAL SETTING**

The subject property contains approximately 2.4 acres and is located west of Green Road and north of Baxter Road in Ann Arbor, Michigan. The subject property is located in Section 23, Township 2 South, Range 6 East, Ann Arbor, Washtenaw County, Michigan.

The Subject Property is developed with four 2-story multi-tenant apartment buildings with associated parking area, play area and landscaping. The adjoining properties are developed with commercial and residential properties.

Refer to Figure 1 for a Site Location Map. See Figure 2 for a Site Map with soil boring locations.

### **2.2 SUBJECT PROPERTY HISTORY AND LAND USE**

Reasonably ascertainable records for the subject property extended back to approximately 1903. Standard historical sources were able to document the first developed use of the subject property occurred prior to 1903. The property was used for residential purposes from at least 1903 to 1937. A farmstead occupied the Subject Property from at least 1937 to the 1970s. The current multi-tenant apartment building was constructed in 1970.

## **2.3 ADJACENT PROPERTY LAND USE**

As determined during ERG's July 2013 Phase I ESA, the current uses of the adjoining properties are as follows:

- North: University of Michigan North Campus Auxiliary Services Complex – 1919 Green Road
- South: Baxter Street with Michigan Municipal League office building – 1675 Green Road
- East: Green Road with apartments beyond
- West: Baxter Park (City of Ann Arbor public park)

## **2.4 PREVIOUS ENVIRONMENTAL INVESTIGATIONS**

No previous environmental investigations were provided to ERG for review.

## **2.5 IDENTIFIED RECOGNIZED ENVIRONMENTAL CONDITIONS (RECS)**

ERG's July 2013 Phase I ESA identified the following RECs:

### **On-Site RECs**

1. A farmstead occupied the Subject Property from at least 1937 to the 1970s. It is likely that pesticides, maintenance materials, and fuels were used and stored at the Subject Property. The probable long-term use and storage of chemicals on the Subject Property is considered an REC.

No other RECs were identified on the Subject Property.

### **Off-Site RECs**

2. The north adjoining property (1919 Green Road) was listed in the environmental databases with two former generators of waste (Bendix Corp and University of Michigan). The records indicated that these operations are not currently generating waste. Although the waste generating activities at those sites are closed, they were located directly up gradient on the north adjoining property. Based on the known historical chemical use and storage in close proximity to the Subject Property, the Bendix Corp. and University of Michigan property is considered an REC.

No other RECs were identified on the adjoining or nearby properties.

### 3.0 PHASE II ENVIRONMENTAL SITE ASSESSMENT ACTIVITIES

#### 3.1 SCOPE OF ASSESSMENT

To evaluate the RECs, ERG conducted a subsurface investigation of the subject property that included: (1) conducting a geophysical survey, (2) advancing seven soil borings, (3) installing three temporary groundwater monitor wells, (4) installing four soil gas collection points, and (5) collecting five soil samples, three groundwater samples and four soil gas samples. The samples were submitted for laboratory analyses of one or more of the following: VOCs, PNAs, PCBs, organochlorine pesticides (pesticides), arsenic and lead.

The following table summarizes each REC, the site investigation activities performed to address each REC, and the laboratory parameters used to address each REC.

**Summary of ERG's Scope of Investigation**

REC #	Environmental Concern	Investigation Activity	Analytical Parameters
REC 1	historical farmstead on site	geophysical survey and SB-1/TW-1, SB-2, SB-3, SB-6 and SB-7	VOCs, VOCs (soil gas) PNAs, pesticides, arsenic and/or lead
REC 2	historical industrial use of north adjoining property	SB-4/TW-2 and SB-5/TW-3	VOCs, VOCs (soil gas) PNAs, PCBs,

##### 3.1.1 Geophysical Survey

ERG retained Fibertec Environmental Services (Fibertec) to conduct a geophysical survey of the central portion of the subject property, where the former farm related operations were located. The purpose of the geophysical survey was to evaluate whether abandoned USTs are present at the subject property. On July 17, 2013, Fibertec conducted a ground penetrating radar (GPR) survey of the central portion of the subject property. Figure 2 depicts the approximate boundary of the geophysical survey area.

The GPR survey was performed using a GSSI Sir 3000 GPR unit with a 400MHz antenna. No anomalies appearing to be a UST were observed.

Refer to Appendix A for a copy of the geophysical survey report.

##### 3.1.2 Soil Evaluation

On July 18, 2013, Fibertec advanced seven soil borings at the subject property. Fibertec used hydraulic drive/direct-push (Geoprobe<sup>®</sup>) and hand auger sampling techniques and followed the drilling procedures outlined in ASTM publication D 6282-98 "Standard Guide for Direct Push Soil Sampling for Environmental Site Characterizations".

Continuous soil samples were collected from the soil borings in five-foot intervals to the maximum depth explored of 20 feet bgs. ERG personnel inspected, field-screened, and logged the samples collected at each soil boring location. Refer to Figure 2 for a site map with soil boring locations. Soil boring logs are provided in Appendix B.

### **3.1.3 Groundwater Evaluation**

Groundwater was encountered groundwater in three of the seven soil borings (SB-1/TW-1, SB-4/TW-2 and SB-5/TW-3) advanced at the subject property. A temporary groundwater monitoring well was installed at each location where groundwater was encountered. One-inch PVC risers with five-foot screens were utilized for the temporary groundwater monitoring wells. Groundwater sample collection methodology was in general accordance with MDEQ low-flow sampling methodologies in that the wells were allowed to stabilize prior to collecting the groundwater samples at a low flow rate. The volume of groundwater in SB-1/TW-1 was only sufficient to collect enough sample to analyze for VOCs. Refer to Figure 2 for a site map with soil boring/temporary well locations.

### **3.1.4 Soil Gas Evaluation**

The soil gas evaluation was conducted to assess potential soil gas associated with the former farmstead operations and potential vapor migration from the north adjoining property.

The soil gas evaluation activities were conducted in general accordance with the guidelines established by the American Society for Testing and Materials (ASTM) in the Standard Practice for Vapor Encroachment Screening on Property Involved in Real Estate Transactions Designation E 2600-10 (ASTM Standard Practice E 2600-10) and MDEQ Remediation and Redevelopment Division Guidance Document for the Vapor Intrusion Pathway, May 2013. This included performing a leak check using a helium chamber and purging three (3) gas point/sand pack volumes at low-flow (200 ml/minute) from each soil gas monitoring point, followed by the collection of one (1) soil gas sample, using vacuum bottle methods, for laboratory analysis of VOCs. The vacuum bottles were regulated with a flow rate of 200 ml/minute, which was pre-set at the laboratory.

Four soil gas samples (SB-4, through SB-7) were collected and submitted for laboratory analysis of VOCs. Refer to Figure 2 for the locations of the sub-surface soil gas points.

## **3.2 QUALITY ASSURANCE/QUALITY CONTROL**

To ensure the accuracy of data collected during on site activities, ERG employed proper quality assurance/quality control (QA/QC) measures. The QA/QC procedures included, but were not limited to, (1) decontamination of sampling equipment before and between sampling events, (2) calibration of field equipment, (3) documentation of field activities, (4) sample preservation techniques, and (5) collection of QA/QC samples.

### **3.2.1 Decontamination of Equipment**

During sample collection, ERG adhered to proper decontamination procedures. Sampling equipment was decontaminated using the following methods to minimize potential cross-contamination of soil samples:

- Steam-cleaning or washing and scrubbing the equipment with non-phosphate detergent
- Rinsing the equipment
- Air-drying the equipment

### **3.2.2 Calibration of Field Equipment**

All field instruments were calibrated prior to first use on-site to ensure accuracy. Field instruments utilized during investigation activities at this subject property included a photoionization detector (PID). During ERG's Phase II ESA, the PID was used to screen all soil samples. The PID was maintained in a calibrated condition using 100 parts per million (ppm) isobutylene span gas prior to the subsurface investigation.

### **3.2.3 Documentation of Activities**

During ERG's Phase II ESA activities, subject property conditions (i.e., soil boring locations, weather conditions) were documented. ERG visually inspected the soil and groundwater samples and prepared a geologic log for each soil boring. The logs include soil characteristics such as (1) color, (2) composition (e.g., sand, clay, or gravel), (3) soil moisture and water table depth, and (4) signs of possible contamination (i.e., stained or discolored soil, odors). All soil samples were delivered to a laboratory under chain-of-custody documentation. See Appendix B for ERG's soil boring logs. See Figure 2 for site map with soil boring locations.

### **3.2.4 Sample Preservation Techniques**

ERG collected soil samples according to USEPA Publication SW-846, "*Test Methods for Evaluating Solid Waste*." Soil samples were collected in laboratory-supplied containers, stored on ice or at approximately four degrees Celsius, and submitted under chain-of-custody documentation. Soil samples collected for VOC analyses were field preserved with methanol in accordance with U.S. EPA Method 5035. Groundwater samples for VOC analysis were preserved with HCL in accordance with U.S. EPA Method 5030B/8260B.

### **3.2.5 QA/QC Sample Collection**

ERG collected QA/QC samples for the groundwater water matrix in accordance with QA/QC sample procedures outlined in DEQ's Sampling Strategies and Statistics Training Materials for Part 201 Cleanup Criteria. ERG's QA/QC samples included a trip blank.

### 3.3 LABORATORY ANALYSES AND METHODS

ERG submitted five soil samples, three groundwater samples and four soil gas samples for laboratory analyses. The following table summarizes the location, depth, matrix, and laboratory analysis for each sample.

**Summary of Laboratory Analyses**

Sample Name/ Depth	Matrix	VOCs	PNAs	PCBs	Arsenic and Lead	Pesticides	VOCs (soil gas)
SB-1 (1.5-2')	Soil	x	x		x	x	
TW-1 (13.5')	GW	x					
SB-2 (8-9')	Soil	x	x		x	x	
SB-3 (9-16')	Soil	x	x		x	x	
SB-4 (7')	Soil Gas						x
TW-2 (12')	GW	x	x	x			
SB-5 (7')	Soil Gas						x
TW-3 (10')	GW	x	x	x			
SB-6 (1-2')	Soil	x	x		x	x	
SB-6 (10')	Soil Gas						x
SB-7 (12-13')	Soil	x	x		x	x	
SB-7 (10')	Soil Gas						x

The laboratory analyzed the soil and/or groundwater samples for: (1) VOCs in accordance with USEPA Method 5035/8260B; (2) PNAs in accordance with USEPA Method 3546/8270C; (3) Arsenic and Lead in accordance with USEPA Method 0200.2-M/6020A; (4) PCBs in accordance with USEPA Method 3535A/8082A; and (5) Pesticides in accordance with USEPA Method 3546/8081B. Soil gas samples were analyzed for VOCs by method TO-15.

## 4.0 EVALUATION AND PRESENTATION OF RESULTS

### 4.1 SUBSURFACE CONDITIONS

#### 4.1.1 Soil and Groundwater Conditions Based on Published Material

The United States Geological Survey Division (U.S.G.S.) 7.5-Minute Topographic Map of the Ann Arbor East, Michigan Quadrangle, 1965 (photo revised in 1983) for the subject property was reviewed. Based on the topographic map, the subject property is located at an elevation of 900 feet above mean sea level. The immediate subject property area appears to slope to the west southwest.

The "Quaternary Geology of Southern Michigan", Department of Geological Sciences, University of Michigan, Ann Arbor, Michigan (1982); and "Bedrock Geology of Southern Michigan", Michigan Department of Natural Resources, Geological Survey Division (1987) were reviewed.

According to the aforementioned sources, in this area of Washtenaw County, quaternary deposits are underlain by bedrock comprised of Coldwater Shale. Bedrock is covered by end moraines of medium textured till. The surface sediments in the area of the site generally consist of end moraines of medium-textured till, gray to reddish brown, non-sorted glacial debris. The matrix is dominantly loam and silty loam texture with variable amounts of cobbles and boulders. It occurs in narrow linear belts of hummocky relief marking former stillstands of ice-sheet margins and includes small areas of ground moraine as well as outwash. The drift thickness is approximately 20 to 30 meters

ERG's research did not identify known groundwater supply or monitoring wells on the subject property.

#### 4.1.2 Soil and Groundwater Conditions Based on Field Observations

During drilling activities, ERG encountered the following soil types:

- FILL from ground surface, or just below the topsoil or concrete, to a depth of two feet bgs in SB-1. In general, the fill consisted of sand, silt, clay and gravel that was moist. Fill was not encountered in SB-2 through SB-7.
- SAND from ground surface to a depth of 3 feet bgs in SB-3 and SB-6.
- SILTS/CLAYS from depths ranging from ground surface, to a depth of 20 bgs, the maximum depth explored.
- Saturated sand layers were encountered from 10-14 feet bgs in SB-1, 10-12.5 feet bgs in SB-4 and 8.5-10 feet bgs in SB-5.

Positive PID readings were recorded at two soil intervals:

- SB-1 at 1.5 feet bgs (185.5 ppm)
- SB-6 at 1 foot bgs (26.4 ppm)

No visual or olfactory evidence of impacts were recorded at any of the boring locations.

See Figure 2 for a site map with soil boring locations. See Appendix B for soil boring logs.

## 4.2 MDEQ RELEVANT EXPOSURE PATHWAYS AND APPLICABLE CRITERIA

### 4.2.1 Relevant Exposure Pathways

As defined in Michigan Public Act 451 Part 201, "relevant pathway" means an exposure pathway that is reasonable and relevant because there is a reasonable potential for exposure to a hazardous substance. The analysis of potential exposure pathways is based on known existing conditions at the subject property. The following subsections identify the relevant exposure pathways based on the subject property conditions observed.

#### Ingestion of Groundwater Pathway

Groundwater was encountered at a depth of approximately 10 feet bgs in three soil borings drilled at the subject property. In the absence of additional deeper soil borings to confirm a confining clay layer, Ingestion of Groundwater is considered a relevant exposure pathway.

#### Groundwater Venting to Surface Water Pathway

Groundwater Venting to Surface Water is not a human exposure pathway, but rather an exposure pathway based on aquatic toxicity. The subject property is located in reasonably close proximity to surface water bodies. Therefore, Groundwater Venting to Surface Water is a relevant exposure pathway.

#### Groundwater Contact Pathway

Groundwater Contact is a relevant exposure pathway.

#### Volatilization to Indoor Air Inhalation Pathway

Volatilization to Indoor Air Inhalation is a relevant exposure pathway.

#### Volatilization to Ambient Air Pathway

Volatilization to Ambient Air is a relevant exposure pathway.

#### Particulate Inhalation Pathway

Particulate Inhalation is a relevant exposure pathway.

#### Direct Contact Pathway

Direct Contact is a relevant exposure pathway.

### 4.2.2 Applicable Criteria

Applicable criterion means a cleanup criterion for a relevant pathway. A criterion is not applicable if the exposure pathway is not relevant. Based on the exposure pathway evaluation, the applicable pathways at the subject property include:

- Drinking Water (DW)/Drinking Water Protection (DWP)
- Groundwater Surfacewater Interface (GSI)/ Groundwater Surfacewater Interface Protection (GSIP)
- Soil Volatilization to Indoor Air Inhalation (SVIAI)/Groundwater Volatilization to Indoor Air Inhalation (GVIAI);
- Infinite Source Volatile Soil Inhalation (VSIC);
- Groundwater Contact Protection (GCP);
- Particulate Soil Inhalation (PSI); and
- Soil Direct Contact (DC)/Groundwater Contact.

#### **4.3 LABORATORY ANALYTICAL RESULTS**

ERG collected soil and groundwater samples for the purpose of determining if the subject property meets the definition of a *facility*. Analytical results were compared with MDEQ GRCC provided in MDEQ RRD's Operational Memorandum No. 1, Tables 1 and 2.

ERG collected soil gas samples for the purpose of evaluating potential vapor sources on and off-site. Soil gas results were compared to MDEQ Vapor Intrusion Deep Soil Gas Screening Levels provided in MDEQ Remediation and Redevelopment Division Guidance Document for the Vapor Intrusion Pathway, May 2013.

##### **4.3.1 Soil Analytical Results**

ERG submitted five soil samples for laboratory analysis of one or more of the following: VOCs, PNAs, PCB, Arsenic, Lead and/or pesticides. Reported concentrations that exceeded relevant and applicable GRCC include:

- One sample SB-1 (1.5-2') had reported concentrations of three PNAs and Arsenic that exceeded one or more GRCC (DC, DWP, GSIP).
- One sample SB-3 (9-10') had a reported concentration of Arsenic that exceeded two GRCC (DWP, GSIP).

No other exceedences of soil GRCC were reported.

Refer to Table 1 for a summary of soil analytical results and a comparison to GRCC. Refer to Appendix C for the analytical laboratory report.

##### **4.3.2 Groundwater Analytical Results**

ERG submitted three groundwater samples for laboratory analysis of one or more of the following: VOCs, PNAs and PCBs. No reported groundwater concentrations exceeded any relevant and applicable GRCC.

Refer to Table 2 for a summary of groundwater analytical results and a comparison to GRCC. Refer to Appendix C for the analytical laboratory report.

#### **4.3.3 Soil Gas Analytical Results**

ERG submitted four soil gas samples for laboratory analysis of VOCs. No reported concentrations of VOCs in soil gas exceeded the applicable MDEQ Vapor Intrusion Deep Soil Gas Screening Levels.

Refer to Table 3 for a summary of soil gas analytical results and a comparison to MDEQ Vapor Intrusion Deep Soil Gas Screening Levels. Refer to Appendix C for the analytical laboratory report.

### **5.0 SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS**

#### **5.1 SUMMARY OF ENVIRONMENTAL CONCERNS**

ERG's July 2013 Phase I ESA identified the following RECs:

##### **On-Site RECs**

1. A farmstead occupied the Subject Property from at least 1937 to the 1970s. It is likely that pesticides, maintenance materials, and fuels were used and stored at the Subject Property. The probable long-term use and storage of chemicals on the Subject Property is considered an REC.

##### **Off-Site RECs**

2. The north adjoining property (1919 Green Road) was listed in the environmental databases with two former generators of waste (Bendix Corp and University of Michigan). The records indicated that these operations are not currently generating waste. Although the waste generating activities at those sites are closed, they were located directly up gradient on the north adjoining property. Based on the known historical chemical use and storage in close proximity to the Subject Property, the Bendix Corp. and University of Michigan property is considered an REC.

#### **5.2 SUMMARY OF SUBSURFACE INVESTIGATION**

To evaluate the RECs, ERG conducted a subsurface investigation of the subject property that included: (1) conducting a geophysical survey, (2) advancing seven soil borings, (3) installing three temporary groundwater monitor wells, (4) installing four soil gas collection points, and (5) collecting five soil samples, three groundwater samples and four soil gas samples. The samples were submitted for laboratory analyses of one or more of the following: VOCs, PNAs, PCBs, organochlorine pesticides (pesticides), arsenic and lead.

### 5.3 CONCLUSIONS

ERG conducted a geophysical survey, and soil/groundwater/soil gas sampling in areas most likely to be impacted by contaminants based on the past use of the subject property. The results of the investigation indicate the following:

- No anomalies appearing to be a UST were observed.
- No reported groundwater concentrations exceeded the relevant and applicable GRCC.
- No reported concentrations of VOCs in soil gas exceeded the applicable MDEQ Vapor Intrusion Deep Soil Gas Screening Levels.
- One sample SB-1 (1.5-2') had reported concentrations of three PNAs and Arsenic that exceeded one or more GRCC (DC, DWP, GSIP).
- One sample SB-3 (9-10') had a reported concentration of Arsenic that exceeded two GRCC (DWP, GSIP).

Based on the results of the Phase II ESA, the subject property meets the definition of a *facility* as defined in Part 201 of the NREPA, Michigan Public Act 451, of 1994, as amended.

### 5.4 RECOMMENDATIONS

ERG recommends future owners/operators prepare a BEA within 45 days after the earlier of the date of purchase, occupancy, or foreclosure and disclose it to MDEQ and subsequent purchaser or transferee.

In addition, because the subject property meets the definition of a *facility*, current and future owners/operators are required to comply with Section 20107(a). ERG recommends conducting a Section 20107(a) Compliance Analysis to assure compliance with Due Care obligations and preparing a Due Care Plan. The Due Care Plan will comply with applicable federal (HOME) requirements and MSHDA standards and will be submitted to DEQ under the Response Activity Plan review process. Due Care obligations include:

- Undertaking measures to prevent exacerbation of existing contamination.
- Exercising due care by undertaking response activities to mitigate unacceptable exposure to hazardous substances, mitigate fire and explosion hazards due to hazardous substances, and allow for the intended use of the subject property in a manner that protects health and safety.
- Taking reasonable precautions against the reasonably foreseeable acts or omissions of a third party and the consequences that could result from those acts or omissions.
- Provide notifications to the MDEQ and others in regard to mitigating fire and explosions hazards, discarded or abandoned containers, contamination migrating beyond property boundaries, as applicable.

The Due Care Plan will include a discussion of the items required for the "No Further Action Report". In addition, the Due Care Plan at a minimum will describe:

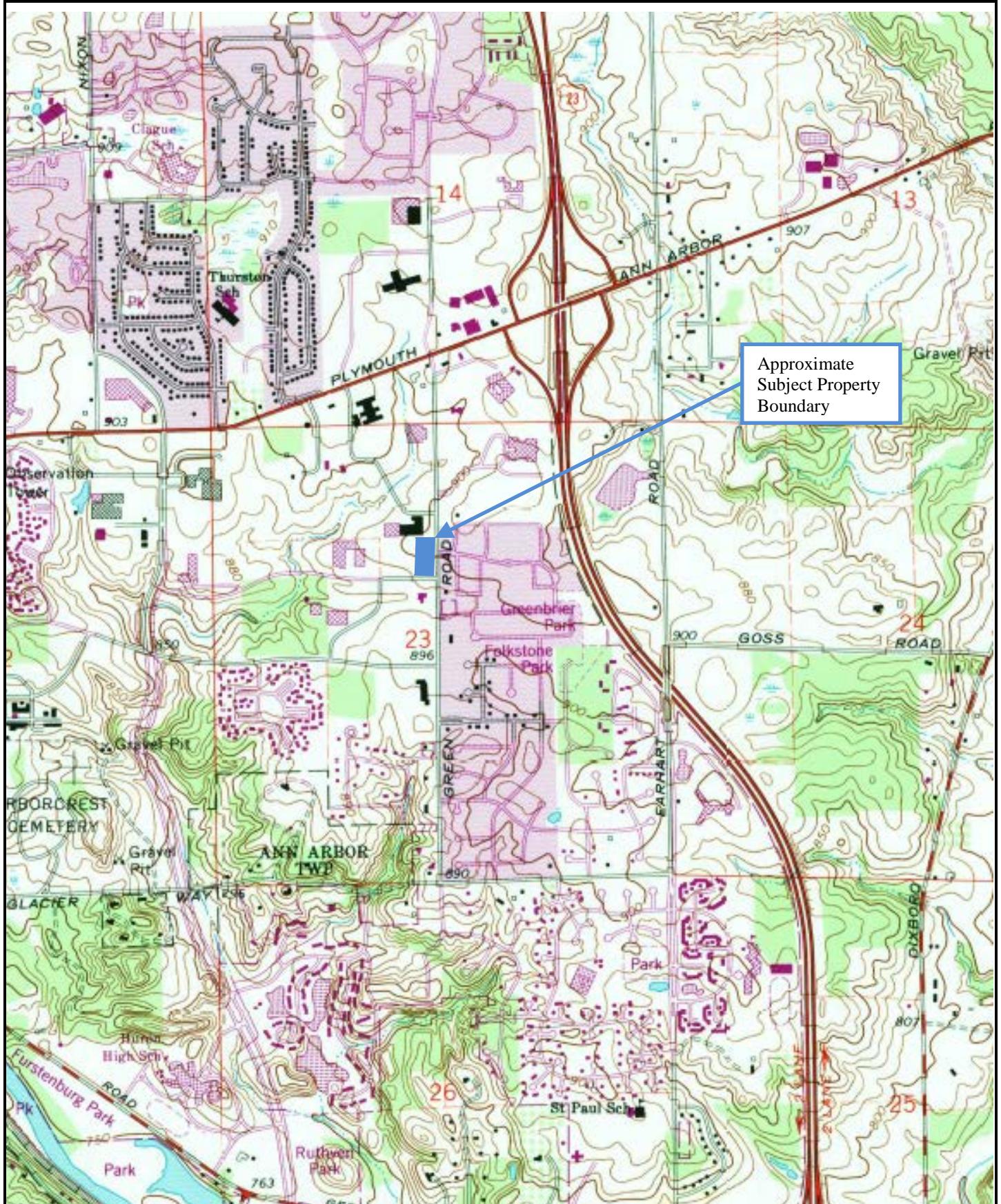
- the components of the Post Closure Plan
- any proposed Restrictions
- if Post Closure Agreement is required,
- if any ongoing Remedial Actions or Monitoring will be necessary

The location where soil concentrations exceed the Direct Contact GRCC is below an existing paved parking lot. Therefore, it is anticipated that the pavement will provide the necessary exposure barrier.

## **FIGURES**

**Figure 1: Location Map**

**Figure 2: Site Map**



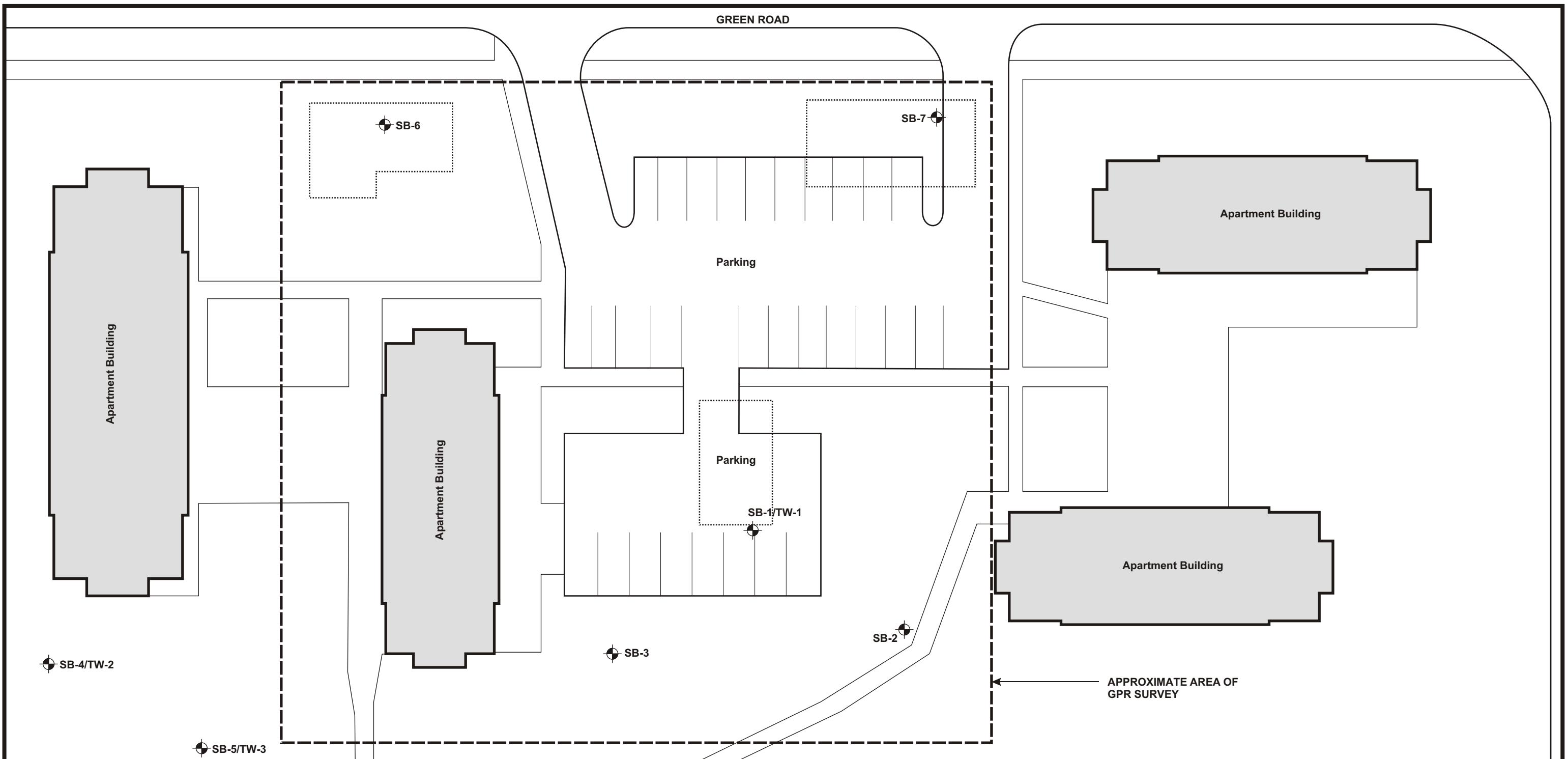
**Figure : 1 - Site Vicinity Map**  
Green Baxter Court  
1701-1747 Green Road, Ann Arbor, Michigan

Source: USGS Ann Arbor East 7.5 Minute Topographic Map



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Legend:

- Soil Boring Location
- TW Temporary Well
- Approximate Historical Structure Location

Note:

1. Source: Gene Terrill Associates Architects "Developmental Plan Site 'N'" dated 3/31/69 and Google Earth image dated 5/9/2010.

Scale:

1-Inch = 30-Feet



Drawn By: S Bennett

Date: 8/12/13

**Soil Boring Diagram**

**Green Baxter**  
1701-1749 Green Road  
Ann Arbor, MI

**ERG** Environmental Resources Group  
18003 Center Oaks Court • Suite 106 • Wixom, MI • 48393  
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## **TABLES**

**Table 1: Soil Analytical Data Summary Table**

**Table 2: Groundwater Analytical Data Summary Table**

**Table 1: Soil Gas Analytical Data Summary Table**

**TABLE 1: SOIL**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number (CAS)	Statewide Default Background Levels	Drinking Water Protection Criteria & RBSLs	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	SB-1 (1.5-2')	SB-2 (8-9')	SB-3 (9-10')	SB-6 (1-2')	SB-7 (12-13')	
				Groundwater Surface Water Interface Protection Criteria & RBSLs	Groundwater Contact Protection Criteria & RBSLs	Soil Volatilization to Indoor Air Inhalation Criteria & RBSLs	Infinite Source Volatile Soil Inhalation Criteria (VSIC) & RBSLs	Direct Contact Criteria & RBSLs	Particulate Soil Inhalation Criteria & RBSLs							
				Depth												
<b>Volatiles</b>																
Acetone (l)	67-64-1	NA	15,000	34,000	1.1E+8 (C)	1.1E+8 (C)	1.3E+8	2.3E+7	1.7E+11		<1,000	<1,000	<1,000	<1,000	<1,000	<1,000
Acrylonitrile (l)	107-13-1	NA	100 (M); 52	100 (M); 40	2.8E+5	6,600	5,000	16,000	5.8E+7		<100	<100	<100	<100	<100	<100
Benzene (l)	71-43-2	NA	100	4,000 (X)	2.2E+5	1,600	13,000	1.8E+5	4.7E+8		<50	<50	<50	<50	<50	<50
Bromobenzene (l)	108-86-1	NA	550	NA	3.6E+5	3.1E+5	4.5E+5	5.4E+5	2.4E+8		<100	<100	<100	<100	<100	<100
Bromochloromethane	74-97-5	NA	18	ID	12,000	3.10E+05	4.50E+05	5.40E+05	5.30E+08		<110	<120	<110	<110	<110	<110
Bromodichloromethane	75-27-4	NA	1,600 (W)	ID	2.8E+5	1,200	9,100	1.1E+5	1.1E+8		<100	<100	<100	<100	<100	<100
Bromoform	75-25-2	NA	1,600 (W)	ID	8.7E+5 (C)	1.5E+5	9.0E+5	8.2E+5	3.6E+9		<110	<120	<110	<110	<110	<110
Bromomethane	74-83-9	NA	200	700	1.4E+6	860	11,000	3.2E+5	1.5E+8		<200	<200	<200	<200	<200	<200
2-Butanone (MEK) (l)	78-93-3	NA	2.6E+5	44,000	2.7E+7 (C)	2.7E+7 (C)	2.9E+7	2.7E+7 (C,DD)	2.9E+10		<750	<750	<750	<750	<750	<750
n-Butylbenzene	104-51-8	NA	1,600	ID	1.2E+5	ID	ID	2.5E+6	8.8E+8		<50	<50	<50	<50	<50	<50
sec-Butylbenzene	135-98-8	NA	1,600	ID	88,000	ID	ID	2.5E+6	1.8E+8		<50	<50	<50	<50	<50	<50
tert-Butylbenzene (l)	98-06-6	NA	1,600	ID	1.8E+5	ID	ID	2.5E+6	2.9E+8		<50	<50	<50	<50	<50	<50
Carbon disulfide (I,R)	75-15-0	NA	16,000	ID	2.8E+5 (C)	76,000	1.3E+6	2.8E+5 (C,DD)	2.1E+10		<250	<250	<250	<250	<250	<250
Carbon tetrachloride	56-23-5	NA	100	900 (X)	92,000	190	3,500	96,000	1.7E+8		<55	<58	<57	<55	<57	<57
Chlorobenzene (l)	108-90-7	NA	2,000	500	2.6E+5 (C)	1.2E+5	7.7E+5	2.6E+5 (C)	2.1E+9		<55	<58	<57	<55	<57	<57
Chloroethane	75-00-3	NA	8,600	22,000 (X)	9.5E+5 (C)	9.5E+5 (C)	3.0E+7	9.5E+5 (C)	2.9E+11		<270	<290	<280	<270	<280	<280
Chloroform	67-66-3	NA	1,600 (W)	7,000	1.5E+6 (C)	7,200	45,000	1.2E+6	1.6E+9		<55	<58	<57	<55	<57	<57
Chloromethane (l)	74-87-3	NA	5,200	ID	1.1E+6 (C)	2,300	40,000	1.1E+6 (C)	2.6E+9		<250	<250	<250	<250	<250	<250
o-Chlorotoluene (l)	95-49-8	NA	3,300	ID	5.0E+5 (C)	2.7E+5	1.2E+6	5.0E+5 (C)	2.1E+9		<50	<50	<50	<50	<50	<50
Dibromochloromethane	124-48-1	NA	1,600 (W)	ID	3.6E+5	3,900	24,000	1.1E+5	1.6E+8		<100	<100	<100	<100	<100	<100
Dibromochloropropane	96-12-8	NA	10 (M); 4.0	ID	1,200 (C)	220	260	1,200 (C)	7.0E+5		<27	<29	<28	<27	<28	<28

Units in µg/kg = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

&lt; = Analyte was not detected at or above the specified reporting limit

Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

**TABLE 1: SOIL**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number (CAS)	Statewide Default Background Levels	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA							Sample Identification	SB-1 (1.5-2')	SB-2 (8-9')	SB-3 (9-10')	SB-6 (1-2')	SB-7 (12-13')	
			Drinking Water Protection Criteria & RBSLs	Groundwater Surface Water Interface Protection Criteria & RBSLs	Groundwater Contact Protection Criteria & RBSLs	Soil Volatilization to Indoor Air Inhalation Criteria & RBSLs	Infinite Source Volatile Soil Inhalation Criteria (VSIC) & RBSLs	Direct Contact Criteria & RBSLs	Particulate Soil Inhalation Criteria & RBSLs							
			Depth	1.5'-2'	8'-9'	9'-10'	1'-2'	12'-13'								
<b>Volatiles</b>																
Dibromomethane	74-95-3	NA	1,600	NA	2.0E+6 (C)	ID	ID	2.0E+6 (C)	ID		<250	<250	<250	<250	<250	<250
1,2-Dichlorobenzene	95-50-1	NA	14,000	280	2.1E+5 (C)	2.1E+5 (C)	3.9E+7	2.1E+5 (C)	4.4E+10		<100	<100	<100	<100	<100	<100
1,3-Dichlorobenzene	541-73-1	NA	170	680	51,000	26,000	79,000	1.7E+5 (C)	8.8E+7		<100	<100	<100	<100	<100	<100
1,4-Dichlorobenzene	106-46-7	NA	1,700	360	1.4E+5	19,000	77,000	4.0E+5	5.7E+8		<100	<100	<100	<100	<100	<100
Dichlorodifluoromethane	75-71-8	NA	95,000	ID	1.0E+6 (C)	9.0E+5	5.3E+7	1.0E+6 (C)	1.5E+12		<250	<250	<250	<250	<250	<250
1,1-Dichloroethane	75-34-3	NA	18,000	15,000	8.9E+5 (C)	2.3E+5	2.1E+6	8.9E+5 (C)	1.5E+10		<55	<58	<57	<55	<57	<57
1,2-Dichloroethane (l)	107-06-2	NA	100	7,200 (X)	3.8E+5	2,100	6,200	91,000	1.5E+8		<55	<58	<57	<55	<57	<57
cis-1,2-Dichloroethylene	156-59-2	NA	1,400	12,000	6.4E+5 (C)	22,000	1.8E+5	6.4E+5 (C)	1.0E+9		<50	<50	<50	<50	<50	<50
trans-1,2-Dichloroethylene	156-60-5	NA	2,000	30,000 (X)	1.4E+6 (C)	23,000	2.8E+5	1.4E+6 (C)	2.1E+9		<50	<50	<50	<50	<50	<50
1,1-Dichloroethylene (l)	75-35-4	NA	140	2,600	2.2E+5	62	1,100	2.0E+5	7.8E+7		<50	<50	<50	<50	<50	<50
1,2-Dichloropropane (l)	78-87-5	NA	100	4,600 (X)	3.2E+5	4,000	25,000	1.4E+5	1.2E+8		<55	<58	<57	<55	<57	<57
cis-1,3-Dichloropropylene	10061-01-5	NA	NA	NA	NA	NA	NA	NA	NA		<55	<58	<57	<55	<57	<57
trans-1,3-Dichloropropylene	10061-02-6	NA	NA	NA	NA	NA	NA	NA	NA		<55	<58	<57	<55	<57	<57
Ethylbenzene (l)	100-41-4	NA	1,500	360	1.4E+5 (C)	87,000	7.2E+5	1.4E+5 (C)	1.3E+10		<50	<50	<50	<50	<50	<50
Ethylene dibromide	106-93-4	NA	20 (M); 1.0	110 (X)	500	670	1,700	92	1.8E+7		<55	<58	<57	<55	<57	<57
2-Hexanone	591-78-6	NA	20,000	ID	2.5E+6 (C)	9.9E+5	1.1E+6	2.5E+6 (C)	1.2E+9		<2,500	<2,500	<2,500	<2,500	<2,500	<2,500
Isopropyl benzene	98-82-8	NA	91,000	3,200	3.9E+5 (C)	3.9E+5 (C)	1.7E+6	3.9E+5 (C)	2.6E+9		<250	<250	<250	<250	<250	<250
Methyl Iodide	74-88-4	NA	NA	NA	NA	NA	NA	NA	NA		<110	<120	<110	<110	<110	<110
4-Methyl-2-pentanone (MIBK) (l)	108-10-1	NA	36,000	ID	2.7E+6 (C)	2.7E+6 (C)	4.5E+7	2.7E+6 (C)	6.0E+10		<2,500	<2,500	<2,500	<2,500	<2,500	<2,500
Methylene chloride	75-09-2	NA	100	30,000 (X)	2.3E+6 (C)	45,000	2.1E+5	1.3E+6	8.3E+9		<100	<100	<100	<100	<100	<100
2-Methylnaphthalene	91-57-6	NA	57,000	4,200	5.5E+6	2.7E+6	1.5E+6	8.1E+6	2.9E+8		<330	<330	<330	<330	<330	<330
Methyl-tert-butyl ether (MTBE)	1634-04-4	NA	800	1.4E+5 (X)	5.9E+6 (C)	5.9E+6 (C)	2.5E+7	1.5E+6	8.8E+10		<250	<250	<250	<250	<250	<250
Naphthalene	91-20-3	NA	35,000	730	2.1E+6	2.5E+5	3.0E+5	1.6E+7	8.8E+7		<330	<330	<330	<330	<330	<330

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**TABLE 1: SOIL**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number (CAS)	Statewide Default Background Levels	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA							Sample Identification	SB-1 (1.5-2')	SB-2 (8-9')	SB-3 (9-10')	SB-6 (1-2')	SB-7 (12-13')					
			Drinking Water Protection Criteria & RBSLs	Groundwater Surface Water Interface Protection Criteria & RBSLs	Groundwater Contact Protection Criteria & RBSLs	Soil Volatilization to Indoor Air Inhalation Criteria & RBSLs	Infinite Source Volatile Soil Inhalation Criteria (VSIC) & RBSLs	Direct Contact Criteria & RBSLs	Particulate Soil Inhalation Criteria & RBSLs											
									Depth											
<i>Volatiles</i>																				
n-Propylbenzene (l)	103-65-1	NA	1,600	ID	3.0E+5	ID	ID	2.5E+6	5.9E+8		<100	<100	<100	<100	<100	<100				
Styrene	100-42-5	NA	2,700	2,100 (X)	2.7E+5	2.5E+5	9.7E+5	4.0E+5	6.9E+9		<50	<50	<50	<50	<50	<50				
1,1,1,2-Tetrachloroethane	630-20-6	NA	1,500	ID	4.4E+5 (C)	6,200	36,000	4.4E+5 (C)	5.3E+8		<100	<100	<100	<100	<100	<100				
1,1,2,2-Tetrachloroethane	79-34-5	NA	170	1,600 (X)	94,000	4,300	10,000	53,000	6.8E+7		<55	<58	<57	<55	<57	<57				
Tetrachloroethylene	127-18-4	NA	100	1,200 (X)	88,000 (C)	11,000	1.7E+5	88,000 (C)	1.2E+9		<50	<50	<50	<50	<50	<50				
Toluene (l)	108-88-3	NA	16,000	5,400	2.5E+5 (C)	2.5E+5 (C)	2.8E+6	2.5E+5 (C)	1.2E+10		<50	<50	<50	<50	<50	<50				
1,2,4-Trichlorobenzene	120-82-1	NA	4,200	5,900 (X)	1.1E+6 (C)	1.1E+6 (C)	2.8E+7	9.9E+5 (DD)	1.1E+10		<330	<330	<330	<330	<330	<330				
1,1,1-Trichloroethane	71-55-6	NA	4,000	1,800	4.6E+5 (C)	2.5E+5	3.8E+6	4.6E+5 (C)	2.9E+10		<55	<58	<57	<55	<57	<57				
1,1,2-Trichloroethane	79-00-5	NA	100	6,600 (X)	4.2E+5	4,600	17,000	1.8E+5	2.5E+8		<55	<58	<57	<55	<57	<57				
Trichloroethylene	79-01-6	NA	100	4,000 (X)	4.4E+5	1,000	11,000	5.0E+5 (C,DD)	5.9E+7		<55	<58	<57	<55	<57	<57				
Trichlorofluoromethane	75-69-4	NA	52,000	NA	5.6E+5 (C)	5.6E+5 (C)	9.2E+7	5.6E+5 (C)	1.7E+12		<100	<100	<100	<100	<100	<100				
1,2,3-Trichloropropane	96-18-4	NA	840	NA	8.3E+5 (C)	4,000	9,200	8.3E+5 (C)	8.8E+6		<110	<120	<110	<110	<110	<110				
1,2,3-Trimethylbenzene	526-73-8	NA	NA	NA	NA	NA	NA	NA	NA		<100	<100	<100	<100	<100	<100				
1,2,4-Trimethylbenzene (l)	95-63-6	NA	2,100	570	1.1E+5 (C)	1.1E+5 (C)	2.1E+7	1.1E+5 (C)	3.6E+10		<100	<100	<100	<100	<100	<100				
1,3,5-Trimethylbenzene (l)	108-67-8	NA	1,800	1,100	94,000 (C)	94,000 (C)	1.6E+7	94,000 (C)	3.6E+10		<100	<100	<100	<100	<100	<100				
Vinyl chloride	75-01-4	NA	40	260 (X)	20,000	270	4,200	3,800	8.9E+8		<40	<40	<40	<40	<40	<40				
Xylenes (l)	1330-20-7	NA	5,600	820	1.5E+5 (C)	1.5E+5 (C)	4.6E+7	1.5E+5 (C)	1.3E+11		<150	<150	<150	<150	<150	<150				

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**TABLE 1: SOIL**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number (CAS)	Statewide Default Background Levels	Drinking Water Protection Criteria & RBSLs	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	SB-1 (1.5-2')	SB-2 (8-9')	SB-3 (9-10')	SB-6 (1-2')	SB-7 (12-13')	
				Groundwater Surface Water Interface Protection Criteria & RBSLs	Groundwater Contact Protection Criteria & RBSLs	Soil Volatilization to Indoor Air Inhalation Criteria & RBSLs	Infinite Source Volatile Soil Inhalation Criteria (VSIC) & RBSLs	Direct Contact Criteria & RBSLs	Particulate Soil Inhalation Criteria & RBSLs							
				Depth												
<b>Semivolatiles</b>																
Acenaphthene	83-32-9	NA	3.0E+5	8,700	9.7E+5	1.9E+8	8.1E+7	4.1E+7	6.2E+9		<330	<330	<330	<330	<330	
Acenaphthylene	208-96-8	NA	5,900	ID	4.4E+5	1.6E+6	2.2E+6	1.6E+6	1.0E+9		<b>2,000</b>	<330	<330	<330	<330	
Anthracene	120-12-7	NA	41,000	ID	41,000	1.0E+9 (D)	1.4E+9	2.3E+8	2.9E+10		<b>1,700</b>	<330	<330	<330	<330	
Benzo(a)anthracene (Q)	56-55-3	NA	NLL	NLL	NLL	NLV	NLV	20,000	ID		<b>5,800</b>	<330	<330	<330	<330	
Benzo(a)pyrene (Q)	50-32-8	NA	NLL	NLL	NLL	NLV	NLV	2,000	1.9E+6		<b>7,800</b>	<330	<330	<330	<330	
Benzo(b)fluoranthene (Q)	205-99-2	NA	NLL	NLL	NLL	ID	ID	20,000	ID		<b>9,000</b>	<330	<330	<330	<330	
Benzo(g,h,i)perylene	191-24-2	NA	NLL	NLL	NLL	NLV	NLV	2.5E+6	3.5E+8		<b>6,500</b>	<330	<330	<330	<330	
Benzo(k)fluoranthene (Q)	207-08-9	NA	NLL	NLL	NLL	NLV	NLV	2.0E+5	ID		<b>3,000</b>	<330	<330	<330	<330	
Chrysene (Q)	218-01-9	NA	NLL	NLL	ID	ID	ID	2.0E+6	ID		<b>4,600</b>	<330	<330	<330	<330	
Dibenzo(a,h)anthracene (Q)	53-70-3	NA	NLL	NLL	NLL	NLV	NLV	2,000	ID		<b>1,900</b>	<330	<330	<330	<330	
Fluoranthene	206-44-0	NA	7.3E+5	<b>5,500</b>	7.3E+5	1.0E+9 (D)	7.4E+8	4.6E+7	4.1E+9		<b>9,400</b>	<330	<330	<330	<330	
Fluorene	86-73-7	NA	3.9E+5	5,300	8.9E+5	5.8E+8	1.3E+8	2.7E+7	4.1E+9		<b>610</b>	<330	<330	<330	<330	
Indeno(1,2,3-cd)pyrene (Q)	193-39-5	NA	NLL	NLL	NLL	NLV	NLV	20,000	ID		<b>6,800</b>	<330	<330	<330	<330	
2-Methylnaphthalene	91-57-6	NA	57,000	4,200	5.5E+6	2.7E+6	1.5E+6	8.1E+6	2.9E+8		<330	<330	<330	<330	<330	
Naphthalene	91-20-3	NA	35,000	730	2.1E+6	2.5E+5	3.0E+5	1.6E+7	8.8E+7		<330	<330	<330	<330	<330	
Phenanthrene	85-01-8	NA	56,000	<b>2,100</b>	1.1E+6	2.8E+6	1.6E+5	1.6E+6	2.9E+6		<b>4,100</b>	<330	<330	<330	<330	
Pyrene	129-00-0	NA	4.8E+5	ID	4.8E+5	1.0E+9 (D)	6.5E+8	2.9E+7	2.9E+9		<b>8,200</b>	<330	<330	<330	<330	
<b>Metals</b>																
Arsenic	7440-38-2	5,800	4,600	4,600	2.0E+6	NLV	NLV	7,600	9.1E+5		<b>7,300</b>	<b>5,600</b>	<b>6,800</b>	<b>5,500</b>	<b>5,100</b>	
Lead (B)	7439-92-1	21,000	7.0E+5	(G,X)	ID	NLV	NLV	4.0E+5	4.4E+7		<b>21,000</b>	<b>6,500</b>	<b>7,200</b>	<b>5,400</b>	<b>7,900</b>	

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**TABLE 1: SOIL**  
**Norstar Development**  
**Phase II ESA**  
**Project Number: 1127.001**  
**Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan**  
**July 2013**

Constituent	Chemical Abstract Service Number (CAS)	Statewide Default Background Levels	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA							Sample Identification	SB-1 (1.5-2')	SB-2 (8-9')	SB-3 (9-10')	SB-6 (1-2')	SB-7 (12-13')					
			Drinking Water Protection Criteria & RBSLs	Groundwater Surface Water Interface Protection Criteria & RBSLs	Groundwater Contact Protection Criteria & RBSLs	Soil Volatilization to Indoor Air Inhalation Criteria & RBSLs	Infinite Source Volatile Soil Inhalation Criteria (VSIC) & RBSLs	Direct Contact Criteria & RBSLs	Particulate Soil Inhalation Criteria & RBSLs											
									Depth											
<i>Pesticides, Chlorinated</i>																				
Aldrin	309-00-2	NA	NLL	NLL	NLL	1.3E+6	58,000	1,000	8.0E+5		<37	<20	<20	<20	<20	<20				
delta-BHC	319-86-8	NA	NA	NA	NA	NA	NA	NA	NA		<37	<20	<20	<20	<20	<20				
Chlordane (J)	57-74-9	NA	NLL	NLL	NLL	1.1E+7	1.2E+6	31,000	2.1E+7		<91	<25	<25	<25	<25	<25				
4-4'-DDD	72-54-8	NA	NLL	NLL	NLL	NLV	NLV	95,000	5.6E+7		<37	<20	<20	<20	<20	<20				
4-4'-DDE	72-55-9	NA	NLL	NLL	NLL	NLV	NLV	45,000	4.0E+7		<37	<20	<20	<20	<20	<20				
4-4'-DDT	50-29-3	NA	NLL	NLL	NLL	NLV	NLV	57,000	4.0E+7		<b>860</b>	<20	<20	<20	<20	<20				
Dieldrin	60-57-1	NA	NLL	NLL	NLL	1.4E+5	19,000	1,100	8.5E+5		<37	<20	<20	<20	<20	<20				
Endosulfan I	959-98-8	NA	NLL	NLL	NLL	NLV	NLV	1,400,000	ID		<37	<20	<20	<20	<20	<20				
Endosulfan II	33213-65-9	NA	NLL	NLL	NLL	NLV	NLV	1,400,000	ID		<37	<20	<20	<20	<20	<20				
Endosulfan Sulfate	1031-07-8	NA	NA	NA	NA	NA	NA	NA	NA		<37	<20	<20	<20	<20	<20				
Endrin	72-20-8	NA	NLL	NLL	NLL	NLV	NLV	65,000	ID		<37	<20	<20	<20	<20	<20				
Endrin Aldehyde	7421-93-4	NA	NA	NA	NA	NA	NA	NA	NA		<37	<20	<20	<20	<20	<20				
Heptachlor	76-44-8	NA	NLL	NLL	NLL	3.5E+5	62,000	5,600	3.0E+6		<37	<20	<20	<20	<20	<20				
Heptachlor epoxide	1024-57-3	NA	NLL	NLL	NLL	NLV	NLV	3,100	1.5E+6		<37	<20	<20	<20	<20	<20				
alpha-Hexachlorocyclohexane	319-84-6	NA	18	ID	2,500	30,000	12,000	2,600	2.1E+6		<37	<20	<20	<20	<20	<20				
beta-Hexachlorocyclohexane	319-85-7	NA	37	ID	5,100	NLV	NLV	5,400	7.4E+6		<37	<20	<20	<20	<20	<20				
Lindane	58-89-9	NA	20 (M); 7.0	20 (M); 1.1	7,100	ID	ID	8,300	ID		<37	<20	<20	<20	<20	<20				
Methoxychlor	72-43-5	NA	16,000	NA	18,000	ID	ID	1.9E+6	ID		<150	<50	<50	<50	<50	<50				
Toxaphene	8001-35-2	NA	24,000	8,200	3.6E+5	NLV	NLV	20,000	1.2E+7		<1,800	<390	<380	<360	<380	<380				

Units in µg/kg = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

< = Analyte was not detected at or above the specified reporting limit

Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

**FOOTNOTES  
for****Part 201 Criteria and Part 213 Risk-Based Screening Levels****Document Release Date: September 28, 2012**

- (A) Criterion is the state of Michigan drinking water standard established pursuant to Section 5 of 1976 PA 399, MCL 325.1005.
- (B) Background, as defined in R 299.5701(b), may be substituted if higher than the calculated cleanup criterion. Background levels may be less than criteria for some inorganic compounds.
- (C) Value presented is a screening level based on the chemical-specific generic soil saturation concentration ( $C_{sat}$ ) since the calculated risk-based criterion is greater than  $C_{sat}$ . Concentrations greater than  $C_{sat}$  are acceptable cleanup criteria for this pathway where a site-specific demonstration indicates that free-phase material containing a hazardous substance is not present.
- (D) Calculated criterion exceeds 100 percent, hence it is reduced to 100 percent or 1.0E+9 parts per billion (ppb).
- (E) Criterion is the aesthetic drinking water value, as required by Section 20120a(5) of the Natural Resources and Environmental Protection Act, 1994 PA 451, as amended (NREPA). A notice of aesthetic impact may be employed as an institutional control mechanism if groundwater concentrations exceed the aesthetic drinking water criterion, but do not exceed the applicable health-based drinking water value provided in the following table:

Hazardous Substance	Chemical Abstract Service Number	Residential Health-Based Drinking Water Value	Non-Residential Health-Based Drinking Water Value
Aluminum	7429905	300	4,100
tertiary Amyl methyl ether	994058	910	2,600
Copper	7440508	1,400	4,000
Diethyl ether	60297	3,700	10,000
Ethylbenzene	100414	700	700
Iron	7439896	2,000	5,600
Manganese	7439965	860	2,500
Methyl-tert-butyl ether (MTBE)	1634044	240	690
Toluene	108883	1,000	1,000
1,2,4-Trimethylbenzene	95636	1,000	2,900
1,3,5-Trimethylbenzene	108678	1,000	2,900
Xylenes	1330207	10,000	10,000

- (F) Criterion is based on adverse impacts to plant life and phytotoxicity.
- (G) Groundwater surface water interface (GSI) criterion depends on the pH or water hardness, or both, of the receiving surface water. The final chronic value (FCV) for the protection of aquatic life shall be calculated based on the pH or hardness of the receiving surface water. Where water hardness exceeds 400 mg CaCO<sub>3</sub>/L, use 400 mg CaCO<sub>3</sub>/L for the FCV calculation. The FCV formula provides values in units of ug/L or ppb. The generic GSI criterion is the lesser of

the calculated FCV, the wildlife value (WV), and the surface water human non-drinking water value (HNDV). The soil GSI protection criteria for these hazardous substances are the greater of the 20 times the GSI criterion or the GSI soil-water partition values using the GSI criteria developed with the procedure described in this footnote.

Hazardous Substance	FCV Formula ug/L	FCV Conversion Factor (CF)	WV ug/L	HNDV ug/L
Acetate	$\text{EXP}(0.2732*(\text{pH}) + 7.0362)$	NA	NA	1.3E+6
Acetic Acid	$\text{EXP}(0.2732*(\text{pH}) + 7.0362)$	NA	NA	1.3E+6
Barium	$\text{EXP}(1.0629*(\text{LnH}) + 1.1869)$	NA	NA	1.6E+5
Beryllium	$\text{EXP}(2.5279*(\text{LnH}) - 10.7689)$	NA	NA	1,200
Cadmium <sup>®</sup>	$(\text{EXP}(0.7852*(\text{LnH}) - 2.715)) * \text{CF}$	$1.101672 - ((\text{LnH}) * (0.041838))$	NA	130
Chromium (III) <sup>®</sup>	$(\text{EXP}(0.819*(\text{LnH}) + 0.6848)) * \text{CF}$	0.86	NA	9,400
Copper	$(\text{EXP}(0.8545*(\text{LnH}) - 1.702)) * \text{CF}$	0.96	NA	38,000
Lead <sup>®</sup>	$(\text{EXP}(0.9859*(\text{LnH}) - 1.270)) * \text{CF}$	$1.46203 - ((\text{LnH}) * (0.14571))$	NA	190
Manganese <sup>®</sup>	$\text{EXP}(0.8784*(\text{LnH}) + 3.5385)$	NA	NA	59,000
Nickel	$(\text{EXP}(0.846*(\text{LnH}) + 0.0584)) * \text{CF}$	0.997	NA	2.1E+5
Pentachlorophenol <sup>®</sup>	$\text{EXP}(1.005*(\text{pH}) - 5.134)$	NA	NA	2.8
Zinc	$(\text{EXP}(0.8473*(\text{LnH}) + 0.884)) * \text{CF}$	0.986	NA	16,000

where,

- $\text{EXP}(x)$  = The base of the natural logarithm raised to power x ( $e^x$ ).
- $\text{LnH}$  = The natural logarithm of water hardness in mg CaCO<sub>3</sub>/L.
- $*$  = The multiplication symbol.
- <sup>®</sup> = The GSI criterion developed here may not be protective for surface water that is used as a drinking water source. Refer to footnote (X) for further guidance.

A spreadsheet that may be used to calculate GSI and GSI protection criteria for (G)-footnoted hazardous substances is available on the Department of Environmental Quality (DEQ) internet web site.

- (H) Valence-specific chromium data (Cr III and Cr VI) shall be compared to the corresponding valence-specific cleanup criteria. If both Cr III and Cr VI are present in groundwater, the total concentration of both cannot exceed the drinking water criterion of 100 ug/L. If analytical data are provided for total chromium only, they shall be compared to the cleanup criteria for Cr VI. Cr III soil cleanup criterion for protection of drinking water can only be used at sites where groundwater is prevented from being used as a public water supply, currently and in the future, through an approved land or resource use restriction.
- (I) Hazardous substance may exhibit the characteristic of ignitability as defined in 40 C.F.R. §261.21 (revised as of July 1, 2001), which is adopted by reference in these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, Remediation and Redevelopment Division (RRD), 525 West Allegan Street, Lansing, Michigan 48933, at cost.

- (J) Hazardous substance may be present in several isomer forms. Isomer-specific concentrations shall be added together for comparison to criteria.
- (K) Hazardous substance may be flammable or explosive, or both.
- (L) Criteria for lead are derived using a biologically based model, as allowed for under Section 20120a(10) of the NREPA, and are not calculated using the algorithms and assumptions specified in pathway-specific rules. The generic residential drinking water criterion of 4 ug/L is linked to the generic residential soil direct contact criterion of 400 mg/kg. A higher concentration in the drinking water, up to the state action level of 15 ug/L, may be allowed as a site-specific remedy and still allow for drinking water use, under Section 20120a(2) of the NREPA if soil concentrations are appropriately lower than 400 mg/kg. If a site-specific criterion is approved based on this subdivision, a notice shall be filed on the deed for all property where the groundwater concentrations will exceed 4 ug/L to provide notice of the potential for unacceptable risk if soil or groundwater concentrations increase. Acceptable combinations of site-specific soil and drinking water concentrations are presented in the following table:

Acceptable Combinations of Lead in Drinking Water and Soil

Drinking Water Concentration (ug/L)	Soil Concentration (mg/kg)
5	386-395
6	376-385
7	376-385
8	366-375
9	356-365
10	346-355
11	336-345
12	336-345
13	326-335
14	316-325
15	306-315

- (M) Calculated criterion is below the analytical target detection limit, therefore, the criterion defaults to the target detection limit.
- (N) The concentrations of all potential sources of nitrate-nitrogen (e.g., ammonia-N, nitrite-N, nitrate-N) in groundwater that is used as a source of drinking water shall not, when added together, exceed the nitrate drinking water criterion of 10,000 ug/L. Where leaching to groundwater is a relevant pathway, soil concentrations of all potential sources of nitrate-nitrogen shall not, when added together, exceed the nitrate drinking water protection criterion of 2.0E+5 ug/kg.
- (O) The concentration of all polychlorinated and polybrominated dibenzodioxin and dibenzofuran isomers present at a facility, expressed as an equivalent concentration of 2,3,7,8-tetrachlorodibenzo-p-dioxin based upon their relative potency, shall be added together and compared to the criteria for 2,3,7,8-tetrachlorodibenzo-p-dioxin. The generic cleanup criteria for 2,3,7,8-tetrachlorodibenzo-p-dioxin are not calculated according to the algorithms presented in R 299.5714 to R 299.5726. The generic cleanup criteria are being held at the values that the DEQ has used since August 1998, in recognition of

- the fact that national efforts to reassess risks posed by dioxin are not yet complete. Until these studies are complete, it is premature to select a revised slope factor and/or reference dose for calculation of generic cleanup criteria.
- (P) Amenable cyanide methods or method OIA-1677 shall be used to quantify cyanide concentrations for compliance with all groundwater criteria. Total cyanide methods or method OIA-1677 shall be used to quantify cyanide concentrations for compliance with soil criteria. Nonresidential direct contact criteria may not be protective of the potential for release of hydrogen cyanide gas. Additional land or resource use restrictions may be necessary to protect for the acute inhalation concerns associated with hydrogen cyanide gas.
- (Q) Criteria for carcinogenic polycyclic aromatic hydrocarbons were developed using relative potential potencies to benzo(a)pyrene.
- (R) Hazardous substance may exhibit the characteristic of reactivity as defined in 40 C.F.R. §261.23 (revised as of July 1, 2001), which is adopted by reference in these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost.
- (S) Criterion defaults to the hazardous substance-specific water solubility limit.
- (T) Refer to the federal Toxic Substances Control Act (TSCA), 40 C.F.R. §761, Subpart D and 40 C.F.R. §761, Subpart G, to determine the applicability of TSCA cleanup standards. Subpart D and Subpart G of 40 C.F.R. §761 (July 1, 2001) are adopted by reference in these rules and are available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulations may be purchased, at a cost as of the time of adoption of these rules of \$55, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401, or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost. Alternatives to compliance with the TSCA standards listed below are possible under 40 C.F.R. §761 Subpart D. New releases may be subject to the standards identified in 40 C.F.R. §761, Subpart G. Use Part 201 soil direct contact cleanup criteria in the following table if TSCA standards are not applicable.

Land Use Category	TSCA, Subpart D Cleanup Standards	Part 201 Soil Direct Contact Cleanup Criteria
Residential	1,000 ppb, or 10,000 ppb if capped	4,000 ppb
Nonresidential	1,000 ppb, or 10,000 ppb if capped	16,000 ppb

- (U) Hazardous substance may exhibit the characteristic of corrosivity as defined in 40 C.F.R. §261.22 (revised as of July 1, 2001), which is adopted by reference in

these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost.

- (V) Criterion is the aesthetic drinking water value as required by Section 20120(a)(5) of the NREPA. Concentrations up to 200 ug/L may be acceptable, and still allow for drinking water use, as part of a site-specific cleanup under Section 20120a(2) of the NREPA.
- (W) Concentrations of trihalomethanes in groundwater shall be added together to determine compliance with the Michigan drinking water standard of 80 ug/L. Concentrations of trihalomethanes in soil shall be added together to determine compliance with the drinking water protection criterion of 1,600 ug/kg.
- (X) The GSI criterion shown in the generic cleanup criteria tables is not protective for surface water that is used as a drinking water source. For a groundwater discharge to the Great Lakes and their connecting waters or discharge in close proximity to a water supply intake in inland surface waters, the generic GSI criterion shall be the surface water human drinking water value (HDV) listed in the table in this footnote, except for those HDV indicated with an asterisk. For HDV with an asterisk, the generic GSI criterion shall be the lowest of the HDV, the WV, and the calculated FCV. See formulas in footnote (G). Soil protection criteria based on the HDV shall be as listed in the table in this footnote, except for those values with an asterisk. Soil GSI protection criteria based on the HDV shall be as listed in the table in this footnote, except for those values with an asterisk. Soil GSI protection criteria for compounds with an asterisk shall be the greater of 20 times the GSI criterion or the GSI soil-water partition values using the GSI criteria developed with the procedure described in this footnote.

Hazardous Substance	Chemical Abstract Service Number	Surface Water Human Drinking Water Values (HDV) (ug/L)	Soil GSI Protection Criteria for HDV (ug/kg)
Acrylamide	79061	0.5 (M); 0.12	10
Alachlor	15972608	3.5	88
Antimony	7440360	2.0 (M); 1.7	1,200
Benzene	71432	12	240
Boron	7440428	4,000	80,000
Bromate	15541454	10 (M); 0.5	200
n-Butanol	71363	3,500	70,000
Butyl benzyl phthalate	85687	6.9	13,000
Cadmium	7440439	2.5*	*
Carbon tetrachloride	56235	5.6	110
Chloride	16887006	50,000	1.0E+6
Chloroethane	75003	170	3,400
Chromium (III)	16065831	120*	*
Cyanazine	21725462	2.0 (M); 0.93	200 (M); 40
1,2-Dichloroethane	107062	6.0	120
trans-1,2-Dichloroethylene	156605	470	9,400
1,2-Dichloropropane	78875	9.1	180
1,3-Dichloropropene	542756	3.3	100 (M); 66
N,N-Dimethylacetamide	127195	700	14,000
1,4-Dioxane	123911	34	680
Ethylene dibromide	106934	0.17	20 (M); 3.4
Ethylene glycol	107211	56,000	1.1E+6
Hexachloroethane	67721	5.3	310
Isophorone	78591	310	6,200
Isopropyl alcohol	67630	28,000	5.6E+5
Lead	7439921	14*	*
Manganese	7439965	1,300*	*
Methanol	67561	14,000	2.8E+5
Methyl-tert-butyl ether (MTBE)	1634044	100	2,000
Methylene chloride	75092	47	940
Molybdenum	7439987	120	2,400
Nitrobenzene	98953	4.7	330 (M); 94
Pentachlorophenol	87865	1.8*	*
Styrene	100425	20	530
1,2,4,5-Tetrachlorobenzene	95943	2.8	3,300
1,1,2,2-Tetrachloroethane	79345	3.2	64
Tetrachloroethylene	127184	11	220
Tetrahydrofuran	109999	350	7,000
Thallium	7440280	2.0 (M); 1.2	1,400
1,2,4-Trichlorobenzene	120821	80	4,700
1,1,2-Trichloroethane	79005	12	240
Trichloroethylene	79016	29	580
Vinyl chloride	75014	1.0 (M); 0.25	40 (M); 20

- (Y) Source size modifiers shown in the following table shall be used to determine soil inhalation criteria for ambient air when the source size is not one-half acre. The modifier shall be multiplied by the generic soil inhalation criteria shown in the

table of generic cleanup criteria to determine the applicable criterion.

Source Size sq. feet or acres	Modifier
400 sq feet	3.17
1000 sq feet	2.2
2000 sq feet	1.76
1/4 acre	1.15
1/2 acre	1
1 acre	0.87
2 acre	0.77
5 acre	0.66
10 acre	0.6
32 acre	0.5
100 acre	0.43

- (Z) Mercury is typically measured as total mercury. The generic cleanup criteria, however, are based on data for different species of mercury. Specifically, data for elemental mercury, chemical abstract service (CAS) number 7439976, serve as the basis for the soil volatilization to indoor air criteria, groundwater volatilization to indoor air, and soil inhalation criteria. Data for methyl mercury, CAS number 22967926, serve as the basis for the GSI criterion; and data for mercuric chloride, CAS number 7487947, serve as the basis for the drinking water, groundwater contact, soil direct contact, and the groundwater protection criteria. Comparison to criteria shall be based on species-specific analytical data only if sufficient facility characterization has been conducted to rule out the presence of other species of mercury.
- (AA) Comparison to these criteria may take into account an evaluation of whether the hazardous substances are adsorbed to particulates rather than dissolved in water and whether filtered groundwater samples were used to evaluate groundwater.
- (BB) The state drinking water standard for asbestos is in units of fibers per milliliter of water (f/mL) longer than 10 millimicrons. Soil concentrations of asbestos are determined by polarized light microscopy.
- (CC) Groundwater: The generic GSI criteria are based on the toxicity of unionized ammonia ( $\text{NH}_3$ ); the criteria are 29 ug/L and 53 ug/L for cold water and warm water surface water, respectively. As a result, the GSI criterion shall be compared to the percent of the total ammonia concentration in the groundwater that will become  $\text{NH}_3$  in the surface water. This percent  $\text{NH}_3$  is a function of the pH and temperature of the receiving surface water and can be estimated using the following table, taken from Emerson, et al., (Journal of the Fisheries Research Board of Canada, Volume 32(12):2382, 1975).

Percent NH<sub>3</sub> in Aqueous Ammonia Solutions for 0-30 °C and pH 6-10

Temp (°F)	Temp (°C)	pH									
		6.0	6.5	7.0	7.5	8.0	8.5	9.0	9.5	10.0	
32.0	0	0.00827	0.0261	0.0826	0.261	0.820	2.55	7.64	20.7	45.3	
33.8	1	0.00899	0.0284	0.0898	0.284	0.891	2.77	8.25	22.1	47.3	
35.6	2	0.00977	0.0309	0.0977	0.308	0.968	3.00	8.90	23.6	49.4	
37.4	3	0.0106	0.0336	0.106	0.335	1.05	3.25	9.60	25.1	51.5	
39.2	4	0.0115	0.0364	0.115	0.363	1.14	3.52	10.3	26.7	53.5	
41.0	5	0.0125	0.0395	0.125	0.394	1.23	3.80	11.1	28.3	55.6	
42.8	6	0.0136	0.0429	0.135	0.427	1.34	4.11	11.9	30.0	57.6	
44.6	7	0.0147	0.0464	0.147	0.462	1.45	4.44	12.8	31.7	59.5	
46.4	8	0.0159	0.0503	0.159	0.501	1.57	4.79	13.7	33.5	61.4	
48.2	9	0.0172	0.0544	0.172	0.542	1.69	5.16	14.7	35.3	63.3	
50.0	10	0.0186	0.0589	0.186	0.586	1.83	5.56	15.7	37.1	65.1	
51.8	11	0.0201	0.0637	0.201	0.633	1.97	5.99	16.8	38.9	66.8	
53.6	12	0.0218	0.0688	0.217	0.684	2.13	6.44	17.9	40.8	68.5	
55.4	13	0.0235	0.0743	0.235	0.738	2.30	6.92	19.0	42.6	70.2	
57.2	14	0.0254	0.0802	0.253	0.796	2.48	7.43	20.2	44.5	71.7	
59.0	15	0.0274	0.0865	0.273	0.859	2.67	7.97	21.5	46.4	73.3	
60.8	16	0.0295	0.0933	0.294	0.925	2.87	8.54	22.8	48.3	74.7	
62.6	17	0.0318	0.101	0.317	0.996	3.08	9.14	24.1	50.2	76.1	
64.4	18	0.0343	0.108	0.342	1.07	3.31	9.78	25.5	52.0	77.4	
66.2	19	0.0369	0.117	0.368	1.15	3.56	10.5	27.0	53.9	78.7	
68.0	20	0.0397	0.125	0.396	1.24	3.82	11.2	28.4	55.7	79.9	
69.8	21	0.0427	0.135	0.425	1.33	4.10	11.9	29.9	57.5	81.0	
71.6	22	0.0459	0.145	0.457	1.43	4.39	12.7	31.5	59.2	82.1	
73.4	23	0.0493	0.156	0.491	1.54	4.70	13.5	33.0	60.9	83.2	
75.2	24	0.0530	0.167	0.527	1.65	5.03	14.4	34.6	62.6	84.1	
77.0	25	0.0569	0.180	0.566	1.77	5.38	15.3	36.3	64.3	85.1	
78.8	26	0.0610	0.193	0.607	1.89	5.75	16.2	37.9	65.9	85.9	
80.6	27	0.0654	0.207	0.651	2.03	6.15	17.2	39.6	67.4	86.8	
82.4	28	0.0701	0.221	0.697	2.17	6.56	18.2	41.2	68.9	87.3	
84.2	29	0.0752	0.237	0.747	2.32	7.00	19.2	42.9	70.4	88.3	
86.0	30	0.0805	0.254	0.799	2.48	7.46	20.3	44.6	71.8	89.0	

The generic approach for estimating NH<sub>3</sub> assumes a default pH of 8 and default temperatures of 68°F and 85°F for cold water and warm water surface water, respectively. The resulting percent NH<sub>3</sub> is 3.8 percent and 7.2 percent for cold water and warm water, respectively. This default percentage shall be multiplied by the total ammonia-nitrogen (NH<sub>3</sub>-N) concentration in the groundwater and the resulting NH<sub>3</sub> concentration compared to the applicable GSI criterion. As an

alternative, the maximum pH and temperature data from the specific receiving surface water can be used to estimate, from the table in this footnote, a lower percent unionized ammonia concentration for comparison to the generic GSI.

- Soil: The generic soil GSI protection criteria for unionized ammonia are 580 ug/kg and 1,100 ug/kg for cold water and warm water surface water, respectively.
- (DD) Hazardous substance causes developmental effects. Residential direct contact criteria are protective of both prenatal and postnatal exposure. Nonresidential direct contact criteria are protective for a pregnant adult receptor.
- (EE) The following are applicable generic GSI criteria as required by Section 20120a(15) of the NREPA.

Hazardous Substance	GSI (ug/L)	Notes
Phosphorus	1,000	Criteria applicable unless receiving water is a surface water that has a phosphorus waste load allocation or is an inland lake. In those cases, contact the department for applicable values.
Total dissolved solids (TDS)	5.0E+5	If TDS data are not available, the TDS criterion may be used a screening level for the sum of the concentrations of the following substances: Calcium, Chlorides, Iron, Magnesium, Potassium, Sodium, Sulfate.
Dissolved Oxygen (DO): Cold receiving waters Warm receiving waters	≥ 7,000 ≥ 5,000	Since a low level of DO can be harmful to aquatic life, the criterion represents a minimum level that on-site samples must exceed. This is in contrast to other criteria which represent "not to exceed" concentrations. DO criteria are not applicable if groundwater Carbonaceous Biochemical Oxygen Demand (CBOD) is less than 10,000 ug/L and groundwater ammonia concentration is less than 2,000 ug/L.

- (FF) The chloride GSI criterion shall be 125 mg/l when the discharge is to surface waters of the state designated as public water supply sources or 50 mg/l when the discharge is to the Great Lakes or connecting waters. Chloride GSI criteria shall not apply for surface waters of the state that are not designated as a public water supply source, however, the total dissolved solids criterion is applicable.
- (GG) Risk-based criteria are not available for methane due to insufficient toxicity data. An acceptable soil gas concentration (presented for both residential and nonresidential land uses) was derived utilizing 25 percent of the lower explosive level for methane. This equates to 1.25 percent or 8.4E+6 ug/m<sup>3</sup>.

"ID" means insufficient data to develop criterion.

"NA" means a criterion or value is not available or, in the case of background and CAS numbers, not applicable.

"NLL" means hazardous substance is not likely to leach under most soil conditions.

"NLV" means hazardous substance is not likely to volatilize under most conditions.

**TABLE 2: GROUNDWATER**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	TW-1	TW-2	TW-3	TRIP BLANK TB001102	
		Drinking Water Criteria & RBSLs	Groundwater Surface Water Interface Criteria & RBSLs	Groundwater Volatilization to Indoor Air Inhalation Criteria & RBSLs	Groundwater Contact Criteria and RBSLs	Flammability and Explosivity Screening Level	Acute Inhalation Screening Level						
							Sample Media		water	water	water	-	
							Collection Date		7/18/2013	7/18/2013	7/18/2013	7/18/2013	
									Depth (ft. bgs)	12'	9'	8'	NA
<b>Volatiles</b>													
Acetone (l)	67-64-1	730	1,700	1.0E+9 (D,S)	3.1E+7	1.5E+7	1.0E+9 (D)		<50	<50	<50	<50	
Acrylonitrile (l)	107-13-1	2.6	2.0 (M); 1.2	34,000	14,000	6.4E+6	ID		<2.0	<2.0	<2.0	<2.0	
Benzene (l)	71-43-2	5.0 (A)	200 (X)	5,600	11,000	68,000	67,000		<1.0	<1.0	<1.0	<1.0	
Bromobenzene (l)	108-86-1	18	NA	1.8E+5	12,000	ID	ID		<1.0	<1.0	<1.0	<1.0	
Bromochloromethane	74-97-5	NA	NA	NA	NA	NA	NA		<1.0	<1.0	<1.0	<1.0	
Bromodichloromethane	75-27-4	80 (A,W)	ID	4,800	14,000	ID	ID		<1.0	<1.0	<1.0	<1.0	
Bromoform	75-25-2	80 (A,W)	ID	4.7E+5	1.4E+5	ID	ID		<1.0	<1.0	<1.0	<1.0	
Bromomethane	74-83-9	10	35	4,000	70,000	ID	ID		<5.0	<5.0	<5.0	<5.0	
2-Butanone (MEK) (l)	78-93-3	13,000	2,200	2.4E+8 (S)	2.4E+8 (S)	ID	2.4E+8 (S)		<25	<25	<25	<25	
n-Butylbenzene	104-51-8	80	ID	ID	5,900	ID	ID		<1.0	<1.0	<1.0	<1.0	
sec-Butylbenzene	135-98-8	80	ID	ID	4,400	ID	ID		<1.0	<1.0	<1.0	<1.0	
tert-Butylbenzene (l)	98-06-6	80	ID	ID	8,900	ID	ID		<1.0	<1.0	<1.0	<1.0	
Carbon disulfide (I,R)	75-15-0	800	ID	2.5E+5	1.2E+6 (S)	13,000	ID		<5.0	<5.0	<5.0	<5.0	
Carbon tetrachloride	56-23-5	5.0 (A)	45 (X)	370	4,600	ID	96,000		<1.0	<1.0	<1.0	<1.0	
Chlorobenzene (l)	108-90-7	100 (A)	25	2.1E+5	86,000	1.6E+5	ID		<1.0	<1.0	<1.0	<1.0	
Chloroethane	75-00-3	430	1,100 (X)	5.7E+6 (S)	4.4E+5	1.1E+5	ID		<5.0	<5.0	<5.0	<5.0	
Chloroform	67-66-3	80 (A,W)	350	28,000	1.5E+5	ID	ID		<1.0	<1.0	<1.0	<1.0	
Chloromethane (l)	74-87-3	260	ID	8,600	4.9E+5	36,000	2.1E+5		<5.0	<5.0	<5.0	<5.0	
o-Chlorotoluene (l)	95-49-8	150	ID	2.2E+5	44,000	ID	ID		<5.0	<5.0	<5.0	<5.0	
Dibromochloromethane	124-48-1	80 (A,W)	ID	14,000	18,000	ID	ID		<5.0	<5.0	<5.0	<5.0	
Dibromochloropropane	96-12-8	0.2 (A)	ID	220	390	NA	ID		<1.0	<1.0	<1.0	<1.0	
Dibromomethane	74-95-3	80	NA	ID	5.3E+5	ID	ID		<5.0	<5.0	<5.0	<5.0	
1,2-Dichlorobenzene	95-50-1	600 (A)	13	1.6E+5 (S)	1.6E+5 (S)	NA	1.6E+5 (S)		<1.0	<1.0	<1.0	<1.0	
1,3-Dichlorobenzene	541-73-1	6.6	28	18,000	2,000	ID	ID		<1.0	<1.0	<1.0	<1.0	
1,4-Dichlorobenzene	106-46-7	75 (A)	17	16,000	6,400	NA	ID		<1.0	<1.0	<1.0	<1.0	

Units in µg/L = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

&lt; = Analyte was not detected at or above the specified reporting limit

Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

**TABLE 2: GROUNDWATER**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	TW-1	TW-2	TW-3	TRIP BLANK TB001102	
		Drinking Water Criteria & RBSLs	Groundwater Surface Water Interface Criteria & RBSLs	Groundwater Volatilization to Indoor Air Inhalation Criteria & RBSLs	Groundwater Contact Criteria and RBSLs	Flammability and Explosivity Screening Level	Acute Inhalation Screening Level						
							Sample Media		water	water	water	-	
							Collection Date		7/18/2013	7/18/2013	7/18/2013	7/18/2013	
								Depth (ft. bgs)	12'	9'	8'	NA	
<b>Volatiles</b>													
Dichlorodifluoromethane	75-71-8	1,700	ID	2.2E+5	3.0E+5 (S)	ID	ID		<5.0	<5.0	<5.0	<5.0	
1,1-Dichloroethane	75-34-3	880	740	1.0E+6	2.4E+6	3.8E+5	ID		<1.0	<1.0	<1.0	<1.0	
1,2-Dichloroethane (l)	107-06-2	5.0 (A)	360 (X)	9,600	19,000	2.5E+6	ID		<1.0	<1.0	<1.0	<1.0	
cis-1,2-Dichloroethylene	156-59-2	70 (A)	620	93,000	2.0E+5	5.3E+5	ID		<1.0	<1.0	<1.0	<1.0	
trans-1,2-Dichloroethylene	156-60-5	100 (A)	1,500 (X)	85,000	2.2E+5	2.3E+5	ID		<1.0	<1.0	<1.0	<1.0	
1,1-Dichloroethylene (l)	75-35-4	7.0 (A)	130	200	11,000	97,000	1.4E+5		<1.0	<1.0	<1.0	<1.0	
1,2-Dichloropropane (l)	78-87-5	5.0 (A)	230 (X)	16,000	16,000	5.5E+5	2.8E+6 (S)		<1.0	<1.0	<1.0	<1.0	
cis-1,3-Dichloropropylene	10061-01-5	NA	NA	NA	NA	NA	NA		<1.0	<1.0	<1.0	<1.0	
trans-1,3-Dichloropropylene	10061-02-6	NA	NA	NA	NA	NA	NA		<1.0	<1.0	<1.0	<1.0	
Ethylbenzene (l)	100-41-4	74 (E)	18	1.1E+5	1.7E+5 (S)	43,000	1.7E+5 (S)		<1.0	<1.0	<1.0	<1.0	
Ethylene dibromide	106-93-4	0.05 (A)	5.7 (X)	2,400	25	ID	ID		<1.0	<1.0	<1.0	<1.0	
2-Hexanone	591-78-6	1,000	ID	4.2E+6	5.2E+6	NA	ID		<50	<50	<50	<50	
Isopropyl benzene	98-82-8	800	28	56,000 (S)	56,000 (S)	29,000	ID		<5.0	<5.0	<5.0	<5.0	
Methyl iodide	74-88-4	NA	NA	NA	NA	NA	NA		<5.0	<5.0	<5.0	<5.0	
4-Methyl-2-pentanone (MIBK) (l)	108-10-1	1,800	ID	2.0E+7 (S)	1.3E+7	ID	2.0E+7 (S)		<50	<50	<50	<50	
Methylene chloride	75-09-2	5.0 (A)	1,500 (X)	2.2E+5	2.2E+5	ID	ID		<5.0	<5.0	<5.0	<5.0	
2-Methylnaphthalene	91-57-6	260	19	25,000 (S)	25,000 (S)	ID	ID		<5.0	<5.0	<5.0	<5.0	
Methyl-tert-butyl ether (MTBE)	1634-04-4	40 (E)	7,100 (X)	4.7E+7 (S)	6.1E+5	ID	ID		<5.0	<5.0	<5.0	<5.0	
Naphthalene	91-20-3	520	11	31,000 (S)	31,000 (S)	NA	31,000 (S)		<5.0	<5.0	<5.0	<5.0	
n-Propylbenzene (l)	103-65-1	80	ID	ID	15,000	ID	ID		<1.0	<1.0	<1.0	<1.0	
Styrene	100-42-5	100 (A)	80 (X)	1.7E+5	9,700	1.4E+5	3.1E+5 (S)		<1.0	<1.0	<1.0	<1.0	
1,1,1,2-Tetrachloroethane	630-20-6	77	ID	15,000	30,000	ID	ID		<1.0	<1.0	<1.0	<1.0	
1,1,2,2-Tetrachloroethane	79-34-5	8.5	78 (X)	12,000	4,700	ID	ID		<1.0	<1.0	<1.0	<1.0	
Tetrachloroethylene	127-18-4	5.0 (A)	60 (X)	25,000	12,000	ID	2.0E+5 (S)		<1.0	<1.0	<1.0	<1.0	

Units in µg/L = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

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Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

**TABLE 2: GROUNDWATER**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	TW-1	TW-2	TW-3	TRIP BLANK TB001102	
		Drinking Water Criteria & RBSLs	Groundwater Surface Water Interface Criteria & RBSLs	Groundwater Volatilization to Indoor Air Inhalation Criteria & RBSLs	Groundwater Contact Criteria and RBSLs	Flammability and Explosivity Screening Level	Acute Inhalation Screening Level						
							Sample Media		water	water	water	-	
							Collection Date		7/18/2013	7/18/2013	7/18/2013	7/18/2013	
								Depth (ft. bgs)	12'	9'	8'	NA	
<b>Volatiles</b>													
Toluene (l)	108-88-3	790 (E)	270	5.3E+5 (S)	5.3E+5 (S)	61,000	ID		<b>4.7</b>	<1.0	<1.0	<1.0	<1.0
1,2,4-Trichlorobenzene	120-82-1	70 (A)	99 (X)	3.0E+5 (S)	19,000	NA	3.0E+5 (S)		<5.0	<5.0	<5.0	<5.0	<5.0
1,1,1-Trichloroethane	71-55-6	200 (A)	89	6.6E+5	1.3E+6 (S)	ID	1.3E+6 (S)		<1.0	<1.0	<1.0	<1.0	<1.0
1,1,2-Trichloroethane	79-00-5	5.0 (A)	330 (X)	17,000	21,000	NA	ID		<1.0	<1.0	<1.0	<1.0	<1.0
Trichloroethylene	79-01-6	5.0 (A)	200 (X)	2,200	22,000	ID	1.1E+6 (S)		<1.0	<1.0	<1.0	<1.0	<1.0
Trichlorofluoromethane	75-69-4	2,600	NA	1.1E+6 (S)	1.1E+6 (S)	ID	1.1E+6 (S)		<1.0	<1.0	<1.0	<1.0	<1.0
1,2,3-Trichloropropane	96-18-4	42	NA	8,300	84,000	NA	ID		<1.0	<1.0	<1.0	<1.0	<1.0
1,2,3-Trimethylbenzene	526-73-8	NA	NA	NA	NA	NA	NA		<1.0	<1.0	<1.0	<1.0	<1.0
1,2,4-Trimethylbenzene (l)	95-63-6	63 (E)	17	56,000 (S)	56,000 (S)	56,000 (S)	ID		<b>1.4</b>	<1.0	<1.0	<1.0	<1.0
1,3,5-Trimethylbenzene (l)	108-67-8	72 (E)	45	61,000 (S)	61,000 (S)	ID	ID		<1.0	<1.0	<1.0	<1.0	<1.0
Vinyl chloride	75-01-4	2.0 (A)	13 (X)	1,100	1,000	33,000	ID		<1.0	<1.0	<1.0	<1.0	<1.0
Xylenes (l)	1330-20-7	280 (E)	41	1.9E+5 (S)	1.9E+5 (S)	70,000	1.9E+5 (S)		<b>5.6</b>	<3.0	<3.0	<3.0	<3.0
<b>Semivolatiles</b>													
Acenaphthene	83-32-9	1,300	38	4,200 (S)	4,200 (S)	ID	ID		NS	<5.0	<5.0	NS	
Acenaphthylene	208-96-8	52	ID	3,900 (S)	3,900 (S)	ID	ID		NS	<5.0	<5.0	NS	
Anthracene	120-12-7	43 (S)	ID	43 (S)	43 (S)	ID	ID		NS	<5.0	<5.0	NS	
Benzo(a)anthracene (Q)	56-55-3	2.1	ID	NLV	9.4 (S,AA)	ID	ID		NS	<1.0	<1.0	NS	
Benzo(a)pyrene (Q)	50-32-8	5.0 (A)	ID	NLV	1.0 (M,AA); 0.64	ID	ID		NS	<1.0	<1.0	NS	
Benzo(b)fluoranthene (Q)	205-99-2	1.5 (S, AA)	ID	ID	1.5 (S,AA)	ID	ID		NS	<1.0	<1.0	NS	
Benzo(g,h,i)perylene	191-24-2	1.0 (M); 0.26 (S)	ID	NLV	1.0 (M,AA); 0.26 (S)	ID	ID		NS	<1.0	<1.0	NS	
Benzo(k)fluoranthene (Q)	207-08-9	1.0 (M); 0.8 (S)	NA	NLV	1.0 (M,AA); 0.8 (S)	ID	ID		NS	<1.0	<1.0	NS	
Chrysene (Q)	218-01-9	1.6 (S)	ID	ID	1.6 (S,AA)	ID	ID		NS	<1.0	<1.0	NS	
Dibenzo(a,h)anthracene (Q)	53-70-3	2.0 (M); 0.21	ID	NLV	2.0 (M,AA); 0.31	ID	ID		NS	<2.0	<2.0	NS	

Units in µg/L = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

&lt; = Analyte was not detected at or above the specified reporting limit

Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

NS = Not Sampled

**TABLE 2: GROUNDWATER**

Norstar Development

Phase II ESA

Project Number: 1127.001

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Constituent	Chemical Abstract Service Number	MDEQ GENERIC RESIDENTIAL CLEANUP CRITERIA						Sample Identification	TW-1	TW-2	TW-3	TRIP BLANK TB001102	
		Drinking Water Criteria & RBSLs	Groundwater Surface Water Interface Criteria & RBSLs	Groundwater Volatilization to Indoor Air Inhalation Criteria & RBSLs	Groundwater Contact Criteria and RBSLs	Flammability and Explosivity Screening Level	Acute Inhalation Screening Level						
							Sample Media		water	water	water	-	
							Collection Date		7/18/2013	7/18/2013	7/18/2013	7/18/2013	
							Depth (ft. bgs)		12'	9'	8'	NA	
<i>Semivolatiles</i>													
Fluoranthene	206-44-0	210 (S)	1.6	210 (S)	210 (S)	ID	ID		NS	<1.0	<1.0	<1.0	NS
Fluorene	86-73-7	880	12	2,000 (S)	2,000 (S)	ID	ID		NS	<5.0	<5.0	<5.0	NS
Indeno(1,2,3-cd)pyrene (Q)	193-39-5	2.0 (M); 0.022 (S)	ID	NLV	2.0 (M, AA); 0.022 (S)	ID	ID		NS	<2.0	<2.0	<2.0	NS
2-Methylnaphthalene	91-57-6	260	19	25,000 (S)	25,000 (S)	ID	ID		<5.0	<5.0	<5.0	<5.0	NS
Naphthalene	91-20-3	520	11	31,000 (S)	31,000 (S)	NA	31,000 (S)		<5.0	<5.0	<5.0	<5.0	NS
Phenanthrene	85-01-8	52	2.0 (M); 1.4	1,000 (S)	1,000 (S)	ID	ID		NS	<2.0	<2.0	<2.0	NS
Pyrene	129-00-0	140 (S)	ID	140 (S)	140 (S)	ID	ID		NS	<5.0	<5.0	<5.0	NS
<i>PCBs</i>													
<i>Polychlorinated biphenyls (PCBs) (J.T)</i>	1336363	0.5 (A)	0.2 (M); 2.6E-5	45 (S)	3.3 (AA)	ID	ID		NS	<0.20	<0.22	<0.22	NS

Units in µg/L = parts per billion (ppb)

See attached for MDEQ Part 201/213 footnotes

< = Analyte was not detected at or above the specified reporting limit

Bolded values = Analyte concentration above reporting limit.

Shaded values = Analyte concentration exceeds applicable Generic Residential Cleanup Criteria

NS = Not Sampled

**FOOTNOTES  
for****Part 201 Criteria and Part 213 Risk-Based Screening Levels****Document Release Date: September 28, 2012**

- (A) Criterion is the state of Michigan drinking water standard established pursuant to Section 5 of 1976 PA 399, MCL 325.1005.
- (B) Background, as defined in R 299.5701(b), may be substituted if higher than the calculated cleanup criterion. Background levels may be less than criteria for some inorganic compounds.
- (C) Value presented is a screening level based on the chemical-specific generic soil saturation concentration ( $C_{sat}$ ) since the calculated risk-based criterion is greater than  $C_{sat}$ . Concentrations greater than  $C_{sat}$  are acceptable cleanup criteria for this pathway where a site-specific demonstration indicates that free-phase material containing a hazardous substance is not present.
- (D) Calculated criterion exceeds 100 percent, hence it is reduced to 100 percent or 1.0E+9 parts per billion (ppb).
- (E) Criterion is the aesthetic drinking water value, as required by Section 20120a(5) of the Natural Resources and Environmental Protection Act, 1994 PA 451, as amended (NREPA). A notice of aesthetic impact may be employed as an institutional control mechanism if groundwater concentrations exceed the aesthetic drinking water criterion, but do not exceed the applicable health-based drinking water value provided in the following table:

Hazardous Substance	Chemical Abstract Service Number	Residential Health-Based Drinking Water Value	Non-Residential Health-Based Drinking Water Value
Aluminum	7429905	300	4,100
tertiary Amyl methyl ether	994058	910	2,600
Copper	7440508	1,400	4,000
Diethyl ether	60297	3,700	10,000
Ethylbenzene	100414	700	700
Iron	7439896	2,000	5,600
Manganese	7439965	860	2,500
Methyl-tert-butyl ether (MTBE)	1634044	240	690
Toluene	108883	1,000	1,000
1,2,4-Trimethylbenzene	95636	1,000	2,900
1,3,5-Trimethylbenzene	108678	1,000	2,900
Xylenes	1330207	10,000	10,000

- (F) Criterion is based on adverse impacts to plant life and phytotoxicity.
- (G) Groundwater surface water interface (GSI) criterion depends on the pH or water hardness, or both, of the receiving surface water. The final chronic value (FCV) for the protection of aquatic life shall be calculated based on the pH or hardness of the receiving surface water. Where water hardness exceeds 400 mg CaCO<sub>3</sub>/L, use 400 mg CaCO<sub>3</sub>/L for the FCV calculation. The FCV formula provides values in units of ug/L or ppb. The generic GSI criterion is the lesser of

the calculated FCV, the wildlife value (WV), and the surface water human non-drinking water value (HNDV). The soil GSI protection criteria for these hazardous substances are the greater of the 20 times the GSI criterion or the GSI soil-water partition values using the GSI criteria developed with the procedure described in this footnote.

Hazardous Substance	FCV Formula ug/L	FCV Conversion Factor (CF)	WV ug/L	HNDV ug/L
Acetate	EXP(0.2732*(pH) + 7.0362)	NA	NA	1.3E+6
Acetic Acid	EXP(0.2732*(pH) + 7.0362)	NA	NA	1.3E+6
Barium	EXP(1.0629*(LnH)+1.1869)	NA	NA	1.6E+5
Beryllium	EXP(2.5279*(LnH)-10.7689)	NA	NA	1,200
Cadmium <sup>®</sup>	(EXP(0.7852*(LnH)-2.715))*CF	1.101672-((LnH)*(0.041838))	NA	130
Chromium (III) <sup>®</sup>	(EXP(0.819*(LnH)+0.6848))*CF	0.86	NA	9,400
Copper	(EXP(0.8545*(LnH)-1.702)) *CF	0.96	NA	38,000
Lead <sup>®</sup>	(EXP(0.9859*(LnH)-1.270))*CF	1.46203-((LnH)*(0.14571))	NA	190
Manganese <sup>®</sup>	EXP(0.8784*(LnH)+3.5385)	NA	NA	59,000
Nickel	(EXP(0.846*(LnH)+0.0584))*CF	0.997	NA	2.1E+5
Pentachlorophenol <sup>®</sup>	EXP(1.005*(pH)-5.134)	NA	NA	2.8
Zinc	(EXP(0.8473*(LnH)+0.884))*CF	0.986	NA	16,000

where,

- EXP(x) = The base of the natural logarithm raised to power x ( $e^x$ ).
- LnH = The natural logarithm of water hardness in mg CaCO<sub>3</sub>/L.
- \* = The multiplication symbol.
- <sup>®</sup> = The GSI criterion developed here may not be protective for surface water that is used as a drinking water source. Refer to footnote (X) for further guidance.

A spreadsheet that may be used to calculate GSI and GSI protection criteria for (G)-footnoted hazardous substances is available on the Department of Environmental Quality (DEQ) internet web site.

- (H) Valence-specific chromium data (Cr III and Cr VI) shall be compared to the corresponding valence-specific cleanup criteria. If both Cr III and Cr VI are present in groundwater, the total concentration of both cannot exceed the drinking water criterion of 100 ug/L. If analytical data are provided for total chromium only, they shall be compared to the cleanup criteria for Cr VI. Cr III soil cleanup criterion for protection of drinking water can only be used at sites where groundwater is prevented from being used as a public water supply, currently and in the future, through an approved land or resource use restriction.
- (I) Hazardous substance may exhibit the characteristic of ignitability as defined in 40 C.F.R. §261.21 (revised as of July 1, 2001), which is adopted by reference in these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, Remediation and Redevelopment Division (RRD), 525 West Allegan Street, Lansing, Michigan 48933, at cost.

- (J) Hazardous substance may be present in several isomer forms. Isomer-specific concentrations shall be added together for comparison to criteria.
- (K) Hazardous substance may be flammable or explosive, or both.
- (L) Criteria for lead are derived using a biologically based model, as allowed for under Section 20120a(10) of the NREPA, and are not calculated using the algorithms and assumptions specified in pathway-specific rules. The generic residential drinking water criterion of 4 ug/L is linked to the generic residential soil direct contact criterion of 400 mg/kg. A higher concentration in the drinking water, up to the state action level of 15 ug/L, may be allowed as a site-specific remedy and still allow for drinking water use, under Section 20120a(2) of the NREPA if soil concentrations are appropriately lower than 400 mg/kg. If a site-specific criterion is approved based on this subdivision, a notice shall be filed on the deed for all property where the groundwater concentrations will exceed 4 ug/L to provide notice of the potential for unacceptable risk if soil or groundwater concentrations increase. Acceptable combinations of site-specific soil and drinking water concentrations are presented in the following table:

Acceptable Combinations of Lead in Drinking Water and Soil

Drinking Water Concentration (ug/L)	Soil Concentration (mg/kg)
5	386-395
6	376-385
7	376-385
8	366-375
9	356-365
10	346-355
11	336-345
12	336-345
13	326-335
14	316-325
15	306-315

- (M) Calculated criterion is below the analytical target detection limit, therefore, the criterion defaults to the target detection limit.
- (N) The concentrations of all potential sources of nitrate-nitrogen (e.g., ammonia-N, nitrite-N, nitrate-N) in groundwater that is used as a source of drinking water shall not, when added together, exceed the nitrate drinking water criterion of 10,000 ug/L. Where leaching to groundwater is a relevant pathway, soil concentrations of all potential sources of nitrate-nitrogen shall not, when added together, exceed the nitrate drinking water protection criterion of 2.0E+5 ug/kg.
- (O) The concentration of all polychlorinated and polybrominated dibenzodioxin and dibenzofuran isomers present at a facility, expressed as an equivalent concentration of 2,3,7,8-tetrachlorodibenzo-p-dioxin based upon their relative potency, shall be added together and compared to the criteria for 2,3,7,8-tetrachlorodibenzo-p-dioxin. The generic cleanup criteria for 2,3,7,8-tetrachlorodibenzo-p-dioxin are not calculated according to the algorithms presented in R 299.5714 to R 299.5726. The generic cleanup criteria are being held at the values that the DEQ has used since August 1998, in recognition of

- the fact that national efforts to reassess risks posed by dioxin are not yet complete. Until these studies are complete, it is premature to select a revised slope factor and/or reference dose for calculation of generic cleanup criteria.
- (P) Amenable cyanide methods or method OIA-1677 shall be used to quantify cyanide concentrations for compliance with all groundwater criteria. Total cyanide methods or method OIA-1677 shall be used to quantify cyanide concentrations for compliance with soil criteria. Nonresidential direct contact criteria may not be protective of the potential for release of hydrogen cyanide gas. Additional land or resource use restrictions may be necessary to protect for the acute inhalation concerns associated with hydrogen cyanide gas.
- (Q) Criteria for carcinogenic polycyclic aromatic hydrocarbons were developed using relative potential potencies to benzo(a)pyrene.
- (R) Hazardous substance may exhibit the characteristic of reactivity as defined in 40 C.F.R. §261.23 (revised as of July 1, 2001), which is adopted by reference in these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost.
- (S) Criterion defaults to the hazardous substance-specific water solubility limit.
- (T) Refer to the federal Toxic Substances Control Act (TSCA), 40 C.F.R. §761, Subpart D and 40 C.F.R. §761, Subpart G, to determine the applicability of TSCA cleanup standards. Subpart D and Subpart G of 40 C.F.R. §761 (July 1, 2001) are adopted by reference in these rules and are available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulations may be purchased, at a cost as of the time of adoption of these rules of \$55, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401, or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost. Alternatives to compliance with the TSCA standards listed below are possible under 40 C.F.R. §761 Subpart D. New releases may be subject to the standards identified in 40 C.F.R. §761, Subpart G. Use Part 201 soil direct contact cleanup criteria in the following table if TSCA standards are not applicable.

Land Use Category	TSCA, Subpart D Cleanup Standards	Part 201 Soil Direct Contact Cleanup Criteria
Residential	1,000 ppb, or 10,000 ppb if capped	4,000 ppb
Nonresidential	1,000 ppb, or 10,000 ppb if capped	16,000 ppb

- (U) Hazardous substance may exhibit the characteristic of corrosivity as defined in 40 C.F.R. §261.22 (revised as of July 1, 2001), which is adopted by reference in

these rules and is available for inspection at the DEQ, 525 West Allegan Street, Lansing, Michigan. Copies of the regulation may be purchased, at a cost as of the time of adoption of these rules of \$45, from the Superintendent of Documents, Government Printing Office, Washington, DC 20401 (stock number 869-044-00155-1), or from the DEQ, RRD, 525 West Allegan Street, Lansing, Michigan 48933, at cost.

- (V) Criterion is the aesthetic drinking water value as required by Section 20120(a)(5) of the NREPA. Concentrations up to 200 ug/L may be acceptable, and still allow for drinking water use, as part of a site-specific cleanup under Section 20120a(2) of the NREPA.
- (W) Concentrations of trihalomethanes in groundwater shall be added together to determine compliance with the Michigan drinking water standard of 80 ug/L. Concentrations of trihalomethanes in soil shall be added together to determine compliance with the drinking water protection criterion of 1,600 ug/kg.
- (X) The GSI criterion shown in the generic cleanup criteria tables is not protective for surface water that is used as a drinking water source. For a groundwater discharge to the Great Lakes and their connecting waters or discharge in close proximity to a water supply intake in inland surface waters, the generic GSI criterion shall be the surface water human drinking water value (HDV) listed in the table in this footnote, except for those HDV indicated with an asterisk. For HDV with an asterisk, the generic GSI criterion shall be the lowest of the HDV, the WV, and the calculated FCV. See formulas in footnote (G). Soil protection criteria based on the HDV shall be as listed in the table in this footnote, except for those values with an asterisk. Soil GSI protection criteria based on the HDV shall be as listed in the table in this footnote, except for those values with an asterisk. Soil GSI protection criteria for compounds with an asterisk shall be the greater of 20 times the GSI criterion or the GSI soil-water partition values using the GSI criteria developed with the procedure described in this footnote.

Hazardous Substance	Chemical Abstract Service Number	Surface Water Human Drinking Water Values (HDV) (ug/L)	Soil GSI Protection Criteria for HDV (ug/kg)
Acrylamide	79061	0.5 (M); 0.12	10
Alachlor	15972608	3.5	88
Antimony	7440360	2.0 (M); 1.7	1,200
Benzene	71432	12	240
Boron	7440428	4,000	80,000
Bromate	15541454	10 (M); 0.5	200
n-Butanol	71363	3,500	70,000
Butyl benzyl phthalate	85687	6.9	13,000
Cadmium	7440439	2.5*	*
Carbon tetrachloride	56235	5.6	110
Chloride	16887006	50,000	1.0E+6
Chloroethane	75003	170	3,400
Chromium (III)	16065831	120*	*
Cyanazine	21725462	2.0 (M); 0.93	200 (M); 40
1,2-Dichloroethane	107062	6.0	120
trans-1,2-Dichloroethylene	156605	470	9,400
1,2-Dichloropropane	78875	9.1	180
1,3-Dichloropropene	542756	3.3	100 (M); 66
N,N-Dimethylacetamide	127195	700	14,000
1,4-Dioxane	123911	34	680
Ethylene dibromide	106934	0.17	20 (M); 3.4
Ethylene glycol	107211	56,000	1.1E+6
Hexachloroethane	67721	5.3	310
Isophorone	78591	310	6,200
Isopropyl alcohol	67630	28,000	5.6E+5
Lead	7439921	14*	*
Manganese	7439965	1,300*	*
Methanol	67561	14,000	2.8E+5
Methyl-tert-butyl ether (MTBE)	1634044	100	2,000
Methylene chloride	75092	47	940
Molybdenum	7439987	120	2,400
Nitrobenzene	98953	4.7	330 (M); 94
Pentachlorophenol	87865	1.8*	*
Styrene	100425	20	530
1,2,4,5-Tetrachlorobenzene	95943	2.8	3,300
1,1,2,2-Tetrachloroethane	79345	3.2	64
Tetrachloroethylene	127184	11	220
Tetrahydrofuran	109999	350	7,000
Thallium	7440280	2.0 (M); 1.2	1,400
1,2,4-Trichlorobenzene	120821	80	4,700
1,1,2-Trichloroethane	79005	12	240
Trichloroethylene	79016	29	580
Vinyl chloride	75014	1.0 (M); 0.25	40 (M); 20

- (Y) Source size modifiers shown in the following table shall be used to determine soil inhalation criteria for ambient air when the source size is not one-half acre. The modifier shall be multiplied by the generic soil inhalation criteria shown in the

table of generic cleanup criteria to determine the applicable criterion.

Source Size sq. feet or acres	Modifier
400 sq feet	3.17
1000 sq feet	2.2
2000 sq feet	1.76
1/4 acre	1.15
1/2 acre	1
1 acre	0.87
2 acre	0.77
5 acre	0.66
10 acre	0.6
32 acre	0.5
100 acre	0.43

- (Z) Mercury is typically measured as total mercury. The generic cleanup criteria, however, are based on data for different species of mercury. Specifically, data for elemental mercury, chemical abstract service (CAS) number 7439976, serve as the basis for the soil volatilization to indoor air criteria, groundwater volatilization to indoor air, and soil inhalation criteria. Data for methyl mercury, CAS number 22967926, serve as the basis for the GSI criterion; and data for mercuric chloride, CAS number 7487947, serve as the basis for the drinking water, groundwater contact, soil direct contact, and the groundwater protection criteria. Comparison to criteria shall be based on species-specific analytical data only if sufficient facility characterization has been conducted to rule out the presence of other species of mercury.
- (AA) Comparison to these criteria may take into account an evaluation of whether the hazardous substances are adsorbed to particulates rather than dissolved in water and whether filtered groundwater samples were used to evaluate groundwater.
- (BB) The state drinking water standard for asbestos is in units of fibers per milliliter of water (f/mL) longer than 10 millimicrons. Soil concentrations of asbestos are determined by polarized light microscopy.
- (CC) Groundwater: The generic GSI criteria are based on the toxicity of unionized ammonia ( $\text{NH}_3$ ); the criteria are 29 ug/L and 53 ug/L for cold water and warm water surface water, respectively. As a result, the GSI criterion shall be compared to the percent of the total ammonia concentration in the groundwater that will become  $\text{NH}_3$  in the surface water. This percent  $\text{NH}_3$  is a function of the pH and temperature of the receiving surface water and can be estimated using the following table, taken from Emerson, et al., (Journal of the Fisheries Research Board of Canada, Volume 32(12):2382, 1975).

Percent NH<sub>3</sub> in Aqueous Ammonia Solutions for 0-30 °C and pH 6-10

Temp (°F)	Temp (°C)	pH									
		6.0	6.5	7.0	7.5	8.0	8.5	9.0	9.5	10.0	
32.0	0	0.00827	0.0261	0.0826	0.261	0.820	2.55	7.64	20.7	45.3	
33.8	1	0.00899	0.0284	0.0898	0.284	0.891	2.77	8.25	22.1	47.3	
35.6	2	0.00977	0.0309	0.0977	0.308	0.968	3.00	8.90	23.6	49.4	
37.4	3	0.0106	0.0336	0.106	0.335	1.05	3.25	9.60	25.1	51.5	
39.2	4	0.0115	0.0364	0.115	0.363	1.14	3.52	10.3	26.7	53.5	
41.0	5	0.0125	0.0395	0.125	0.394	1.23	3.80	11.1	28.3	55.6	
42.8	6	0.0136	0.0429	0.135	0.427	1.34	4.11	11.9	30.0	57.6	
44.6	7	0.0147	0.0464	0.147	0.462	1.45	4.44	12.8	31.7	59.5	
46.4	8	0.0159	0.0503	0.159	0.501	1.57	4.79	13.7	33.5	61.4	
48.2	9	0.0172	0.0544	0.172	0.542	1.69	5.16	14.7	35.3	63.3	
50.0	10	0.0186	0.0589	0.186	0.586	1.83	5.56	15.7	37.1	65.1	
51.8	11	0.0201	0.0637	0.201	0.633	1.97	5.99	16.8	38.9	66.8	
53.6	12	0.0218	0.0688	0.217	0.684	2.13	6.44	17.9	40.8	68.5	
55.4	13	0.0235	0.0743	0.235	0.738	2.30	6.92	19.0	42.6	70.2	
57.2	14	0.0254	0.0802	0.253	0.796	2.48	7.43	20.2	44.5	71.7	
59.0	15	0.0274	0.0865	0.273	0.859	2.67	7.97	21.5	46.4	73.3	
60.8	16	0.0295	0.0933	0.294	0.925	2.87	8.54	22.8	48.3	74.7	
62.6	17	0.0318	0.101	0.317	0.996	3.08	9.14	24.1	50.2	76.1	
64.4	18	0.0343	0.108	0.342	1.07	3.31	9.78	25.5	52.0	77.4	
66.2	19	0.0369	0.117	0.368	1.15	3.56	10.5	27.0	53.9	78.7	
68.0	20	0.0397	0.125	0.396	1.24	3.82	11.2	28.4	55.7	79.9	
69.8	21	0.0427	0.135	0.425	1.33	4.10	11.9	29.9	57.5	81.0	
71.6	22	0.0459	0.145	0.457	1.43	4.39	12.7	31.5	59.2	82.1	
73.4	23	0.0493	0.156	0.491	1.54	4.70	13.5	33.0	60.9	83.2	
75.2	24	0.0530	0.167	0.527	1.65	5.03	14.4	34.6	62.6	84.1	
77.0	25	0.0569	0.180	0.566	1.77	5.38	15.3	36.3	64.3	85.1	
78.8	26	0.0610	0.193	0.607	1.89	5.75	16.2	37.9	65.9	85.9	
80.6	27	0.0654	0.207	0.651	2.03	6.15	17.2	39.6	67.4	86.8	
82.4	28	0.0701	0.221	0.697	2.17	6.56	18.2	41.2	68.9	87.3	
84.2	29	0.0752	0.237	0.747	2.32	7.00	19.2	42.9	70.4	88.3	
86.0	30	0.0805	0.254	0.799	2.48	7.46	20.3	44.6	71.8	89.0	

The generic approach for estimating NH<sub>3</sub> assumes a default pH of 8 and default temperatures of 68°F and 85°F for cold water and warm water surface water, respectively. The resulting percent NH<sub>3</sub> is 3.8 percent and 7.2 percent for cold water and warm water, respectively. This default percentage shall be multiplied by the total ammonia-nitrogen (NH<sub>3</sub>-N) concentration in the groundwater and the resulting NH<sub>3</sub> concentration compared to the applicable GSI criterion. As an

alternative, the maximum pH and temperature data from the specific receiving surface water can be used to estimate, from the table in this footnote, a lower percent unionized ammonia concentration for comparison to the generic GSI.

- Soil: The generic soil GSI protection criteria for unionized ammonia are 580 ug/kg and 1,100 ug/kg for cold water and warm water surface water, respectively.
- (DD) Hazardous substance causes developmental effects. Residential direct contact criteria are protective of both prenatal and postnatal exposure. Nonresidential direct contact criteria are protective for a pregnant adult receptor.
- (EE) The following are applicable generic GSI criteria as required by Section 20120a(15) of the NREPA.

Hazardous Substance	GSI (ug/L)	Notes
Phosphorus	1,000	Criteria applicable unless receiving water is a surface water that has a phosphorus waste load allocation or is an inland lake. In those cases, contact the department for applicable values.
Total dissolved solids (TDS)	5.0E+5	If TDS data are not available, the TDS criterion may be used a screening level for the sum of the concentrations of the following substances: Calcium, Chlorides, Iron, Magnesium, Potassium, Sodium, Sulfate.
Dissolved Oxygen (DO): Cold receiving waters Warm receiving waters	≥ 7,000 ≥ 5,000	Since a low level of DO can be harmful to aquatic life, the criterion represents a minimum level that on-site samples must exceed. This is in contrast to other criteria which represent "not to exceed" concentrations. DO criteria are not applicable if groundwater Carbonaceous Biochemical Oxygen Demand (CBOD) is less than 10,000 ug/L and groundwater ammonia concentration is less than 2,000 ug/L.

- (FF) The chloride GSI criterion shall be 125 mg/l when the discharge is to surface waters of the state designated as public water supply sources or 50 mg/l when the discharge is to the Great Lakes or connecting waters. Chloride GSI criteria shall not apply for surface waters of the state that are not designated as a public water supply source, however, the total dissolved solids criterion is applicable.
- (GG) Risk-based criteria are not available for methane due to insufficient toxicity data. An acceptable soil gas concentration (presented for both residential and nonresidential land uses) was derived utilizing 25 percent of the lower explosive level for methane. This equates to 1.25 percent or 8.4E+6 ug/m<sup>3</sup>.

"ID" means insufficient data to develop criterion.

"NA" means a criterion or value is not available or, in the case of background and CAS numbers, not applicable.

"NLL" means hazardous substance is not likely to leach under most soil conditions.

"NLV" means hazardous substance is not likely to volatilize under most conditions.

**TABLE 3: SOIL GAS**

Norstar Development

Phase II ESA

Project Number: 1127.002

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Parameters	Chemical Abstract Service Number	Vapor Intrusion Deep Soil Gas Screening Levels (ppbv)	Sample Identification	SB-4	SB-5	SB-6	SB-7
			Depth	7'	7'	10'	10'
			Collection Date	7/18/2013	7/18/2013	7/18/2013	7/18/2013
<b>Volatile Organic Compounds (VOCs)</b>							
Acetone	67-64-1	8.20E+05		<b>130</b>	<84	<b>150</b>	<b>500</b>
Benzene	71-43-2	3.20E+02		<b>76</b>	<b>39</b>	<b>53</b>	<b>77</b>
Benzyl Chloride	100-44-7	3.40E+01		<12	<12	<12	<25
Bromodichloromethane	75-27-4	7.10E+01		<1.4	<1.4	<1.4	<2.8
Bromoform	75-25-2	7.60E+02		<1.4	<1.4	<1.4	<2.8
Bromomethane	74-83-9	4.30E+02		<2.6	<2.6	<2.6	<5.3
1,3-Butadiene	106-99-0	NA		<14	<14	<14	<28
2-Butanone	78-93-3	5.60E+05		<b>9.5</b>	<6.9	<b>9.6</b>	<b>20</b>
Carbon Disulfide	75-15-0	7.40E+04		<b>60</b>	<b>22</b>	<13	<26
Carbon Tetrachloride	56-23-5	2.30E+02		<1.3	<1.3	<1.3	<2.7
Chlorobenzene	108-90-7	5.00E+03		<1.4	<1.4	<1.4	<2.8
Chloroethane	75-00-3	1.30E+06		<13	<13	<13	<26
Chloroform	67-66-3	7.30E+02		<1.4	<1.4	<1.4	<2.8
Chloromethane	74-87-3	6.50E+03		<2.6	<2.6	<2.6	<5.1
Cyclohexane	110-82-7	5.80E+05		<b>4.1</b>	<b>2.7</b>	<b>5.9</b>	<5.2
Dibromochloromethane	124-48-1	4.10E+01		<1.4	<1.4	<1.4	<2.9
1,2-Dichlorobenzene	95-50-1	1.60E+04		<13	<13	<13	<26
1,3-Dichlorobenzene	541-73-1	1.60E+02		<1.4	<1.4	<1.4	<2.8
1,4-Dichlorobenzene	106-46-7	2.10E+02		<1.4	<1.4	<1.4	<2.8
Dichlorodifluoromethane	75-71-8	3.30E+06		<1.3	<1.3	<1.3	<3.0
1,1-Dichloroethane	75-34-3	4.10E+04		<1.4	<1.4	<1.4	<2.8
1,2-Dichloroethane	107-06-2	8.20E+01		<2.6	<2.6	<2.6	<5.2
1,1-Dichloroethene	75-35-4	1.70E+04		<1.4	<1.4	<1.4	<2.8
cis-1,2-Dichloroethene	156-59-2	5.80E+02		<1.4	<1.4	<1.4	<2.8

Units = ppbv = parts per billion by volume

< = Analyte not detected at or above the reporting limit

Bolded values = analyte concentrations above reporting limit

Shaded boxes = analyte exceeds applicable screening criteria

See Attached MDEQ Vapor Intrusion Screening Criteria footnotes

NS = Not Sampled

NA = Criteria Not Available

**TABLE 3: SOIL GAS**

Norstar Development

Phase II ESA

Project Number: 1127.002

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Parameters	Chemical Abstract Service Number	Vapor Intrusion Deep Soil Gas Screening Levels (ppbv)	Sample Identification	SB-4	SB-5	SB-6	SB-7
			Depth	7'	7'	10'	10'
			Collection Date	7/18/2013	7/18/2013	7/18/2013	7/18/2013
<b>Volatile Organic Compounds (VOCs)</b>							
trans-1,2-Dichloroethene	156-60-5	5.80E+03		<1.4	<1.4	<1.4	<2.8
1,2-Dichloropropane	78-87-5	2.90E+02		<1.4	<1.4	<1.4	<2.8
cis-1,3-Dichloropropene	10061-01-5	NA		<1.4	<1.4	<1.4	<2.8
trans-1,3-Dichloropropene	10061-02-6	NA		<2.4	<2.4	<2.4	<4.8
1,4-Dioxane	123-91-1	NA		<2.6	<2.6	<2.6	<5.1
Ethyl Acetate	141-78-6	2.90E+05		<13	<13	<13	<27
Ethylbenzene	100-41-4	6.40E+03		<b>48</b>	<b>20</b>	<b>57</b>	<b>53</b>
Ethylene Dibromide	106-93-4	1.90E+00		<1.4	<1.4	<1.4	<2.8
4-Ethyltoluene	622-96-8	NA		<b>18</b>	<13	<b>93</b>	<26
n-Heptane	142-82-5	2.80E+05		<b>18</b>	<b>12</b>	<b>10</b>	<b>11</b>
Hexachlorobutadiene	87-68-3	3.70E+01		<2.6	<2.6	<2.6	<5.1
n-Hexane	110-54-3	6.60E+04		<b>29</b>	<b>18</b>	<b>29</b>	<b>26</b>
2-Hexanone	591-78-6	2.50E+03		<14	<14	<14	<29
Isopropanol	67-63-0	NA		<14	<14	<14	<29
Methylene Chloride	75-09-2	8.80E+03		<13	<13	<13	<26
2-Methylnaphthalene	91-57-6	5.70E+02		<24	<24	<24	<49
4-Methyl-2-pentanone	108-10-1	2.40E+05		<b>7.2</b>	<b>4.0</b>	<b>8.5</b>	<b>17</b>
MTBE	1634-04-4	2.70E+05		<1.4	<1.4	<1.4	<2.8
Naphthalene	91-20-3	1.40E+02		<14	<14	<14	<28
Propylene	115-07-1	NA		<b>420</b>	<b>82</b>	<b>470</b>	<b>520</b>
Styrene	100-42-5	3.50E+03		<2.6	<2.6	<2.6	<5.2
1,1,2,2-Tetrachloroethane	79-34-5	2.20E+01		<2.7	<2.7	<2.7	<5.3
Tetrachloroethene	127-18-4	1.70E+03		<b>6.3</b>	<b>5.6</b>	<b>2.0</b>	<2.8
Tetrahydrofuran	109-99-9	2.00E+03		<2.9	<2.9	<2.9	<5.8
Toluene	108-88-3	4.40E+05		<b>250</b>	<b>140</b>	<b>150</b>	<b>210</b>
1,2,4-Trichlorobenzene	120-82-1	1.80E+02		<2.5	<2.5	<2.5	<5.0

Units = ppbv = parts per billion by volume

< = Analyte not detected at or above the reporting limit

Bolded values = analyte concentrations above reporting limit

Shaded boxes = analyte exceeds applicable screening criteria

See Attached MDEQ Vapor Intrusion Screening Criteria footnotes

NS = Not Sampled

NA = Criteria Not Available

**TABLE 3: SOIL GAS**

Norstar Development

Phase II ESA

Project Number: 1127.002

Green Baxter, 1701-1749 Green Road, Ann Arbor, Michigan

July 2013

Parameters	Chemical Abstract Service Number	Vapor Intrusion Deep Soil Gas Screening Levels (ppbv)	Sample Identification	SB-4	SB-5	SB-6	SB-7
			Depth	7'	7'	10'	10'
			Collection Date	7/18/2013	7/18/2013	7/18/2013	7/18/2013
<b>Volatile Organic Compounds (VOCs)</b>							
1,1,1-Trichloroethane	71-55-6	3.60E+05		<1.4	<1.4	<1.4	<2.8
1,1,2-Trichloroethane	79-00-5	9.80E+01		<1.4	<1.4	<1.4	<2.8
Trichloroethene	79-01-6	1.20E+02		<1.4	<1.4	<1.4	<2.8
Trichlorofluoromethane	75-69-4	3.30E+06		<1.4	<1.4	<1.4	<2.8
1,1,2-Trichlorotrifluoroethane	76-13-1	8.40E+05		<1.4	<1.4	<1.4	<2.8
1,2,4-Trimethylbenzene	95-63-6	1.50E+04		<b>52</b>	<b>32</b>	<b>130</b>	<b>88</b>
1,3,5-Trimethylbenzene	108-67-8	1.50E+04		<b>18</b>	<13	<b>100</b>	<26
Vinyl Acetate	108-05-4	1.90E+04		<13	<13	<13	<27
Vinyl Chloride	75-01-4	2.10E+02		<6.6	<6.6	<6.6	<13
Xylenes	1330207	NA		<b>259</b>	<b>109</b>	<b>324</b>	<b>309</b>
m&p-Xylene	136777-61-2	NA		<b>200</b>	<b>86</b>	<b>240</b>	<b>240</b>
o-Xylene	95-47-6	7.60E+03		<b>59</b>	<b>23</b>	<b>84</b>	<b>69</b>

Units = ppbv = parts per billion by volume

< = Analyte not detected at or above the reporting limit

Bolded values = analyte concentrations above reporting limit

Shaded boxes = analyte exceeds applicable screening criteria

See Attached MDEQ Vapor Intrusion Screening Criteria footnotes

NS = Not Sampled

NA = Criteria Not Available

**APPENDIX A**

**Geophysical Survey Report**

July 25, 2013

Andy Foerg  
Environmental Resources Group LLC  
28003 Center Oaks Suite 106  
Wixom, MI 48393

Tel: (248) 763-3639  
Fax: (248) 363-0461  
Email: andy.foerg@ergrp.net

Dear Andy:

On July 17, 2013 Jason Hurshman, representing Fibertec performed a Ground Penetrating Radar (GPR) scan at 1737 Greed Road in Ann Arbor, Michigan focusing on the identification of buried underground storage tanks (UST).

Using a GSSI Sir 3000 GPR unit with a 400MHz antenna an area approximately 195' x 235' was scanned and no anomalies appearing to be a UST were observed. A rough map of the area scanned is provided below. In addition the scan files from the GPR have also been provided. The file names correspond to the numbered grid on the map.

Fibertec appreciates your business and looks forward to future opportunities to partner with Environmental Resources Group LLC. Please do not hesitate to call with any questions.

Sincerely,



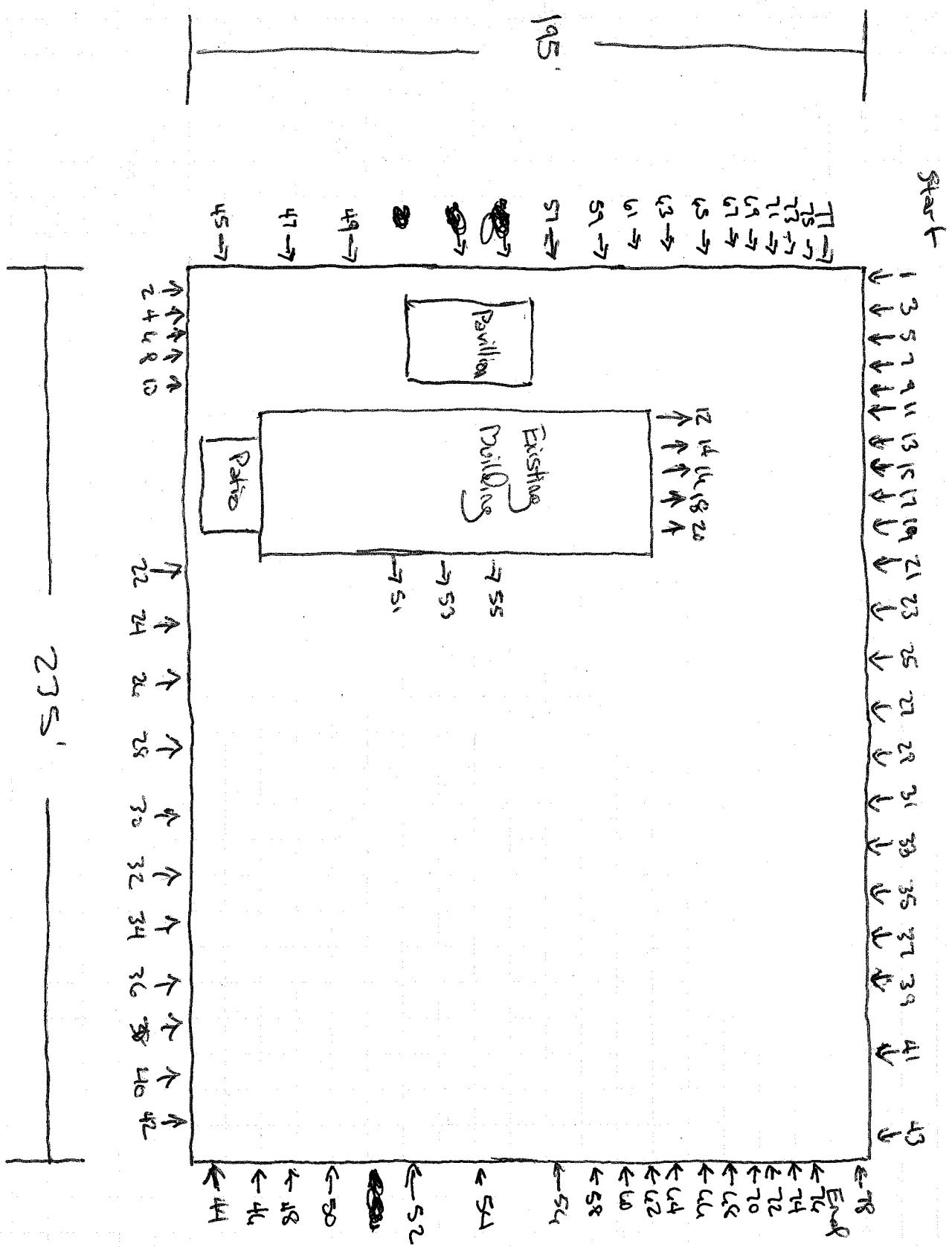
Michael C. McCourtie  
Client Services Coordinator  
Fibertec Environmental Service  
Subsurface Investigations

**Notes & Disclaimer:** Fibertec utilizes standards of care consistent with ordinary standards of professional care of others using similar tools. Client recognizes that Fibertec's investigation may not identify all underground structures. The information upon which Fibertec relies may contain errors or be incomplete, and the limitations of Fibertec's equipment may not identify all structures buried on the surveyed property. Therefore, the client agrees to the fullest extent permitted by law, to waive all claims and causes of action against the Fibertec and anyone to whom Fibertec may be liable, for damages to underground utilities and structures, resulting from subsurface penetrations in locations established by Fibertec as not having detectable underground structures or utilities.

Green R.R

← N

Start  
↓  
1 3 5 7 9 11 13 15 17 19 21 23 25 27 29 31 33 35 37 39 41 43  
↓  
End  
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Measurements  
and 5' grid  
are approx.

## **APPENDIX B**

### **Soil Boring Logs**

# SOIL BORING LOG

Page 1 of 1



Environmental Resources Group  
28003 Center Oaks Court • Suite 205 • Novi, MI • 48375  
Phone 248-773-7986 • Fax 248-924-3108

Project: Green Baxter Phase II		Project Number: 1127.001	Client: Norstar Development	Boring No.: SB-1/TW-1
Address, City, State: 1701-1749 Green Road, Ann Arbor, MI		Drilling Contractor: Fibertec	Drill Rig Type: 6620 DT geoprobe	
Logged By: L. Lambert		Started: 7/18/2013	Bit Type: N/A	Hole Diameter: N/A
Drill Crew: Ryan		Completed: 7/18/2013	Hammer Type: N/A	Well Diameter: N/A
Comments: sampled soil 11:15 / water @ 11:20		Backfilled: natural cuttings & bentonite granules	Hammer Weight: N/A	Hammer Drop: N/A
		Groundwater Depth: N/A	Elevation: N/A	Total Depth of Boring: 15 feet
Depth (feet)	Sample Type	Recovery (feet)	Graphic Log	Lithology
				<u>Soil Group Name:</u> modifier, color, moisture, density/consistency, grain size, other descriptors
				<u>Rock Description:</u> modifier, color, hardness/degree of concentration, bedding and joint characteristics, solutions, void conditions.
2	☒	2.5'		4" Asphalt FILL, including Sand, Silt, Clay and Gravel, moist.
4				Brown, Clay, soft, some Gravel, moist. Brown, Clay, medium stiff, few Gravel, damp.
6				S. A., damp.
8				Brown, Silt, some Clay, some fine Sand, few Gravel, moist.
10				Brown Silt and Clay, few Gravel, damp.
12	☒	4'		S. A., interbedded fine Sand, damp.
14	☒			Brown, fine Sand and Clay, soft, moist. Brown, fine to medium Sand, few Gravel, moist.
				S. A., wet.
				Gray, Clay, some Silt, soft, moist.
				EOB
Standard Penetration Split Spoon Sampler (SPT) Direct Push Sampler Water Sample Hand Auger		Soil gas sample Stabilized Ground water Groundwater At time of Drilling Screen Interval		

# SOIL BORING LOG

Page 1 of 1



Environmental Resources Group  
28003 Center Oaks Court • Suite 106 • Wixom, MI • 48393  
Phone 248-773-7986 • Fax 248-924-3108

Project: Green Baxter Phase II		Project Number: 1127.001		Client: Norstar Development	Boring No.: SB-2		
Address, City, State: 1701-1749 Green Road, Ann Arbor, MI				Drilling Contractor: Fibertec	Drill Rig Type: 6620 DT geoprobe		
Logged By: L. Lambert		Date	Started: 7/18/2013	Bit Type: N/A	Hole Diameter: N/A		
Drill Crew: Ryan			Completed: 7/18/2013	Hammer Type: N/A	Well Diameter: N/A		
Comments: sampled soil @ 12:15			Backfilled: natural cuttings & bentonite granules	Hammer Weight: N/A	Hammer Drop: N/A		
		Groundwater Depth: N/A	Elevation: N/A	Total Depth of Boring: 10 feet			
Depth (feet)	Sample Type	Recovery (feet)	Graphic Log	Lithology	PID	Well Construction	Soil Gas Sample Screen Interval
				<b>Soil Group Name:</b> modifier, color, moisture, density/consistency, grain size, other descriptors			
				<b>Rock Description:</b> modifier, color, hardness/degree of concentration, bedding and joint characteristics, solutions, void conditions.			
				Brown, fine Sand and Clay and Silt, some Organics, moist.	0.0		
				Brown, Clay, some fine Sand and Silt, some Organics, moist.	0.0		
2	HA			S. A. A., few Gravel, few Organics, moist.	0.0		
				Brown, Clay, moist.	0.0		
				S. A. A., mottled white, trace Gravel, damp.	0.0		
4				Brown, fine to medium Sand and Clay, wet.	0.0		
				Brown, Silt and Clay, damp.	0.0		
6				S. A. A., sand lens @ 8.75'-9'	0.0		
				Gray, Clay, damp.	0.0		
8	X			EOB			
10							
Standard Penetration Split Spoon Sampler (SPT) Direct Push Sampler Water Sample HA Hand Auger				Soil gas sample Stabilized Ground water Groundwater At time of Drilling Screen Interval			

# SOIL BORING LOG

Page 1 of 1



Environmental Resources Group  
28003 Center Oaks Court • Suite 106 • Wixom, MI • 48393  
Phone 248-773-7986 • Fax 248-924-3108

Project: Green Baxter Phase II		Project Number: 1127.001		Client: Norstar Development	Boring No.: SB-3		
Address, City, State: 1701-1749 Green Road, Ann Arbor, MI		Drilling Contractor: Fibertec		Drill Rig Type: 6620 DT geoprobe			
Logged By: L. Lambert		Date	Started: 7/18/2013	Bit Type: N/A	Hole Diameter: N/A		
Drill Crew: Ryan			Completed: 7/18/2013	Hammer Type: N/A	Well Diameter: N/A		
Comments: sampled soil @ 12:50			Backfilled: natural cuttings & bentonite granules	Hammer Weight: N/A	Hammer Drop: N/A		
		Groundwater Depth: N/A		Elevation: N/A	Total Depth of Boring: 10 feet		
Depth (feet)	Sample Type	Recovery (feet)	Graphic Log	Lithology	PID	Well Construction	Soil Gas Sample Screen Interval
				<b>Soil Group Name:</b> modifier, color, moisture, density/consistency, grain size, other descriptors			
				<b>Rock Description:</b> modifier, color, hardness/degree of concentration, bedding and joint characteristics, solutions, void conditions.			
				Brown, fine Sand and Silt, some Clay, some Organics, trace Gravel, damp.	0.0		
2				Brown, fine Sand and Silt, few Organics, trace Clay, damp.	0.0		
	HA			Light brown, Clay and Silt, trace fine Sand, damp.	0.0		
4				Brown, Silt, trace coarse Sand, trace Gravel, damp.	0.0		
6					0.0		
8					0.0		
10				EOB	0.0		
Standard Penetration Split Spoon Sampler (SPT) Direct Push Sampler Water Sample Hand Auger		Soil gas sample Stabilized Ground water Groundwater At time of Drilling Screen Interval					

# SOIL BORING LOG

Page 1 of 1



Environmental Resources Group  
28023 Center Oaks Court • Suite 106 • Wixom, MI • 48393  
Phone 248-773-7986 • Fax 248-724-3208

Project: Green Baxter Phase II		Project Number: 1127.001		Client: Norstar Development	Boring No.: SB-4/TW-2
Address, City, State: 1701-1749 Green Road, Ann Arbor, MI		Drilling Contractor: Fibertec		Drill Rig Type: 6620 DT geoprobe	
Logged By: L. Lambert		Started: 7/18/2013		Bit Type: N/A	Hole Diameter: N/A
Drill Crew: Ryan		Completed: 7/18/2013		Hammer Type: N/A	Well Diameter: N/A
Comments: sampled water @ 14:00		Backfilled: natural cuttings & bentonite granules		Hammer Weight: N/A	Hammer Drop: N/A
		Groundwater Depth: N/A		Elevation: N/A	Total Depth of Boring: 20 feet
Depth (feet)	Sample Type	Recovery (feet)	Graphic Log	Lithology	
				Soil Group Name: modifier, color, moisture, density/consistency, grain size, other descriptors	PID
				Rock Description: modifier, color, hardness/degree of concentration, bedding and joint characteristics, solutions, void conditions.	Well Construction
2				Brown, Clay and Silt, trace Organics, moist.	
4				Brown, Clay, some fine Sand, trace Organics, moist.	
6				Brown, Clay, some Silt, moist.	
8				Brown, Clay and Silt, very moist.	
10				Brown, fine Sand, Clay and Silt, trace Gravel, very moist.	
12				S. A. A., wet.	
14				Gray, Clay, trace Gravel, moist.	
16				Brown, fine to medium Sand with Gray Clay, wet.	
18				Brown, fine to medium Sand, saturated.	
20				Gray, Clay, soft, some fine to medium Sand, few Gravel, very moist.	
				S. A. A., trace Gravel, very moist.	
				Brown, fine to coarse Sand, few Gravel, interbedded Clay lens, moist.	
				EOB	
Standard Penetration Split Spoon Sampler (SPT) Direct Push Sampler Water Sample HA Hand Auger					
Soil gas sample Stabilized Ground water Groundwater At time of Drilling Screen Interval					

# SOIL BORING LOG



**Environmental Resources Group**  
2809 Center Oaks Court • Suite 125 • Winona, MN 58383  
Phone: 218-722-7285 • Fax: 218-722-7208

# **SOIL BORING LOG**



**Environmental Resources Group**  
26003 Centre Oaks Court • Suite 205 • Austin, TX • 78733  
Phone 210.775.7955 • Fax 210.722.7259

Project: Green Baxter Phase II		Project Number: 1127.001	Client: Norstar Development	Boring No.: SB-6
Address, City, State: 1701-1749 Green Road, Ann Arbor, MI		Drilling Contractor: Fibertec	Drill Rig Type: 6620 DT geoprobe	Hole Diameter: N/A
Logged By: L. Lambert		Started: 7/18/2013	Bit Type: N/A	Hole Diameter: N/A
Drill Crew: Ryan		Completed: 7/18/2013	Hammer Type: N/A	Well Diameter: N/A
Comments: sampled soil @ 16:40		Backfilled: natural cuttings & bentonite granules	Hammer Weight: N/A	Hammer Drop: N/A
		Groundwater Depth: N/A	Elevation: N/A	Total Depth of Boring: 20 feet
Depth (feet)	Sample Type	Recovery (feet)	Graphic Log	Lithology
				<b>Soil Group Name:</b> modifier, color, moisture, density/consistency, grain size, other descriptors
				<b>Rock Description:</b> modifier, color, hardness/degree of concentration, bedding and joint characteristics, solutions, void conditions.
2				Brown, fine to medium Sand, trace Organics, moist. S. A. A., moist.
4				S. A. A., trace Gravel, moist.
6				Brown, fine to medium Sand and Silt, trace Gravel, damp.
8				S. A. A., damp.
10				Brown, Silt, few Gravel, damp.
12				Brown, Silt and Clay, mottled white, trace Gravel, damp.
14				Brown, fine to medium Sand and Clay, interbedded fine to coarse Sand lenses to 15', wet.
16				Gray, Clay, soft, few Gravel, moist.
18				
20				EOB
				PID
				Well Construction
				Soil Gas Sample Screen Interval

# SOIL BORING LOG



**Environmental Resources Group**  
2803 Center Oaks Court • Suite 206 • Wixom, MI • 48393  
Phone 248-773-7966 • Fax 248-924-3108

**APPENDIX C**

**Analytical Data Reports**

Thursday, July 25, 2013

Fibertec Project Number: 57032  
Project Identification: Green Baxter /1127.001  
Submittal Date: 07/18/2013

Ms. Laura Lambert  
Environmental Resources Group  
28003 Center Oaks Court Suite 106  
Wixom, MI 48393

Dear Ms. Lambert,

Thank you for selecting Fibertec Environmental Services as your analytical laboratory. The samples you submitted have been analyzed in accordance with NELAC standards and the results compiled in the attached report. Any exceptions to NELAC compliance are noted in the report. These results apply only to those samples submitted. Please note samples will be disposed of 30 days after reporting date.

If you have any questions regarding these results or if we may be of further assistance to you, please contact me at (517) 699-0345.

Sincerely,



Daryl P. Strandbergh  
Laboratory Director

DPS/kc

Enclosures

---

1914 Holloway Drive  
11766 E. Grand River  
8660 S. Mackinaw Trail

Holt, MI 48842  
Brighton, MI 48116  
Cadillac, MI 49601

T: (517) 699-0345  
T: (810) 220-3300  
T: (231) 775-8368

F: (517) 699-0388  
F: (810) 220-3311  
F: (231) 775-8584

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-1 (1.5'-2')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>1</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>11:15</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Dry Weight Determination (ASTM D 2974-87)</b>						<b>Aliquot ID: 57032-001A</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: BMG</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Percent Moisture (Water Content) (NN)	<b>8.9</b>		%	0.1	1.0	07/22/13	MC130722	07/23/13	MC130722
<b>Trace Elements by ICP/MS (EPA 0200.2-M/EPA 6020A)</b>									
Parameter(s) Result Q Units Reporting Limit Dilution Prep Date Prep Batch Analysis Date Analysis Batch									
1. Arsenic	<b>7300</b>		µg/kg	100	20	07/24/13	PT13G24D	07/24/13	T213G24A
2. Lead	<b>21000</b>		µg/kg	1000	20	07/24/13	PT13G24D	07/24/13	T213G24A
<b>Organochlorine Pesticides (EPA 3546/EPA 8081B)</b>									
Parameter(s) Result Q Units Reporting Limit Dilution Prep Date Prep Batch Analysis Date Analysis Batch									
1. Aldrin	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
2. alpha-BHC (NN)	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
3. beta-BHC (NN)	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
4. delta-BHC	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
5. gamma-BHC (NN)	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
6. Chlordane (NN)	<b>U</b>		µg/kg	91	25	07/23/13	PS13G23A	07/24/13	SC13G23B
7. 4,4'-DDD	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
8. 4,4'-DDE	<b>U</b>	J,L-	µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
9. 4,4'-DDT	<b>860</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
10. Dieldrin	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
11. Endosulfan I	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
12. Endosulfan II	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
13. Endosulfan Sulfate	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
14. Endrin	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
15. Endrin Aldehyde	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
16. Heptachlor	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
17. Heptachlor Epoxide	<b>U</b>		µg/kg	37	25	07/23/13	PS13G23A	07/24/13	SC13G23B
18. Methoxychlor	<b>U</b>		µg/kg	150	25	07/23/13	PS13G23A	07/24/13	SC13G23B
19. Toxaphene (NN)	<b>U</b>		µg/kg	1800	25	07/23/13	PS13G23A	07/24/13	SC13G24A

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>						<b>Aliquot ID: 57032-001</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	<b>U</b>		µg/kg	1000	1.0	07/19/13	V913G19A	07/19/13	V913G19A
2. Acrylonitrile	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
3. Benzene	<b>U</b>		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
4. Bromobenzene	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
5. Bromochloromethane	<b>U</b>		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
6. Bromodichloromethane	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A

1914 Holloway Drive  
11766 E. Grand River  
8660 S. Mackinaw Trail

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T: (231) 775-8368

F: (517) 699-0388  
F: (810) 220-3311  
F: (231) 775-8584

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-1 (1.5'-2')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>1</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>11:15</b>

Sample Comments: **Soil results have been calculated and reported on a dry weight basis unless otherwise noted.**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-001</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
7. Bromoform	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
8. Bromomethane	U		µg/kg	200	1.0	07/19/13	V913G19A	07/19/13	V913G19A
9. 2-Butanone	U		µg/kg	750	1.0	07/19/13	V913G19A	07/19/13	V913G19A
10. n-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
11. sec-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
12. tert-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
13. Carbon Disulfide	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
14. Carbon Tetrachloride	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
15. Chlorobenzene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
16. Chloroethane	U		µg/kg	270	1.0	07/19/13	V913G19A	07/19/13	V913G19A
17. Chloroform	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
18. Chloromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
19. 2-Chlorotoluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
20. Dibromochloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/kg	27	1.0	07/19/13	V913G19A	07/19/13	V913G19A
22. Dibromomethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
23. 1,2-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
24. 1,3-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
25. 1,4-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
26. Dichlorodifluoromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
27. 1,1-Dichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
28. 1,2-Dichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
29. 1,1-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
30. cis-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
31. trans-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
32. 1,2-Dichloropropane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
33. cis-1,3-Dichloropropene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
34. trans-1,3-Dichloropropene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
35. Ethylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
36. Ethylene Dibromide	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
37. 2-Hexanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
38. Isopropylbenzene	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
39. Methyl Iodide	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
40. Methylene Chloride	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
41. 4-Methyl-2-pentanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
42. MTBE	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
43. Naphthalene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
44. n-Propylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
45. Styrene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-1 (1.5'-2')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>1</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>11:15</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-001</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
46. 1,1,2-Tetrachloroethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
47. 1,1,2,2-Tetrachloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
48. Tetrachloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
49. Toluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
50. 1,2,4-Trichlorobenzene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
51. 1,1,1-Trichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
52. 1,1,2-Trichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
53. Trichloroethene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
54. Trichlorofluoromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
55. 1,2,3-Trichloropropane	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
57. 1,2,4-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
58. 1,3,5-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
59. Vinyl Chloride	U		µg/kg	40	1.0	07/19/13	V913G19A	07/19/13	V913G19A
60. Xylenes	U		µg/kg	150	1.0	07/19/13	V913G19A	07/19/13	V913G19A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3546/EPA 8270C)</b>				<b>Aliquot ID: 57032-001A</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: TMC</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acenaphthene (SIM)	U		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
2. Acenaphthylene (SIM)	2000		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
3. Anthracene (SIM)	1700		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
4. Benzo(a)anthracene (SIM)	5800		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
5. Benzo(a)pyrene (SIM)	7800		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
6. Benzo(b)fluoranthene (SIM)	9000		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
7. Benzo(ghi)perylene (SIM)	6500		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
8. Benzo(k)fluoranthene (SIM)	3000		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
9. Chrysene (SIM)	4600		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
10. Dibenzo(a,h)anthracene (SIM)	1900		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
11. Fluoranthene (SIM)	9400		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
12. Fluorene (SIM)	610		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
13. Indeno(1,2,3-cd)pyrene (SIM)	6800		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
14. 2-Methylnaphthalene (SIM)	U		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
15. Phenanthrene (SIM)	4100		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A
16. Pyrene (SIM)	8200		µg/kg	330	20	07/23/13	PS13G23A	07/24/13	S513G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-1</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>2</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>11:20</b>

**Sample Comments:**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>						<b>Aliquot ID: 57032-002</b>	<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
1. Acetone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
2. Acrylonitrile	U		µg/L	2.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
3. Benzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
4. Bromobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
5. Bromochloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
6. Bromodichloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
7. Bromoform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
8. Bromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
9. 2-Butanone	U		µg/L	25	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
10. n-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
11. sec-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
12. tert-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
13. Carbon Disulfide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
14. Carbon Tetrachloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
15. Chlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
16. Chloroethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
17. Chloroform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
18. Chloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
19. 2-Chlorotoluene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
20. Dibromochloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
22. Dibromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
23. 1,2-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
24. 1,3-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
25. 1,4-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
26. Dichlorodifluoromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
27. 1,1-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
28. 1,2-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
29. 1,1-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
30. cis-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
31. trans-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
32. 1,2-Dichloropropane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
33. cis-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
34. trans-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
35. Ethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
36. Ethylene Dibromide	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
37. 2-Hexanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
38. Isopropylbenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
39. Methyl Iodide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-1</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>2</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>11:20</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>				<b>Aliquot ID: 57032-002</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>		
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
40. Methylene Chloride	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
41. 2-Methylnaphthalene (NN)	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
42. 4-Methyl-2-pentanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
43. MTBE	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
44. Naphthalene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
45. n-Propylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
46. Styrene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
47. 1,1,1,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
48. 1,1,2,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
49. Tetrachloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
50. Toluene	4.7		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
51. 1,2,4-Trichlorobenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
52. 1,1,1-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
53. 1,1,2-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
54. Trichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
55. Trichlorofluoromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
56. 1,2,3-Trichloropropane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
57. 1,2,3-Trimethylbenzene (NN)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
58. 1,2,4-Trimethylbenzene	1.4		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
59. 1,3,5-Trimethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
60. Vinyl Chloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
61. Xylenes	5.6		µg/L	3.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-2 (8-9')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>3</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>12:15</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Dry Weight Determination (ASTM D 2974-87)</b>						<b>Aliquot ID: 57032-003A</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: BMG</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Percent Moisture (Water Content) (NN)	<b>14</b>		%	0.1	1.0	07/22/13	MC130722	07/23/13	MC130722
<b>Trace Elements by ICP/MS (EPA 0200.2-M/EPA 6020A)</b>									
<b>Aliquot ID: 57032-003A</b>									
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Arsenic	<b>5600</b>		µg/kg	100	20	07/24/13	PT13G24D	07/24/13	T213G24A
2 Lead	<b>6500</b>		µg/kg	1000	20	07/24/13	PT13G24D	07/24/13	T213G24A
<b>Organochlorine Pesticides (EPA 3546/EPA 8081B)</b>									
<b>Aliquot ID: 57032-003A</b>									
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Aldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
2 alpha-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
3. beta-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
4. delta-BHC	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
5. gamma-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
6. Chlordane (NN)	<b>U</b>		µg/kg	25	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
7. 4,4'-DDD	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
8. 4,4'-DDE	<b>U</b>	J,L-	µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
9. 4,4'-DDT	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
10. Dieldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
11. Endosulfan I	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
12 Endosulfan II	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
13. Endosulfan Sulfate	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
14. Endrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
15. Endrin Aldehyde	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
16. Heptachlor	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
17. Heptachlor Epoxide	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
18. Methoxychlor	<b>U</b>		µg/kg	50	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
19. Toxaphene (NN)	<b>U</b>		µg/kg	390	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>						<b>Aliquot ID: 57032-003</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	<b>U</b>		µg/kg	1000	1.0	07/19/13	V913G19A	07/19/13	V913G19A
2. Acrylonitrile	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
3. Benzene	<b>U</b>		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
4. Bromobenzene	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
5. Bromochloromethane	<b>U</b>		µg/kg	120	1.0	07/19/13	V913G19A	07/19/13	V913G19A
6. Bromodichloromethane	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-2 (8-9')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>3</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>12:15</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
7. Bromoform	U		µg/kg	120	1.0	07/19/13	V913G19A	07/19/13	V913G19A
8. Bromomethane	U		µg/kg	200	1.0	07/19/13	V913G19A	07/19/13	V913G19A
9. 2-Butanone	U		µg/kg	750	1.0	07/19/13	V913G19A	07/19/13	V913G19A
10. n-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
11. sec-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
12. tert-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
13. Carbon Disulfide	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
14. Carbon Tetrachloride	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
15. Chlorobenzene	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
16. Chloroethane	U		µg/kg	290	1.0	07/19/13	V913G19A	07/19/13	V913G19A
17. Chloroform	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
18. Chloromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
19. 2-Chlorotoluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
20. Dibromochloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/kg	29	1.0	07/19/13	V913G19A	07/19/13	V913G19A
22. Dibromomethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
23. 1,2-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
24. 1,3-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
25. 1,4-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
26. Dichlorodifluoromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
27. 1,1-Dichloroethane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
28. 1,2-Dichloroethane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
29. 1,1-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
30. cis-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
31. trans-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
32. 1,2-Dichloropropane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
33. cis-1,3-Dichloropropene	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
34. trans-1,3-Dichloropropene	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
35. Ethylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
36. Ethylene Dibromide	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
37. 2-Hexanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
38. Isopropylbenzene	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
39. Methyl Iodide	U		µg/kg	120	1.0	07/19/13	V913G19A	07/19/13	V913G19A
40. Methylene Chloride	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
41. 4-Methyl-2-pentanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
42. MTBE	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
43. Naphthalene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
44. n-Propylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
45. Styrene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-2 (8-9')</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>3</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>12:15</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-003</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
46. 1,1,2-Tetrachloroethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
47. 1,1,2,2-Tetrachloroethane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
48. Tetrachloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
49. Toluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
50. 1,2,4-Trichlorobenzene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
51. 1,1,1-Trichloroethane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
52. 1,1,2-Trichloroethane	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
53. Trichloroethene	U		µg/kg	58	1.0	07/19/13	V913G19A	07/19/13	V913G19A
54. Trichlorofluoromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
55. 1,2,3-Trichloropropane	U		µg/kg	120	1.0	07/19/13	V913G19A	07/19/13	V913G19A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
57. 1,2,4-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
58. 1,3,5-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
59. Vinyl Chloride	U		µg/kg	40	1.0	07/19/13	V913G19A	07/19/13	V913G19A
60. Xylenes	U		µg/kg	150	1.0	07/19/13	V913G19A	07/19/13	V913G19A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3546/EPA 8270C)</b>				<b>Aliquot ID: 57032-003A</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: TMC</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acenaphthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
2. Acenaphthylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
3. Anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
4. Benzo(a)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
5. Benzo(a)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
6. Benzo(b)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
7. Benzo(ghi)perylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
8. Benzo(k)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
9. Chrysene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
10. Dibenzo(a,h)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
11. Fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
12. Fluorene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
14. 2-Methylnaphthalene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
15. Phenanthrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
16. Pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-3 (9-10)</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>4</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>12:50</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

**Dry Weight Determination (ASTM D 2974-87)**      **Aliquot ID: 57032-004A**      **Matrix: Soil/Solid**      **Analyst: BMG**

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Percent Moisture (Water Content) (NN)	<b>12</b>		%	0.1	1.0	07/22/13	MC130722	07/23/13	MC130722

**Trace Elements by ICP/MS (EPA 0200.2-M/EPA 6020A)**      **Aliquot ID: 57032-004A**      **Matrix: Soil/Solid**      **Analyst: JLH**

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Arsenic	<b>6800</b>		µg/kg	100	20	07/24/13	PT13G24D	07/24/13	T213G24A
2 Lead	<b>7200</b>		µg/kg	1000	20	07/24/13	PT13G24D	07/24/13	T213G24A

**Organochlorine Pesticides (EPA 3546/EPA 8081B)**      **Aliquot ID: 57032-004A**      **Matrix: Soil/Solid**      **Analyst: GAN**

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Aldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
2 alpha-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
3. beta-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
4. delta-BHC	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
5. gamma-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
6. Chlordane (NN)	<b>U</b>		µg/kg	25	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
7. 4,4'-DDD	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
8. 4,4'-DDE	<b>U</b>	J,L-	µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
9. 4,4'-DDT	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
10. Dieldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
11. Endosulfan I	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
12. Endosulfan II	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
13. Endosulfan Sulfate	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
14. Endrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
15. Endrin Aldehyde	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
16. Heptachlor	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
17. Heptachlor Epoxide	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
18. Methoxychlor	<b>U</b>		µg/kg	50	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
19. Toxaphene (NN)	<b>U</b>		µg/kg	380	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B

**Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)**      **Aliquot ID: 57032-004**      **Matrix: Soil/Solid**      **Analyst: CCD**

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	<b>U</b>		µg/kg	1000	1.0	07/19/13	V913G19A	07/19/13	V913G19A
2 Acrylonitrile	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
3. Benzene	<b>U</b>		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
4. Bromobenzene	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
5. Bromochloromethane	<b>U</b>		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
6. Bromodichloromethane	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-3 (9-10)</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>4</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>12:50</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions:	Q: Qualifier (see definitions at end of report)	NA: Not Applicable	NN: Parameter not included in NELAC Scope of Analysis.		

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-004</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
7. Bromoform	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
8. Bromomethane	U		µg/kg	200	1.0	07/19/13	V913G19A	07/19/13	V913G19A
9. 2-Butanone	U		µg/kg	750	1.0	07/19/13	V913G19A	07/19/13	V913G19A
10. n-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
11. sec-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
12. tert-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
13. Carbon Disulfide	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
14. Carbon Tetrachloride	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
15. Chlorobenzene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
16. Chloroethane	U		µg/kg	280	1.0	07/19/13	V913G19A	07/19/13	V913G19A
17. Chloroform	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
18. Chloromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
19. 2-Chlorotoluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
20. Dibromochloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/kg	28	1.0	07/19/13	V913G19A	07/19/13	V913G19A
22. Dibromomethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
23. 1,2-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
24. 1,3-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
25. 1,4-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
26. Dichlorodifluoromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
27. 1,1-Dichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
28. 1,2-Dichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
29. 1,1-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
30. cis-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
31. trans-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
32. 1,2-Dichloropropane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
33. cis-1,3-Dichloropropene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
34. trans-1,3-Dichloropropene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
35. Ethylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
36. Ethylene Dibromide	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
37. 2-Hexanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
38. Isopropylbenzene	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
39. Methyl Iodide	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
40. Methylene Chloride	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
41. 4-Methyl-2-pentanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
42. MTBE	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
43. Naphthalene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
44. n-Propylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
45. Styrene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	Environmental Resources Group	Sample Description:	SB-3 (9-10)	Chain of Custody:	124238
Client Project Name:	Green Baxter	Sample No:	4	Collect Date:	07/18/13
Client Project No:	1127.001	Sample Matrix:	Soil/Solid	Collect Time:	12:50
Sample Comments: Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)					Aliquot ID: 57032-004		Matrix: Soil/Solid		Analyst: CCD
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
46.1,1,1,2-Tetrachloroethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
47.1,1,2,2-Tetrachloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
48. Tetrachloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
49. Toluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
50. 1,2,4-Trichlorobenzene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
51. 1,1,1-Trichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
52. 1,1,2-Trichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
53. Trichloroethene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
54. Trichlorofluoromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
55. 1,2,3-Trichloroproppane	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
57. 1,2,4-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
58. 1,3,5-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
59. Vinyl Chloride	U		µg/kg	40	1.0	07/19/13	V913G19A	07/19/13	V913G19A
60. Xylenes	U		µg/kg	150	1.0	07/19/13	V913G19A	07/19/13	V913G19A

Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3546/EPA 8270C)					Aliquot ID: 57032-004A		Matrix: Soil/Solid		Analyst: TMC
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acenaphthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
2. Acenaphthylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
3. Anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
4. Benzo(a)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
5. Benzo(a)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
6. Benzo(b)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
7. Benzo(ghi)perylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
8. Benzo(k)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
9. Chrysene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
10. Dibenzo(a,h)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
11. Fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
12. Fluorene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
14. 2-Methylnaphthalene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
15. Phenanthrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
16. Pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-2</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>5</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:00</b>

**Sample Comments:**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Polychlorinated Biphenyls (PCBs) (EPA 3535A/EPA 8082A)</b>				<b>Aliquot ID: 57032-005A</b>		<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>		
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Aroclor-1016	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
2. Aroclor-1221	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
3. Aroclor-1232	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
4. Aroclor-1242	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
5. Aroclor-1248	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
6. Aroclor-1254	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
7. Aroclor-1260	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
8. Aroclor-1262 (NN)	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A
9. Aroclor-1268 (NN)	U		µg/L	0.20	1.0	07/23/13	PS13G23C	07/23/13	SB13G23A

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>				<b>Aliquot ID: 57032-005</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>		
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
2. Acrylonitrile	U		µg/L	2.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
3. Benzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
4. Bromobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
5. Bromochloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
6. Bromodichloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
7. Bromoform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
8. Bromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
9. 2-Butanone	U		µg/L	25	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
10. n-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
11. sec-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
12. tert-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
13. Carbon Disulfide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
14. Carbon Tetrachloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
15. Chlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
16. Chloroethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
17. Chloroform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
18. Chloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
19. 2-Chlorotoluene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
20. Dibromochloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
22. Dibromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
23. 1,2-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
24. 1,3-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
25. 1,4-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
26. Dichlorodifluoromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-2</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>5</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:00</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>				<b>Aliquot ID: 57032-005</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>		
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
27. 1,1-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
28. 1,2-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
29. 1,1-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
30. cis-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
31. trans-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
32. 1,2-Dichloropropane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
33. cis-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
34. trans-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
35. Ethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
36. Ethylene Dibromide	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
37. 2-Hexanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
38. Isopropylbenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
39. Methyl Iodide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
40. Methylene Chloride	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
41. 4-Methyl-2-pentanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
42. MTBE	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
43. Naphthalene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
44. n-Propylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
45. Styrene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
46. 1,1,1,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
47. 1,1,2,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
48. Tetrachloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
49. Toluene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
50. 1,2,4-Trichlorobenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
51. 1,1,1-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
52. 1,1,2-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
53. Trichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
54. Trichlorofluoromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
55. 1,2,3-Trichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
57. 1,2,4-Trimethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
58. 1,3,5-Trimethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
59. Vinyl Chloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
60. Xylenes	U		µg/L	3.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3535A/EPA 8270C)</b>				<b>Aliquot ID: 57032-005A</b>		<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>		
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
1. Acenaphthene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-2</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>5</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:00</b>

**Sample Comments:**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3535A/EPA 8270C)</b>					<b>Aliquot ID: 57032-005A</b>		<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
2. Acenaphthylene (SIM)	U		µg/L	5.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
3. Anthracene (SIM)	U		µg/L	5.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
4. Benzo(a)anthracene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
5. Benzo(a)pyrene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
6. Benzo(b)fluoranthene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
7. Benzo(ghi)perylene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
8. Benzo(k)fluoranthene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
9. Chrysene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
10. Dibenzo(a,h)anthracene (SIM)	U		µg/L	2.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
11. Fluoranthene (SIM)	U		µg/L	1.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
12. Fluorene (SIM)	U		µg/L	5.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/L	2.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
14. 2-Methylnaphthalene (SIM)	U		µg/L	5.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
15. Phenanthrene (SIM)	U		µg/L	2.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B
16. Pyrene (SIM)	U		µg/L	5.0	1:1	07/23/13	PS13G23B	07/23/13	S513G23B

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-3</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>6</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:55</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Polychlorinated Biphenyls (PCBs) (EPA 3535A/EPA 8082A)</b>				<b>Aliquot ID: 57032-006A</b>		<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>		
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Aroclor-1016	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
2 Aroclor-1221	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
3. Aroclor-1232	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
4. Aroclor-1242	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
5. Aroclor-1248	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
6. Aroclor-1254	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
7. Aroclor-1260	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
8. Aroclor-1262 (NN)	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A
9. Aroclor-1268 (NN)	U		µg/L	0.22	1.1	07/23/13	PS13G23C	07/23/13	SB13G23A

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>				<b>Aliquot ID: 57032-006</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>		
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
2 Acrylonitrile	U		µg/L	2.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
3. Benzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
4. Bromobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
5. Bromochloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
6. Bromodichloromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
7. Bromoform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
8. Bromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
9. 2-Butanone	U		µg/L	25	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
10. n-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
11. sec-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
12. tert-Butylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
13. Carbon Disulfide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
14. Carbon Tetrachloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
15. Chlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
16. Chloroethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
17. Chloroform	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
18. Chloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
19. 2-Chlorotoluene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
20. Dibromochloromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
22. Dibromomethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
23. 1,2-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
24. 1,3-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
25. 1,4-Dichlorobenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
26. Dichlorodifluoromethane	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-3</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>6</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:55</b>

**Sample Comments:**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>				<b>Aliquot ID: 57032-006</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>		
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
27. 1,1-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
28. 1,2-Dichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
29. 1,1-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
30. cis-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
31. trans-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
32. 1,2-Dichloropropane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
33. cis-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
34. trans-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
35. Ethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
36. Ethylene Dibromide	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
37. 2-Hexanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
38. Isopropylbenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
39. Methyl Iodide	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
40. Methylene Chloride	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
41. 4-Methyl-2-pentanone	U		µg/L	50	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
42. MTBE	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
43. Naphthalene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
44. n-Propylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
45. Styrene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
46. 1,1,1,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
47. 1,1,2,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
48. Tetrachloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
49. Toluene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
50. 1,2,4-Trichlorobenzene	U		µg/L	5.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
51. 1,1,1-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
52. 1,1,2-Trichloroethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
53. Trichloroethene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
54. Trichlorofluoromethane	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
55. 1,2,3-Trichloropropene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
57. 1,2,4-Trimethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
58. 1,3,5-Trimethylbenzene	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
59. Vinyl Chloride	U		µg/L	1.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A
60. Xylenes	U		µg/L	3.0	1.0	07/24/13	VB13G24A	07/24/13	VB13G24A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3535A/EPA 8270C)</b>				<b>Aliquot ID: 57032-006A</b>		<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>		
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
1. Acenaphthene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TW-3</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>6</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>14:55</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3535A/EPA 8270C)</b>						<b>Aliquot ID: 57032-006A</b>	<b>Matrix: Ground Water</b>	<b>Analyst: TMC</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
2. Acenaphthylene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
3. Anthracene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
4. Benzo(a)anthracene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
5. Benzo(a)pyrene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
6. Benzo(b)fluoranthene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
7. Benzo(ghi)perylene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
8. Benzo(k)fluoranthene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
9. Chrysene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
10. Dibenzo(a,h)anthracene (SIM)	U		µg/L	2.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
11. Fluoranthene (SIM)	U		µg/L	1.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
12. Fluorene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/L	2.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
14. 2-Methylnaphthalene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
15. Phenanthrene (SIM)	U		µg/L	2.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B
16. Pyrene (SIM)	U		µg/L	5.0	1.1	07/23/13	PS13G23B	07/23/13	S513G23B

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Client Identification:	Environmental Resources Group	Sample Description:	SB-6 (1-2'	Chain of Custody:	124238
Client Project Name:	Green Baxter	Sample No:	7	Collect Date:	07/18/13
Client Project No:	1127.001	Sample Matrix:	Soil/Solid	Collect Time:	16:40
Sample Comments: Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

Dry Weight Determination (ASTM D 2974-87)						Aliquot ID: 57032-007A	Matrix: Soil/Solid	Analyst: BMG	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Percent Moisture (Water Content) (NN)	8.3		%	0.1	1.0	07/22/13	MC130722	07/23/13	MC130722
<b>Trace Elements by ICP/MS (EPA 0200.2-M/EPA 6020A)</b>									
Parameter(s) Result Q Units Reporting Limit Dilution Prep Date Prep Batch Analysis Date Analysis Batch									
1. Arsenic	5500		µg/kg	100	20	07/24/13	PT13G24D	07/24/13	T213G24A
2. Lead	5400		µg/kg	1000	20	07/24/13	PT13G24D	07/24/13	T213G24A
<b>Organochlorine Pesticides (EPA 3546/EPA 8081B)</b>									
Parameter(s) Result Q Units Reporting Limit Dilution Prep Date Prep Batch Analysis Date Analysis Batch									
1. Aldrin	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
2. alpha-BHC (NN)	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
3. beta-BHC (NN)	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
4. delta-BHC	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
5. gamma-BHC (NN)	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
6. Chlordane (NN)	U		µg/kg	25	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
7. 4,4'-DDD	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
8. 4,4'-DDE	U	J,L-	µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
9. 4,4'-DDT	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
10. Dieldrin	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
11. Endosulfan I	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
12. Endosulfan II	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
13. Endosulfan Sulfate	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
14. Endrin	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
15. Endrin Aldehyde	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
16. Heptachlor	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
17. Heptachlor Epoxide	U		µg/kg	20	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
18. Methoxychlor	U		µg/kg	50	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B
19. Toxaphene (NN)	U		µg/kg	360	5.0	07/23/13	PS13G23A	07/23/13	SC13G23B

Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)						Aliquot ID: 57032-007	Matrix: Soil/Solid	Analyst: CCD	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	U		µg/kg	1000	1.0	07/19/13	V913G19A	07/19/13	V913G19A
2. Acrylonitrile	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
3. Benzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
4. Bromobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
5. Bromochloromethane	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
6. Bromodichloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-6 (1-2'</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>7</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>16:40</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Matrix: Soil/Solid	Analyst: CCD
7. Bromoform	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
8. Bromomethane	U		µg/kg	200	1.0	07/19/13	V913G19A	07/19/13	V913G19A
9. 2-Butanone	U		µg/kg	750	1.0	07/19/13	V913G19A	07/19/13	V913G19A
10. n-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
11. sec-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
12. tert-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
13. Carbon Disulfide	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
14. Carbon Tetrachloride	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
15. Chlorobenzene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
16. Chloroethane	U		µg/kg	270	1.0	07/19/13	V913G19A	07/19/13	V913G19A
17. Chloroform	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
18. Chloromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
19. 2-Chlorotoluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
20. Dibromochloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/kg	27	1.0	07/19/13	V913G19A	07/19/13	V913G19A
22. Dibromomethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
23. 1,2-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
24. 1,3-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
25. 1,4-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
26. Dichlorodifluoromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
27. 1,1-Dichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
28. 1,2-Dichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
29. 1,1-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
30. cis-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
31. trans-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
32. 1,2-Dichloropropane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
33. cis-1,3-Dichloropropene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
34. trans-1,3-Dichloropropene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
35. Ethylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
36. Ethylene Dibromide	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
37. 2-Hexanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
38. Isopropylbenzene	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
39. Methyl Iodide	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
40. Methylene Chloride	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
41. 4-Methyl-2-pentanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
42. MTBE	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
43. Naphthalene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
44. n-Propylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
45. Styrene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-6 (1-2'</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>7</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>16:40</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>			<b>Aliquot ID: 57032-007</b>			<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
46. 1,1,2-Tetrachloroethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
47. 1,1,2,2-Tetrachloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
48. Tetrachloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
49. Toluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
50. 1,2,4-Trichlorobenzene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
51. 1,1,1-Trichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
52. 1,1,2-Trichloroethane	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
53. Trichloroethene	U		µg/kg	55	1.0	07/19/13	V913G19A	07/19/13	V913G19A
54. Trichlorofluoromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
55. 1,2,3-Trichloropropane	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
57. 1,2,4-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
58. 1,3,5-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
59. Vinyl Chloride	U		µg/kg	40	1.0	07/19/13	V913G19A	07/19/13	V913G19A
60. Xylenes	U		µg/kg	150	1.0	07/19/13	V913G19A	07/19/13	V913G19A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3546/EPA 8270C)</b>			<b>Aliquot ID: 57032-007A</b>			<b>Matrix: Soil/Solid</b>		<b>Analyst: TMC</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acenaphthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
2. Acenaphthylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
3. Anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
4. Benzo(a)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
5. Benzo(a)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
6. Benzo(b)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
7. Benzo(ghi)perylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
8. Benzo(k)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
9. Chrysene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
10. Dibenzo(a,h)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
11. Fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
12. Fluorene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
14. 2-Methylnaphthalene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
15. Phenanthrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
16. Pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-7 (12-13)'</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>8</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>17:25</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Dry Weight Determination (ASTM D 2974-87)</b>						<b>Aliquot ID: 57032-008A</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: BMG</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Percent Moisture (Water Content) (NN)	<b>12</b>		%	0.1	1.0	07/22/13	MC130722	07/23/13	MC130722
<b>Trace Elements by ICP/MS (EPA 0200.2-M/EPA 6020A)</b>									
<b>Parameter(s)</b> <b>Result</b> <b>Q</b> <b>Units</b> <b>Reporting Limit</b> <b>Dilution</b> <b>Prep Date</b> <b>Prep Batch</b> <b>Analysis Date</b> <b>Analysis Batch</b>									
1. Arsenic	<b>5100</b>		µg/kg	100	20	07/24/13	PT13G24D	07/24/13	T213G24A
2. Lead	<b>7900</b>		µg/kg	1000	20	07/24/13	PT13G24D	07/24/13	T213G24A
<b>Organochlorine Pesticides (EPA 3546/EPA 8081B)</b>									
<b>Parameter(s)</b> <b>Result</b> <b>Q</b> <b>Units</b> <b>Reporting Limit</b> <b>Dilution</b> <b>Prep Date</b> <b>Prep Batch</b> <b>Analysis Date</b> <b>Analysis Batch</b>									
1. Aldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
2. alpha-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
3. beta-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
4. delta-BHC	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
5. gamma-BHC (NN)	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
6. Chlordane (NN)	<b>U</b>		µg/kg	25	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
7. 4,4'-DDD	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
8. 4,4'-DDE	<b>U</b>	J,L-	µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
9. 4,4'-DDT	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
10. Dieldrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
11. Endosulfan I	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
12. Endosulfan II	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
13. Endosulfan Sulfate	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
14. Endrin	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
15. Endrin Aldehyde	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
16. Heptachlor	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
17. Heptachlor Epoxide	<b>U</b>		µg/kg	20	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
18. Methoxychlor	<b>U</b>		µg/kg	50	5.0	07/23/13	PS13G23A	07/24/13	SC13G23B
19. Toxaphene (NN)	<b>U</b>		µg/kg	380	5.0	07/23/13	PS13G23A	07/24/13	SC13G24A

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>						<b>Aliquot ID: 57032-008</b>	<b>Matrix: Soil/Solid</b>	<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone	<b>U</b>		µg/kg	1000	1.0	07/19/13	V913G19A	07/19/13	V913G19A
2. Acrylonitrile	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
3. Benzene	<b>U</b>		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
4. Bromobenzene	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
5. Bromochloromethane	<b>U</b>		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
6. Bromodichloromethane	<b>U</b>		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-7 (12-13)'</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>8</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>17:25</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
<b>Definitions:</b> Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-008</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
7. Bromoform	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
8. Bromomethane	U		µg/kg	200	1.0	07/19/13	V913G19A	07/19/13	V913G19A
9. 2-Butanone	U		µg/kg	750	1.0	07/19/13	V913G19A	07/19/13	V913G19A
10. n-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
11. sec-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
12. tert-Butylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
13. Carbon Disulfide	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
14. Carbon Tetrachloride	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
15. Chlorobenzene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
16. Chloroethane	U		µg/kg	280	1.0	07/19/13	V913G19A	07/19/13	V913G19A
17. Chloroform	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
18. Chloromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
19. 2-Chlorotoluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
20. Dibromochloromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/kg	28	1.0	07/19/13	V913G19A	07/19/13	V913G19A
22. Dibromomethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
23. 1,2-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
24. 1,3-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
25. 1,4-Dichlorobenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
26. Dichlorodifluoromethane	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
27. 1,1-Dichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
28. 1,2-Dichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
29. 1,1-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
30. cis-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
31. trans-1,2-Dichloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
32. 1,2-Dichloropropane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
33. cis-1,3-Dichloropropene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
34. trans-1,3-Dichloropropene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
35. Ethylbenzene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
36. Ethylene Dibromide	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
37. 2-Hexanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
38. Isopropylbenzene	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
39. Methyl Iodide	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
40. Methylene Chloride	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
41. 4-Methyl-2-pentanone	U		µg/kg	2500	1.0	07/19/13	V913G19A	07/19/13	V913G19A
42. MTBE	U		µg/kg	250	1.0	07/19/13	V913G19A	07/19/13	V913G19A
43. Naphthalene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
44. n-Propylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
45. Styrene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-7 (12-13)'</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>8</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Soil/Solid</b>	Collect Time:	<b>17:25</b>
<b>Sample Comments:</b> Soil results have been calculated and reported on a dry weight basis unless otherwise noted.					
Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.					

<b>Volatile Organic Compounds (VOCs) by GC/MS, 5035 (EPA 5035/EPA 8260B)</b>				<b>Aliquot ID: 57032-008</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: CCD</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
46.1,1,2-Tetrachloroethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
47.1,1,2,2-Tetrachloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
48. Tetrachloroethene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
49. Toluene	U		µg/kg	50	1.0	07/19/13	V913G19A	07/19/13	V913G19A
50. 1,2,4-Trichlorobenzene	U		µg/kg	330	1.0	07/19/13	V913G19A	07/19/13	V913G19A
51. 1,1,1-Trichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
52. 1,1,2-Trichloroethane	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
53. Trichloroethene	U		µg/kg	57	1.0	07/19/13	V913G19A	07/19/13	V913G19A
54. Trichlorofluoromethane	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
55. 1,2,3-Trichloropropane	U		µg/kg	110	1.0	07/19/13	V913G19A	07/19/13	V913G19A
56. 1,2,3-Trimethylbenzene (NN)	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
57. 1,2,4-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
58. 1,3,5-Trimethylbenzene	U		µg/kg	100	1.0	07/19/13	V913G19A	07/19/13	V913G19A
59. Vinyl Chloride	U		µg/kg	40	1.0	07/19/13	V913G19A	07/19/13	V913G19A
60. Xylenes	U		µg/kg	150	1.0	07/19/13	V913G19A	07/19/13	V913G19A

<b>Polynuclear Aromatic Hydrocarbons (PNAs) (EPA 3546/EPA 8270C)</b>				<b>Aliquot ID: 57032-008A</b>		<b>Matrix: Soil/Solid</b>		<b>Analyst: TMC</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acenaphthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
2. Acenaphthylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
3. Anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
4. Benzo(a)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
5. Benzo(a)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
6. Benzo(b)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
7. Benzo(ghi)perylene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
8. Benzo(k)fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
9. Chrysene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
10. Dibenzo(a,h)anthracene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
11. Fluoranthene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
12. Fluorene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
13. Indeno(1,2,3-cd)pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
14. 2-Methylnaphthalene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
15. Phenanthrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A
16. Pyrene (SIM)	U		µg/kg	330	1.0	07/23/13	PS13G23A	07/24/13	S513G24A

Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TRIP BLANK TB001102</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>9</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>NA</b>

**Sample Comments:**

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>Volatile Organic Compounds (VOCs) by GC/MS (EPA 5030B/EPA 8260B)</b>					<b>Aliquot ID: 57032-009</b>		<b>Matrix: Ground Water</b>	<b>Analyst: JPL</b>	
<b>Parameter(s)</b>	<b>Result</b>	<b>Q</b>	<b>Units</b>	<b>Reporting Limit</b>	<b>Dilution</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analysis Date</b>	<b>Analysis Batch</b>
1. Acetone	U		µg/L	50	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
2. Acrylonitrile	U		µg/L	2.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
3. Benzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
4. Bromobenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
5. Bromochloromethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
6. Bromodichloromethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
7. Bromoform	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
8. Bromomethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
9. 2-Butanone	U		µg/L	25	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
10. n-Butylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
11. sec-Butylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
12. tert-Butylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
13. Carbon Disulfide	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
14. Carbon Tetrachloride	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
15. Chlorobenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
16. Chloroethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
17. Chloroform	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
18. Chloromethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
19. 2-Chlorotoluene	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
20. Dibromochloromethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
21. 1,2-Dibromo-3-chloropropane (SIM) (N)	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
22. Dibromomethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
23. 1,2-Dichlorobenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
24. 1,3-Dichlorobenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
25. 1,4-Dichlorobenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
26. Dichlorodifluoromethane	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
27. 1,1-Dichloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
28. 1,2-Dichloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
29. 1,1-Dichloroethene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
30. cis-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
31. trans-1,2-Dichloroethene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
32. 1,2-Dichloropropane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
33. cis-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
34. trans-1,3-Dichloropropene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
35. Ethylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
36. Ethylene Dibromide	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
37. 2-Hexanone	U		µg/L	50	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
38. Isopropylbenzene	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
39. Methyl Iodide	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>TRIP BLANK TB001102</b>	Chain of Custody:	<b>124238</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>9</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Ground Water</b>	Collect Time:	<b>NA</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Matrix: Ground Water	Analyst: JPL
40. Methylene Chloride	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
41. 2-Methylnaphthalene (NN)	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
42. 4-Methyl-2-pentanone	U		µg/L	50	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
43. MTBE	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
44. Naphthalene	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
45. n-Propylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
46. Styrene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
47. 1,1,1,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
48. 1,1,2,2-Tetrachloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
49. Tetrachloroethene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
50. Toluene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
51. 1,2,4-Trichlorobenzene	U		µg/L	5.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
52. 1,1,1-Trichloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
53. 1,1,2-Trichloroethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
54. Trichloroethene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
55. Trichlorofluoromethane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
56. 1,2,3-Trichloroproppane	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
57. 1,2,3-Trimethylbenzene (NN)	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
58. 1,2,4-Trimethylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
59. 1,3,5-Trimethylbenzene	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
60. Vinyl Chloride	U		µg/L	1.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A
61. Xylenes	U		µg/L	3.0	1.0	07/23/13	VH13G23A	07/23/13	VH13G23A

**Definitions/ Qualifiers:**

- A: Spike recovery or precision unusable due to dilution.
- B: The analyte was detected in the associated method blank.
- E: The analyte was detected at a concentration greater than the calibration range, therefore the result is estimated.
- J: The concentration is an estimated value.
- M: Modified Method
- U: The analyte was not detected at or above the reporting limit.
- X: Matrix Interference has resulted in a raised reporting limit or distorted result.
- W: Results reported on a wet-weight basis.
- \*: Value reported is outside QA limits

**Exception Summary:**

- L- : Recovery in the associated laboratory sample (LCS) exceeds the lower control limit. Results may be biased low.



**E-10395**

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Friday, July 26, 2013

Fibertec Project Number: 57028  
Project Identification: Green Baxter /1127.001  
Submittal Date: 07/18/2013

Ms. Laura Lambert  
Environmental Resources Group  
28003 Center Oaks Court Suite 106  
Wixom, MI 48393

Dear Ms. Lambert,

Thank you for selecting Fibertec Environmental Services as your analytical laboratory. The samples you submitted have been analyzed in accordance with NELAC standards and the results compiled in the attached report. Any exceptions to NELAC compliance are noted in the report. These results apply only to those samples submitted. Please note samples will be disposed of 30 days after reporting date.

If you have any questions regarding these results or if we may be of further assistance to you, please contact me at (517) 699-0345.

Sincerely,



Daryl P. Strandbergh  
Laboratory Director

DPS/kc

Enclosures

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-7</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>1</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:44</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

TO-15 (TO-15)	Aliquot ID: 57028-001						Matrix: Air	Analyst: RDK	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone (NN)	<b>500</b>		ppbv	28	10	07/25/13	V413G25C	07/26/13	V413G25C
2. Benzene (NN)	<b>77</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
3. Benzyl Chloride (NN)	<b>U</b>		ppbv	25	10	07/25/13	V413G25C	07/26/13	V413G25C
4. Bromodichloromethane (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
5. Bromoform (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
6. Bromomethane (NN)	<b>U</b>		ppbv	5.3	10	07/25/13	V413G25C	07/26/13	V413G25C
7. 1,3-Butadiene (NN)	<b>U</b>		ppbv	28	10	07/25/13	V413G25C	07/26/13	V413G25C
8. 2-Butanone (NN)	<b>20</b>		ppbv	14	10	07/25/13	V413G25C	07/26/13	V413G25C
9. Carbon Disulfide (NN)	<b>U</b>		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
10. Carbon Tetrachloride (NN)	<b>U</b>		ppbv	2.7	10	07/25/13	V413G25C	07/26/13	V413G25C
11. Chlorobenzene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
12. Chloroethane (NN)	<b>U</b>		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
13. Chloroform (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
14. Chloromethane (NN)	<b>U</b>		ppbv	5.1	10	07/25/13	V413G25C	07/26/13	V413G25C
15. Cyclohexane (NN)	<b>U</b>		ppbv	5.2	10	07/25/13	V413G25C	07/26/13	V413G25C
16. Dibromochloromethane (NN)	<b>U</b>		ppbv	2.9	10	07/25/13	V413G25C	07/26/13	V413G25C
17. 1,2-Dichlorobenzene (NN)	<b>U</b>		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
18. 1,3-Dichlorobenzene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
19. 1,4-Dichlorobenzene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
20. Dichlorodifluoromethane (NN)	<b>U</b>		ppbv	3.0	10	07/25/13	V413G25C	07/26/13	V413G25C
21. 1,1-Dichloroethane (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
22. 1,2-Dichloroethane (NN)	<b>U</b>		ppbv	5.2	10	07/25/13	V413G25C	07/26/13	V413G25C
23. 1,1-Dichloroethene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
24. cis-1,2-Dichloroethene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
25. trans-1,2-Dichloroethene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
26. 1,2-Dichloropropane (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
27. cis-1,3-Dichloropropene (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
28. trans-1,3-Dichloropropene (NN)	<b>U</b>		ppbv	4.8	10	07/25/13	V413G25C	07/26/13	V413G25C
29. 1,4-Dioxane (NN)	<b>U</b>		ppbv	5.1	10	07/25/13	V413G25C	07/26/13	V413G25C
30. Ethyl Acetate (NN)	<b>U</b>		ppbv	27	10	07/25/13	V413G25C	07/26/13	V413G25C
31. Ethylbenzene (NN)	<b>53</b>	J,L+	ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
32. Ethylene Dibromide (NN)	<b>U</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
33. 4-Ethyltoluene (NN)	<b>U</b>		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
34. n-Heptane (NN)	<b>11</b>		ppbv	5.3	10	07/25/13	V413G25C	07/26/13	V413G25C
35. Hexachlorobutadiene (NN)	<b>U</b>		ppbv	5.1	10	07/25/13	V413G25C	07/26/13	V413G25C
36. n-Hexane (NN)	<b>26</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
37. 2-Hexanone (NN)	<b>U</b>		ppbv	29	10	07/25/13	V413G25C	07/26/13	V413G25C
38. Isopropanol (NN)	<b>U</b>		ppbv	29	10	07/25/13	V413G25C	07/26/13	V413G25C
39. Methylene Chloride (NN)	<b>U</b>		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-7</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>1</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:44</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>TO-15 (TO-15)</b>		<b>Aliquot ID: 57028-001</b>					<b>Matrix: Air</b>	<b>Analyst: RDK</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
40. 2-Methylnaphthalene (NN)	U		ppbv	49	10	07/25/13	V413G25C	07/26/13	V413G25C
41. 4-Methyl-2-pentanone (NN)	<b>17</b>		ppbv	5.1	10	07/25/13	V413G25C	07/26/13	V413G25C
42. MTBE (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
43. Naphthalene (NN)	U		ppbv	28	10	07/25/13	V413G25C	07/26/13	V413G25C
44. Propylene (NN)	<b>520</b> J,L+		ppbv	5.5	10	07/25/13	V413G25C	07/26/13	V413G25C
45. Styrene (NN)	U		ppbv	5.2	10	07/25/13	V413G25C	07/26/13	V413G25C
46. 1,1,2,2-Tetrachloroethane (NN)	U		ppbv	5.3	10	07/25/13	V413G25C	07/26/13	V413G25C
47. Tetrachloroethene (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
48. Tetrahydrofuran (NN)	U		ppbv	5.8	10	07/25/13	V413G25C	07/26/13	V413G25C
49. Toluene (NN)	<b>210</b>		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
50. 1,2,4-Trichlorobenzene (NN)	U		ppbv	5.0	10	07/25/13	V413G25C	07/26/13	V413G25C
51. 1,1,1-Trichloroethane (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
52. 1,1,2-Trichloroethane (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
53. Trichloroethene (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
54. Trichlorofluoromethane (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
55. 1,1,2-Trichlorotrifluoroethane (NN)	U		ppbv	2.8	10	07/25/13	V413G25C	07/26/13	V413G25C
56. 1,2,4-Trimethylbenzene (NN)	<b>88</b> J,L+		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
57. 1,3,5-Trimethylbenzene (NN)	U		ppbv	26	10	07/25/13	V413G25C	07/26/13	V413G25C
58. Vinyl Acetate (NN)	U		ppbv	27	10	07/25/13	V413G25C	07/26/13	V413G25C
59. Vinyl Chloride (NN)	U		ppbv	13	10	07/25/13	V413G25C	07/26/13	V413G25C
60. m&p-Xylene (NN)	<b>240</b> J,L+		ppbv	5.5	10	07/25/13	V413G25C	07/26/13	V413G25C
61. o-Xylene (NN)	<b>69</b> J,L+		ppbv	5.3	10	07/25/13	V413G25C	07/26/13	V413G25C
62. Xylenes (NN)	<b>310</b> J,L+		ppbv	11	10	07/25/13	V413G25C	07/26/13	V413G25C

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-6</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>2</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:51</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

TO-15 (TO-15)	Aliquot ID: 57028-002						Matrix: Air	Analyst: RDK	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone (NN)	<b>150</b>		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
2. Benzene (NN)	<b>53</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
3. Benzyl Chloride (NN)	U		ppbv	12	5.0	07/24/13	V413G24A	07/25/13	V413G24A
4. Bromodichloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
5. Bromoform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
6. Bromomethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
7. 1,3-Butadiene (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
8. 2-Butanone (NN)	<b>9.6</b>		ppbv	6.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
9. Carbon Disulfide (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
10. Carbon Tetrachloride (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
11. Chlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
12. Chloroethane (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
13. Chloroform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
14. Chloromethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
15. Cyclohexane (NN)	<b>5.9</b>		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
16. Dibromochloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
17. 1,2-Dichlorobenzene (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
18. 1,3-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
19. 1,4-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
20. Dichlorodifluoromethane (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
21. 1,1-Dichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
22. 1,2-Dichloroethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
23. 1,1-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
24. cis-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
25. trans-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
26. 1,2-Dichloropropane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
27. cis-1,3-Dichloropropene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
28. trans-1,3-Dichloropropene (NN)	U		ppbv	2.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
29. 1,4-Dioxane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
30. Ethyl Acetate (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
31. Ethylbenzene (NN)	<b>57</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
32. Ethylene Dibromide (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
33. 4-Ethyltoluene (NN)	<b>93</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
34. n-Heptane (NN)	<b>10</b>		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
35. Hexachlorobutadiene (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
36. n-Hexane (NN)	<b>29</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
37. 2-Hexanone (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
38. Isopropanol (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
39. Methylene Chloride (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-6</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>2</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:51</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>TO-15 (TO-15)</b>		<b>Aliquot ID: 57028-002</b>					<b>Matrix: Air</b>	<b>Analyst: RDK</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
40. 2-Methylnaphthalene (NN)	U		ppbv	24	5.0	07/24/13	V413G24A	07/25/13	V413G24A
41. 4-Methyl-2-pentanone (NN)	<b>8.5</b>		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
42. MTBE (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
43. Naphthalene (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
44. Propylene (NN)	<b>470</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
45. Styrene (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
46. 1,1,2,2-Tetrachloroethane (NN)	U		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
47. Tetrachloroethene (NN)	<b>2.0</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
48. Tetrahydrofuran (NN)	U		ppbv	2.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
49. Toluene (NN)	<b>150</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
50. 1,2,4-Trichlorobenzene (NN)	U		ppbv	2.5	5.0	07/24/13	V413G24A	07/25/13	V413G24A
51. 1,1,1-Trichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
52. 1,1,2-Trichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
53. Trichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
54. Trichlorofluoromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
55. 1,1,2-Trichlorotrifluoroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
56. 1,2,4-Trimethylbenzene (NN)	<b>130</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
57. 1,3,5-Trimethylbenzene (NN)	<b>100</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
58. Vinyl Acetate (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
59. Vinyl Chloride (NN)	U		ppbv	6.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
60. m&p-Xylene (NN)	<b>240</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
61. o-Xylene (NN)	<b>84</b>		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
62. Xylenes (NN)	<b>320</b>		ppbv	5.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-5</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>3</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:59</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

TO-15 (TO-15)	Aliquot ID: 57028-003						Matrix: Air	Analyst: RDK	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone (NN)	U		ppbv	84	5.0	07/24/13	V413G24A	07/25/13	V413G24A
2. Benzene (NN)	39		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
3. Benzyl Chloride (NN)	U		ppbv	12	5.0	07/24/13	V413G24A	07/25/13	V413G24A
4. Bromodichloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
5. Bromoform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
6. Bromomethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
7. 1,3-Butadiene (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
8. 2-Butanone (NN)	U		ppbv	6.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
9. Carbon Disulfide (NN)	22		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
10. Carbon Tetrachloride (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
11. Chlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
12. Chloroethane (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
13. Chloroform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
14. Chloromethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
15. Cyclohexane (NN)	2.7		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
16. Dibromochloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
17. 1,2-Dichlorobenzene (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
18. 1,3-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
19. 1,4-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
20. Dichlorodifluoromethane (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
21. 1,1-Dichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
22. 1,2-Dichloroethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
23. 1,1-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
24. cis-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
25. trans-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
26. 1,2-Dichloropropane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
27. cis-1,3-Dichloropropene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
28. trans-1,3-Dichloropropene (NN)	U		ppbv	2.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
29. 1,4-Dioxane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
30. Ethyl Acetate (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
31. Ethylbenzene (NN)	20		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
32. Ethylene Dibromide (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
33. 4-Ethyltoluene (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
34. n-Heptane (NN)	12		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
35. Hexachlorobutadiene (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
36. n-Hexane (NN)	18		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
37. 2-Hexanone (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
38. Isopropanol (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
39. Methylene Chloride (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-5</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>3</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>18:59</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>TO-15 (TO-15)</b>		<b>Aliquot ID: 57028-003</b>					<b>Matrix: Air</b>	<b>Analyst: RDK</b>		
Parameter(s)		Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
40. 2-Methylnaphthalene (NN)		U		ppbv	24	5.0	07/24/13	V413G24A	07/25/13	V413G24A
41. 4-Methyl-2-pentanone (NN)		<b>4.0</b>		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
42. MTBE (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
43. Naphthalene (NN)		U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
44. Propylene (NN)		<b>82</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
45. Styrene (NN)		U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
46. 1,1,2,2-Tetrachloroethane (NN)		U		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
47. Tetrachloroethene (NN)		<b>5.6</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
48. Tetrahydrofuran (NN)		U		ppbv	2.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
49. Toluene (NN)		<b>140</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
50. 1,2,4-Trichlorobenzene (NN)		U		ppbv	2.5	5.0	07/24/13	V413G24A	07/25/13	V413G24A
51. 1,1,1-Trichloroethane (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
52. 1,1,2-Trichloroethane (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
53. Trichloroethene (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
54. Trichlorofluoromethane (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
55. 1,1,2-Trichlorotrifluoroethane (NN)		U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
56. 1,2,4-Trimethylbenzene (NN)		<b>32</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
57. 1,3,5-Trimethylbenzene (NN)		U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
58. Vinyl Acetate (NN)		U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
59. Vinyl Chloride (NN)		U		ppbv	<b>6.6</b>	5.0	07/24/13	V413G24A	07/25/13	V413G24A
60. m&p-Xylene (NN)		<b>86</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
61. o-Xylene (NN)		<b>23</b>		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
62. Xylenes (NN)		<b>110</b>		ppbv	5.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-4</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>4</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>19:05</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

TO-15 (TO-15)	Aliquot ID: 57028-004						Matrix: Air	Analyst: RDK	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
1. Acetone (NN)	<b>130</b>		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
2. Benzene (NN)	<b>76</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
3. Benzyl Chloride (NN)	U		ppbv	12	5.0	07/24/13	V413G24A	07/25/13	V413G24A
4. Bromodichloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
5. Bromoform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
6. Bromomethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
7. 1,3-Butadiene (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
8. 2-Butanone (NN)	<b>9.5</b>		ppbv	6.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
9. Carbon Disulfide (NN)	<b>60</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
10. Carbon Tetrachloride (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
11. Chlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
12. Chloroethane (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
13. Chloroform (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
14. Chloromethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
15. Cyclohexane (NN)	<b>4.1</b>		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
16. Dibromochloromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
17. 1,2-Dichlorobenzene (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
18. 1,3-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
19. 1,4-Dichlorobenzene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
20. Dichlorodifluoromethane (NN)	U		ppbv	1.3	5.0	07/24/13	V413G24A	07/25/13	V413G24A
21. 1,1-Dichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
22. 1,2-Dichloroethane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
23. 1,1-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
24. cis-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
25. trans-1,2-Dichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
26. 1,2-Dichloropropane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
27. cis-1,3-Dichloropropene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
28. trans-1,3-Dichloropropene (NN)	U		ppbv	2.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
29. 1,4-Dioxane (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
30. Ethyl Acetate (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
31. Ethylbenzene (NN)	<b>48</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
32. Ethylene Dibromide (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
33. 4-Ethyltoluene (NN)	<b>18</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
34. n-Heptane (NN)	<b>18</b>		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
35. Hexachlorobutadiene (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
36. n-Hexane (NN)	<b>29</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
37. 2-Hexanone (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
38. Isopropanol (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
39. Methylene Chloride (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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Client Identification:	<b>Environmental Resources Group</b>	Sample Description:	<b>SB-4</b>	Chain of Custody:	<b>124237</b>
Client Project Name:	<b>Green Baxter</b>	Sample No:	<b>4</b>	Collect Date:	<b>07/18/13</b>
Client Project No:	<b>1127.001</b>	Sample Matrix:	<b>Air</b>	Collect Time:	<b>19:05</b>

Sample Comments:

Definitions: Q: Qualifier (see definitions at end of report) NA: Not Applicable NN: Parameter not included in NELAC Scope of Analysis.

<b>TO-15 (TO-15)</b>		<b>Aliquot ID: 57028-004</b>					<b>Matrix: Air</b>	<b>Analyst: RDK</b>	
Parameter(s)	Result	Q	Units	Reporting Limit	Dilution	Prep Date	Prep Batch	Analysis Date	Analysis Batch
40. 2-Methylnaphthalene (NN)	U		ppbv	24	5.0	07/24/13	V413G24A	07/25/13	V413G24A
41. 4-Methyl-2-pentanone (NN)	<b>7.2</b>		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
42. MTBE (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
43. Naphthalene (NN)	U		ppbv	14	5.0	07/24/13	V413G24A	07/25/13	V413G24A
44. Propylene (NN)	<b>420</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
45. Styrene (NN)	U		ppbv	2.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
46. 1,1,2,2-Tetrachloroethane (NN)	U		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
47. Tetrachloroethene (NN)	<b>6.3</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
48. Tetrahydrofuran (NN)	U		ppbv	2.9	5.0	07/24/13	V413G24A	07/25/13	V413G24A
49. Toluene (NN)	<b>250</b>		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
50. 1,2,4-Trichlorobenzene (NN)	U		ppbv	2.5	5.0	07/24/13	V413G24A	07/25/13	V413G24A
51. 1,1,1-Trichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
52. 1,1,2-Trichloroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
53. Trichloroethene (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
54. Trichlorofluoromethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
55. 1,1,2-Trichlorotrifluoroethane (NN)	U		ppbv	1.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A
56. 1,2,4-Trimethylbenzene (NN)	<b>52</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
57. 1,3,5-Trimethylbenzene (NN)	<b>18</b>		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
58. Vinyl Acetate (NN)	U		ppbv	13	5.0	07/24/13	V413G24A	07/25/13	V413G24A
59. Vinyl Chloride (NN)	U		ppbv	6.6	5.0	07/24/13	V413G24A	07/25/13	V413G24A
60. m&p-Xylene (NN)	<b>200</b>		ppbv	2.8	5.0	07/24/13	V413G24A	07/25/13	V413G24A
61. o-Xylene (NN)	<b>59</b>		ppbv	2.7	5.0	07/24/13	V413G24A	07/25/13	V413G24A
62. Xylenes (NN)	<b>260</b>		ppbv	5.4	5.0	07/24/13	V413G24A	07/25/13	V413G24A

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**Definitions/ Qualifiers:**

- A: Spike recovery or precision unusable due to dilution.
- B: The analyte was detected in the associated method blank.
- E: The analyte was detected at a concentration greater than the calibration range, therefore the result is estimated.
- J: The concentration is an estimated value.
- M: Modified Method
- U: The analyte was not detected at or above the reporting limit.
- X: Matrix Interference has resulted in a raised reporting limit or distorted result.
- W: Results reported on a wet-weight basis.
- \*: Value reported is outside QA limits

**Exception Summary:**

L+ : Recovery in the associated laboratory sample (LCS) exceeds the upper control limit. Results may be biased high.



Accreditation Number:

**E-10395**

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Chain of Custody #  
**124237**