

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: P3440

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, NJ	Project Number :	D3779922
Laboratory Sample ID	(s) : <u>P3440</u>	Sampling Date(s) :	8/01/2024

List DKQP Methods Used (e.g., 8260,8270, et Cetra) 6010D,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?	$\mathbf{\nabla}$	Yes		No	
1B	EPH Method: Was the EPH method conducted without significant modifications (see Section 11.3 of respective DKQ methods)		Yes		No	✓ N/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?	Ø	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?	$\mathbf{\nabla}$	Yes		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."



Client Sample Number

Cover Page

- Order ID : P3440
- Project ID : Former Schlumberger Site Princeton NJ
 - **Client :** JACOBS Engineering Group, Inc.

Lab Sample Number

P3440-01923-K1-WS-080124P3440-02923-K1-WS-080124MSP3440-03923-K1-WS-080124MSDP3440-04922-K1-WS-080124P3440-05TB-01-080124

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 :

N. N. Pandya

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 # P3440 Test Name: VOCMS Group6

A. Number of Samples and Date of Receipt:

5 Water samples were received on 08/01/2024.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOC-SIMGroup1, 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 MS recoveries met the requirements for all compounds .

The MSD recoveries met the acceptable requirements .

The RPD met criteria.

The Blank Spike 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:

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:



Signature_

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.

N. N. Pandya

APPROVED

By Nimisha Pandya QA/QC Supervisor at 10:49 am, Aug 16, 2024

2.1



CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3440 Test Name: SVOCMS Group3

A. Number of Samples and Date of Receipt:

5 Water samples were received on 08/01/2024.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOC-SIMGroup1, 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 except for 923-K1-WS-080124MSD [Terphenyl-d14 - 134%] and 922-K1-WS-080124 [Terphenyl-d14 - 148%] these compounds did not meet the NJDKQP criteria but met the in-house criteria.

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

The MS {P3440-02MS} with File ID: BN033232.D recoveries met the requirements for all compounds except for 2-Methylnaphthalene[68%], Phenanthrene[136%] these compounds did not meet the NJDKQP criteria and in-house criteria due to matrix interference.

The MSD {P3440-03MSD} with File ID: BN033233.D recoveries met the acceptable requirements except for Phenanthrene[152%] this compound did not meet the NJDKQP criteria and in-house criteria due to matrix interference.

The RPD met criteria . The Blank Spike met requirements for all samples .



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

The % RSD is greater than 20% in the Initial Calibration method (Method 8270Sim-BN080524.M) for 1,4-Dioxane, this compound is passing on Linear Regression.

The Continuous Calibration File ID BN033228.D met the requirements except for Benzo(b)fluoranthene is failing marginally low and and 2,4,6-Tribromophenol, failure surrogate is not associated with the client list, as per criteria affected surrogates were passing, therefore no corrective action was taken.

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.

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.

N. N. Pandya

Signature_

APPROVED

By Nimisha Pandya QA/QC Supervisor at 10:50 am, Aug 16, 2024



CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3440 Test Name: SVOCMS Group6

A. Number of Samples and Date of Receipt:

5 Water samples were received on 08/01/2024.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOC-SIMGroup1, 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_F using GC Column DB-UI 8270D which is 20 meters, 0.18 mm ID, 0.36 um dfThe 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. The Internal Standards Areas met the acceptable requirements. The Retention Times were acceptable for all samples.

The MS {P3440-02MS} with File ID: BF138838.D recoveries met the requirements for all compounds except for Benzaldehyde[0%] this compound did not meet the NJDKQP criteria and in-house criteria due to matrix interference.

The MSD {P3440-03MSD} with File ID: BF138839.D recoveries met the acceptable requirements except for Benzaldehyde[0%] this compound did not meet the NJDKQP criteria and in-house criteria due to matrix interference.

The RPD met criteria . The Blank Spike 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 BF138834.D met the requirements except for Pentachlorophenol but no positive hit in associated samples therefore no corrective action taken.



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 8 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.

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.

N. N. Pandya Signature_

APPROVED

By Nimisha Pandya QA/QC Supervisor at 10:50 am, Aug 16, 2024



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

CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3440 Test Name: Metals Group4,Mercury

A. Number of Samples and Date of Receipt:

5 Water samples were received on 08/01/2024.

B. Parameters:

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOC-SIMGroup1, 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 6010D, 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 analysis met criteria for all samples.

The Matrix Spike (923-K1-WS-080124MS) analysis met criteria for all samples except for Strontium due to Chemical interference during digestion Process.

The Matrix Spike Duplicate (923-K1-WS-080124MSD) analysis met criteria for all samples except for Strontium 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 met the acceptable 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_

N. N. Pandya

APPROVED

By Nimisha Pandya QA/QC Supervisor at 10:51 am, Aug 16, 2024

24



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

CASE NARRATIVE

JACOBS Engineering Group, Inc. Project Name: Former Schlumberger Site Princeton NJ Project # N/A Chemtech Project # P3440 Test Name: Hexavalent Chromium

A. Number of Samples and Date of Receipt:

5 Water samples were received on 08/01/2024.

B. Parameters:

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, Mercury, Metals Group4, SVOC-SIMGroup1, 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.

N. N. Pandya

Signature_

APPROVED By Nimisha Pandya QA/QC Supervisor at 10:51 am, Aug 16, 2024

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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	Method qualifiers"P"for ICP instrument"PM"for ICP when Microwave Digestion is used"CV"for Manual Cold Vapor AA"AV"for automated Cold Vapor AA"CA"for MIDI-Distillation Spectrophotometric"AS"for Semi – Automated Spectrophotometric"C"for Manual Spectrophotometric"T"for Titrimetric"NR"for analyte not required to be analyzedIndicates the analyte's concentration exceeds the calibrated range of the instrument for that specific analysis.					
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	 Indicates an estimated value. This flag is used: (1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.) (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.
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 #: P3440

Completed

4

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	<u> </u>
Do requested analyses on Chain of Custody agree with the log-in page	<u> </u>
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	<u> </u>
All manual calculations and /or hand notations verified	<u>✓</u>

1st Level QA Review Signature:

SOHIL JODHANI

N. N. Pandya

Date: 08/16/2024

APPROVED By Nimisha Pandya QA/QC Supervisor at 10:51 am, Aug 16, 2024

2nd Level QA Review Signature:

P3440



Hit Summary Sheet SW-846

 SDG No.:
 P3440

 Client:
 JACOBS Engineering Group, Inc.

Sample ID	Client ID	Matrix	Parameter	Concentration	С	MDL	RDL	Units
Client ID:	923-K1-WS-080124							
P3440-01	923-K1-WS-080124 Wa	ater	Acetone	4.90	J	1.40	5.00	ug/L
P3440-01	923-K1-WS-080124 Wa	ater	Toluene	0.71	J	0.18	1.00	ug/L
			Total Voc :	5.6	1			
			Total Concentration:	5.6	1			
Client ID:	922-K1-WS-080124							
P 3440-04	922-K1-WS-080124 Wa	ater	Toluene	1.10		0.18	1.00	ug/L
			Total Voc :	1.1	D			
			Total Concentration:	1.10)			

В

С

5

P3440





5

A B C D



Date Collected:

Date Received:

Level :

Report of Analysis

uL

08/01/24

08/01/24

P3440

Water

5000

LOW

1.00

ug/L

VOCMS Group6

0

A B C

5

Client Sample ID:	923-K1-WS-080124	4	SDG No.:
Lab Sample ID:	P3440-01		Matrix:
Analytical Method:	SW8260		% Solid:
Sample Wt/Vol:	5 Units:	mL	Final Vol:
Soil Aliquot Vol:		uL	Test:

ID: 0.25

JACOBS Engineering Group, Inc.

RXI-624

Former Schlumberger Site Princeton NJ

GC Column: Prep Method :

Client:

Project:

File ID/Qc Batch:	Dilution:	Prep Date		Date Analyzed	Prep Batch ID	
VN083251.D	1			08/13/24 12:30	VN081324	
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	4.90	J	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	0.71	J	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
05 47 6		0.1.4		0.1.4	1.00	17

95-47-6

o-Xylene

U

0.14

0.14



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	923-K1-WS-080124	SDG No.:	P3440
Lab Sample ID:	P3440-01	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	
VN083251.D	1			08/13/24 12:30	VN081324	
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	56.1		70 (74) - 130 (125)	112%	SPK: 50
1868-53-7	Dibromofluoromethane	52.5		70 (75) - 130 (124)	105%	SPK: 50
2037-26-5	Toluene-d8	53.3		70 (86) - 130 (113)	107%	SPK: 50
460-00-4	4-Bromofluorobenzene	56.2		70 (77) - 130 (121)	112%	SPK: 50
INTERNAL STAN	DARDS					
363-72-4	Pentafluorobenzene	149000	8.224			
540-36-3	1,4-Difluorobenzene	288000	9.1			
3114-55-4	Chlorobenzene-d5	294000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	131000	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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B



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	922-K1-WS-080124	SDG No.:	P3440
Lab Sample ID:	P3440-04	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	
VN083261.D	1			08/13/24 16:31	VN081324	
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	1.10		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

B C

D

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P3440



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	922-K1-WS-080124	SDG No.:	P3440
Lab Sample ID:	P3440-04	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	
VN083261.D	1			08/13/24 16:31	VN081324	
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	55.0		70 (74) - 130 (125)	110%	SPK: 50
1868-53-7	Dibromofluoromethane	52.4		70 (75) - 130 (124)	105%	SPK: 50
2037-26-5	Toluene-d8	52.8		70 (86) - 130 (113)	106%	SPK: 50
460-00-4	4-Bromofluorobenzene	54.8		70 (77) - 130 (121)	110%	SPK: 50
INTERNAL STAN	DARDS					
363-72-4	Pentafluorobenzene	132000	8.224			
540-36-3	1,4-Difluorobenzene	256000	9.106			
3114-55-4	Chlorobenzene-d5	256000	11.865			
3855-82-1	1,4-Dichlorobenzene-d4	112000	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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B



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	TB-01-080124	SDG No.:	P3440
Lab Sample ID:	P3440-05	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	
VN083262.D	1			08/13/24 16:56	VN081324	
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	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

D

5



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	TB-01-080124	SDG No.:	P3440
Lab Sample ID:	P3440-05	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	
VN083262.D	1			08/13/24 16:56	VN081324	
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	56.6		70 (74) - 130 (125)	113%	SPK: 50
1868-53-7	Dibromofluoromethane	52.9		70 (75) - 130 (124)	106%	SPK: 50
2037-26-5	Toluene-d8	53.3		70 (86) - 130 (113)	107%	SPK: 50
460-00-4	4-Bromofluorobenzene	57.1		70 (77) - 130 (121)	114%	SPK: 50
INTERNAL STAN	DARDS					
363-72-4	Pentafluorobenzene	143000	8.229			
540-36-3	1,4-Difluorobenzene	276000	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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С



5

С

D

LAB CHRONICLE

OrderID: Client: Contact:	P3440 JACOBS Engineering Group, Ir Mary I. Murphy	ıc.		OrderDate: Project: Location:	8/1/2024 12:28 Former Schlum D31,VOA Ref. ;	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3440-01	923-K1-WS-080124	Water			08/01/24			08/01/24
P3440-04	922-K1-WS-080124	Water	VOCMS Group6	8260-Low	08/01/24		08/13/24	08/01/24
			VOCMS Group6	8260-Low	,		08/13/24	,
P3440-05	TB-01-080124	Water	VOCMS Group6	8260-Low	08/01/24		08/13/24	08/01/24



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Hit Summary Sheet SW-846

SDG No.: P3440

Client:

JACOBS Engineering Group, Inc.

Sample ID Client ID :	Client ID 922-K1-WS-080124		Parameter	Concentration	С	MDL	RDL	Units
P3440-04	922-K1-WS-080124	WATER	Phenanthrene	0.040	J	0.02	0.11	ug/L
P3440-04	922-K1-WS-080124	WATER	Fluoranthene	0.040	J	0.02	0.11	ug/L
P3440-04	922-K1-WS-080124	WATER	Pyrene	0.030	J	0.02	0.11	ug/L
			Total Svoc :		0.	11		
			Total Concentration:		0	.11		

6

B C

D





6

A B C D



6

C D

		Repo	rt of Anal	ysis		
Client:	JACOBS Engineer	ing Group, Inc.		Date Collected	l: 08/01/24	ł
Project:	Former Schlumberg	ger Site Princeton N	1J	Date Received	: 08/01/24	Į.
Client Sample I				SDG No.:	P3440	
-				Matrix:	Water	
Lab Sample ID						
Analytical Met	hod: SW8270SIM			% Solid:	0	
Sample Wt/Vol	: 910 Units:	mL		Final Vol:	1000	uL
Soil Aliquot Vo	1:	uL		Test:	SVOCM	IS Group3
Extraction Type	e:	Deca	inted : N	Level :	LOW	
Injection Volum	ne :	GPC Factor :	1.0	GPC Cleanup	: N	PH :
Prep Method :	SW3510C					
File ID/Qc Batch	: Dilution:	Prep Date		Date Analyzed	Prep Batch	ID
BN033231.D	1	08/02/24		08/03/24 10:24	PB162464	
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Unit
TARGETS	NT 1.4 1	0.020		0.020	0.11	/*
91-20-3	Naphthalene	0.030	U	0.030	0.11	ug/L
91-57-6	2-Methylnaphthalene	0.030	U	0.030	0.11	ug/L
208-96-8 83-32-9	Acenaphthylene	0.020	U	0.020	0.11 0.11	ug/L
83-32-9 86-73-7	Acenaphthene Fluorene	0.020 0.020	U U	0.020 0.020	0.11	ug/L
80-73-7 85-01-8	Phenanthrene	0.020	U U	0.020	0.11	ug/L ug/L
120-12-7	Anthracene	0.020	U U	0.020	0.11	ug/L ug/L
206-44-0	Fluoranthene	0.030	U U	0.030	0.11	ug/L ug/L
200-44-0 129-00-0	Pyrene	0.030	U	0.030	0.11	ug/L ug/L
56-55-3	Benzo(a)anthracene	0.020	U	0.020	0.11	ug/L ug/L
218-01-9	Chrysene	0.020	U	0.020	0.11	ug/L ug/L
205-99-2	Benzo(b)fluoranthene	0.040	U	0.040	0.11	ug/L
207-08-9	Benzo(k)fluoranthene	0.040	U	0.040	0.11	ug/L
50-32-8	Benzo(a)pyrene	0.040	U	0.060	0.11	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	0.040	U	0.040	0.11	ug/L
53-70-3	Dibenzo(a,h)anthracene	0.040	U	0.040	0.11	ug/L
191-24-2	Benzo(g,h,i)perylene	0.040	U	0.040	0.11	ug/L
123-91-1	1,4-Dioxane	0.070	U	0.070	0.22	ug/L
SURROGATES						
7297-45-2	2-Methylnaphthalene-d10	0.24		30 (30) - 150 (150)	61%	SPK: 0.
93951-69-0	Fluoranthene-d10	0.31		30 (30) - 150 (150)	77%	SPK: 0.
4165-60-0	Nitrobenzene-d5	0.24		30 (11) - 130 (175)	60%	SPK: 0.
321-60-8	2-Fluorobiphenyl	0.28		30 (10) - 130 (175)	71%	SPK: 0.
1718-51-0	Terphenyl-d14	0.36		30 (54) - 130 (171)	89%	SPK: 0.
NTERNAL STAN		2040	7 504			
3855-82-1	1,4-Dichlorobenzene-d4	2960 0520	7.524			
1146-65-2	Naphthalene-d8	9520 5300	10.297			
15067-26-2	Acenaphthene-d10	5300	14.144			

1517-22-2

Phenanthrene-d10

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16.915

10000



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С

				Report	t of Analy	sis				
Client:	JACOBS E	ngineeri	ng Grouj	o, Inc.			Date Collected:	C	08/01/24	
Project:	Former Sch	nlumberg	er Site P	rinceton NJ			Date Received:	C	08/01/24	
Client Sample ID:	923-K1-W	S-080124	4				SDG No.:	F	P 3440	
Lab Sample ID:	P3440-01						Matrix:	١	Water	
Analytical Method:	SW8270SI	М					% Solid:	C)	
Sample Wt/Vol:	910	Units:	mL				Final Vol:	1	000	uL
Soil Aliquot Vol:			uL				Test:	S	SVOCMS (Group3
Extraction Type :				Decan	ted : N		Level :	Ι	LOW	
Injection Volume :			GP	C Factor :	1.0		GPC Cleanup :	Ν	P	H :
Prep Method :	SW3510C									
File ID/Qc Batch:	Dilution:			Prep Date		Date A	nalyzed	Prep	Batch ID	
BN033231.D	1			08/02/24 09	9:25	08/03/	24 10:24	PB1	62464	
CAS Number Pa	arameter			Conc.	Qualifier	MDL		LOQ /	CRQL	Units
1719-03-5 C	hrysene-d12			8930	21.131					
1520-96-3 Ре	erylene-d12			10300	23.307					

- 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

P3440

- 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



6

A B C D

24
24
uL
MS Group3
PH :
ı ID
L Units
ug/L
SPK: 0.4

P3440

1517-22-2

Phenanthrene-d10

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16.915

10600



6
U

С

			Repor	t of Anal	ysis				
Client:	JACOBS E	ngineering	Group, Inc.			Date Collected:		08/01/24	
Project:	Former Sch	lumberger	Site Princeton N.	J		Date Received:		08/01/24	
Client Sample ID:	922-K1-WS	S-080124				SDG No.:		P3440	
Lab Sample ID:	P3440-04					Matrix:		Water	
Analytical Metho	d: SW8270SI	М				% Solid:		0	
Sample Wt/Vol:	930	Units: 1	mL			Final Vol:		1000	uL
Soil Aliquot Vol:		I	uL			Test:		SVOCMS	Group3
Extraction Type :			Decar	nted : N		Level :		LOW	
Injection Volume	:		GPC Factor :	1.0		GPC Cleanup :	Ν	F	РН:
Prep Method :	SW3510C								
File ID/Qc Batch:	Dilution:		Prep Date		Date A	nalyzed	Pr	ep Batch ID	
BN033234.D	1		08/02/24 0	9:25	08/03/	24 12:12	PI	3162464	
CAS Number	Parameter		Conc.	Qualifier	MDL		LOQ	/ CRQL	Units
1719-03-5	Chrysene-d12		7010	21.131					
1520-96-3	Perylene-d12		6900	23.312					

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- LOD = Limit of Detection
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- 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

6

LAB CHRONICLE

OrderID: Client: Contact:	P3440 JACOBS Engineering Group, Ir Mary I. Murphy	COBS Engineering Group, Inc.		OrderDate: Project: Location:	8/1/2024 12:28 Former Schlum D31,VOA Ref. :	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3440-01	923-K1-WS-080124	Water			08/01/24			08/01/24
			SVOCMS Group3	8270-Modifie d		08/02/24	08/03/24	
			SVOCMS Group6	8270E		08/02/24	08/07/24	
P3440-04	922-K1-WS-080124	Water			08/01/24			08/01/24
			SVOCMS Group3	8270-Modifie d		08/02/24	08/03/24	
			SVOCMS Group6	8270E		08/02/24	08/07/24	



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В	

7

Hit Summary Sheet SW-846

SDG No.:	P3440				
Client:	JACOBS Engine	eering Group, Inc.			
Sample ID Client ID :	Client ID	Client ID Matrix Parameter		Concentration C MDL	RDL Units
			Total Svoc :	0.00	
			Total Concentration:	0.00	





7

A B C D



7

		Report	t of Ana	lysis			
Client:	JACOBS Engineerin	ng Group, Inc.			Date Collected:	08/01/24	
Project:	Former Schlumberg	er Site Princeton NJ			Date Received:	08/01/24	
Client Sample II	c c				SDG No.:	P3440	
Lab Sample ID:		T			Matrix:	Water	
Analytical Meth	od: SW8270				% Solid:	0	
Sample Wt/Vol:	880 Units:	mL			Final Vol:	1000	uL
Soil Aliquot Vol	:	uL			Test:	SVOCM	S Group6
Extraction Type	:	Decan	ted : N	N	Level :	LOW	
Injection Volume	e :	GPC Factor :	1.0		GPC Cleanup :	Ν	PH :
Prep Method :	SW3510C				ľ		
-	Dibtion	Draw Data		Dett	Analyzad	Duce Date 1	
File ID/Qc Batch:		Prep Date			Analyzed	Prep Batch	D
BF138837.D	1	08/02/24 09	9:23	08/07/24 12:34		PB162463	
CAS Number	Parameter	Conc.	Qualifier	MDL		LOQ / CRQL	Units
FARGETS 110-86-1	Pyridine	1.80	U	1.80		5.70	ug/L
100-52-7	Benzaldehyde	4.50	U	4.50		11.4	ug/L ug/L
95-48-7	2-Methylphenol	1.30	U	1.30		5.70	ug/L
98-86-2	Acetophenone	1.30	U	1.30		5.70	ug/L
65794-96-9	3+4-Methylphenols	1.30	U	1.30		11.4	ug/L ug/L
98-95-3	Nitrobenzene	1.40	U	1.40		5.70	ug/L ug/L
20-83-2	2,4-Dichlorophenol	1.00	U	1.00		5.70	ug/L
91-20-3	Naphthalene	1.20	U	1.20		5.70	ug/L
37-68-3	Hexachlorobutadiene	1.40	U	1.40		5.70	ug/L
91-57-6	2-Methylnaphthalene	1.30	Ū	1.30		5.70	ug/L
88-06-2	2,4,6-Trichlorophenol	1.00	U	1.00		5.70	ug/L
95-95-4	2,4,5-Trichlorophenol	1.10	U	1.10		5.70	ug/L
208-96-8	Acenaphthylene	1.20	U	1.20		5.70	ug/L
33-32-9	Acenaphthene	0.92	U	0.92		5.70	ug/L
132-64-9	Dibenzofuran	1.10	U	1.10		5.70	ug/L
6-73-7	Fluorene	1.10	U	1.10		5.70	ug/L
18-74-1	Hexachlorobenzene	1.30	U	1.30		5.70	ug/L
37-86-5	Pentachlorophenol	2.10	U	2.10		11.4	ug/L
5-01-8	Phenanthrene	1.00	U	1.00		5.70	ug/L
6-74-8	Carbazole	1.30	U	1.30		5.70	ug/L
4-74-2	Di-n-butylphthalate	1.70	U	1.70		5.70	ug/L ug/L
06-44-0	Fluoranthene	1.50	U	1.50		5.70	ug/L ug/L
29-00-0	Pyrene	1.20	U	1.20		5.70	ug/L
56-55-3	Benzo(a)anthracene	1.10	U	1.10		5.70	ug/L
218-01-9	Chrysene	0.98	U	0.98		5.70	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	2.10	U	2.10		5.70	ug/L
205-99-2	Benzo(b)fluoranthene	1.30	U	1.30		5.70	ug/L
207-08-9	Benzo(k)fluoranthene	1.40	U	1.40		5.70	ug/L ug/L
50 22 8	Benzo(a)pureno	1.40	U	1.40		5.70	ug/L

50-32-8

Benzo(a)pyrene

U

1.90

5.70

ug/L

1.90



7

Report of Analysis Client: JACOBS Engineering Group, Inc. Date Collected: 08/01/24 Project: Former Schlumberger Site Princeton NJ Date Received: 08/01/24 Client Sample ID: 923-K1-WS-080124 SDG No .: P3440 Lab Sample ID: P3440-01 Matrix: Water Analytical Method: SW8270 % Solid: 0 Final Vol: 1000 uL Sample Wt/Vol: 880 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 : Dilution: File ID/Qc Batch: Prep Date Date Analyzed Prep Batch ID BF138837.D 1 08/02/24 09:23 08/07/24 12:34 PB162463 MDL Units CAS Number Conc. Qualifier LOQ / CRQL Parameter 193-39-5 1.20 U 1.20 5.70 ug/L Indeno(1,2,3-cd)pyrene 53-70-3 Dibenzo(a h)anthracene 1 30 H 1 30 5 70 11**σ**/Ι

53-70-3	Dibenzo(a,h)anthracene	1.30	U	1.30	5.70	ug/L
191-24-2	Benzo(g,h,i)perylene	1.30	U	1.30	5.70	ug/L
123-91-1	1,4-Dioxane	1.40	U	1.40	5.70	ug/L
90-12-0	1-Methylnaphthalene	0.98	U	0.98	5.70	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	63.6		15 (10) - 110 (139)	42%	SPK: 150
13127-88-3	Phenol-d6	38.2		15 (10) - 110 (134)	25%	SPK: 150
4165-60-0	Nitrobenzene-d5	94.8		30 (49) - 130 (133)	95%	SPK: 100
321-60-8	2-Fluorobiphenyl	98.3		30 (52) - 130 (132)	98%	SPK: 100
118-79-6	2,4,6-Tribromophenol	154		15 (32) - 110 (145)	102%	SPK: 150
1718-51-0	Terphenyl-d14	114		30 (36) - 130 (145)	114%	SPK: 100
INTERNAL STAN	DARDS					
3855-82-1	1,4-Dichlorobenzene-d4	45200	6.84			
1146-65-2	Naphthalene-d8	184000	8.116			
15067-26-2	Acenaphthene-d10	98300	9.869			
1517-22-2	Phenanthrene-d10	162000	11.357			
1719-03-5	Chrysene-d12	84400	13.992			
1520-96-3	Perylene-d12	89900	15.457			

- 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



7

		Report	t of Ana	lysis				
Client:	JACOBS Engineerin			Date Collected:	08/01/24			
Project:	Former Schlumberger Site Princeton NJ				Date Received:	08/01/24		
Client Sample II	-				SDG No.:	P3440		
Lab Sample ID:		r			Matrix:	Water		
Analytical Meth	od: SW8270				% Solid:	0		
Sample Wt/Vol:	890 Units:	mL			Final Vol:	1000	uL	
Soil Aliquot Vol:		uL			Test:	SVOCM	S Group6	
Extraction Type	Extraction Type :		ted : N	N	Level :	LOW		
Injection Volume	e :	GPC Factor :	1.0		GPC Cleanup :	Ν	PH :	
Prep Method :	SW3510C							
File ID/Os Databa	Dilution	Dran Data		Dat-	Analyzad	Dron Databi		
	File ID/Qc Batch: Dilution:		Prep Date		Date Analyzed		Prep Batch ID	
BF138840.D 1		08/02/24 09	9:23	08/07/24 14:05		PB162463		
CAS Number	Parameter	Conc.	Qualifier	· MDL		LOQ / CRQL	Units	
DODTO								
FARGETS 110-86-1	Pyridine	1.70	U	1.70		5.60	ug/L	
.00-52-7	Benzaldehyde	4.50	U	4.50		11.2	ug/L ug/L	
95-48-7	2-Methylphenol	1.30	U	1.30		5.60	ug/L ug/L	
98-86-2	Acetophenone	1.20	U	1.20		5.60	ug/L ug/L	
55794-96-9	3+4-Methylphenols	1.20	U	1.20		11.2	ug/L ug/L	
)8-95-3	Nitrobenzene	1.40	U	1.30		5.60	ug/L ug/L	
120-83-2	2,4-Dichlorophenol	0.99	U	0.99		5.60	ug/L ug/L	
1-20-3	Naphthalene	1.10	U	1.10		5.60	ug/L ug/L	
37-68-3	Hexachlorobutadiene	1.40	U	1.40		5.60	ug/L ug/L	
91-57-6	2-Methylnaphthalene	1.30	U	1.30		5.60	ug/L	
38-06-2	2,4,6-Trichlorophenol	1.00	U	1.00		5.60	ug/L	
95-95-4	2,4,5-Trichlorophenol	1.10	U	1.10		5.60	ug/L ug/L	
208-96-8	Acenaphthylene	1.20	U	1.20		5.60	ug/L ug/L	
83-32-9	Acenaphthene	0.91	U	0.91		5.60	ug/L ug/L	
32-64-9	Dibenzofuran	1.00	U	1.00		5.60	ug/L ug/L	
6-73-7	Fluorene	1.10	U	1.10		5.60	ug/L ug/L	
18-74-1	Hexachlorobenzene	1.30	U	1.30		5.60	ug/L ug/L	
37-86-5	Pentachlorophenol	2.10	U	2.10		11.2	ug/L ug/L	
35-01-8	Phenanthrene	1.00	U	1.00		5.60	ug/L ug/L	
36-74-8	Carbazole	1.30	U	1.30		5.60	ug/L ug/L	
4-74-2	Di-n-butylphthalate	1.70	U	1.70		5.60	ug/L ug/L	
06-44-0	Fluoranthene	1.40	U	1.70		5.60	ug/L ug/L	
29-00-0	Pyrene	1.40	U	1.40		5.60	ug/L ug/L	
6-55-3	Benzo(a)anthracene	1.10	U	1.10		5.60	ug/L ug/L	
218-01-9	Chrysene	0.97	U	0.97		5.60	ug/L ug/L	
117-81-7	Bis(2-ethylhexyl)phthalate	2.10	U	2.10		5.60	ug/L	
205-99-2	Benzo(b)fluoranthene	1.30	U	1.30		5.60	ug/L ug/L	
207-08-9	Benzo(k)fluoranthene	1.30	U	1.30		5.60	ug/L ug/L	
50 22 8	Benzo(a)pureno	1.00	U	1.00		5.60	ug/L	

50-32-8

Benzo(a)pyrene

U

1.90

5.60

ug/L

1.90



7

Report of Analysis Client: JACOBS Engineering Group, Inc. Date Collected: 08/01/24 Project: Former Schlumberger Site Princeton NJ Date Received: 08/01/24 Client Sample ID: 922-K1-WS-080124 SDG No .: P3440 Lab Sample ID: P3440-04 Matrix: Water Analytical Method: SW8270 % Solid: 0 890 Final Vol: 1000 uL Sample Wt/Vol: 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 BF138840.D 1 08/02/24 09:23 08/07/24 14:05 PB162463 MDL Units CAS Number Conc. Qualifier LOQ / CRQL Parameter

193-39-5	Indeno(1,2,3-cd)pyrene	1.10	U	1.10	5.60	ug/L
53-70-3	Dibenzo(a,h)anthracene	1.30	U	1.30	5.60	ug/L
191-24-2	Benzo(g,h,i)perylene	1.30	U	1.30	5.60	ug/L
123-91-1	1,4-Dioxane	1.40	U	1.40	5.60	ug/L
90-12-0	1-Methylnaphthalene	0.97	U	0.97	5.60	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	57.9		15 (10) - 110 (139)	39%	SPK: 150
13127-88-3	Phenol-d6	34.5		15 (10) - 110 (134)	23%	SPK: 150
4165-60-0	Nitrobenzene-d5	92.4		30 (49) - 130 (133)	92%	SPK: 100
321-60-8	2-Fluorobiphenyl	96.8		30 (52) - 130 (132)	97%	SPK: 100
118-79-6	2,4,6-Tribromophenol	147		15 (32) - 110 (145)	98%	SPK: 150
1718-51-0	Terphenyl-d14	108		30 (36) - 130 (145)	108%	SPK: 100
INTERNAL STA	NDARDS					
3855-82-1	1,4-Dichlorobenzene-d4	41800	6.84			
1146-65-2	Naphthalene-d8	170000	8.116			
15067-26-2	Acenaphthene-d10	89900	9.869			
1517-22-2	Phenanthrene-d10	146000	11.357			
1719-03-5	Chrysene-d12	76100	13.998			
1520-96-3	Perylene-d12	79900	15.457			

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



LAB CHRONICLE

OrderID: Client: Contact:	P3440 JACOBS Engineering Group, Ir Mary I. Murphy	OrderDate: Project: Location:	8/1/2024 12:28:00 PM Former Schlumberger Site Princeton NJ D31,VOA Ref. #3 Water					
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3440-01	923-K1-WS-080124	Water			08/01/24			08/01/24
			SVOCMS Group3	8270-Modifie d		08/02/24	08/03/24	
			SVOCMS Group6	8270E		08/02/24	08/07/24	
P3440-04	922-K1-WS-080124	Water			08/01/24			08/01/24
			SVOCMS Group3	8270-Modifie d		08/02/24	08/03/24	
			SVOCMS Group6	8270E		08/02/24	08/07/24	



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

Hit Summary Sheet SW-846

SDG No.:	P3440			Order ID:		P3440		
Client:	JACOBS Engineering Grou	up, Inc.		Project ID):	Former Schlumbe	erger Site Princetor	n NJ
Sample ID	Client ID	Matrix	Parameter	Concentration	С	MDL	RDL	Units
Client ID :	923-K1-WS-080124							
P3440-01	923-K1-WS-080124	Water	Aluminum	50.9		28.3	50.0	ug/L
P3440-01	923-K1-WS-080124	Water	Barium	58.9		6.28	50.0	ug/L
P3440-01	923-K1-WS-080124	Water	Boron	47.6	J	9.95	50.0	ug/L
P3440-01	923-K1-WS-080124	Water	Calcium	19400		33.0	1000	ug/L
P3440-01	923-K1-WS-080124	Water	Chromium	0.76	J	0.66	5.00	ug/L
P3440-01	923-K1-WS-080124	Water	Iron	2490		18.5	50.0	ug/L
P3440-01	923-K1-WS-080124	Water	Magnesium	3680		39.4	1000	ug/L
P3440-01	923-K1-WS-080124	Water	Manganese	777		1.46	10.0	ug/L
P3440-01	923-K1-WS-080124	Water	Nickel	1.18	J	0.85	20.0	ug/L
P3440-01	923-K1-WS-080124	Water	Potassium	3290		685	1000	ug/L
P3440-01	923-K1-WS-080124	Water	Sodium	79700		237	1000	ug/L
P3440-01	923-K1-WS-080124	Water	Strontium	129		2.32	10.0	ug/L
P3440-01	923-K1-WS-080124	Water	Zinc	10.1	J	1.75	20.0	ug/L
Client ID :	922-K1-WS-080124							
P3440-04	922-K1-WS-080124	Water	Aluminum	45.7	J	28.3	50.0	ug/L
P3440-04	922-K1-WS-080124	Water	Barium	82.2		6.28	50.0	ug/L
P3440-04	922-K1-WS-080124	Water	Boron	58.2		9.95	50.0	ug/L
P3440-04	922-K1-WS-080124	Water	Calcium	35100		33.0	1000	ug/L
P3440-04	922-K1-WS-080124	Water	Cobalt	4.73	J	0.50	15.0	ug/L
P3440-04	922-K1-WS-080124	Water	Iron	4080		18.5	50.0	ug/L
P3440-04	922-K1-WS-080124	Water	Magnesium	8070		39.4	1000	ug/L
P3440-04	922-K1-WS-080124	Water	Manganese	1020		1.46	10.0	ug/L
P3440-04	922-K1-WS-080124	Water	Nickel	0.95	J	0.85	20.0	ug/L
P3440-04	922-K1-WS-080124	Water	Potassium	6330		685	1000	ug/L
P3440-04	922-K1-WS-080124	Water	Sodium	187000		237	1000	ug/L
P3440-04	922-K1-WS-080124	Water	Strontium	243		2.32	10.0	ug/L
P3440-04	922-K1-WS-080124	Water	Zinc	22.7		1.75	20.0	ug/L

B C

D





8

A B C D



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	923-K1-WS-080124	SDG No.:	P3440
Lab Sample ID:	P3440-01	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	50.9		1	28.3	50.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-36-0	Antimony	2.06	U	1	2.06	25.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-38-2	Arsenic	3.48	U	1	3.48	10.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-39-3	Barium	58.9		1	6.28	50.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-41-7	Beryllium	0.13	U	1	0.13	3.00	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-42-8	Boron	47.6	J	1	9.95	50.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-43-9	Cadmium	0.094	U	1	0.094	3.00	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-70-2	Calcium	19400		1	33.0	1000	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-47-3	Chromium	0.76	J	1	0.66	5.00	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-48-4	Cobalt	0.50	U	1	0.50	15.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-50-8	Copper	7.07	U	1	7.07	10.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7439-89-6	Iron	2490		1	18.5	50.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7439-92-1	Lead	3.51	U	1	3.51	6.00	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7439-95-4	Magnesium	3680		1	39.4	1000	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7439-96-5	Manganese	777		1	1.46	10.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7439-97-6	Mercury	0.081	U	1	0.081	0.20	ug/L	08/12/24 16:13	08/13/24 10:23	SW7470A	
7439-98-7	Molybdenum	3.67	U	1	3.67	100	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-02-0	Nickel	1.18	J	1	0.85	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-09-7	Potassium	3290		1	685	1000	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7782-49-2	Selenium	5.88	U	1	5.88	10.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-22-4	Silver	0.58	U	1	0.58	5.00	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-23-5	Sodium	79700		1	237	1000	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-24-6	Strontium	129	Ν	1	2.32	10.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-28-0	Thallium	2.32	U	1	2.32	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-31-5	Tin	1.89	U	1	1.89	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-32-6	Titanium	2.35	U	1	2.35	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-62-2	Vanadium	3.06	U	1	3.06	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010
7440-66-6	Zinc	10.1	J	1	1.75	20.0	ug/L	08/05/24 09:15	08/07/24 17:00	SW6010	SW3010

Color Before:	Colorless	Clarity Before:	Clear	Texture:
Color After:	Colorless	Clarity After:	Clear	Artifacts:
Comments:	Mercury			
MDL = Methodologiest MDL = Limit of D = Dilution	of Quantitation od Detection Limit	requirements		J = Estimated Value B = Analyte Found in Associated Method Blank * = indicates the duplicate analysis is not within control limits. E = Indicates the reported value is estimated because of the presence of interference. OR = Over Range N =Spiked sample recovery not within control limits

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B C

D



Client:	JACOBS Engineering Group, Inc.	Date Collected:	08/01/24
Project:	Former Schlumberger Site Princeton NJ	Date Received:	08/01/24
Client Sample ID:	922-K1-WS-080124	SDG No.:	P3440
Lab Sample ID:	P3440-04	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	45.7	J	1	28.3	50.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-36-0	Antimony	2.06	U	1	2.06	25.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-38-2	Arsenic	3.48	U	1	3.48	10.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-39-3	Barium	82.2		1	6.28	50.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-41-7	Beryllium	0.13	U	1	0.13	3.00	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-42-8	Boron	58.2		1	9.95	50.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-43-9	Cadmium	0.094	U	1	0.094	3.00	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-70-2	Calcium	35100		1	33.0	1000	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-47-3	Chromium	0.66	U	1	0.66	5.00	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-48-4	Cobalt	4.73	J	1	0.50	15.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-50-8	Copper	7.07	U	1	7.07	10.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7439-89-6	Iron	4080		1	18.5	50.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7439-92-1	Lead	3.51	U	1	3.51	6.00	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7439-95-4	Magnesium	8070		1	39.4	1000	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7439-96-5	Manganese	1020		1	1.46	10.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7439-97-6	Mercury	0.081	U	1	0.081	0.20	ug/L	08/12/24 16:13	08/13/24 10:39	SW7470A	L
7439-98-7	Molybdenum	3.67	U	1	3.67	100	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-02-0	Nickel	0.95	J	1	0.85	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-09-7	Potassium	6330		1	685	1000	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7782-49-2	Selenium	5.88	U	1	5.88	10.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-22-4	Silver	0.58	U	1	0.58	5.00	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-23-5	Sodium	187000		1	237	1000	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-24-6	Strontium	243	Ν	1	2.32	10.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-28-0	Thallium	2.32	U	1	2.32	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-31-5	Tin	1.89	U	1	1.89	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-32-6	Titanium	2.35	U	1	2.35	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-62-2	Vanadium	3.06	U	1	3.06	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010
7440-66-6	Zinc	22.7		1	1.75	20.0	ug/L	08/05/24 09:15	08/07/24 18:38	SW6010	SW3010

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

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8

B C

D



Α

D

8

LAB CHRONICLE

OrderID: Client: Contact:	P3440 JACOBS Engineering Group, Ir Mary I. Murphy	ıc.		OrderDate: Project: Location:	8/1/2024 12:28 Former Schlum D31,VOA Ref. 3	berger Site Pri	nceton NJ	
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
P3440-01	923-K1-WS-080124	Water			08/01/24			08/01/24
			Mercury	7470A		08/12/24	08/13/24	
			Metals Group4	6010D		08/05/24	08/07/24	
P3440-04	922-K1-WS-080124	Water			08/01/24			08/01/24
			Mercury	7470A		08/12/24	08/13/24	
			Metals Group4	6010D		08/05/24	08/07/24	





9

В



Client:	JACOBS Engineering Group,	Inc.	Date Col	lected: 08/01	/24 09:15
Project:	Former Schlumberger Site Pri	nceton NJ	Date Rec	eived: 08/01	/24
Client Sample ID:	923-K1-WS-080124		SDG No.	: P3440)
Lab Sample ID:	P3440-01		Matrix:	WATE	ER
			% Solid:	0	
Parameter	Conc. Qua. DF MDL	LOQ / CRQL	Units Prej	o Date Date Ana.	Ana Met.
Dissolved Hexavalent Chromium	0.0030 U 1 0.0030	0.010	mg/L	08/01/24 1	5:34 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
- * = indicates the duplicate analysis is not within control limits.
- E = Indicates the reported value is estimated because of the presence of interference.
- OR = Over Range
- N =Spiked sample recovery not within control limits



Client: Project: Client Sample ID: Lab Sample ID:		ngineering Group, lumberger Site Pri -080124		I	Date Collected: Date Received: SDG No.: Matrix:	08/01/24 11:10 08/01/24 P3440 WATER		
				9	% Solid:	0		
Parameter Dissolved Hexavalent Chromium	Conc. Qua. 0.0030 U	DF MDL 1 0.0030	LOQ / CRQL 0.010	Units mg/L	Prep Date	Date Ana. 08/01/24 15:3	Ana Met. 8 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
- * = indicates the duplicate analysis is not within control limits.
- E = Indicates the reported value is estimated because of the presence of interference.
- OR = Over Range
- N =Spiked sample recovery not within control limits



С

Q

LAB CHRONICLE

OrderID: Client: Contact:	P3440 JACOBS Engineering Group, Ir Mary I. Murphy	nc.		OrderDate: Project: Location:		4 12:28:00 PM Schlumberger Site Princeton NJ DA Ref. #3 Water					
LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received			
P3440-01	923-K1-WS-080124	WATER			08/01/24 09:15			08/01/24			
			Hexavalent Chromium	7196A			08/01/24 15:34				
P3440-04	922-K1-WS-080124	WATER			08/01/24 11:10			08/01/24			
			Hexavalent Chromium	7196A			08/01/24 15:38				



<u>SHIPPING</u> DOCUMENTS

10

	CUSTODY RECORD	(908) 789-8900 • Fax (908) 789-8922										(1 1		
No. of the Party	CLIENT PROJECT INFORMATION														ORMATION	1. 11 July 1	-		
COMPANY: J		PROJECT NAME: STC PTC BILL TO: Mary										lary	Murp	hy		PO#:			
ADDRESS: 4	12 Mt Kemble Ave Suik 4/00	PROJE	CT N	0.: D	377997	Z LOC	ATION:	Prince	ton J	inchan	ADD	RESS:			/				
CITY Mor	ISTATE: NJ ZIP: 07960					hig Mu					CITY					STA	TE:	ZIP;	
ATTENTION: -	J.hn Yutank	e-mail:	Ma	wy.	Morphi	Ja Jacol	5.00	m			ATTE	NTION	e e e e e e e e e e e e e e e e e e e			PHO	ONE:		
PHONE (28)	ATA TURNAROUND INFORMATION		12	01)9	136-058	6 F/	AX:								AN	ALYSI	11		
FAX (RUSH) HARDCOPY (DA EDD: *TO BE APPROV STANDARD HAR	Leve	el 1 (Re el 2 (Re el 3 (Re aw Da	esults esults esults ta)	Only) □ + QC) □ + QC □	RABLE IN Level 4 (QC NJ Reduce NYS ASP 4 Other	C + Full d D U	Raw Dati S EPA C	ĹP 🌍	2601 F	All SITU Of all I	to the	11944			3 9				
CHEMTECH			1 mm	/IPLE		MPLE	LES	4/		1		SERVA	TIVES	1		•		MMENTS	
SAMPLE ID	PROJECT SAMPLE IDENTIFICATION	SAMPLE MATRIX		GRAB B	DATE	ECTION TIME	# OF BOTTLES	A/E	E	B/E 3	E 4	5	6	7	8	9	A-HCI B-HN03 C-H2SO4	D-NaOH E-ICE F-OTHER	
1.	923-K1-WS-573+080124	ws		X	8/1/24	0915	18	6	6	3	3				-	1	MS/M		1
2.	922- KI-WS-080124	WS		1	8/124	110	6	2	2	1	Ĩ				-		- /	<i>v</i>	1
3.	16-01-08012	DI		-	8/1/24	1700	1	1		1			1		1	1			1
4.	7				1111	1,200													1
5.								1			1				1	1			-
6.															-				-
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8.										1									-
9.										1					-				-
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p. he	SAMPLE CUSTODY MUST BE DOCI	JMENTEI	D BEI	LOW	EACH TI	L ME SAMP	LES C	L HANGE	POSS	SESSIO		L UDING			I ELIVEF	RY .	1	Elson inter	
RELINQUISHED BY 1. RELINQUISHED BY 2.	AMPLER: DATE/TIME: 1220 RECRIVED BY: 8/1/2 AMPLER: DATE/TIME: RECRIVED BY: 2.	\bigcirc	_	220		ions of bottles nts: See nd ECC	and the second second		_	COMPLIAN	T D NO	N COMPLI	ANT D	COOLER 1	TEMP		, ECD-S	°C V9Cs ,	
RELINDUIGHED BY	BIRY 3.				Page	a f		CLIEN		Hand D								t Complete	1
DarP3440	WHITE - CHEMTER	CH COPY FO	B RET	URN TO	Page	of	51 CHE	CHEMT			ed Up - SAMPLE		eld Sam	oling			C YES	O NO	1



Laboratory Certification

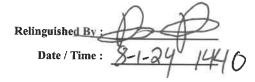
Certified By	License No.
CAS EPA CLP Contract	68HERH20D0011
Connecticut	PH-0830
DOD ELAP (L-A-B)	L2219
Maine	2022022
Maryland	296
New Hampshire	255423
New Jersey	20012
New York	11376
Pennsylvania	68-00548
Soil Permit	P330-21-00137
Texas	T104704488



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

LOGIN REPORT/SAMPLE TRANSFER

Client Contact : Invoice Name :		 P3440 JACO05 JACOBS Engineering Grou Mary I. Murphy JACOBS Engineering Grou Mary I. Murphy 			Project Name		8/1/2024 12:28:00 PM Former Schlumberger Site 1 8/1/2024 2:00:00 PM		Project Mgr : Report Type : EDD Type : Hard Copy Date : Date Signoff :			
LAB ID	CLIEN	T ID		MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD		FAX DATE	DUE DATES
P3440-01	92	23-K1-WS-0	80124	Water	08/01/2024	09:15						
							VOCMS Group6		8260-Low	10 Bus. Days		
P3440-02		P3440-011	MS	Water	08/01/2024	09:15						
P3440-03		P3440-01M	ISD	Water	08/01/2024	09:15	VOCMS Group6		8260-Low	10 Bus. Days		
							VOCMS Group6		8260-Low	10 Bus. Days		
P3440-04	92	2-K1-WS-08	80124	Water	08/01/2024	11:10						
							VOCMS Group6		8260-Low	10 Bus. Days		
P3440-05		TB-01-0801	24	Water	08/01/2024	12:00						
							VOCMS Group6		8260-Low	10 Bus. Days		



14.20 Rg # 4 M **Received By :** Date / Time : 8 (124

Storage Area : VOA Refridgerator Room

Page 1 of 1 51 of 51