

1983 Marcus Ave., Suite 109  
Lake Success, New York 11042  
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Fax (516) 328-1381

### LETTER OF TRANSMITTAL

Date:	04/14/06	Job No.	25001
Attention: <b>Mr. Carl Hoffman</b>			
Re:	<b>Katonah Quarterly Water Monitoring</b>		

**TO:**

**NYSDEC**  
**625 Broadway**  
**Albany, NY 12233-7013**

WE ARE SENDING YOU:  Included  Under separate cover via \_\_\_\_\_ the following items:

Shop Drawings  Prints  Plans  Qualifications  Specifications  
 Copy of Letter  Report  \_\_\_\_\_

COPIES	DATE	NO.	
1	4/14/06		<i>Katonah Quarterly Water Monitoring Report - 4th Quarter</i>

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REMARKS

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If there are any questions, please call me.

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COPY TO File

SIGNED



James Hahn  
James J. Hahn Engineering  
Putnam Business Park  
1689 Route 22  
Brewster, NY 10509

April 10, 2006

Dear Mr. Hahn:

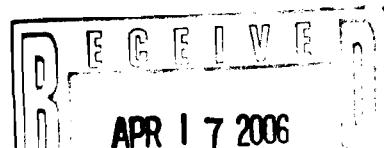
Enclosed please find the quarterly monitoring report for the end of the 4<sup>th</sup> quarter of 2005 for the Katonah Municipal Well, Town of Bedford, Westchester County, New York (NYSDEC Site ID # 3-60-007).

Please call me with any questions.

Sincerely,

Francesco Portelos  
Project Engineer

cc: Kenneth Caffrey, PE, NYSDOH  
Carl Hoffman, NYSDEC  
William Nixon, Town of Bedford  
Paul Kutzy, Westchester County DOH  
Damian Duda, USEPA Region 2



**GROUNDWATER QUALITY MONITORING  
QUARTERLY REPORT  
DECEMBER 2005  
KATONAH MUNICIPAL WELL  
TOWN OF BEDFORD  
WESTCHESTER, NEW YORK  
NYSDEC Site ID # 3-60-007**

**PREPARED FOR:**

**James J. Hahn Engineering  
Millbrook Office Center  
Route 22 & Milltown Road  
Brewster, New York 10509**

**PREPARED BY:**

**Environmental Planning & Management, Inc.  
1983 Marcus Avenue, Suite 109  
Lake Success, New York 11042**

*RECEIVED*  
*APR 17 2006*

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## **1.0 INTRODUCTION**

This quarterly groundwater sampling and analysis report has been prepared for the Katonah Municipal Well Site in Katonah, Town of Bedford, New York. This submittal is in accordance with the groundwater monitoring requirements of the New York State Department of Health (NYSDOH) and the U.S. Environmental Protection Agency (USEPA). This report includes the data collection and analysis results of the remedial system operation, for the end of the 4<sup>th</sup> quarter of 2005. Sampling of the remedial system was conducted on December 20<sup>th</sup>, 2005.

## **2.0 SAMPLE COLLECTION**

Environmental Planning & Management, Inc., collected samples on December 20<sup>th</sup>, 2005. Three sample sets were collected from sampling taps; the raw water sampling tap (RW), the stripper number two effluent sampling tap (STEFF), and the distribution sampling tap (DIST). One field duplicate sample (DUP) was collected on December 20<sup>th</sup>, 2005 of the RW sampling tap. Sample locations are shown on Figure 1 - Sampling Tap Location Schematic. Sampling was conducted in accordance with the approved Project Operation Plan.

Samples were labeled at the field location and placed into transport coolers containing ice. A trip blank and chain-of-custody documentation accompanied the samples to the laboratory for analysis. The samples were analyzed by Chemtech , in accordance with CLP methods, for volatile organics (Principal Organic Contaminants), by method 524.2, revision number 3.

### 3.0 FINDINGS

Table 1 provides a summary of the analytical results for the quarterly water quality monitoring, as well as the applicable NYSDOH Drinking Water Standards and the U.S. EPA clean-up requirement for Tetrachloroethene. As indicated by the laboratory analysis, the treatment system effluent meets the NYSDOH drinking water standards and the USEPA clean-up level of less than one part per billion (ppb) (or non-detectable) for Tetrachloroethene and meets the levels of less than 100 parts per billion for Trihalomethanes.

Tetrachloroethene was detected in the raw water (untreated) sample, RW, at a concentration of 30ug/l (ppb), exceeding the NYSDOH drinking water standard for that compound.

One VOC, Methylene Chloride was detected in the treated (stripper number 2) water sample, STEFF at a concentration of 1.0ppb. This value is well below the NYSDOH drinking water standards.

Two VOC's, Dibromochloromethane and Bromodichloromethane were found in the distribution water sample, DIST, at concentrations of 4.0ppb and 1.7ppb respectively. These values are well below the NYSDOH drinking water standards.

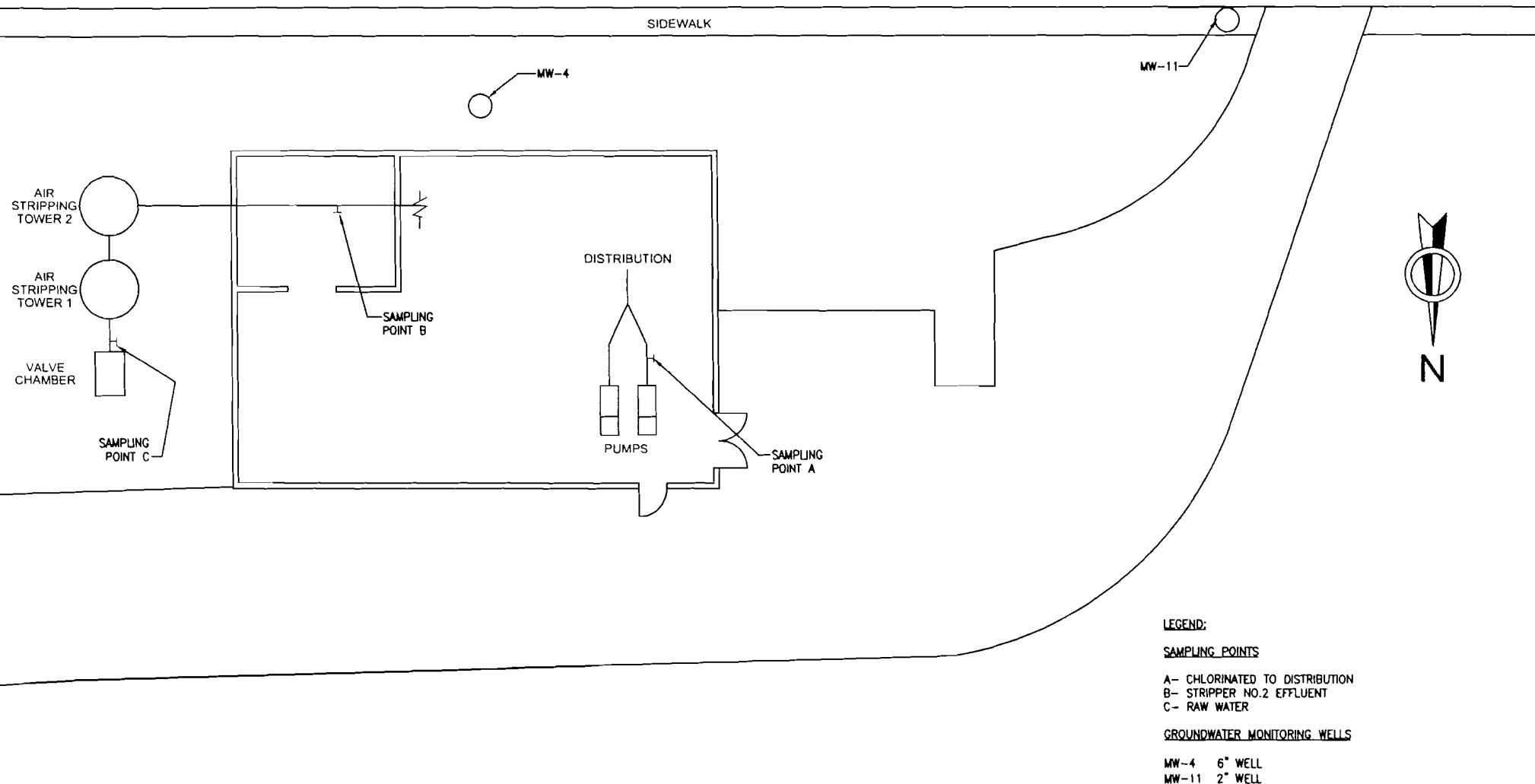
No VOC's were detected in the trip blank water sample, TB.

Analytical results found in DUP, a duplicate sample of the Raw Water sample, RW, are similar.

Refer to Table 1 for a summary of the groundwater analysis results for volatile organic compounds (VOC's). Table 1 reflects the detectable concentration values which have been qualified as a result of data validation. Refer to Appendix A for the data validation report which details the changes in the detectable concentration values discussed above.

The PCE concentration in the Influent (raw water) has increased over the last sampling event (see Figure 2). To date, the PCE level in the raw water samples is not of significant concern, since the treated water and distribution water samples continue to exhibit non-detectable or insignificant concentrations of PCE. However, changes in PCE levels will continue to be closely monitored.

# JAY STREET



**Table 1 - SUMMARY OF QUARTERLY ANALYTICAL RESULTS  
KATONAH MUNICIPAL WELL  
October 2005**

Date Collected	12/20/2005					
Sample Location	Raw Water (Influent)	RW DUP	STEFF (Treated Water)	DIST (Distribution Water)	TB (Trip Blank)	NYSDOH\\ USEPA Standard
<i>Volatile Organic Compounds (ppb)</i>						
Tetrachloroethene	30.00J	29.00J	.16U	.16U	.16U	5/1*
Trichloroethene	.9J	.8J	.15U	.15U	.15U	5
cis-1,2-Dichloroethene	.9J	.8J	.12U	.12U	.12U	5
Methylene Chloride	.27U	.27U	1.00	.27U	.27U	5
Dibromochloromethane	.17U	.17U	.17U	4.00	.17U	50
Bromodichloromethane	.17U	.17U	.17U	1.70	.17U	50

\* 1 ppb is the USEPA cleanup standard for the site

1 - Determined undetect following data validation

Level exceeds the USEPA/NYSDOH standard

U Denotes detection limit/not detected

J Denotes an estimated value

N Presumptive evidence of a compound

R Determined unusable following data validation

NS No standard

B Denotes Detection in the Field Blank as well.



#### **4.0 FUTURE ACTIONS**

Water quality monitoring will continue to be conducted quarterly at the treatment system influent, stripper number 2 effluent, and distribution entry point. Groundwater monitoring well samples will be collected bi-annually. EPM will communicate with the Town of Bedford Water Department to schedule a date when all the taps are available for sampling.

The next sampling event, the end of the first quarterly event for year fifteen, is tentatively scheduled for March 2006.

## **APPENDIX A**

### **Katonah Municipal Well Site Data Validation Groundwater Quality Monitoring Quarterly Report - December 2005**

**Samples Collected by Environmental Planning & Management, Inc.  
Samples Analyzed by Chemtech**

**Data Validation Performed by:**

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**Julie Smith  
Environmental Chemist**

## PROJECT DESCRIPTION

**Report Prepared by:** Julie Smith, Environmental Chemist

**Date of Validation Report:** March 23, 2006

**EPM Project Name/No.** 25001-Katonah 4th Quarter

**Laboratory:** Chemtech

**Laboratory Project Name:** T6267

**Deliverable Format:** NYSDEC ASP B

**Sample Date:** December 20, 2005

<b>Samples Validated:</b>	EPM Sample ID	Laboratory Sample ID
RW		T6267-01
RW-MS		T6267-02
RW-MSD		T6267-03
DIST		T6267-04
STEFF		T6267-05
DUP		T6267-06
TB		T6267-07

### Validation Protocols/

**References:** U.S. Environmental Protection Agency (USEPA) Standard Operating Procedure for the Validation of Organic Data Acquired Using Method 524.2, October 2001.

U.S. Environmental Protection Agency (USEPA) Measurement of Purgeable Organic Compounds in Water by Capillary Column Gas Chromatography/Mass Spectrometry, Method 524.2, Methods for Chemical Analysis of Water and Wastes, 1995.

U.S. Environmental Protection Agency (USEPA) National Functional Guidelines for Organic Data Review, 1999.

## INTRODUCTION

Data qualification provides guidance regarding data usability. As part of the environmental laboratory analytical reporting process under most environmental methods of analysis, the laboratory is required to append data qualifiers to reported analytical observations to account for minor, acceptable QC deficiencies that arise during the course of standard operations. As part of the analytical data validation process, additional data qualifiers may be applied. These qualifiers are applied for other QC deficiencies that impact data quality but that may not have been identified by the laboratory or that may not be part of the reporting requirement of the applied analytical method. In many cases, the laboratory may be compliant with the requirements of the applied analytical methods but may not be compliant with the data validation review protocols.

In general, the data qualifiers are intended to assist the data user with the overall data interpretation by serving as descriptive indicators of the data quality of the associated analytical observations. There are a number of other data qualifiers that describe the representativeness of the associated data and also serve to provide information about the quality of the associated control data.

- U** The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- UJ** The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- J** The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample. When data are qualified as estimated (qualified "J"), there generally is no information on the quantitative impact on the associated result although there may be useful information on the direction of bias of the result
- R** The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet QC criteria. The presence or absence of the analyte cannot be confirmed. In some cases, sample data are qualified as unusable and rejected (qualified "R") due to major method non-compliance or extreme deficiencies in associated QC data. In these cases, there is no information as to the presence or absence of the rejected analyte in the affected sample.

## VALIDATION SUMMARY

The analytical data has been reviewed in accordance with the appropriate regulatory guidelines and/or associated analytical methodology. If required, the data has been qualified, negated, or rejected according to applicable validation protocols and professional judgment. The analytical validation was performed based upon the following parameters:

- \* Completeness of data package
- \* Blank Contamination
- \* Hold Times
- \* GC/MS Performance Check (Tuning) Summaries
- \* System Monitoring Compound (Surrogate) Recoveries
- \* Internal Standard Area Performance
- \* Initial and Continuing Calibration Results  
Matrix Spike (MS) and Matrix Spike Duplicate (MSD) Summaries
- \* Laboratory Control Sample
- \* Target Compound Identification and Quantitation

\* All criteria were met for this parameter

## OVERALL DATA ASSESSMENT

The volatile organics data was validated for compliance with the requirements set forth in EPA Method 524.2 and as described by the Standard Operating Procedure for the Validation of Organic Data Acquired Using Method 524.2. Overall, the data quality is acceptable. The data validation review has identified aspects of the analytical data that require qualification. The laboratory analytical data contained herein are deemed usable and in compliance with the New York ASP B Deliverable Format requirements.

## VOLATILE ORGANIC RESULTS

### General Comments

Documentation as required by the project sample analyses was included in the data package. Additionally, there were no discrepancies found between the reviewed raw data and summary forms.

### Blank Contamination

Laboratory method blanks (instrument blanks) are clean liquid matrix samples prepared by the laboratory and analyzed in the same manner as the investigative samples. Laboratory method blanks are used to identify whether investigative samples have been contaminated during the sample preparation, sample analysis or from a previous sample (instrument carry-over).

Trip-blanks are carbon-free deionized water samples that accompany volatile investigative samples during all stages of shipment, storage and analysis. The trip-blanks are used to assess the potential for artificial introduction of volatile compounds into the investigative samples during the transportation and sample handling processes.

- The VOA target compounds, acetone and 2-butanone, were detected in the method blank VBLK01. No qualifier is required since acetone and 2-butanone was not detected in the associated project samples.

### **Hold Times**

Technical hold times were assessed by comparing the sample dates with that of the preparation dates and/or analysis dates.

- All volatile analyses performed on the associated project samples were within the required hold times. The sample cooler temperature (4.0°C) upon verified time of sample receipt (VTSR) in the laboratory fell within the 4°C ( $\pm 2^{\circ}\text{C}$ ) requirement.

### **Internal Standard Area Performance**

Internal standards are analytes that are added to the investigative samples prior to analysis to ensure that GC/MS sensitivity and responses remain stable. Internal standards are reported with the volatile analyses.

- The volatile internal standard area counts and retention times fell within control limits for the associated project samples. No qualifier is required.

### **Matrix Spike/Matrix Spike Duplicate**

Matrix spikes are samples spiked with known concentrations of analytes of interest. The MS/MSD percent recoveries and duplicate results are used to assess extraction efficiencies, possible matrix effects, and overall analytical accuracy and precision.

- A matrix spike/matrix spike duplicate (MS/MSD) was performed on EPM Sample RW. The volatile percent recoveries (%R), for the most part, fell within control limits in both the MS and MSD samples. However, the MS recovery and RPD for target compound, tetrachloroethene, fell outside control limits (low). Therefore, the positive tetrachloroethene results in the associated project samples are regarded as estimated values and are flagged with a (J).
- The relative percent differences (RPD) fell within control limits. No qualifier is required.

### **Laboratory Control Sample**

The laboratory control sample (LCS) and/or blank spike (BS) are blank samples fortified (spiked) with known concentrations of analytes of interest. The percent recoveries of the LCS and/or BS are used to assess overall analytical accuracy and precision.

- The volatile LCS results fell within acceptable control limits with the exception of methylene chloride (low) and acetone (low). The non-detected methylene chloride and acetone results reported for the associated project samples are qualified (UJ).

### **System Monitoring Compounds (Surrogates)**

System monitoring compounds are those compounds that are not expected to be detected in the investigative samples but that are chemically similar to analytes of interest. Surrogate compound percent recoveries are used to assess extraction efficiencies, possible matrix effects and overall analytical accuracy.

- The recoveries of the volatile surrogates, 4-bromofluorobenzene (BFB) and 1,2-dichlorobenzene-d4 (DCB), fell within control limits for the reviewed project samples.

### **Initial Calibration and Continuing Calibration Results**

Control limits for initial and continuing instrument calibrations are established to ensure that the instrument is capable of producing accurate quantitative data at the beginning and throughout each of the analyses.

- The volatile initial and continuing calibration response factors (RRF), percent relative standard deviations (%RSD), percent differences (%D) and/or correlation coefficients fell within acceptable control limits. No qualifier is required.

### **GC/MS Performance Check (Tuning) Summaries**

Gas chromatograph/mass spectrometer (GC/MS) instrument tuning and performance checks are performed to ensure the instrument's ability to provide appropriate mass-resolution, identification and sensitivity.

- The bromofluorobenzene (BFB) tuning compound mass-ion abundance criteria for the volatile organic compound analyses were reported within control limits. The samples were analyzed within twelve hours of BFB injection. No qualifier is required.

### **Compound Identification and Quantitation**

The laboratory calculations are verified and compound identifications are reviewed and assessed by the data reviewer.

- Sample DUP was collected and submitted as a blind field duplicate of project sample RW. The reproducibility of the associated analyses is good providing a positive indication of the accuracy and precision associated with these samples.

### **Tentatively Identified Compound**

Area not examined, validation not requested.

**Volatiles Method 524.2**

Chemtech Project #T6267

**Hold Time**

Samples collected on 12/20/05 and analyzed on 12/27/05.

**Instrument Performance Check (BFB)**

12/23/2005	12:44 PM	Instr# MSVOAF	meets QC requirements	initial calibration	/
12/27/2005	2:25 PM		meets QC requirements	sample analysis	

**Initial Calibration**

12/23/2005 Instr# MSVOAF

Compound	1 RRF	2 RRF		20 RRF		30 RRF	Mean	STDEV	%RSD
Fluorobenzene (IS)	544844		530210		542416		594603		
tetrachloroethene	156610	0.287	227801	0.215	2655827	0.245	3967465	0.222	0.242
Bromoform	57034	0.105	96603	0.091	1129955	0.104	1734983	0.097	0.099
Trichloroethene	197710	0.363	279511	0.264	3336857	0.308	4979225	0.279	0.303
BFB	258415	0.474	266534	0.503	273500	0.504	282777	0.476	0.489

control limit RSD &lt; 20%

correlation coefficient &gt;0.990

**Continuing Calibration** 10 ppb ccc 12/27/2005 Instr# MSVOAF

Compound	10 RRF	% Difference	Status
Fluorobenzene (IS)	860743		
tetrachloroethene	1835335	0.2132	12.0
Bromoform	693178	0.0805	18.9
Trichloroethene	2494541	0.2898	4.4
BFB	403333	0.4686	4.2

control limit D &lt; 30%

**Surrogate Recovery**

Recoveries of surrogates, BFB and 1,2-dichloroethane -d4, fell within control limits for the reviewed samples.

**Internal Standard Summary**

Internal standard areas and retention times fell within control limits for the reviewed samples.

Blanks	Conc.	Compound
VBLK01	6.9 ug/L	acetone
	1.7 ug/L	2-butanone
Trip Blank	ND	

Not detected in samples  
Not detected in samples

QC	Sample RW	RW	%R	RW	MSD	%R	RPD	QC limits
Trichloroethene	0.9	6.8	118	6.5	112	-5	70-130	15
Tetrachloroethene	30	32	40	32	43	0	70-130	15 outside control limits
Bromodichloromethane	ND	5.5	110	5.8	116	5	70-130	15
cis-1,2-dichloroethene	0.9	6.2	106	6.3	108	2	70-130	15

Field Duplicate Results	RW	DUP	RPD
c-1,2-dichloroethene	0.9 J	0.8 J	0
trichloroethene	0.9 J	0.8 J	0
tetrachloroethene	30	29	3.4

## Sample Results

Sample ID	DIST	(area of IS)(RRT)	Raw Result	Dilution	Final Result	Reported Result
Lab ID T6267-04						
124986		1	= <u>124986</u>			
558047		0.463	= <u>258375.76</u>			
Bromodichloromethane						
321652		1	= <u>321652</u>			
558047		0.348	= <u>194200.36</u>			
Dibromochloromethane						
444769		1	= <u>444769</u>			
558047		0.201	= <u>112167.45</u>			
Bromoform						
199920		1	= <u>199920</u>			
558047		0.101	= <u>56362.75</u>			
Sample ID RW						
Tetrachloroethene						
Lab ID T6267-01						
3903930		1	= <u>3903930</u>			
540674		0.242	= <u>131044.71</u>			
Trichloroethene						
149918		1	= <u>149918</u>			
540674		0.303	= <u>163984.55</u>			

**APPENDIX B**  
**LABORATORY ANALYSIS SUMMARY REPORT**







284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	RW	SDG No.:	T6267
Lab Sample ID:	T6267-01	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000557.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
106-43-4	4-Chlorotoluene	0.15	U	1.0	0.15	ug/L
98-06-6	tert-Butylbenzene	0.15	U	1.0	0.15	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.15	U	1.0	0.15	ug/L
135-98-8	Sec-butylbenzene	0.14	U	1.0	0.14	ug/L
99-87-6	p-Isopropyltoluene	0.14	U	1.0	0.14	ug/L
541-73-1	1,3-Dichlorobenzene	0.15	U	1.0	0.15	ug/L
106-46-7	1,4-Dichlorobenzene	0.17	U	1.0	0.17	ug/L
104-51-8	n-Butylbenzene	0.12	U	1.0	0.12	ug/L
95-50-1	1,2-Dichlorobenzene	0.16	U	1.0	0.16	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	0.19	U	1.0	0.19	ug/L
120-82-1	1,2,4-Trichlorobenzene	0.11	U	1.0	0.11	ug/L
87-68-3	Hexachlorobutadiene	0.13	U	1.0	0.13	ug/L
91-20-3	Naphthalene	0.14	U	1.0	0.14	ug/L
87-61-6	1,2,3-Trichlorobenzene	0.16	U	1.0	0.16	ug/L

**SURROGATES**

2199-69-1	1,2-Dichlorobenzene-d4	1.15	115 %	80 - 120	SPK: 1
460-00-4	4-Bromofluorobenzene	1.03	103 %	80 - 120	SPK: 1

**INTERNAL STANDARDS**

462-06-6	Fluorobenzene	540674	9.12
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U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound





284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	DIST	SDG No.:	T6267
Lab Sample ID:	T6267-04	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID			
VF000558.D	1	12/27/2005	VF122305			

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
78-87-5	1,2-Dichloropropane	0.14	U	1.0	0.14	ug/L
126-98-7	Methacrylonitrile	0.62	U	1.0	0.62	ug/L
109-99-9	Tetrahydrofuran	0.45	U	2.4	0.45	ug/L
109-69-3	1-Chlorobutane	0.17	U	1.0	0.17	ug/L
74-95-3	Dibromomethane	0.19	U	1.0	0.19	ug/L
75-27-4	Bromodichloromethane	1.7		1.0	0.17	ug/L
108-10-1	4-Methyl-2-Pentanone	0.90	U	5.0	0.90	ug/L
80-62-6	Methyl methacrylate	0.32	U	2.0	0.32	ug/L
97-63-2	Ethyl methacrylate	0.16	U	1.0	0.16	ug/L
108-88-3	Toluene	0.13	U	1.0	0.13	ug/L
10061-02-6	t-1,3-Dichloropropene	0.14	U	1.0	0.14	ug/L
10061-01-5	cis-1,3-Dichloropropene	0.13	U	1.0	0.13	ug/L
79-00-5	1,1,2-Trichloroethane	0.18	U	1.0	0.18	ug/L
142-28-9	1,3-Dichloropropane	0.14	U	1.0	0.14	ug/L
591-78-6	2-Hexanone	0.81	U	5.0	0.81	ug/L
124-48-1	Dibromochloromethane	4.0		1.0	0.17	ug/L
106-93-4	1,2-Dibromoethane	0.17	U	1.0	0.17	ug/L
127-18-4	Tetrachloroethene	0.16	U	1.0	0.16	ug/L
108-90-7	Chlorobenzene	0.13	U	1.0	0.13	ug/L
630-20-6	1,1,1,2-Tetrachloroethane	0.17	U	1.0	0.17	ug/L
67-72-1	Hexachloroethane	0.17	U	1.0	0.17	ug/L
100-41-4	Ethyl Benzene	0.14	U	1.0	0.14	ug/L
126777-61-2	m/p-Xylenes	0.29	U	1.0	0.29	ug/L
95-47-6	o-Xylene	0.15	U	1.0	0.15	ug/L
100-42-5	Styrene	0.14	U	1.0	0.14	ug/L
75-25-2	Bromoform	3.6		1.0	0.17	ug/L
108-86-1	Bromobenzene	0.14	U	1.0	0.14	ug/L
98-82-8	Isopropylbenzene	0.14	U	1.0	0.14	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	0.18	U	1.0	0.18	ug/L
96-18-4	1,2,3-Trichloropropane	0.20	U	1.0	0.20	ug/L
103-65-1	N-propylbenzene	0.14	U	1.0	0.14	ug/L
95-49-8	2-Chlorotoluene	0.11	U	1.0	0.11	ug/L
108-67-8	1,3,5-Trimethylbenzene	0.15	U	1.0	0.15	ug/L

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	DIST	SDG No.:	T6267
Lab Sample ID:	T6267-04	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000558.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
106-43-4	4-Chlorotoluene	0.15	U	1.0	0.15	ug/L
98-06-6	tert-Butylbenzene	0.15	U	1.0	0.15	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.15	U	1.0	0.15	ug/L
135-98-8	Sec-butylbenzene	0.14	U	1.0	0.14	ug/L
99-87-6	p-Isopropyltoluene	0.14	U	1.0	0.14	ug/L
541-73-1	1,3-Dichlorobenzene	0.15	U	1.0	0.15	ug/L
106-46-7	1,4-Dichlorobenzene	0.17	U	1.0	0.17	ug/L
104-51-8	n-Butylbenzene	0.12	U	1.0	0.12	ug/L
95-50-1	1,2-Dichlorobenzene	0.16	U	1.0	0.16	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	0.19	U	1.0	0.19	ug/L
120-82-1	1,2,4-Trichlorobenzene	0.11	U	1.0	0.11	ug/L
87-68-3	Hexachlorobutadiene	0.13	U	1.0	0.13	ug/L
91-20-3	Naphthalene	0.14	U	1.0	0.14	ug/L
87-61-6	1,2,3-Trichlorobenzene	0.16	U	1.0	0.16	ug/L

### SURROGATES

2199-69-1	1,2-Dichlorobenzene-d4	1.03	103 %	80 - 120	SPK: 1
460-00-4	4-Bromofluorobenzene	1.01	101 %	80 - 120	SPK: 1

### INTERNAL STANDARDS

462-06-6	Fluorobenzene	558047	9.12
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U = Not Detected

J = Estimated Value

RL = Reporting Limit

B = Analyte Found in Associated Method Blank

MDL = Method Detection Limit

N = Presumptive Evidence of a Compound

E = Value Exceeds Calibration Range



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	STEFF	SDG No.:	T6267
Lab Sample ID:	T6267-05	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000559.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
<b>TARGETS</b>						
75-71-8	Dichlorodifluoromethane	0.06	U	1.0	0.06	ug/L
74-87-3	Chloromethane	0.07	U	1.0	0.07	ug/L
75-01-4	Vinyl Chloride	0.07	U	1.0	0.07	ug/L
74-83-9	Bromomethane	0.23	U	1.0	0.23	ug/L
75-00-3	Chloroethane	0.17	U	1.0	0.17	ug/L
75-69-4	Trichlorofluoromethane	0.09	U	1.0	0.09	ug/L
75-65-0	tert-Butyl Alcohol	2.9	U	10	2.9	ug/L
60-29-7	Diethyl Ether	0.16	U	1.0	0.16	ug/L
75-35-4	1,1-Dichloroethene	0.14	U	1.0	0.14	ug/L
74-88-4	Iodomethane	0.08	U	1.0	0.08	ug/L
107-5-1	Allyl Chloride	0.15	U	1.0	0.15	ug/L
107-13-1	Acrylonitrile	0.46	U	2.0	0.46	ug/L
67-64-1	Acetone	1.1	U	5.8	1.1	ug/L
75-15-0	Carbon disulfide	0.14	U	1.0	0.14	ug/L
1634-04-4	Methyl tert-butyl Ether	0.15	U	1.0	0.15	ug/L
79-20-9	Methyl acrylate	0.16	U	1.0	0.16	ug/L
75-09-2	Methylene Chloride	1.0		1.0	0.27	ug/L
156-60-5	trans-1,2-Dichloroethene	0.14	U	1.0	0.14	ug/L
75-34-3	1,1-Dichloroethane	0.16	U	1.0	0.16	ug/L
78-93-3	2-Butanone	0.99	U	5.0	0.99	ug/L
56-23-5	Carbon Tetrachloride	0.15	U	1.0	0.15	ug/L
594-20-7	2,2-Dichloropropane	0.19	U	1.0	0.19	ug/L
156-59-2	cis-1,2-Dichloroethene	0.12	U	1.0	0.12	ug/L
67-66-3	Chloroform	0.16	U	1.0	0.16	ug/L
71-55-6	1,1,1-Trichloroethane	0.14	U	1.0	0.14	ug/L
110-57-6	t-1,4-Dichloro-2-butene	0.45	U	2.0	0.45	ug/L
563-58-6	1,1-Dichloropropene	0.16	U	1.0	0.16	ug/L
108-20-3	Isopropyl Ether	0.18	U	1.0	0.18	ug/L
107-12-0	Propionitrile	1.7	U	10	1.7	ug/L
71-43-2	Benzene	0.14	U	1.0	0.14	ug/L
107-06-2	1,2-Dichloroethane	0.21	U	1.0	0.21	ug/L
79-01-6	Trichloroethene	0.15	U	1.0	0.15	ug/L

U = Not Detected

J = Estimated Value

RL = Reporting Limit

B = Analyte Found in Associated Method Blank

MDL = Method Detection Limit

N = Presumptive Evidence of a Compound

E = Value Exceeds Calibration Range



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	STEFF	SDG No.:	T6267
Lab Sample ID:	T6267-05	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000559.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
78-87-5	1,2-Dichloropropane	0.14	U	1.0	0.14	ug/L
126-98-7	Methacrylonitrile	0.62	U	1.0	0.62	ug/L
109-99-9	Tetrahydrofuran	0.45	U	2.4	0.45	ug/L
109-69-3	1-Chlorobutane	0.17	U	1.0	0.17	ug/L
74-95-3	Dibromomethane	0.19	U	1.0	0.19	ug/L
75-27-4	Bromodichloromethane	0.17	U	1.0	0.17	ug/L
108-10-1	4-Methyl-2-Pentanone	0.90	U	5.0	0.90	ug/L
80-62-6	Methyl methacrylate	0.32	U	2.0	0.32	ug/L
97-63-2	Ethyl methacrylate	0.16	U	1.0	0.16	ug/L
108-88-3	Toluene	0.13	U	1.0	0.13	ug/L
10061-02-6	t-1,3-Dichloropropene	0.14	U	1.0	0.14	ug/L
10061-01-5	cis-1,3-Dichloropropene	0.13	U	1.0	0.13	ug/L
79-00-5	1,1,2-Trichloroethane	0.18	U	1.0	0.18	ug/L
142-28-9	1,3-Dichloropropane	0.14	U	1.0	0.14	ug/L
591-78-6	2-Hexanone	0.81	U	5.0	0.81	ug/L
124-48-1	Dibromochloromethane	0.17	U	1.0	0.17	ug/L
106-93-4	1,2-Dibromoethane	0.17	U	1.0	0.17	ug/L
127-18-4	Tetrachloroethene	0.16	U	1.0	0.16	ug/L
108-90-7	Chlorobenzene	0.13	U	1.0	0.13	ug/L
630-20-6	1,1,1,2-Tetrachloroethane	0.17	U	1.0	0.17	ug/L
67-72-1	Hexachloroethane	0.17	U	1.0	0.17	ug/L
100-41-4	Ethyl Benzene	0.14	U	1.0	0.14	ug/L
126777-61-2	m/p-Xylenes	0.29	U	1.0	0.29	ug/L
95-47-6	o-Xylene	0.15	U	1.0	0.15	ug/L
100-42-5	Styrene	0.14	U	1.0	0.14	ug/L
75-25-2	Bromoform	0.17	U	1.0	0.17	ug/L
108-86-1	Bromobenzene	0.14	U	1.0	0.14	ug/L
98-82-8	Isopropylbenzene	0.14	U	1.0	0.14	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	0.18	U	1.0	0.18	ug/L
96-18-4	1,2,3-Trichloropropane	0.20	U	1.0	0.20	ug/L
103-65-1	N-propylbenzene	0.14	U	1.0	0.14	ug/L
95-49-8	2-Chlorotoluene	0.11	U	1.0	0.11	ug/L
108-67-8	1,3,5-Trimethylbenzene	0.15	U	1.0	0.15	ug/L

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	STEFF	SDG No.:	T6267
Lab Sample ID:	T6267-05	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000559.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
106-43-4	4-Chlorotoluene	0.15	U	1.0	0.15	ug/L
98-06-6	tert-Butylbenzene	0.15	U	1.0	0.15	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.15	U	1.0	0.15	ug/L
135-98-8	Sec-butylbenzene	0.14	U	1.0	0.14	ug/L
99-87-6	p-Isopropyltoluene	0.14	U	1.0	0.14	ug/L
541-73-1	1,3-Dichlorobenzene	0.15	U	1.0	0.15	ug/L
106-46-7	1,4-Dichlorobenzene	0.17	U	1.0	0.17	ug/L
104-51-8	n-Butylbenzene	0.12	U	1.0	0.12	ug/L
95-50-1	1,2-Dichlorobenzene	0.16	U	1.0	0.16	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	0.19	U	1.0	0.19	ug/L
120-82-1	1,2,4-Trichlorobenzene	0.11	U	1.0	0.11	ug/L
87-68-3	Hexachlorobutadiene	0.13	U	1.0	0.13	ug/L
91-20-3	Naphthalene	0.14	U	1.0	0.14	ug/L
87-61-6	1,2,3-Trichlorobenzene	0.16	U	1.0	0.16	ug/L
<b>SURROGATES</b>						
2199-69-1	1,2-Dichlorobenzene-d4	1.1	110 %	80 - 120	SPK:	1
460-00-4	4-Bromofluorobenzene	0.96	96 %	80 - 120	SPK:	1
<b>INTERNAL STANDARDS</b>						
462-06-6	Fluorobenzene		554298	9.11		

U = Not Detected

J = Estimated Value

RL = Reporting Limit

B = Analyte Found in Associated Method Blank

MDL = Method Detection Limit

N = Presumptive Evidence of a Compound

E = Value Exceeds Calibration Range



284 Sheffield Street, Mountainside, NJ 07042 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	DUP	SDG No.:	T6267
Lab Sample ID:	T6267-06	Matrix:	WATER
Analytical Method:	S24.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000560.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
<b>TARGETS</b>						
75-71-8	Dichlorodifluoromethane	0.06	U	1.0	0.06	ug/L
74-87-3	Chloromethane	0.07	U	1.0	0.07	ug/L
75-01-4	Vinyl Chloride	0.07	U	1.0	0.07	ug/L
74-83-9	Bromomethane	0.23	U	1.0	0.23	ug/L
75-00-3	Chloroethane	0.17	U	1.0	0.17	ug/L
75-69-4	Trichlorofluoromethane	0.09	U	1.0	0.09	ug/L
75-65-0	tert-Butyl Alcohol	2.9	U	10	2.9	ug/L
60-29-7	Diethyl Ether	0.16	U	1.0	0.16	ug/L
75-35-4	1,1-Dichloroethene	0.14	U	1.0	0.14	ug/L
74-88-4	Iodomethane	0.08	U	1.0	0.08	ug/L
107-5-1	Allyl Chloride	0.15	U	1.0	0.15	ug/L
107-13-1	Acrylonitrile	0.46	U	2.0	0.46	ug/L
67-64-1	Acetone	1.1	U	5.8	1.1	ug/L
75-15-0	Carbon disulfide	0.14	U	1.0	0.14	ug/L
1634-04-4	Methyl tert-butyl Ether	0.15	U	1.0	0.15	ug/L
79-20-9	Methyl acrylate	0.16	U	1.0	0.16	ug/L
75-09-2	Methylene Chloride	0.27	U	1.0	0.27	ug/L
156-60-5	trans-1,2-Dichloroethene	0.14	U	1.0	0.14	ug/L
75-34-3	1,1-Dichloroethane	0.16	U	1.0	0.16	ug/L
78-93-3	2-Butanone	0.99	U	5.0	0.99	ug/L
56-23-5	Carbon Tetrachloride	0.15	U	1.0	0.15	ug/L
594-20-7	2,2-Dichloropropane	0.19	U	1.0	0.19	ug/L
156-59-2	cis-1,2-Dichloroethene	0.8	J	1.0	0.12	ug/L
67-66-3	Chloroform	0.16	U	1.0	0.16	ug/L
71-55-6	1,1,1-Trichloroethane	0.14	U	1.0	0.14	ug/L
110-57-6	t-1,4-Dichloro-2-butene	0.45	U	2.0	0.45	ug/L
563-58-6	1,1-Dichloropropene	0.16	U	1.0	0.16	ug/L
108-20-3	Isopropyl Ether	0.18	U	1.0	0.18	ug/L
107-12-0	Propionitrile	1.7	U	10	1.7	ug/L
71-43-2	Benzene	0.14	U	1.0	0.14	ug/L
107-06-2	1,2-Dichloroethane	0.21	U	1.0	0.21	ug/L
79-01-6	Trichloroethene	0.8	J	1.0	0.15	ug/L

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	DUP	SDG No.:	T6267
Lab Sample ID:	T6267-06	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000560.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
78-87-5	1,2-Dichloropropane	0.14	U	1.0	0.14	ug/L
126-98-7	Methacrylonitrile	0.62	U	1.0	0.62	ug/L
109-99-9	Tetrahydrofuran	0.45	U	2.4	0.45	ug/L
109-69-3	1-Chlorobutane	0.17	U	1.0	0.17	ug/L
74-95-3	Dibromomethane	0.19	U	1.0	0.19	ug/L
75-27-4	Bromodichloromethane	0.17	U	1.0	0.17	ug/L
108-10-1	4-Methyl-2-Pentanone	0.90	U	5.0	0.90	ug/L
80-62-6	Methyl methacrylate	0.32	U	2.0	0.32	ug/L
97-63-2	Ethyl methacrylate	0.16	U	1.0	0.16	ug/L
108-88-3	Toluene	0.13	U	1.0	0.13	ug/L
10061-02-6	t-1,3-Dichloropropene	0.14	U	1.0	0.14	ug/L
10061-01-5	cis-1,3-Dichloropropene	0.13	U	1.0	0.13	ug/L
79-00-5	1,1,2-Trichloroethane	0.18	U	1.0	0.18	ug/L
142-28-9	1,3-Dichloropropane	0.14	U	1.0	0.14	ug/L
591-78-6	2-Hexanone	0.81	U	5.0	0.81	ug/L
124-48-1	Dibromochloromethane	0.17	U	1.0	0.17	ug/L
106-93-4	1,2-Dibromoethane	0.17	U	1.0	0.17	ug/L
127-18-4	Tetrachloroethene	29		1.0	0.16	ug/L
108-90-7	Chlorobenzene	0.13	U	1.0	0.13	ug/L
630-20-6	1,1,1,2-Tetrachloroethane	0.17	U	1.0	0.17	ug/L
67-72-1	Hexachloroethane	0.17	U	1.0	0.17	ug/L
100-41-4	Ethyl Benzene	0.14	U	1.0	0.14	ug/L
126777-61-2	m/p-Xylenes	0.29	U	1.0	0.29	ug/L
95-47-6	o-Xylene	0.15	U	1.0	0.15	ug/L
100-42-5	Styrene	0.14	U	1.0	0.14	ug/L
75-25-2	Bromoform	0.17	U	1.0	0.17	ug/L
108-86-1	Bromobenzene	0.14	U	1.0	0.14	ug/L
98-82-8	Isopropylbenzene	0.14	U	1.0	0.14	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	0.18	U	1.0	0.18	ug/L
96-18-4	1,2,3-Trichloropropane	0.20	U	1.0	0.20	ug/L
103-65-1	N-propylbenzene	0.14	U	1.0	0.14	ug/L
95-49-8	2-Chlorotoluene	0.11	U	1.0	0.11	ug/L
108-67-8	1,3,5-Trimethylbenzene	0.15	U	1.0	0.15	ug/L

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound



284 Sheffield Street, Mountainside, NJ 07092 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	DUP	SDG No.:	T6267
Lab Sample ID:	T6267-06	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000560.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
106-43-4	4-Chlorotoluene	0.15	U	1.0	0.15	ug/L
98-06-6	tert-Butylbenzene	0.15	U	1.0	0.15	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.15	U	1.0	0.15	ug/L
135-98-8	Sec-butylbenzene	0.14	U	1.0	0.14	ug/L
99-87-6	p-Isopropyltoluene	0.14	U	1.0	0.14	ug/L
541-73-1	1,3-Dichlorobenzene	0.15	U	1.0	0.15	ug/L
106-46-7	1,4-Dichlorobenzene	0.17	U	1.0	0.17	ug/L
104-51-8	n-Butylbenzene	0.12	U	1.0	0.12	ug/L
95-50-1	1,2-Dichlorobenzene	0.16	U	1.0	0.16	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	0.19	U	1.0	0.19	ug/L
120-82-1	1,2,4-Trichlorobenzene	0.11	U	1.0	0.11	ug/L
87-68-3	Hexachlorobutadiene	0.13	U	1.0	0.13	ug/L
91-20-3	Naphthalene	0.14	U	1.0	0.14	ug/L
87-61-6	1,2,3-Trichlorobenzene	0.16	U	1.0	0.16	ug/L

**SURROGATES**

2199-69-1	1,2-Dichlorobenzene-d4	1.07	107 %	80 - 120	SPK: 1
460-00-4	4-Bromofluorobenzene	0.97	97 %	80 - 120	SPK: 1

**INTERNAL STANDARDS**

462-06-6	Fluorobenzene	446780	9.11
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U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound







284 Sheffield Street, Mountainside, NJ 07042 Phone: 908-789-8900 Fax: 908-789-8922

## Report of Analysis

Client:	EPM, INC.	Date Collected:	12/20/2005
Project:	Katonah	Date Received:	12/21/2005
Client Sample ID:	TB	SDG No.:	T6267
Lab Sample ID:	T6267-07	Matrix:	WATER
Analytical Method:	524.2 Rev3	% Moisture:	100
Sample Wt/Wt:	25.0 Units: mL	Soil Extract Vol:	uL
Soil Aliquot Vol:	uL		

File ID:	Dilution:	Date Analyzed	Analytical Batch ID
VF000556.D	1	12/27/2005	VF122305

CAS Number	Parameter	Conc.	Qualifier	RL	MDL	Units
106-43-4	4-Chlorotoluene	0.15	U	1.0	0.15	ug/L
98-06-6	tert-Butylbenzene	0.15	U	1.0	0.15	ug/L
95-63-6	1,2,4-Trimethylbenzene	0.15	U	1.0	0.15	ug/L
135-98-8	Sec-butylbenzene	0.14	U	1.0	0.14	ug/L
99-87-6	p-Isopropyltoluene	0.14	U	1.0	0.14	ug/L
541-73-1	1,3-Dichlorobenzene	0.15	U	1.0	0.15	ug/L
106-46-7	1,4-Dichlorobenzene	0.17	U	1.0	0.17	ug/L
104-51-8	n-Butylbenzene	0.12	U	1.0	0.12	ug/L
95-50-1	1,2-Dichlorobenzene	0.16	U	1.0	0.16	ug/L
96-12-8	1,2-Dibromo-3-Chloropropane	0.19	U	1.0	0.19	ug/L
120-82-1	1,2,4-Trichlorobenzene	0.11	U	1.0	0.11	ug/L
87-68-3	Hexachlorobutadiene	0.13	U	1.0	0.13	ug/L
91-20-3	Naphthalene	0.14	U	1.0	0.14	ug/L
87-61-6	1,2,3-Trichlorobenzene	0.16	U	1.0	0.16	ug/L

### SURROGATES

2199-69-1	1,2-Dichlorobenzene-d4	1.09	109 %	80 - 120	SPK: 1
460-00-4	4-Bromofluorobenzene	1.01	101 %	80 - 120	SPK: 1

### INTERNAL STANDARDS

462-06-6	Fluorobenzene	579806	9.12
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