

# **ANALYTICAL RESULTS SUMMARY**

VOLATILE ORGANICS GENERAL CHEMISTRY METALS SEMI-VOLATILE ORGANICS

### **PROJECT NAME : FORMER SCHLUMBERGER SITE PRINCETON NJ**

# JACOBS ENGINEERING GROUP, INC.

412 Mt. Kemble Ave

**Downtown Building** 

Morristown, NJ - 07960

Phone No: 9732670555

ORDER ID : P3657 ATTENTION : Mary I. Murphy



Laboratory Certification ID # 20012







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# DATA OF KNOWN QUALITY CONFORMANCE/NON-CONFORMANCE SUMMARY QUESTIONNAIRE

1

Laboratory Name :	Alliance Technical Group LLC	Client :	JACOBS Engineering Group, Inc.
Project Location :	Princeton Junction, NJ	Project Number :	D3779922
Laboratory Sample ID(	s) : <u>P3657</u>	Sampling Date(s) :	8/16/2024

List DKQP Methods Used (e.g., 8260,8270, et Cetra) 6020B,7196A,7470A,8260-Low,8270-Modified,8270E

1	For each analytical method referenced in this laboratory report package, were all specified QA/QC performance criteria followed, including the requirement to explain any criteria falling outside of acceptable guidelines, as specified in the NJDEP Data of Known Quality performance standards?	V	Yes		No	
1A	Were the method specified handling, preservation, and holding time requirements met?	V	Yes		No	
1B	EPH Method: Was the EPH method conducted without significant modifications (see Section 11.3 of respective DKQ methods)		Yes		No	<b>M</b> /A
2	Were all samples received by the laboratory in a condition consistent with that described on the associated chain-of-custody document(s)?	$\mathbf{\nabla}$	Yes		No	
3	Were samples received at an appropriate temperature (4±2° C)?	Ø	Yes		No	□ N/A
4	Were all QA/QC performance criteria specified in the NJDEP DKQP standards achieved?		Yes	V	No	
5	a)Were reporting limits specified or referenced on the chain-of-custody or communicated to the laboratory prior to sample receipt?	Q	Yes		No	
	b)Were these reporting limits met?	$\square$	Yes		No	□ N/A
6	For each analytical method referenced in this laboratory report package, were results reported for all constituents identified in the method-specific analyte lists presented in the DKQP documents and/or site-specific QAPP?	V	Yes		No	
7	Are project-specific matrix spikes and/or laboratory duplicates included in this data set?		Yes	V	No	

Notes: For all questions to which the response was "No" (with the exception of question #7), additional information should be provided in an attached narrative. If the answer to question #1, #1A, or #1B is "No", the data package does not meet the requirements for "Data of Known Quality."



# **Cover Page**

- Order ID : P3657
- Project ID : Former Schlumberger Site Princeton NJ
  - **Client :** JACOBS Engineering Group, Inc.

#### Lab Sample Number

Client Sample Number

P3657-01 P3657-02 917-J-WS-081624 TB-01-081624

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hard copy data package has been authorized by the laboratory manager or his designee, as verified by the following signature.

Signature :

Date: 9/5/2024

NYDOH CERTIFICATION NO - 11376

NJDEP CERTIFICATION NO - 20012



# CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3657 Test Name: VOCMS Group6

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 08/16/2024.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOCMS Group3, SVOCMS Group6 and VOCMS Group6. This data package contains results for VOCMS Group6.

#### **C. Analytical Techniques:**

The analysis performed on instrument MSVOA\_N were done using GC column RXI-624SIL MS 30m 0.25mm 1.4 um. Cat#13868.The analysis of VOCMS Group6 was based on method 8260D.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria.

The Internal Standards Areas met the acceptable requirements.

The Retention Times were acceptable for all samples.

The RPD met criteria.

The Blank Spike met requirements for all samples .

The Blank Spike Duplicate met requirements for all samples .

The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements .

The Continuous Calibration met the requirements .

The Tuning criteria met requirements.

#### **E. Additional Comments:**

Samples for MS/MSD for VOC analysis were not provided with this set of samples. The Blank Spike Duplicate is reported with the data.

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <15% for the Initial Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is > 15% for the Initial Calibration curve for SW-846 analysis.

#### **F. Manual Integration Comments:**



Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. The laboratory manager or his designee, as verified by the following signature has authorized release of the data contained in this hard copy data package.

Signature\_\_\_\_\_

2.1



# CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3657 Test Name: SVOCMS Group3

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 08/16/2024.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOCMS Group3, SVOCMS Group6 and VOCMS Group6. This data package contains results for SVOCMS Group3.

#### **C. Analytical Techniques:**

The samples were analyzed on instrument BNA\_N using GC Column ZB-SemiVolatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGAThe analysis of SVOCMS Group3 was based on method 8270-Modified and extraction was done based on method 3510.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria.

The Internal Standards Areas met the acceptable requirements.

The Retention Times were acceptable for all samples.

The RPD met criteria.

The Blank Spike met requirements for all samples .

The Blank Spike Duplicate met requirements for all samples .

The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements .

The Continuous Calibration met the requirements .

The Tuning criteria met requirements.

#### **E. Additional Comments:**

The Form 6 is not included in the data package because the Initial Calibration was performed using 7 points.

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <20% for the Initial Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is > 20% for the Initial Calibration curve for SW-846 analysis.



For sample # 917-J-WS-081624 some compounds below Method detection limits, therefore it is not reported as Hit in Form-1.

#### **F. Manual Integration Comments:**

Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

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Signature\_\_\_\_\_

2.2



# CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3657 Test Name: SVOCMS Group6

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 08/16/2024.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOCMS Group3, SVOCMS Group6 and VOCMS Group6. This data package contains results for SVOCMS Group6.

#### **C. Analytical Techniques:**

The samples were analyzed on instrument BNA\_P using GC Column ZB-SemiVolatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGAThe analysis of SVOCMS Group6 was based on method 8270E and extraction was done based on method 3510.

#### **D. QA/ QC Samples:**

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria except for 917-J-WS-081624 [2,4 and6-Tribromophenol - 119%]. these compound did not meet the NJDKQP criteria but met the in-house criteria, Therefor no corrective action was required.

The Internal Standards Areas met the acceptable requirements. The Retention Times were acceptable for all samples.

The RPD for {PB162822BSD} with File ID: BP021569.D met criteria except for Benzaldehyde[24%], Due to result difference between BS and BSD, therefor no corrective action was required.

The Blank Spike met requirements for all samples . The Blank Spike Duplicate met requirements for all samples . The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements.



The Continuous Calibration File ID BP021566.D met the requirements except for Pentachlorophenol, Failed high side and a sample does not have hit for this compound, therefor no corrective action was required.

The Continuous Calibration File ID BP021575.D met the requirements except for Pentachlorophenol . Failed high side and samples does not have hit for this compound, Therefor no corrective action was required. The Tuning criteria met requirements.

#### **E. Additional Comments:**

The Form 6 is not included in the data package because the Initial Calibration was performed using 7 points.

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <15% for the Initial Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is > 15% for the Initial Calibration curve for SW-846 analysis.

#### **F. Manual Integration Comments:**

Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

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Signature\_\_\_\_\_



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

# CASE NARRATIVE

24

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3657 Test Name: Metals Group4,Mercury

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 08/16/2024.

#### **B.** Parameters:

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOCMS Group3, SVOCMS Group6 and VOCMS Group6. This data package contains results for Metals Group4,Mercury.

#### **C. Analytical Techniques:**

The analysis of Metals Group4 was based on method 6020B, digestion based on method 3010 (waters). The analysis and digestion of Mercury was based on method 7470A.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Blank Spike met requirements for all samples.

The Duplicate (918-J-WS-081324-FDDUP) analysis met criteria for all samples except for Arsenic due to sample matrix interference.

The Matrix Spike (1027MS) analysis met criteria for all samples except for Mercury due to sample matrix interference. The Matrix Spike (918-J-WS-081324-FDMS) analysis met criteria for all samples except for Molybdenum and Silver due to Chemical Interference during Digestion Process.

The Matrix Spike Duplicate (1027MSD) analysis met criteria for all samples except for Mercury due to sample matrix interference. The Matrix Spike Duplicate (918-J-WS-081324-FDMSD) analysis met criteria for all samples except for Molybdenum and Silver due to Chemical Interference during Digestion Process.

The Blank analysis did not indicate the presence of lab contamination.

The Calibration met the requirements.

The Serial Dilution (918-J-WS-081324-FDL) met criteria for all samples except for Aluminum, Iron, and Manganese due to sample matrix interference.

#### **E. Additional Comments:**

Collision cell is being used to remove potential interferences. The analytes Na, Mg, Al, K, V, Cr, Mn, Fe, Co, Ni, Cu, Zn, As are being analyzed with collision cell and analytes Be, B, Ca, Ti, Se, Sr, Zr, Mo, Ag, Cd, Sn, Sb, Ba, Tl, Pb, U are being analyzed with Non-Collision Cell. Helium gas is used for the Collision Cell analysis.



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Signature\_\_\_\_\_



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

# CASE NARRATIVE

25

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3657 Test Name: Hexavalent Chromium

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 08/16/2024.

#### **B.** Parameters:

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOCMS Group3, SVOCMS Group6 and VOCMS Group6. This data package contains results for Hexavalent Chromium.

#### **C. Analytical Techniques:**

The analysis of Hexavalent Chromium was based on method 7196A.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis. The Blank Spike met requirements for all samples. The Duplicate analysis met criteria for all samples. The Matrix Spike analysis met criteria for all samples. The Matrix Spike Duplicate analysis met criteria for all samples. The Blank analysis did not indicate the presence of lab contamination. The Calibration met the requirements.

#### **E. Additional Comments:**

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. The laboratory manager or his designee, as verified by the following signature has authorized release of the data contained in this hard copy data package.

Signature\_\_\_\_\_



# DATA REPORTING QUALIFIERS- INORGANIC

For reporting results, the following " Results Qualifiers" are used:

J	Indicates the reported value was obtained from a reading that was less than the Contract Required Detection Limit (CRDL), but greater than or equal to the Instrument Detection Limit (IDL).
U	Indicates the analyte was analyzed for, but not detected.
ND	Indicates the analyte was analyzed for, but not detected
Ε	Indicates the reported value is estimated because of the presence of interference
Μ	Indicates Duplicate injection precision not met.
Ν	Indicates the spiked sample recovery is not within control limits.
S	Indicates the reported value was determined by the Method of Standard Addition (MSA).
*	Indicates that the duplicate analysis is not within control limits.
+	Indicates the correlation coefficient for the MSA is less than 0.995.
D	Indicates the reported value is from a secondary analysis with a dilution factor. The original analysis exceeded the calibration range.
M OR	<ul> <li>Method qualifiers</li> <li>"P" for ICP instrument</li> <li>"PM" for ICP when Microwave Digestion is used</li> <li>"CV" for Manual Cold Vapor AA</li> <li>"AV" for automated Cold Vapor AA</li> <li>"CA" for MIDI-Distillation Spectrophotometric</li> <li>"AS" for Semi – Automated Spectrophotometric</li> <li>"C" for Manual Spectrophotometric</li> <li>"T" for Titrimetric</li> <li>"NR" for analyte not required to be analyzed</li> <li>Indicates the analyte's concentration exceeds the calibrated range of the instrument for that specific analysis.</li> </ul>
Q	Indicates the LCS did not meet the control limits requirements
Н	Sample Analysis Out Of Hold Time



# DATA REPORTING QUALIFIERS- ORGANIC

For reporting results, the following " Results Qualifiers" are used:

Value	If the result is a value greater than or equal to the detection limit, report the value
U	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. "10 U". This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
ND	Indicates the analyte was analyzed for, but not detected
J	<ul> <li>Indicates an estimated value. This flag is used:</li> <li>(1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.)</li> <li>(2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3 ug/L was calculated report as 3 J. This is flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others.</li> </ul>
В	Indicates the analyte was found in the blank as well as the sample report as "12 B".
Ε	Indicates the analyte 's concentration exceeds the calibrated range of the instrument for that specific analysis.
D	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
Р	This flag is used for Pesticide/PCB target analyte when there is >25% difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form 1 and flagged with a "P".
Ν	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
Α	This flag indicates that a Tentatively Identified Compound is a suspected aldol- condensation product.
Q	Indicates the LCS did not meet the control limits requirements



#### APPENDIX A

#### **QA REVIEW GENERAL DOCUMENTATION**

Project #: P3657

Completed

For thorough review, the report must have the following:	
GENERAL:	
Are all original paperwork present (chain of custody, record of communication,airbill, sample management lab chronicle, login page)	<u> </u>
Check chain-of-custody for proper relinquish/return of samples	<u> </u>
Is the chain of custody signed and complete	✓ ✓ ✓ ✓
Check internal chain-of-custody for proper relinquish/return of samples /sample extracts	<u>✓</u>
Collect information for each project id from server. Were all requirements followed	<u>✓</u>
COVER PAGE:	
Do numbers of samples correspond to the number of samples in the Chain of Custody on login page	<u>✓</u>
Do lab numbers and client Ids on cover page agree with the Chain of Custody	<u>✓</u>
CHAIN OF CUSTODY:	
Do requested analyses on Chain of Custody agree with form I results	✓
Do requested analyses on Chain of Custody agree with the log-in page	✓
Were the correct method log-in for analysis according to the Analytical Request and Chain of Castody	✓ ✓ ✓
Were the samples received within hold time	<u>✓</u>
Were any problems found with the samples at arrival recorded in the Sample Management Laboratory Chronicle	<u> </u>
ANALYTICAL:	
Was method requirement followed?	<u>✓</u>
Was client requirement followed?	<u> </u>
Does the case narrative summarize all QC failure?	✓ ✓ ✓ ✓ ✓
All runlogs and manual integration are reviewed for requirements	✓
All manual calculations and /or hand notations verified	<u> </u>

1st Level QA Review Signature:

SOHIL JODHANI

Date: 09/05/2024

2nd Level QA Review Signature:

Date:



#### Hit Summary Sheet SW-846

				5 W-040				
SDG No.:	P3657							В
Client:	JACOBS Engineer	ring Group, In	c.					С
								D
Sample ID	Client ID	Matrix	Parameter	Concentration	C MDL	RDL	Units	
Client ID:	917-J-WS-081624							
P3657-01	917-J-WS-081624	Water	Acetone	7.30	1.40	5.00	ug/L	
			Total Voc :	7.30				

Total Concentration: 7.30







A B C D



Date Collected:

Test:

Level :

08/16/24

VOCMS Group6

LOW

uL

# 5

Former Schlumberger Site Princeton NJ	Date Received:	08/16/24
917-J-WS-081624	SDG No.:	P3657
P3657-01	Matrix:	Water
SW8260	% Solid:	0
5 Units: mL	Final Vol:	5000

**Report of Analysis** 

JACOBS Engineering Group, Inc.

RXI-624

uL

ID: 0.25

GC Column: Prep Method :

Soil Aliquot Vol:

Client:

Project:

Client Sample ID: Lab Sample ID: Analytical Method: Sample Wt/Vol:

File ID/Qc Batch:	Dilution:	Prep Date		Date Analyzed	Prep Batch ID	
VN083379.D	1			08/19/24 17:53	VN081924	
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	0.21	U	0.21	1.00	ug/L
74-87-3	Chloromethane	0.35	U	0.35	1.00	ug/L
75-01-4	Vinyl Chloride	0.34	U	0.34	1.00	ug/L
74-83-9	Bromomethane	1.40	U	1.40	5.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	0.25	1.00	ug/L
67-64-1	Acetone	7.30		1.40	5.00	ug/L
75-15-0	Carbon Disulfide	0.32	U	0.32	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	0.16	U	0.16	1.00	ug/L
75-09-2	Methylene Chloride	0.32	U	0.32	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	0.25	U	0.25	1.00	ug/L
110-82-7	Cyclohexane	1.60	U	1.60	5.00	ug/L
78-93-3	2-Butanone	1.30	U	1.30	5.00	ug/L
56-23-5	Carbon Tetrachloride	0.25	U	0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	0.25	U	0.25	1.00	ug/L
67-66-3	Chloroform	0.26	U	0.26	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	0.19	U	0.19	1.00	ug/L
108-87-2	Methylcyclohexane	0.19	U	0.19	1.00	ug/L
71-43-2	Benzene	0.16	U	0.16	1.00	ug/L
107-06-2	1,2-Dichloroethane	0.24	U	0.24	1.00	ug/L
79-01-6	Trichloroethene	0.32	U	0.32	1.00	ug/L
75-27-4	Bromodichloromethane	0.24	U	0.24	1.00	ug/L
108-88-3	Toluene	0.18	U	0.18	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	0.21	U	0.21	1.00	ug/L
124-48-1	Dibromochloromethane	0.18	U	0.18	1.00	ug/L
127-18-4	Tetrachloroethene	0.25	U	0.25	1.00	ug/L
108-90-7	Chlorobenzene	0.13	U	0.13	1.00	ug/L
100-41-4	Ethyl Benzene	0.16	U	0.16	1.00	ug/L
179601-23-1	m/p-Xylenes	0.31	U	0.31	2.00	ug/L
1330-20-7	Total Xylenes	0.45	U	0.45	3.00	ug/L
95-47-6	o-Xylene	0.14	U	0.14	1.00	ug/L



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/16/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/16/24
Client Sample ID:	917-J-WS-081624	SDG No.:	P3657
Lab Sample ID:	P3657-01	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group6
GC Column:	RXI-624 ID: 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date		Date Analyzed	Prep Batch ID	
VN083379.D	1			08/19/24 17:53	VN081924	
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
98-82-8	Isopropylbenzene	0.13	U	0.13	1.00	ug/L
106-46-7	1,4-Dichlorobenzene	0.27	U	0.27	1.00	ug/L
95-50-1	1,2-Dichlorobenzene	0.19	U	0.19	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	53.3		70 (74) - 130 (125)	107%	SPK: 50
1868-53-7	Dibromofluoromethane	48.9		70 (75) - 130 (124)	98%	SPK: 50
2037-26-5	Toluene-d8	48.9		70 (86) - 130 (113)	98%	SPK: 50
460-00-4	4-Bromofluorobenzene	51.3		70 (77) - 130 (121)	103%	SPK: 50
INTERNAL STAN	DARDS					
363-72-4	Pentafluorobenzene	139000	8.224			
540-36-3	1,4-Difluorobenzene	273000	9.106			
3114-55-4	Chlorobenzene-d5	284000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	127000	13.794			

U = Not Detected

- LOQ = Limit of Quantitation
- MDL = Method Detection Limit
- LOD = Limit of Detection
- E = Value Exceeds Calibration Range
- Q = indicates LCS control criteria did not meet requirements
- M = MS/MSD acceptance criteria did not meet requirements

- J = Estimated Value
- B = Analyte Found in Associated Method Blank
- N = Presumptive Evidence of a Compound
- \* = Values outside of QC limits
- D = Dilution
- () = Laboratory InHouse Limit
- A = Aldol-Condensation Reaction Products

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Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/16/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/16/24
Client Sample ID:	TB-01-081624	SDG No.:	P3657
Lab Sample ID:	P3657-02	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group6
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Batch: Dilution: Prep Date Date Analyzed Prep Bate		Prep Batch ID			
VN083378.D	1			08/19/24 17:29	VN081924	
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	0.21	U	0.21	1.00	ug/L
74-87-3	Chloromethane	0.35	U	0.35	1.00	ug/L
75-01-4	Vinyl Chloride	0.34	U	0.34	1.00	ug/L
74-83-9	Bromomethane	1.40	U	1.40	5.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	0.25	1.00	ug/L
67-64-1	Acetone	1.40	U	1.40	5.00	ug/L
75-15-0	Carbon Disulfide	0.32	U	0.32	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	0.16	U	0.16	1.00	ug/L
75-09-2	Methylene Chloride	0.32	U	0.32	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	0.25	U	0.25	1.00	ug/L
110-82-7	Cyclohexane	1.60	U	1.60	5.00	ug/L
78-93-3	2-Butanone	1.30	U	1.30	5.00	ug/L
56-23-5	Carbon Tetrachloride	0.25	U	0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	0.25	U	0.25	1.00	ug/L
67-66-3	Chloroform	0.26	U	0.26	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	0.19	U	0.19	1.00	ug/L
108-87-2	Methylcyclohexane	0.19	U	0.19	1.00	ug/L
71-43-2	Benzene	0.16	U	0.16	1.00	ug/L
107-06-2	1,2-Dichloroethane	0.24	U	0.24	1.00	ug/L
79-01-6	Trichloroethene	0.32	U	0.32	1.00	ug/L
75-27-4	Bromodichloromethane	0.24	U	0.24	1.00	ug/L
108-88-3	Toluene	0.18	U	0.18	1.00	ug/L
79-00-5	1,1,2-Trichloroethane	0.21	U	0.21	1.00	ug/L
124-48-1	Dibromochloromethane	0.18	U	0.18	1.00	ug/L
127-18-4	Tetrachloroethene	0.25	U	0.25	1.00	ug/L
108-90-7	Chlorobenzene	0.13	Ū	0.13	1.00	ug/L
100-41-4	Ethyl Benzene	0.16	U	0.16	1.00	ug/L
179601-23-1	m/p-Xylenes	0.31	Ŭ	0.31	2.00	ug/L
1330-20-7	Total Xylenes	0.45	Ŭ	0.45	3.00	ug/L
95-47-6	o-Xylene	0.14	U	0.14	1.00	ug/L

B C

D

5

P3657



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/16/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/16/24
Client Sample ID:	TB-01-081624	SDG No.:	P3657
Lab Sample ID:	P3657-02	Matrix:	Water
Analytical Method:	SW8260	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group6
GC Column:	RXI-624 ID: 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date		Date Analyzed	Prep Batch ID	
VN083378.D	1			08/19/24 17:29	VN081924	
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
98-82-8	Isopropylbenzene	0.13	U	0.13	1.00	ug/L
106-46-7	1,4-Dichlorobenzene	0.27	U	0.27	1.00	ug/L
95-50-1	1,2-Dichlorobenzene	0.19	U	0.19	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	53.1		70 (74) - 130 (125)	106%	SPK: 50
1868-53-7	Dibromofluoromethane	48.6		70 (75) - 130 (124)	97%	SPK: 50
2037-26-5	Toluene-d8	46.7		70 (86) - 130 (113)	93%	SPK: 50
460-00-4	4-Bromofluorobenzene	49.7		70 (77) - 130 (121)	99%	SPK: 50
INTERNAL STAN	DARDS					
363-72-4	Pentafluorobenzene	134000	8.224			
540-36-3	1,4-Difluorobenzene	268000	9.106			
3114-55-4	Chlorobenzene-d5	272000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	120000	13.794			

U = Not Detected

- LOQ = Limit of Quantitation
- MDL = Method Detection Limit
- LOD = Limit of Detection
- E = Value Exceeds Calibration Range
- Q = indicates LCS control criteria did not meet requirements
- M = MS/MSD acceptance criteria did not meet requirements

- J = Estimated Value
- B = Analyte Found in Associated Method Blank
- N = Presumptive Evidence of a Compound
- \* = Values outside of QC limits
- D = Dilution
- () = Laboratory InHouse Limit
- A = Aldol-Condensation Reaction Products

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С



D

# LAB CHRONICLE

OrderID: Client: Contact:	Client: JACOBS Engineering Group, Inc.				8/16/2024 2:45: Former Schlum G11,VOA Ref. #	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3657-01	917-J-WS-081624	Water		0260 1	08/16/24		00/10/24	08/16/24
P3657-02	TB-01-081624	Water	VOCMS Group6	8260-Low 8260-Low	08/16/24		08/19/24 08/19/24	08/16/24



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#### Hit Summary Sheet SW-846

**SDG No.:** P3657

Client: JACOBS Engineering Group, Inc.

Sample ID	Client ID	Parameter		Concentration	С	MDL	RDL	Units
Client ID :	917-J-WS-081624							
P3657-01	917-J-WS-081624	WATER	Fluorene	0.070	J	0.02	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Phenanthrene	0.050	J	0.02	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Fluoranthene	0.090	J	0.02	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Pyrene	0.060	J	0.02	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Benzo(a)anthracene	0.050	J	0.02	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Chrysene	0.080	J	0.03	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Benzo(b)fluoranthene	0.070	J	0.03	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Benzo(k)fluoranthene	0.050	J	0.04	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Indeno(1,2,3-cd)pyrene	0.050	J	0.04	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Dibenzo(a,h)anthracene	0.040	J	0.04	0.1	ug/L
P3657-01	917-J-WS-081624	WATER	Benzo(g,h,i)perylene	0.050	J	0.04	0.1	ug/L
			Total Svoc :		0.	66		
			<b>Total Concentration:</b>		0	.66		

6

B C

D





A B C D



		•		•			
Client:	JACOBS Engineer	ng Group, Inc.		]	Date Collected:	08/16/24	ŀ
Project:	Former Schlumberg	ger Site Princeton NJ		]	Date Received:	08/16/24	Ļ
Client Sample I	ID: 917-J-WS-081624			5	SDG No.:	P3657	
Lab Sample ID	: P3657-01			1	Matrix:	Water	
Analytical Met					% Solid:	0	
-		-					
Sample Wt/Vol	: 960 Units:	mL		]	Final Vol:	1000	uL
Soil Aliquot Vo	l:	uL			Test:	SVOCM	IS Group3
Extraction Type	2:	Decant	ted : N	[ ]	Level :	LOW	
Injection Volun	ne :	GPC Factor :	1.0	(	GPC Cleanup :	Ν	PH :
Prep Method :	SW3510C						
Trep Method .	5.455100						
File ID/Qc Batch	Dilution:	Prep Date		Date Ana	alyzed	Prep Batch	ID
BN033501.D	1	08/19/24 09	9:50	08/20/24	12:11	PB162821	
CAS Number	Parameter	Conc.	Qualifier	MDL		LOQ / CRQL	Units
TARGETS 91-20-3	Naphthalene	0.030	ΤT	0.020		0.10	11~/T
91-20-3 91-57-6	2-Methylnaphthalene	0.030	U U	0.030 0.030		0.10	ug/L ug/L
208-96-8	Acenaphthylene	0.030	U U	0.030		0.10	ug/L ug/L
83-32-9	Acenaphthene	0.020	U	0.020		0.10	ug/L ug/L
86-73-7	Fluorene	0.020	J	0.020		0.10	ug/L ug/L
85-01-8	Phenanthrene	0.070	J	0.020		0.10	ug/L ug/L
120-12-7	Anthracene	0.030	J U	0.020		0.10	ug/L ug/L
206-44-0	Fluoranthene	0.090	J	0.020		0.10	ug/L ug/L
129-00-0	Pyrene	0.060	J	0.020		0.10	ug/L
56-55-3	Benzo(a)anthracene	0.050	J	0.020		0.10	ug/L
218-01-9	Chrysene	0.080	J	0.020		0.10	ug/L
205-99-2	Benzo(b)fluoranthene	0.070	J	0.030		0.10	ug/L
207-08-9	Benzo(k)fluoranthene	0.050	J	0.040		0.10	ug/L
50-32-8	Benzo(a)pyrene	0.060	Ŭ	0.060		0.10	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	0.050	J	0.040		0.10	ug/L
53-70-3	Dibenzo(a,h)anthracene	0.040	J	0.040		0.10	ug/L
191-24-2	Benzo(g,h,i)perylene	0.050	J	0.040		0.10	ug/L
123-91-1	1,4-Dioxane	0.070	U	0.070		0.21	ug/L
SURROGATES							
7297-45-2	2-Methylnaphthalene-d10	0.31		30 (20) - 150	) (139)	77%	SPK: 0.4
93951-69-0	Fluoranthene-d10	0.37		30 (30) - 150	0 (150)	93%	SPK: 0.4
4165-60-0	Nitrobenzene-d5	0.29		30 (27) - 130	0 (123)	73%	SPK: 0.4
321-60-8	2-Fluorobiphenyl	0.31		30 (34) - 130	) (132)	78%	SPK: 0.4
1718-51-0	Terphenyl-d14	0.46		30 (35) - 130	) (157)	115%	SPK: 0.4
INTERNAL STAN							
3855-82-1	1,4-Dichlorobenzene-d4	7400	7.552				
1146-65-2	Naphthalene-d8	19900	10.314				
15067-26-2	Acenaphthene-d10	10400	14.189				
1517 22 2	Dhamanthrona d10	21000	16 042				

1517-22-2

Phenanthrene-d10

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21000

16.942



6
U

Report of Analysis										
Client:	lient: JACOBS Engineering Group, Inc. Dat								08/16/24	
Project:	Former S	chlumberg	ger Site	Princeton NJ	ſ		Date Received:		08/16/24	
Client Sample ID:	917-J-WS	5-081624					SDG No.:		P3657	
Lab Sample ID:	P3657-01						Matrix:		Water	
Analytical Metho	d: SW82705	SIM					% Solid:		0	
Sample Wt/Vol:	960	Units:	mL				Final Vol:		1000	uL
Soil Aliquot Vol:		uL	uL			Test:		SVOCMS	Group3	
Extraction Type :				Decan	ited : N		Level :		LOW	
Injection Volume	:		G	PC Factor :	1.0		GPC Cleanup :	Ν		PH :
Prep Method :	SW35100	2								
File ID/Qc Batch:	Dilution:			Prep Date		Date A	nalyzed	P	rep Batch II	)
BN033501.D	1			08/19/24 09	9:50	08/20/	24 12:11	P	B162821	
CAS Number	Parameter			Conc.	Qualifier	MDL		LOQ	/ CRQL	Units
1719-03-5 1520-96-3	Chrysene-d12 Perylene-d12			13100 12700	21.148 23.323					

U = Not Detected

- LOQ = Limit of Quantitation
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- LOD = Limit of Detection
- E = Value Exceeds Calibration Range
- Q = indicates LCS control criteria did not meet requirements
- M = MS/MSD acceptance criteria did not meet requirements

- J = Estimated Value
- B = Analyte Found in Associated Method Blank
- N = Presumptive Evidence of a Compound
- \* = Values outside of QC limits
- D = Dilution
- () = Laboratory InHouse Limit
- A = Aldol-Condensation Reaction Products



# A B

D

6

# LAB CHRONICLE

OrderID: Client: Contact:	P3657 JACOBS Engineering Group, I Mary I. Murphy	OrderDate: Project: Location:	8/16/2024 2:45 Former Schlum G11,VOA Ref. #	berger Site Pri	nceton NJ			
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3657-01	917-J-WS-081624	Water			08/16/24			08/16/24
			SVOCMS Group3	8270-Modifie		08/19/24	08/20/24	
			SVOCMS Group6	d 8270E		08/19/24	08/21/24	



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в	

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#### Hit Summary Sheet SW-846

SDG No.:	P3657				
Client:	JACOBS Enginee	ering Group, Inc.			
Sample ID Client ID :	Client ID	Matrix	Parameter	Concentration C MDL	RDL Units
				0.000	
			Total Svoc :	0.00	
			<b>Total Concentration:</b>	0.00	





A B C D



7

		Repor	t of An	alysis			
Client:	JACOBS Engineeri	Date Collected:	08/16/24	1			
Project:	Former Schlumberg	08/16/24	ł				
Client Sample II	D: 917-J-WS-081624				SDG No.:	P3657	
Lab Sample ID:	P3657-01				Matrix:	Water	
-							
Analytical Meth	od: SW8270				% Solid:	0	
Sample Wt/Vol:	960 Units:	mL			Final Vol:	1000	uL
Soil Aliquot Vol	:	uL			Test:	SVOCM	IS Group6
Extraction Type	:	Decan	nted :	Ν	Level :	LOW	
Injection Volume	e :	GPC Factor :	1.0		GPC Cleanup :	N	PH :
Prep Method :	SW3510C						
File ID/Qc Batch:	Dilution:	Prep Date		Data	Analyzed	Prep Batch	ID
		-	0.40				
BP021589.D	1	08/19/24 09	9:40	08/2	1/24 02:44	PB162822	
CAS Number	Parameter	Conc.	Qualifi	er MDL		LOQ / CRQL	Units
FARGETS 110-86-1	Pyridine	1.60	U	1.60		5.20	ug/L
00-52-7	Benzaldehyde	4.20	U	4.20		10.4	ug/L ug/L
95-48-7	2-Methylphenol	1.20	U	1.20		5.20	ug/L ug/L
98-86-2	Acetophenone	1.10	U	1.10		5.20	ug/L
65794-96-9	3+4-Methylphenols	1.20	U	1.20		10.4	ug/L ug/L
98-95-3	Nitrobenzene	1.30	U	1.30		5.20	ug/L ug/L
120-83-2	2,4-Dichlorophenol	0.92	U	0.92		5.20	ug/L ug/L
91-20-3	Naphthalene	1.10	U	1.10		5.20	ug/L
37-68-3	Hexachlorobutadiene	1.30	U	1.30		5.20	ug/L
91-57-6	2-Methylnaphthalene	1.20	U	1.20		5.20	ug/L
38-06-2	2,4,6-Trichlorophenol	0.93	Ŭ	0.93		5.20	ug/L
95-95-4	2,4,5-Trichlorophenol	1.10	U	1.10		5.20	ug/L
208-96-8	Acenaphthylene	1.10	U	1.10		5.20	ug/L
33-32-9	Acenaphthene	0.84	U	0.84		5.20	ug/L
32-64-9	Dibenzofuran	0.97	Ŭ	0.97		5.20	ug/L
36-73-7	Fluorene	1.00	U	1.00		5.20	ug/L
18-74-1	Hexachlorobenzene	1.20	U	1.20		5.20	ug/L
37-86-5	Pentachlorophenol	1.90	U	1.90		10.4	ug/L
5-01-8	Phenanthrene	0.93	U	0.93		5.20	ug/L
6-74-8	Carbazole	1.20	U	1.20		5.20	ug/L
4-74-2	Di-n-butylphthalate	1.50	U	1.50		5.20	ug/L ug/L
06-44-0	Fluoranthene	1.30	U	1.30		5.20	ug/L
29-00-0	Pyrene	1.10	U	1.10		5.20	ug/L
6-55-3	Benzo(a)anthracene	0.98	U	0.98		5.20	ug/L
218-01-9	Chrysene	0.90	U	0.90		5.20	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	2.00	U	2.00		5.20	ug/L ug/L
205-99-2	Benzo(b)fluoranthene	1.20	U	1.20		5.20	ug/L ug/L
207-08-9	Benzo(k)fluoranthene	1.20	U	1.20		5.20	ug/L ug/L
207-08-9		1.20	U	1.20		5.20	ug/L

50-32-8

Benzo(a)pyrene

U

1.70

5.20

ug/L

1.70



7

#### **Report of Analysis** Client: JACOBS Engineering Group, Inc. Date Collected: 08/16/24 Project: Former Schlumberger Site Princeton NJ Date Received: 08/16/24 Client Sample ID: 917-J-WS-081624 SDG No .: P3657 Lab Sample ID: P3657-01 Matrix: Water Analytical Method: SW8270 % Solid: 0 Final Vol: 1000 uL Sample Wt/Vol: 960 Units: mL Soil Aliquot Vol: uL Test: SVOCMS Group6 Extraction Type : Decanted : Ν Level : LOW Injection Volume : GPC Factor : 1.0 GPC Cleanup : Ν PH : SW3510C Prep Method : File ID/Qc Batch: Dilution: Prep Date Date Analyzed Prep Batch ID BP021589.D 1 08/19/24 09:40 08/21/24 02:44 PB162822 Conc. MDL Units Qualifier LOQ / CRQL **CAS Number** Parameter

			-			
193-39-5	Indeno(1,2,3-cd)pyrene	1.10	U	1.10	5.20	ug/L
53-70-3	Dibenzo(a,h)anthracene	1.20	U	1.20	5.20	ug/L
191-24-2	Benzo(g,h,i)perylene	1.20	U	1.20	5.20	ug/L
123-91-1	1,4-Dioxane	1.30	U	1.30	5.20	ug/L
90-12-0	1-Methylnaphthalene	0.90	U	0.90	5.20	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	52.1		15 (10) - 110 (139)	35%	SPK: 150
13127-88-3	Phenol-d6	32.9		15 (10) - 110 (134)	22%	SPK: 150
4165-60-0	Nitrobenzene-d5	82.0		30 (49) - 130 (133)	82%	SPK: 100
321-60-8	2-Fluorobiphenyl	76.3		30 (52) - 130 (132)	76%	SPK: 100
118-79-6	2,4,6-Tribromophenol	179	*	15 (44) - 110 (137)	119%	SPK: 150
1718-51-0	Terphenyl-d14	96.0		30 (48) - 130 (125)	96%	SPK: 100
INTERNAL STAN	DARDS					
3855-82-1	1,4-Dichlorobenzene-d4	353000	7.805			
1146-65-2	Naphthalene-d8	1460000	10.599			
15067-26-2	Acenaphthene-d10	957000	14.457			
1517-22-2	Phenanthrene-d10	2100000	17.269			
1719-03-5	Chrysene-d12	1960000	21.727			
1520-96-3	Perylene-d12	2260000	25.168			

- LOQ = Limit of Quantitation
- MDL = Method Detection Limit
- LOD = Limit of Detection
- E = Value Exceeds Calibration Range
- Q = indicates LCS control criteria did not meet requirements
- M = MS/MSD acceptance criteria did not meet requirements

- J = Estimated Value
- B = Analyte Found in Associated Method Blank
- N = Presumptive Evidence of a Compound
- \* = Values outside of QC limits
- D = Dilution
- () = Laboratory InHouse Limit
- A = Aldol-Condensation Reaction Products



# A B C D

# LAB CHRONICLE

OrderID: Client: Contact:	P3657 JACOBS Engineering Group, Inc. Mary I. Murphy				8/16/2024 2:45 Former Schlum G11,VOA Ref. :	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3657-01	917-J-WS-081624	Water			08/16/24			08/16/24
			SVOCMS Group6	8270E		08/19/24	08/21/24	



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#### Hit Summary Sheet SW-846

SDG No.:	P3657			Order ID:		P3657		
Client:	JACOBS Engineering Group	p, Inc.		Project ID	:	Former Schlumb	erger Site Princetor	ı NJ
Sample ID	Client ID	Matrix	Parameter	Concentration	С	MDL	RDL	Units
Client ID :	917-J-WS-081624							
P3657-01	917-J-WS-081624	Water	Aluminum	92.1		1.98	20.0	ug/L
P3657-01	917-J-WS-081624	Water	Antimony	0.23	J	0.11	2.00	ug/L
P3657-01	917-J-WS-081624	Water	Arsenic	1.85		0.090	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Barium	58.0		0.30	10.0	ug/L
P3657-01	917-J-WS-081624	Water	Calcium	19400		62.5	500	ug/L
P3657-01	917-J-WS-081624	Water	Chromium	1.48	J	0.40	2.00	ug/L
P3657-01	917-J-WS-081624	Water	Cobalt	0.60	J	0.062	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Copper	2.68		0.40	2.00	ug/L
P3657-01	917-J-WS-081624	Water	Iron	3070		9.60	50.0	ug/L
P3657-01	917-J-WS-081624	Water	Lead	1.67		0.11	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Magnesium	3500		26.6	500	ug/L
P3657-01	917-J-WS-081624	Water	Manganese	362		0.24	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Nickel	2.83		0.18	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Potassium	2760		46.1	500	ug/L
P3657-01	917-J-WS-081624	Water	Tin	0.23	J	0.12	5.00	ug/L
P3657-01	917-J-WS-081624	Water	Sodium	72200		85.8	500	ug/L
P3657-01	917-J-WS-081624	Water	Vanadium	0.94	J	0.072	5.00	ug/L
P3657-01	917-J-WS-081624	Water	Zinc	11.7		0.56	5.00	ug/L
P3657-01	917-J-WS-081624	Water	Strontium	133		0.35	1.00	ug/L
P3657-01	917-J-WS-081624	Water	Titanium	1.91	J	0.26	5.00	ug/L

8

A B C

D





A B C D



B C D

# **Report of Analysis**

Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/16/24	
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/16/24	
Client Sample ID:	917-J-WS-081624	SDG No.:	P3657	
Lab Sample ID:	P3657-01	Matrix:	Water	
Level (low/med):	low	% Solid:	0	

Cas	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Prep Date	Date Ana.	Ana Met.	Prep Met.
7429-90-5	Aluminum	92.1		1	1.98	20.0	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-36-0	Antimony	0.23	J	1	0.11	2.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-38-2	Arsenic	1.85	*	1	0.090	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-39-3	Barium	58.0		1	0.30	10.0	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-41-7	Beryllium	0.16	U	1	0.16	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-43-9	Cadmium	0.30	U	1	0.30	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-70-2	Calcium	19400		1	62.5	500	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-47-3	Chromium	1.48	J	1	0.40	2.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-48-4	Cobalt	0.60	J	1	0.062	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-50-8	Copper	2.68		1	0.40	2.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7439-89-6	Iron	3070		1	9.60	50.0	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7439-92-1	Lead	1.67		1	0.11	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7439-95-4	Magnesium	3500		1	26.6	500	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7439-96-5	Manganese	362		1	0.24	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7439-97-6	Mercury	0.081	UN	1	0.081	0.20	ug/L	08/21/24 15:15	08/22/24 15:53	SW7470A	
7439-98-7	Molybdenum	0.93	UN	1	0.93	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-02-0	Nickel	2.83		1	0.18	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-09-7	Potassium	2760		1	46.1	500	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7782-49-2	Selenium	1.38	U	1	1.38	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-22-4	Silver	0.077	UN	1	0.077	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-23-5	Sodium	72200		1	85.8	500	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-24-6	Strontium	133		1	0.35	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-28-0	Thallium	0.085	U	1	0.085	1.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-31-5	Tin	0.23	J	1	0.12	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-32-6	Titanium	1.91	J	1	0.26	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-62-2	Vanadium	0.94	J	1	0.072	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A
7440-66-6	Zinc	11.7		1	0.56	5.00	ug/L	09/04/24 12:30	09/04/24 16:16	SW6020	3010A

Color Before:	Colorless	Clarity Before:	Clear	Texture:				
Color After:	Colorless	Clarity After:	Clear	Artifacts:				
Comments:	Mercury							
MDL = MetholLOD = Limit ofD = Dilution	of Quantitation d Detection Limit	et requirements		<ul> <li>J = Estimated Value</li> <li>B = Analyte Found in Associated Method Blank</li> <li>* = indicates the duplicate analysis is not within control limits.</li> <li>E = Indicates the reported value is estimated because of the presence of interference.</li> <li>OR = Over Range</li> <li>N =Spiked sample recovery not within control limits</li> </ul>				
P3657				of 48				



## Α

B C

D

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OrderID: Client: Contact:	P3657 JACOBS Engineering Group, II Mary I. Murphy	nc.		OrderDate: Project: Location:	8/16/2024 2:45: Former Schlum G11,VOA Ref. #	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3657-01	917-J-WS-081624	Water			08/16/24			08/16/24
			Mercury Metals Group4	7470A 6020B		08/21/24 09/04/24	08/22/24 09/04/24	





9

В



#### **Report of Analysis**

Client:	JACOBS Engineering Group,	Inc.	Date Coll	ected: 08/16/24	4 09:30
Project:	Former Schlumberger Site Pri	nceton NJ	Date Reco	eived: 08/16/24	ł
Client Sample ID:	917-J-WS-081624		SDG No.:	P3657	
Lab Sample ID:	P3657-01		Matrix:	WATER	
			% Solid:	0	
Parameter	Conc. Qua. DF MDL	LOQ / CRQL	Units Prep	Date Date Ana.	Ana Met.
Dissolved Hexavalent Chromium	0.0030 U 1 0.0030	0.010	mg/L	08/16/24 17:3	34 7196A

Comments:

- U = Not Detected
- LOQ = Limit of Quantitation
- MDL = Method Detection Limit
- LOD = Limit of Detection
- D = Dilution
- Q = indicates LCS control criteria did not meet requirements
- H = Sample Analysis Out Of Hold Time

- J = Estimated Value
- B = Analyte Found in Associated Method Blank

- E = Indicates the reported value is estimated because of the presence of interference.
- OR = Over Range
- N =Spiked sample recovery not within control limits

<sup>\* =</sup> indicates the duplicate analysis is not within control limits.



С

#### LAB CHRONICLE

OrderID: Client: Contact:	P3657 JACOBS Engineering Group, I Mary I. Murphy	nc.		OrderDate: Project: Location:	8/16/2024 2:45 Former Schlum G11,VOA Ref. <del>3</del>	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3657-01	917-J-WS-081624	WATER			08/16/24 09:30			08/16/24
			Hexavalent Chromium	7196A			08/16/24 17:34	



## <u>SHIPPING</u> DOCUMENTS

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#### 284 Sheffield Street, Mountainside, NJ 07092 (908) 789-8900 • Fax (908) 789-8922 www.chemtech.net

CHEMTECH PROJECT NO. P3657	
GOUTENO.	
COC Number 2041316	— <u>10</u>
2041316	10.1

10.1

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ATTENTION:	John Yut	ante		e-mail:	No	wy.	Murp	hy@J.	achs	com	0		ATTE	TION:				РНС	NE:	
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1.				1.10	8	-			**	1	2	3	4	5	6	7	8	9	C-H2SO4	F-OTHER
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Method	Analyte	CAS Number	Units	Higher of PQL and Ground Water Ouality	Fresh Surface Water Chronic NJDEP
				Criterion <sup>a</sup>	Ecological Criterion <sup>b</sup>
ECO-SVOCs					
SW8270E	1,4-Dioxane	123-91-1	µg/L	0.4	1
SW8270E	1-Methyinaphthalene	90-12-0	hg/L	1	
SW8270E	2,4,5-Trichlorophenol	95-95-4	hg/L	700	
SW8270E	2,4,6-Trichlorophenol	88-06-2	hg/L	20	
SW8270E	2,4-Dinitrotoluene	121-14-2	hg/L	10	
SW8270E	2-Methyinaphthalene	91-57-6	hg/L	30	
SW8270E	2-Methylphenol	95-48-7	hg/L	ł	
SW8270E	3 & 4-Methylphenol (m,p-Cresols)	65794-96-9	µg/L	1	
SW8270E	Acenaphthene	83-32-9	hg/L	400	
SW8270E	Acenaphthylene	208-96-8	hg/L	I	
SW8270E	Anthracene	120-12-7	Hg/L	2000	
SW8270E	Benzaldehyde	100-52-7	µg/L	1	
SW8270E	Benzo(a)anthracene	56-55-3	hg/L	ł	
SW8270E	Benzo(a)pyrene	50-32-8	hg/L	0.1	
SW8270E	Benzo(b)fluoranthene	205-99-2	µg/L	0.5	
SW8270E	Benzo(g,h,i)perylene	191-24-2	µg/L	ł	
SW8270E	Benzo(k)fluoranthene	207-08-9	µg/L	0.5	
SW8270E	Bis (2-ethylhexyl) phthalate	117-81-7	µg/L	I	
SW8270E	Carbazole	86-74-8	µg/L	I	
SW8270E	Chrysene	218-01-9	µg/L	5	
SW8270E	Dibenzo(a,h)anthracene	53-70-3	µg/L	0.3	
SW8270E	Dibenzofuran	132-64-9	μg/L	ł	
SW8270E	Di-N-Butylphthalate	84-74-2	μg/L	1	
SW8270E	Fluoranthene	206-44-0	µg/L	300	
SW8270E	Fluorene	86-73-7	Hg/L	300	
SW8270E	Hexachlorobenzene	118-74-1	µg/L	0.02	
SW8270E	Hexachlorobutadiene	87-68-3	µg/L	t	
SW8270E	Hexachloroethane	67-72-1	µg/l	7	
SW8270E	Indeno(1,2,3-Cd)Pyrene	193-39-5	µg/L	0.2	
SW8270E	Naphthalene	91-20-3	µg/L	300	
SW8270E	Nitrobenzene	98-95-3	µg/L	9	
SW8270E	Pentachlorophenol	87-86-5	µg/L	0.3	
SW8270E	Phenanthrene	85-01-8	Hg/L	Ŧ	
SW8270E	Pyrene	129-00-0	µg/L	200	
SW8270E	Pyridine	110-86-1	µg/L	1	
ECO-VOCs					
SWR260D	1 1 1 Trickloroothana	71 55 5	-	000	ſ

P3657

Table 3. Surface Water Target Analytes, Methods, Action Levels, and Control Limits

Site Sampling Plan for Ecological Evaluation

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Princeton Technology Center, West Windsor Township, New Jersey

Method	Analyte	CAS Number	Units	Higher of PQL and Ground Water	Surface Water Chronic
				Quality Criterion <sup>a</sup>	NJDEP Ecological Criterion <sup>b</sup>
SW8260D	1,1,2-Trichloroethane	79-00-5	µg/L	m	500
SW8260D	1,1-Dichloroethane	75-34-3	µg/L	50	1
SW8260D	1,1-Dichloroethene	75-35-4	µg/L	۲	65
SW8260D	1,2-Dichlorobenzene	95-50-1	µg/L	600	14
SW8260D	1,2-Dichloroethane	107-06-2	hg/L	2	910
SW8260D	1,2-Dichloroethene (Total)	540-59-0	Hg/L		
SW8260D	1,4-Dichlorobenzene	106-46-7	µg/L	75	9.4
SW8260D	2-Butanone	78-93-3	μg/L	300	1
SW8260D	Acetone	67-64-1	Hg/L	6000	1
SW8260D	Benzene	71-43-2	hg/L		114
SW8260D	Bromodichloromethane	75-27-4	µg/L	1	1
SW8260D	Bromomethane	74-83-9	hg/L	10	1
SW8260D	Carbon disulfide	75-15-0	µg/L	700	ł
SW8260D	Carbon tetrachloride	56-23-5	µg/L	1	240
SW8260D	Chlorobenzene	108-90-7	µg/L	50	47
SW8260D	Chloroethane	75-00-3	µg/L	I	1
SW8260D	Chloroform	67-66-3	μg/L	70	140
SW8260D	Chloromethane	74-87-3	µg/L	ł	ł
SW8260D	cis-1,2-Dichloroethene	156-59-2	μg/L	70	I
SW8260D	Cyclohexane	110-82-7	µg/L	ſ	I
SW8260D	Dibromochloromethane	124-48-1	µg/L	1	ł
SW8260D	Dichlorodifluoromethane	75-71-8	µg/L	1000	I
SW8260D	Ethylbenzene	100-41-4	µg/L	700	14
SW8260D	Freon TF	76-13-1	µg/i	20000	I
SW8260D	Isopropylbenzene	98-82-8	µg/L	700	1
SW8260D	m&p-Xylene	179601-23-1	µg/L	1000	27
SW8260D	Methylcyclohexane	108-87-2	µg/L	1	ł
SW8260D	Methylene Chloride	75-09-2	µg/L	£	940
SW8260D	MTBE	1634-04-4	µg/L	70	51000
SW8260D	o-Xylene	95-47-6	µg/L	1000	27
SW8260D	Tetrachloroethene	127-18-4	µg/L	1	45
SW8260D	Toluene	108-88-3	Hg/L	600	253
SW8260D	trans-1,2-Dichloroethene	156-60-5	µg/L	100	970
SW8260D	Trichloroethene	79-01-6	µg/L	1	47
SW8260D	Vinyl chloride	75-01-4	µg/L	÷	930
SW8260D	Xylenes, Total	1330-20-7	µg/L		
ECO-PAHS					
SW8270E SIM	1,4-Dioxane	123-91-1	ug/L	0.4	

Table 3. Surface Water Target Analytes, Methods, Action Levels, and Control Limits Site Sampling Plan for Ecological Evaluation

Princeton Technology Center, West Windsor Township, New Jersey

Method         Analyte         CAS           Werthod         Analyte         Numbe           SW8270E SIM         Acenaphthene         91-57-           SW8270E SIM         Acenaphthene         91-57-           SW8270E SIM         Acenaphthene         83-32-           SW8270E SIM         Acenaphthene         83-32-           SW8270E SIM         Acenaphthene         83-32-           SW8270E SIM         Anthracene         120-12           SW8270E SIM         Benzo(a)pyrene         56-55-           SW8270E SIM         Benzo(a)pyrene         205-49           SW8270E SIM         Benzo(b)fluoroanthene         207-08           SW8270E SIM         Benzo(c)hyrene         207-08           SW8270E SIM         Benzo(c)hyrene         207-08           SW8270E SIM         Benzo(c)hyrene         207-08           SW8270E SIM         Benzo(c)hyrene         207-08           SW8270E SIM         Naphthalene         207-08			Higher of PQL and	Fresh Surface Water
thod Aaalyte Aaalyte SiM 2-Methylnaphthalene SiM Acenaphthene SiM Acenaphthene SiM Acenaphthene SiM Acenaphthene SiM Acenaphthene SiM Benzo(a)anthracene SiM Benzo(a)anthracene SiM Benzo(a)anthracene SiM Benzo(a)anthracene SiM Benzo(a)apyrene SiM Indeno[1,2,3-cd]apyrene SiM Indeno[1,2,3-cd]apyrene SiM Naphthalene SiM Naphthalene SiM Naphthalene SiM Naphthalene SiM Prena Arrinony Anathracene SiM Anathracene SiM Prena Arrinony Anathracene SiM Prena Arrinony Anathracene SiM Anathracene SiM Prena Arrinony Anathracene SiM Prena Arrino Arrinony Anathracene SiM Prena Arrino Arrinony Anathracene SiM Prena Arrino Arrinony Arrino Arrino Arrino Arrino Arrino Arrino Arri			PQL and	Water
thod Analyte Analyte Analyte Analyte Acenaphthalene SIM Acenaphthalene SIM Acenaphthalene SIM Acenaphthalene SIM Anthracene SIM Benzo(a)anthracene SIM Benzo(a)anthracene SIM Benzo(a)anthracene SIM Benzo(a)anthracene SIM Benzo(b)fluoroanthene SIM Benzo(b)fluoroanthene SIM Benzo(b)fluoroanthene SIM Benzo(b)fluoroanthene SIM Benzo(b)fluoroanthene SIM Pluorenthene SIM Indeno[1,2,3,cd]pyrene SIM Naphthalene SIM Naphthalene SIM Naphthalene SIM Naphthalene SIM Naphthalene SIM Naphthalene SIM Antimony Arsenic Cadmium Antimony Coper Coper Iron Cadmium Barium Barium Barium Barium Barium Antimony Arsenic Cadmium Antimony Cadmium Cadm			Crossed	
SIM       2-Methylnaphthalene         SIM       Acenaphthalene         SIM       Acenaphthalene         SIM       Acenaphthylene         SIM       Acenaphthylene         SIM       Benzo(a)anthracene         SIM       Benzo(a)pyrene         SIM       Benzo(a)pyrene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(k)fluoroanthene         SIM       Benzo(k)fluoroanthene         SIM       Dibenz(a,h)anthracene         SIM       Pitenz(a,h)anthracene         SIM       Pitenz(a,h)anthracene         SIM       Pitenz(a,h)anthracene         SIM       Naphthalene         SIM       Pitenz(a,h)anthracene         SIM       Naphthalene         SIM       Pyrene         SIM       Naphthalene         SIM       Pyrene         SIM       Pyrene         SIM       Pyrene         SIM       Pyrene         SIM       Pyrene         SIM       Pyrene         SIM       Pyr	2	Units	Water	Chronic
SIM       2-Methylnaphthalene         SIM       Acenaphthalene         SIM       Acenaphthylene         SIM       Anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b,j)perylene         SIM       Benzo(k)fluoroanthene         SIM       Rhuoroanthene         SIM       Pinenz(a,h)anthracene         SIM       Pinenz(a,h)anthracene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Phenanthrene         SIM       Pyrene         SIM <th></th> <th></th> <th>Quality</th> <th>NJDEP</th>			Quality	NJDEP
SIM       2-Methylinaphthalene         SIM       Acenaphthene         SIM       Acenaphthylene         SIM       Acenaphthylene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(b)fluoroanthene         SIM       Chrysene         SIM       Chrysene         SIM       Dibenz(a,h)anthracene         SIM       Chrysene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Pyrene         Aluminum       Aluminum         V1196A       Hexavalent Chromium         V1196A       Hexavalent Chromium         V1190M			Criterion <sup>a</sup>	Ecological Criterion <sup>b</sup>
SIM       Acenaphthene         SIM       Acenaphthylene         SIM       Benzo(a)anthracene         SIM       Benzo(b)fluoroanthene         SIM       Chrysene         SIM       Chrysene         SIM       Chrysene         SIM       Chrysene         SIM       Chrysene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Phenanthrene         SIM <td>91-57-6</td> <td>µg/L</td> <td>30</td> <td>330</td>	91-57-6	µg/L	30	330
SIM       Acenaphthylene         SIM       Anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(k)fluoroanthene         SIM       Diberz(a,h)anthracene         SIM       Diberz(a,h)anthracene         SIM       Ploroanthene         SIM       Ploroanthene         SIM       Ploroanthene         SIM       Naphthalene         SIM       Naphthalene         SIM       Phenanthrene         Aluminum       Aluminum         Aluminum       Aluminum         Aluminum       Aluminum         Aluminum       Aluminum </td <td>83-32-9</td> <td>µg/L</td> <td>400</td> <td>38</td>	83-32-9	µg/L	400	38
SIM       Anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)anthracene         SIM       Benzo(a)pyrene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(k)fluoroanthene         SIM       Benzo(k)fluoroanthene         SIM       Dibenz(a,h)anthracene         SIM       Dibenz(a,h)anthracene         SIM       Pitoroanthene         SIM       Fluoroanthene         SIM       Pitorene         SIM       Potene         SIM       Phenanthracene         Altimony       Altimony         Altimony </td <td>208-96-8</td> <td>µg/L</td> <td>ł</td> <td>4840</td>	208-96-8	µg/L	ł	4840
SIM       Benzo(a)anthracene         SIM       Benzo(a)pyrene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(b)fluoroanthene         SIM       Benzo(g,h,i)perylene         SIM       Benzo(g,h,i)perylene         SIM       Chrysene         SIM       Chrysene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Pluoroanthene         SIM       Pluorene         SIM       Pluorene         SIM       Phenanthracene         SIM       Phenanthrene         SIM	120-12-7	µg/L	2000	0.035
SIM       Benzo(a)pyrene         SIM       Benzo(g,hj)perylene         SIM       Benzo(g,hj)perylene         SIM       Benzo(g,hj)perylene         SIM       Benzo(g,hj)perylene         SIM       Chrysene         SIM       Chrysene         SIM       Pluoroanthene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Pluoroanthene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Phenanthracene         Aluminum       Aluminum         Aluminum       Aluminum         Aluminum       Aluminum         Aluminum       Arsenic         Aluminum       Arsenic         Coper       Cadmium         Coper       Coper         Coper       Coper	56-55-3	Hg/L	0.1	0.025
SIM       Benzo(g,h,i)perylene         SIM       Benzo(g,h,i)perylene         SIM       Benzo(g,h,i)perylene         SIM       Chrysene         SIM       Chrysene         SIM       Chrysene         SIM       Pluoroanthene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Naphthalene         SIM       Phenanthrene         SIM	50-32-8	hg/L	0.1	0.014
SIM       Benzo(g,h,i)perylene         SIM       Benzo(k)filuoroanthene         SIM       Chrysene         SIM       Dibenz(a,h)anthracene         SIM       Dibenz(a,h)anthracene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Fluorene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Naphthalene         SIM       Phenanthrene         SIM       Pyrene         Aluminum       Aluminum         Artimony       Artimony         Artimony       Artimony         Artimony <t< td=""><td>205-99-2</td><td>µg/L</td><td>0.2</td><td>9.07</td></t<>	205-99-2	µg/L	0.2	9.07
SIM       Benzo(k)fluoroanthene         SIM       Chrysene         SIM       Diberz(a,h)anthracene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Naphthalene         SIM       Phenanthrene         SIM       Pyrene         Antimony       Antimony         Antimony       Antimony         Antimony       Antimony         Antimony       Antimony         Antimony       Antimony         Antimony       Antimony         Antimony       Antinum         Antimony	191-24-2	µg/L	1	7.64
SIM Chrysene SIM Dibenz(a,h)anthracene SIM Fluoroanthene SIM Fluoroanthene SIM Indeno[1,2,3-cd]pyrene SIM Naphthalene SIM Phenanthrene SIM Pyrene SIM Phenanthrene SIM Pyrene SIM Phenanthrene SIM Pyrene SIM Phenanthrene SIM Pyrene SIM Phenanthrene SIM Phenanthrene Phenanthrene Phenanthrene SIM Phenanthrene SIM Phenanthrene Phenanthrene SIM Phenanthrene SIM Phenanthrene Phenanthrene SIM Phenanthrene SIM Phenanthrene Phenanthrene SIM Phenanthrene SIM Phenanthrene SIM Phenanthrene Phenanthrene Phenanthrene Phenanthrene SIM Phenanthrene Phenanthrene Phenanthrene SIM Phenanthrene Phenanthr	207-08-9	µg/L	0.5	1
SIM       Dibenz(a,h)anthracene         SIM       Fluoroanthene         SIM       Fluoroanthene         SIM       Indeno[1,2,3-cd]pyrene         SIM       Naphthalene         SIM       Phenanthree         SIM       Phenanthree         SIM       Phenanthree         SIM       Pyrene         Antimony       Antimony         Antimony       Antinun         Aresenic       Bery	218-01-9	µg/L	5	1
SIM Fluoroanthene SIM Indeno[1,2,3-cd]pyrene SIM Naphthalene SIM Naphthalene SIM Phenanthrene SIM Pyrene SIM P	53-70-3	µg/L	0.3	1
SIM Fluorene SIM Indeno[1,2,3-cd]pyrene SIM Naphthalene SIM Phenanthrene SIM Pyrene SIM Pyrene SIM Pyrene SIM Pyrene Alumium Alumium Alumium Alumium Alumium Alumium Alumium Alumium Alumium Alumium Alumium Calonium Cadmium Barium Barium Barium Cadmium Cad	206-44-0	µg/L	300	1.9
SiM Indeno[1,2,3-cd]pyrene SIM Naphthalene SIM Phenanthrene SIM Pyrene SIM Pyrene Menanthrene Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Arsenic Beryllium Cadmium Cadmium Chomiu	86-73-7	µg/L	300	19
SIM     Naphthalene       SIM     Phenanthrene       SIM     Pyrene       Aluminum     Mercury       Arsenic     Antimony       Arsenic     Arsenic       Arsenic     Cadmium       Barium     Barium       Cadmium     Cadmium       Copper     Copper       Icon     Lead       Magnesium     Nickel       Nickel     Nickel	193-39-5	µg/L	0.2	4.31
SIM Phenanthrene SIM Pyrene SIM Pyrene SIM Pyrene SIM Pyrene SIM Pyrene SIM Pyrene Mercury Mercury Antimony Antimony Antimony Antimony Antimony Antimony Antimony Antimony Antimony Calcium Beryllium Beryllium Calciu	91-20-3	µg/L	300	13
SIM Pyrene als Pyrene Altanony Mercury Aluminum Altimony Antimony Arsenic Barium Barium Barium Cadmium Cadmium Cadmium Chrom	85-01-8	µg/L	1	3.6
Alternation V1366 Hexavalent Chromium Mercury Altimony Antimony Arsenic Barium Barium Barium Cadmium C	129-00-0	µg/L	200	0.3
V71366 Hexavalent Chromium Mercury Mercury Aluminum Antimony Arsenic Barium Baryllium Baryllium Cadmium Cadmium Cadmium Cadmium Cadmium Cadmium Chromium Chr				
Mercury Aluminum Antimony Arsenic Barium Baryllium Cadmium Cadmium Cadmium Cadmium Cadmium Chromium Copper Iron Lead Magnesium Magnesium Nickel Selenium	18540-29-9	hg/L	1	10
Aluminum Antimony Arsenic Barium Beryllium Cadmium Cadmium Cadmium Chromium Chromium Chromium Copper Iron Lead Magnesium Magnesium Nickel Selenium	7439-97-6	μg/L	2	0.77
Antimony Arsenic Barium Beryllium Cadmium Calcium Chromium Chromium Chromium Chromium Copper Iron Iron Iron Iron Iron Iron Bagnesium Magnesium Selenium Selenium	7429-90-5	hg/L	1	1
Arsenic Barium Barium Cadmium Calcium Calcium Cobalt Copper Iron Iron Iron Iron Magnesium Magnesium Selenium Selenium	7440-36-0	µg/L	9	80
Barium Beryllium Calcium Chromium Chromium Chromium Copper Iron Lead Magnesium Magnesium Nickel Selenium	7440-38-2	µg/L	m	150
Beryllium Cadmium Calcium Chromium Cobalt Copper Iron Iron Lead Magnesium Magnesium Selenium Selenium	7440-39-3	µg/L	6000	220
Cadmium Calcium Chromium Cobalt Cobalt Iron Iron Lead Magnesium Magnesium Nickel Selenium	7440-41-7	µg/L	Ţ	3.6
Calcium Chromium Cobalt Copper Iron Lead Magnesium Magnesium Nickel Potassium Selenium	7440-43-9	µg/L	4	1
Chromium Cobalt Copper Iron Lead Magnesium Magnese Nickel Potassium Selenium	7440-70-2	µg/L	I	I
Cobalt Copper Iron Lead Magnesium Manganese Nickel Potassium Selenium	7440-47-3	µg/L	1	42
Copper Iron Lead Magnesium Manganese Nickel Potassium Selenium	7440-48-4	µg/L	100	24
Iron Lead Marganese Nickel Potassium Selenium	7440-50-8	µg/L	1300	1
Lead Magnesium Manganese Nickel Potassium Selenium	7439-89-6	µg/L	I	1
Magnesium Manganese Nickel Potassium Selenium	7439-92-1	µg/L	5	5.4
Manganese Nickel Potassium Selenium	7439-95-4	µg/L	1	I
Nickel Potassium Selenium	7439-96-5	µg/L	1	1
Potassium Selenium	7440-02-0	µg/L	100	1
Selenium	7440-09-7	µg/L	1	1
COCUMUN	7782-49-2	µg/L	40	ъ
EPA 200.7 Silica 7631-86	7631-86-9	µg/L	ł	1

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10.1

Princeton Technology Center, West Windsor Township, New Jersey

Fresh Surface Water Chronic NJDEP Ecological	Criterion
Higher of PQL and Ground Water Quality Criterion <sup>a</sup>	
Units	
CAS Number	
Analyte	
Method	

					Criterion <sup>b</sup>	
SW6020B	Silver	7440-22-4	µg/L	40	40 0.12	
SW6020B	Sodium	7440-23-5	µg/L	1	:	
SW6020B	Thallium	7440-28-0	hg/L	1	10	
SW6020B	Vanadium	7440-62-2	hg/L	1	12	
SW6020B	Zinc	7440-66-6 µg/L	µg/L	2000	1	
Notes:						

<sup>a</sup> New Jersey Department of Environmental Protection (NJDEP) Ground Water Quality Standards - Class IIA by

Constituent. May 2021. New Jersey Administration Code 7:9C-1.4: Remediation Standards.

<sup>b</sup> NJDEP Ground Water Quality Standards - Class IIA by Constituent. May 2021. New Jersey Administration Code 7:9C-1.4: Remediation Standards. NJDEP Ecological Surface Water SSLs. March 2009.

Bold = MDL and RL exceed screening criteria.

-- = not available (no standard)

 $\mu g/L = microgram(s)$  per liter

CAS = Chemical Abstracts Service

Freon TF = 1,1,2-Trichloro-1,2,2-trifluoroethane

MDL = method detection limit

MTBE = methyl tert butyl ether

NJDEP = New Jersey Department of Environmental Protection

PAH = polycyclic aromatic hydrocarbon

PQL = Practical Quantitation Level as defined in N.J.A.C. 7:9C-1.4

RL = reporting limit

SIM = selected ion method

SVOC = semivolatile organic compound

VOC = volatile organic compound

10.1



### 10 10.2

#### Laboratory Certification

Certified By	License No.
CAS EPA CLP Contract	68HERH20D0011
Connecticut	PH-0830
DOD ELAP (L-A-B)	L2219
Maine	2024021
Maryland	296
New Hampshire	255423
New Jersey	20012
New York	11376
Pennsylvania	68-00548
Soil Permit	525-24-234-08441
Texas	T104704488



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

#### LOGIN REPORT/SAMPLE TRANSFER

	Order ID: P3657	P3657 JACO05		Order Date :		8/16/2024 2:45:00 PM	Project Mgr :				
Client Name : JACOBS Engineering Grou		Project Name :		Former Schlumberger Site	E	<b>Report Type :</b> L	evel 4				
Client Contact : Mary I. Murphy Invoice Name : JACOBS Engineering Grou		Receive DateTime : Purchase Order :		8/16/2024 1 <del>2:00:00 AM</del> 12 = 45	EDD Type : CH2MHILL						
					Hard Copy Date :						
Invoice Contact : Mary I. Murphy		urphy						Date Signoff :			
LAB ID	CLIENT ID		MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD		FAX DATE	DUE DATES
P3657-01	<mark>717-J-WS-0</mark> 917	081624	Water	08/16/2024	09:30						
						VOCMS Group6		8260-Low	10 Bus. Days		
P3657-02	TB-01-08	1624	Water	08/16/2024	10:55						
						VOCMS Group6		8260-Low	10 Bus. Days		

**Relinguished By :** Date / Time : 08-16-24 1524

em **Received By :** 15:27 Rg+ 7 4 16/24 Date / Time : 💋

Storage Area : VOA Refridgerator Room