

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 : P3451

ATTENTION : Mary I. Murphy



Laboratory Certification ID # 20012



| | |
|------------------------------------|----|
| 1) Signature Page | 3 |
| 2) Case Narrative | 5 |
| 2.1) VOCMS Group6- Case Narrative | 5 |
| 2.2) SVOCMS Group3- Case Narrative | 7 |
| 2.3) SVOCMS Group6- Case Narrative | 9 |
| 2.4) Metals-MS- Case Narrative | 11 |
| 2.5) Genchem- Case Narrative | 13 |
| 3) Qualifier Page | 14 |
| 4) QA Checklist | 16 |
| 5) VOCMS Group6 Data | 17 |
| 6) SVOCMS Group3 Data | 24 |
| 7) SVOCMS Group6 Data | 29 |
| 8) Metals-MS Data | 34 |
| 9) Genchem Data | 38 |
| 10) Shipping Document | 41 |
| 10.1) CHAIN OF CUSTODY | 42 |
| 10.2) Lab Certificate | 43 |
| 10.3) Internal COC | 44 |

| |
|----|
| 1 |
| 2 |
| 3 |
| 4 |
| 5 |
| 6 |
| 7 |
| 8 |
| 9 |
| 10 |

Cover Page

Order ID : P3451

Project ID : Former Schlumberger Site Princeton NJ

Client : JACOBS Engineering Group, Inc.

Lab Sample Number

P3451-01
P3451-02

Client Sample Number

921-J-WS-080124
TB-03-080124

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 : _____

Date: 8/27/2024

NYDOH CERTIFICATION NO - 11376

NJDEP CERTIFICATION NO - 20012

DATA OF KNOWN QUALITY CONFORMANCE/NON-CONFORMANCE SUMMARY QUESTIONNAIRE

2

Laboratory Name : Alliance Technical Group LLC Client : JACOBS Engineering Group, Inc.
 Project Location : Princeton, NJ Project Number : D3779922
 Laboratory Sample ID(s) : P3451 Sampling Date(s) : 8/01/2024
 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? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
| 1A | Were the method specified handling, preservation, and holding time requirements met? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
| 1B | EPH Method: Was the EPH method conducted without significant modifications (see Section 11.3 of respective DKQ methods) | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> 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)? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
| 3 | Were samples received at an appropriate temperature (4±2° C)? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A |
| 4 | Were all QA/QC performance criteria specified in the NJDEP DKQP standards achieved? | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No |
| 5 | a)Were reporting limits specified or referenced on the chain-of-custody or communicated to the laboratory prior to sample receipt? b)Were these reporting limits met? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> 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? | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
| 7 | Are project-specific matrix spikes and/or laboratory duplicates included in this data set? | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> 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."

CASE NARRATIVE

JACOBS Engineering Group, Inc.

Project Name: Former Schlumberger Site Princeton NJ

Project # N/A

Chemtech Project # P3451

Test Name: VOCMS Group6

A. Number of Samples and Date of Receipt:

2 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, 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:

Samples for MS/MSD for VOC analysis were not provided with this set of samples therefore lab used from another project.

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.



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

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_____

CASE NARRATIVE

JACOBS Engineering Group, Inc.

Project Name: Former Schlumberger Site Princeton NJ

Project # N/A

Chemtech Project # P3451

Test Name: SVOCMS Group3

A. Number of Samples and Date of Receipt:

2 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, 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%] this compound 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 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.

Signature_____

CASE NARRATIVE

JACOBS Engineering Group, Inc.

Project Name: Former Schlumberger Site Princeton NJ

Project # N/A

Chemtech Project # P3451

Test Name: SVOCMS Group6

A. Number of Samples and Date of Receipt:

2 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, 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 sample 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.

Signature_____

CASE NARRATIVE

JACOBS Engineering Group, Inc.

Project Name: Former Schlumberger Site Princeton NJ

Project # N/A

Chemtech Project # P3451

Test Name: Metals Group4,Mercury

A. Number of Samples and Date of Receipt:

2 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, 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 analysis met criteria for all samples.

The Matrix Spike (923-K1-WS-080124MS) analysis met criteria for all samples except for Silver, Strontium, Titanium due to matrix interference.

The Matrix Spike Duplicate (923-K1-WS-080124MSD) analysis met criteria for all samples except for Silver, Strontium, Titanium due to matrix interference.

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:

The data package has been revised due to the Analytical Method changed for Metals as per client request.

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.

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_____



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 # P3451

Test Name: Hexavalent Chromium

A. Number of Samples and Date of Receipt:

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

B. Parameters:

According to the Chain of Custody document, the following analyses were requested: Hexavalent Chromium, 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 |
| E | Indicates the reported value is estimated because of the presence of interference |
| M | Indicates Duplicate injection precision not met. |
| N | 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 | 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 analyzed |
| OR | Indicates 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 |
| H | 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: <ol style="list-style-type: none"> (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 flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others. |
| B | Indicates the analyte was found in the blank as well as the sample report as “12 B”. |
| E | 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. |
| P | 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”. |
| N | 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. |
| A | 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 #: P3451

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)

✓

Check chain-of-custody for proper relinquish/return of samples

✓

Is the chain of custody signed and complete

✓

Check internal chain-of-custody for proper relinquish/return of samples /sample extracts

✓

Collect information for each project id from server. Were all requirements followed

✓

COVER PAGE:

Do numbers of samples correspond to the number of samples in the Chain of Custody on login page

✓

Do lab numbers and client Ids on cover page agree with the Chain of Custody

✓

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 Custody

✓

Were the samples received within hold time

✓

Were any problems found with the samples at arrival recorded in the Sample Management Laboratory Chronicle

✓

ANALYTICAL:

Was method requirement followed?

✓

Was client requirement followed?

✓

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

✓

1st Level QA Review Signature: SOHIL JODHANI

Date: 08/27/2024

2nd Level QA Review Signature: _____

Date: _____

Hit Summary Sheet
SW-846

SDG No.: P3451

Client: JACOBS Engineering Group, Inc.

| Sample ID | Client ID | Matrix | Parameter | Concentration | C | MDL | RDL | Units |
|-------------------|------------------------|--------|-----------------------------|---------------|---|------|------|-------|
| Client ID: | 921-J-WS-080124 | | | | | | | |
| P3451-01 | 921-J-WS-080124 | Water | Acetone | 10.7 | | 1.40 | 5.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Toluene | 2.00 | | 0.18 | 1.00 | ug/L |
| | | | Total Voc : | | | 12.7 | | |
| | | | Total Concentration: | | | 12.7 | | |

A

B

C

D



SAMPLE DATA

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | 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 |
| VN083283.D | 1 | | 08/14/24 07:16 | 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 | 10.7 | | 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 | 2.00 | | 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 |

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | 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 |
| VN083283.D | 1 | | 08/14/24 07:16 | 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 | 57.0 | | 70 (74) - 130 (125) | 114% | SPK: 50 |
| 1868-53-7 | Dibromofluoromethane | 52.6 | | 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 | 57.9 | | 70 (77) - 130 (121) | 116% | SPK: 50 |
| INTERNAL STANDARDS | | | | | | |
| 363-72-4 | Pentafluorobenzene | 131000 | 8.224 | | | |
| 540-36-3 | 1,4-Difluorobenzene | 258000 | 9.1 | | | |
| 3114-55-4 | Chlorobenzene-d5 | 270000 | 11.865 | | | |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 123000 | 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

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: | TB-03-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-02 | 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 |
| VN083284.D | 1 | | 08/14/24 07:40 | 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 |

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: | TB-03-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-02 | 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 |
| VN083284.D | 1 | | 08/14/24 07:40 | 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 | 56.8 | | 70 (74) - 130 (125) | 114% | SPK: 50 |
| 1868-53-7 | Dibromofluoromethane | 52.3 | | 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 | 55.8 | | 70 (77) - 130 (121) | 112% | SPK: 50 |
| INTERNAL STANDARDS | | | | | | |
| 363-72-4 | Pentafluorobenzene | 139000 | 8.224 | | | |
| 540-36-3 | 1,4-Difluorobenzene | 272000 | 9.1 | | | |
| 3114-55-4 | Chlorobenzene-d5 | 277000 | 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

LAB CHRONICLE

| | | | |
|-----------------|--------------------------------|-------------------|---------------------------------------|
| OrderID: | P3451 | OrderDate: | 8/1/2024 4:38:00 PM |
| Client: | JACOBS Engineering Group, Inc. | Project: | Former Schlumberger Site Princeton NJ |
| Contact: | Mary I. Murphy | Location: | D31,VOA Ref. #3 Water |

| LabID | ClientID | Matrix | Test | Method | Sample Date | Prep Date | Anal Date | Received |
|----------|-----------------|--------|--------------|----------|-------------|-----------|-----------|----------|
| P3451-01 | 921-J-WS-080124 | Water | VOCMS Group6 | 8260-Low | 08/01/24 | | 08/14/24 | 08/01/24 |
| P3451-02 | TB-03-080124 | Water | VOCMS Group6 | 8260-Low | 08/01/24 | | 08/14/24 | 08/01/24 |



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

Hit Summary Sheet
SW-846

SDG No.: P3451
Client: JACOBS Engineering Group, Inc.

| Sample ID | Client ID | | Parameter | Concentration | C | MDL | RDL | Units |
|------------------------------------|-----------------|-------|----------------------|---------------|---|-------------|-----|-------|
| Client ID : 921-J-WS-080124 | | | | | | | | |
| P3451-01 | 921-J-WS-080124 | WATER | Phenanthrene | 0.050 | J | 0.02 | 0.1 | ug/L |
| P3451-01 | 921-J-WS-080124 | WATER | Fluoranthene | 0.060 | J | 0.02 | 0.1 | ug/L |
| P3451-01 | 921-J-WS-080124 | WATER | Pyrene | 0.040 | J | 0.02 | 0.1 | ug/L |
| P3451-01 | 921-J-WS-080124 | WATER | Benzo(b)fluoranthene | 0.030 | J | 0.03 | 0.1 | ug/L |
| Total Svoc : | | | | | | 0.18 | | |
| Total Concentration: | | | | | | 0.18 | | |



SAMPLE DATA

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | Matrix: | Water |
| Analytical Method: | SW8270SIM | % Solid: | 0 |
| Sample Wt/Vol: | 980 Units: mL | Final Vol: | 1000 uL |
| Soil Aliquot Vol: | uL | Test: | SVOCMS Group3 |
| Extraction Type : | Decanted : N | Level : | LOW |
| Injection Volume : | GPC Factor : 1.0 | GPC Cleanup : | N PH : |
| Prep Method : | SW3510C | | |

| | | | | |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed | Prep Batch ID |
| BN033235.D | 1 | 08/02/24 09:25 | 08/03/24 12:48 | PB162464 |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|---------------------------|-------------------------|-------|-----------|---------------------|------------|----------|
| TARGETS | | | | | | |
| 91-20-3 | Naphthalene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 91-57-6 | 2-Methylnaphthalene | 0.030 | U | 0.030 | 0.10 | ug/L |
| 208-96-8 | Acenaphthylene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 83-32-9 | Acenaphthene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 86-73-7 | Fluorene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 85-01-8 | Phenanthrene | 0.050 | J | 0.020 | 0.10 | ug/L |
| 120-12-7 | Anthracene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 206-44-0 | Fluoranthene | 0.060 | J | 0.020 | 0.10 | ug/L |
| 129-00-0 | Pyrene | 0.040 | J | 0.020 | 0.10 | ug/L |
| 56-55-3 | Benzo(a)anthracene | 0.020 | U | 0.020 | 0.10 | ug/L |
| 218-01-9 | Chrysene | 0.030 | U | 0.030 | 0.10 | ug/L |
| 205-99-2 | Benzo(b)fluoranthene | 0.030 | J | 0.030 | 0.10 | ug/L |
| 207-08-9 | Benzo(k)fluoranthene | 0.030 | U | 0.030 | 0.10 | ug/L |
| 50-32-8 | Benzo(a)pyrene | 0.060 | U | 0.060 | 0.10 | ug/L |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.040 | U | 0.040 | 0.10 | ug/L |
| 53-70-3 | Dibenzo(a,h)anthracene | 0.040 | U | 0.040 | 0.10 | ug/L |
| 191-24-2 | Benzo(g,h,i)perylene | 0.040 | U | 0.040 | 0.10 | ug/L |
| 123-91-1 | 1,4-Dioxane | 0.070 | U | 0.070 | 0.20 | ug/L |
| SURROGATES | | | | | | |
| 7297-45-2 | 2-Methylnaphthalene-d10 | 0.25 | | 30 (30) - 150 (150) | 63% | SPK: 0.4 |
| 93951-69-0 | Fluoranthene-d10 | 0.32 | | 30 (30) - 150 (150) | 79% | SPK: 0.4 |
| 4165-60-0 | Nitrobenzene-d5 | 0.28 | | 30 (11) - 130 (175) | 70% | SPK: 0.4 |
| 321-60-8 | 2-Fluorobiphenyl | 0.31 | | 30 (10) - 130 (175) | 78% | SPK: 0.4 |
| 1718-51-0 | Terphenyl-d14 | 0.39 | | 30 (54) - 130 (171) | 98% | SPK: 0.4 |
| INTERNAL STANDARDS | | | | | | |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 2720 | 7.546 | | | |
| 1146-65-2 | Naphthalene-d8 | 9390 | 10.276 | | | |
| 15067-26-2 | Acenaphthene-d10 | 4760 | 14.137 | | | |
| 1517-22-2 | Phenanthrene-d10 | 9240 | 16.908 | | | |

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | Matrix: | Water |
| Analytical Method: | SW8270SIM | % Solid: | 0 |
| Sample Wt/Vol: | 980 Units: mL | Final Vol: | 1000 uL |
| Soil Aliquot Vol: | uL | Test: | SVOCMS Group3 |
| Extraction Type : | Decanted : N | Level : | LOW |
| Injection Volume : | GPC Factor : 1.0 | GPC Cleanup : | N PH : |
| Prep Method : | SW3510C | | |

| | | | | |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed | Prep Batch ID |
| BN033235.D | 1 | 08/02/24 09:25 | 08/03/24 12:48 | PB162464 |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|--------------|-------|-----------|-----|------------|-------|
| 1719-03-5 | Chrysene-d12 | 7130 | 21.125 | | | |
| 1520-96-3 | Perylene-d12 | 7840 | 23.306 | | | |

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: | P3451 | OrderDate: | 8/1/2024 4:38:00 PM |
| Client: | JACOBS Engineering Group, Inc. | Project: | Former Schlumberger Site Princeton NJ |
| Contact: | Mary I. Murphy | Location: | D31,VOA Ref. #3 Water |

| LabID | ClientID | Matrix | Test | Method | Sample Date | Prep Date | Anal Date | Received |
|----------|-----------------|--------|---------------|---------------|-------------|-----------|-----------|----------|
| P3451-01 | 921-J-WS-080124 | Water | SVOCMS Group3 | 8270-Modified | 08/01/24 | 08/02/24 | 08/03/24 | 08/01/24 |



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

Hit Summary Sheet SW-846

SDG No.: P3451
Client: JACOBS Engineering Group, Inc.

| Sample ID | Client ID | Matrix | Parameter | Concentration | C | MDL | RDL | Units |
|-------------|-----------|--------|----------------------|---------------|---|------|-----|-------|
| Client ID : | | | | | | | | |
| | | | | 0.000 | | | | |
| | | | Total Svoc : | | | 0.00 | | |
| | | | Total Concentration: | | | 0.00 | | |



SAMPLE DATA

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | Matrix: | Water |
| Analytical Method: | SW8270 | % Solid: | 0 |
| Sample Wt/Vol: | 980 Units: mL | Final Vol: | 1000 uL |
| Soil Aliquot Vol: | uL | Test: | SVOCMS Group6 |
| Extraction Type : | Decanted : N | Level : | LOW |
| Injection Volume : | GPC Factor : 1.0 | GPC Cleanup : | N PH : |
| Prep Method : | SW3510C | | |

| | | | | |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed | Prep Batch ID |
| BF138844.D | 1 | 08/02/24 09:23 | 08/07/24 16:05 | PB162463 |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|----------------|----------------------------|-------|-----------|------|------------|-------|
| TARGETS | | | | | | |
| 110-86-1 | Pyridine | 1.60 | U | 1.60 | 5.10 | ug/L |
| 100-52-7 | Benzaldehyde | 4.10 | U | 4.10 | 10.2 | ug/L |
| 95-48-7 | 2-Methylphenol | 1.20 | U | 1.20 | 5.10 | ug/L |
| 98-86-2 | Acetophenone | 1.10 | U | 1.10 | 5.10 | ug/L |
| 65794-96-9 | 3+4-Methylphenols | 1.20 | U | 1.20 | 10.2 | ug/L |
| 98-95-3 | Nitrobenzene | 1.30 | U | 1.30 | 5.10 | ug/L |
| 120-83-2 | 2,4-Dichlorophenol | 0.90 | U | 0.90 | 5.10 | ug/L |
| 91-20-3 | Naphthalene | 1.00 | U | 1.00 | 5.10 | ug/L |
| 87-68-3 | Hexachlorobutadiene | 1.30 | U | 1.30 | 5.10 | ug/L |
| 91-57-6 | 2-Methylnaphthalene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.91 | U | 0.91 | 5.10 | ug/L |
| 95-95-4 | 2,4,5-Trichlorophenol | 1.00 | U | 1.00 | 5.10 | ug/L |
| 208-96-8 | Acenaphthylene | 1.10 | U | 1.10 | 5.10 | ug/L |
| 83-32-9 | Acenaphthene | 0.83 | U | 0.83 | 5.10 | ug/L |
| 132-64-9 | Dibenzofuran | 0.95 | U | 0.95 | 5.10 | ug/L |
| 86-73-7 | Fluorene | 0.98 | U | 0.98 | 5.10 | ug/L |
| 118-74-1 | Hexachlorobenzene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 87-86-5 | Pentachlorophenol | 1.90 | U | 1.90 | 10.2 | ug/L |
| 85-01-8 | Phenanthrene | 0.91 | U | 0.91 | 5.10 | ug/L |
| 86-74-8 | Carbazole | 1.20 | U | 1.20 | 5.10 | ug/L |
| 84-74-2 | Di-n-butylphthalate | 1.50 | U | 1.50 | 5.10 | ug/L |
| 206-44-0 | Fluoranthene | 1.30 | U | 1.30 | 5.10 | ug/L |
| 129-00-0 | Pyrene | 1.10 | U | 1.10 | 5.10 | ug/L |
| 56-55-3 | Benzo(a)anthracene | 0.96 | U | 0.96 | 5.10 | ug/L |
| 218-01-9 | Chrysene | 0.88 | U | 0.88 | 5.10 | ug/L |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 1.90 | U | 1.90 | 5.10 | ug/L |
| 205-99-2 | Benzo(b)fluoranthene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 207-08-9 | Benzo(k)fluoranthene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 50-32-8 | Benzo(a)pyrene | 1.70 | U | 1.70 | 5.10 | ug/L |

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | Matrix: | Water |
| Analytical Method: | SW8270 | % Solid: | 0 |
| Sample Wt/Vol: | 980 Units: mL | Final Vol: | 1000 uL |
| Soil Aliquot Vol: | uL | Test: | SVOCMS Group6 |
| Extraction Type : | Decanted : N | Level : | LOW |
| Injection Volume : | GPC Factor : 1.0 | GPC Cleanup : | N PH : |
| Prep Method : | SW3510C | | |

| | | | | |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed | Prep Batch ID |
| BF138844.D | 1 | 08/02/24 09:23 | 08/07/24 16:05 | PB162463 |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|---------------------------|------------------------|--------|-----------|---------------------|------------|----------|
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.00 | U | 1.00 | 5.10 | ug/L |
| 53-70-3 | Dibenzo(a,h)anthracene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 191-24-2 | Benzo(g,h,i)perylene | 1.20 | U | 1.20 | 5.10 | ug/L |
| 123-91-1 | 1,4-Dioxane | 1.30 | U | 1.30 | 5.10 | ug/L |
| 90-12-0 | 1-Methylnaphthalene | 0.88 | U | 0.88 | 5.10 | ug/L |
| SURROGATES | | | | | | |
| 367-12-4 | 2-Fluorophenol | 49.2 | | 15 (10) - 110 (139) | 33% | SPK: 150 |
| 13127-88-3 | Phenol-d6 | 29.6 | | 15 (10) - 110 (134) | 20% | SPK: 150 |
| 4165-60-0 | Nitrobenzene-d5 | 78.6 | | 30 (49) - 130 (133) | 79% | SPK: 100 |
| 321-60-8 | 2-Fluorobiphenyl | 87.4 | | 30 (52) - 130 (132) | 87% | SPK: 100 |
| 118-79-6 | 2,4,6-Tribromophenol | 128 | | 15 (32) - 110 (145) | 86% | SPK: 150 |
| 1718-51-0 | Terphenyl-d14 | 95.3 | | 30 (36) - 130 (145) | 95% | SPK: 100 |
| INTERNAL STANDARDS | | | | | | |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 44300 | 6.839 | | | |
| 1146-65-2 | Naphthalene-d8 | 184000 | 8.122 | | | |
| 15067-26-2 | Acenaphthene-d10 | 95700 | 9.869 | | | |
| 1517-22-2 | Phenanthrene-d10 | 153000 | 11.357 | | | |
| 1719-03-5 | Chrysene-d12 | 77100 | 13.998 | | | |
| 1520-96-3 | Perylene-d12 | 83600 | 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: | P3451 | OrderDate: | 8/1/2024 4:38:00 PM |
| Client: | JACOBS Engineering Group, Inc. | Project: | Former Schlumberger Site Princeton NJ |
| Contact: | Mary I. Murphy | Location: | D31,VOA Ref. #3 Water |

| LabID | ClientID | Matrix | Test | Method | Sample Date | Prep Date | Anal Date | Received |
|----------|-----------------|--------|---------------|---------------|-------------|-----------|-----------|----------|
| P3451-01 | 921-J-WS-080124 | Water | | | 08/01/24 | | | 08/01/24 |
| | | | SVOCMS Group3 | 8270-Modified | | 08/02/24 | 08/03/24 | |
| | | | SVOCMS Group6 | 8270E | | 08/02/24 | 08/07/24 | |

Hit Summary Sheet
SW-846

| | | | |
|-----------------|--------------------------------|--------------------|---------------------------------------|
| SDG No.: | P3451 | Order ID: | P3451 |
| Client: | JACOBS Engineering Group, Inc. | Project ID: | Former Schlumberger Site Princeton NJ |

| Sample ID | Client ID | Matrix | Parameter | Concentration | C | MDL | RDL | Units |
|------------------------------------|-----------------|--------|-----------|---------------|---|-------|------|-------|
| Client ID : 921-J-WS-080124 | | | | | | | | |
| P3451-01 | 921-J-WS-080124 | Water | Aluminum | 151 | | 1.98 | 20.0 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Antimony | 0.23 | J | 0.11 | 2.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Arsenic | 2.11 | | 0.090 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Barium | 48.7 | | 0.30 | 10.0 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Calcium | 14700 | | 62.5 | 500 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Cobalt | 0.46 | J | 0.062 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Copper | 2.72 | | 0.40 | 2.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Iron | 3270 | | 9.60 | 50.0 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Lead | 2.52 | | 0.11 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Magnesium | 2500 | | 26.6 | 500 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Manganese | 352 | | 0.24 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Nickel | 0.67 | J | 0.18 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Potassium | 2540 | | 46.1 | 500 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Tin | 0.31 | J | 0.12 | 5.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Sodium | 51300 | | 85.8 | 500 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Vanadium | 1.61 | J | 0.072 | 5.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Zinc | 7.82 | | 0.56 | 5.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Strontium | 107 | | 0.35 | 1.00 | ug/L |
| P3451-01 | 921-J-WS-080124 | Water | Titanium | 2.85 | J | 0.26 | 5.00 | ug/L |



SAMPLE DATA

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: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | 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 | 151 | | 1 | 1.98 | 20.0 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-36-0 | Antimony | 0.23 | J | 1 | 0.11 | 2.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-38-2 | Arsenic | 2.11 | | 1 | 0.090 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-39-3 | Barium | 48.7 | | 1 | 0.30 | 10.0 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-41-7 | Beryllium | 0.16 | U | 1 | 0.16 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-43-9 | Cadmium | 0.30 | U | 1 | 0.30 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-70-2 | Calcium | 14700 | | 1 | 62.5 | 500 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-47-3 | Chromium | 0.40 | U | 1 | 0.40 | 2.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-48-4 | Cobalt | 0.46 | J | 1 | 0.062 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-50-8 | Copper | 2.72 | | 1 | 0.40 | 2.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7439-89-6 | Iron | 3270 | | 1 | 9.60 | 50.0 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7439-92-1 | Lead | 2.52 | | 1 | 0.11 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7439-95-4 | Magnesium | 2500 | | 1 | 26.6 | 500 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7439-96-5 | Manganese | 352 | | 1 | 0.24 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7439-97-6 | Mercury | 0.081 | U | 1 | 0.081 | 0.20 | ug/L | 08/12/24 16:13 | 08/13/24 10:42 | SW7470A | |
| 7439-98-7 | Molybdenum | 0.93 | U | 1 | 0.93 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-02-0 | Nickel | 0.67 | J | 1 | 0.18 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-09-7 | Potassium | 2540 | | 1 | 46.1 | 500 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7782-49-2 | Selenium | 1.38 | U | 1 | 1.38 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-22-4 | Silver | 0.077 | UN | 1 | 0.077 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-23-5 | Sodium | 51300 | | 1 | 85.8 | 500 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-24-6 | Strontium | 107 | N | 1 | 0.35 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-28-0 | Thallium | 0.085 | U | 1 | 0.085 | 1.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-31-5 | Tin | 0.31 | J | 1 | 0.12 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-32-6 | Titanium | 2.85 | JN | 1 | 0.26 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-62-2 | Vanadium | 1.61 | J | 1 | 0.072 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |
| 7440-66-6 | Zinc | 7.82 | | 1 | 0.56 | 5.00 | ug/L | 08/23/24 15:00 | 08/25/24 17:57 | SW6020 | 3010A |

| | | | | | |
|---------------|-----------|-----------------|-------|------------|--------|
| Color Before: | Colorless | Clarity Before: | Clear | Texture: | Medium |
| Color After: | Colorless | Clarity After: | N/A | Artifacts: | N/A |
| Comments: | Mercury | | | | |

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

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

LAB CHRONICLE

| | | | |
|----------|--------------------------------|------------|---------------------------------------|
| OrderID: | P3451 | OrderDate: | 8/1/2024 4:38:00 PM |
| Client: | JACOBS Engineering Group, Inc. | Project: | Former Schlumberger Site Princeton NJ |
| Contact: | Mary I. Murphy | Location: | D31,VOA Ref. #3 Water |

| LabID | ClientID | Matrix | Test | Method | Sample Date | Prep Date | Anal Date | Received |
|----------|-----------------|--------|---------------|--------|-------------|-----------|-----------|----------|
| P3451-01 | 921-J-WS-080124 | Water | | | 08/01/24 | | | 08/01/24 |
| | | | Mercury | 7470A | | 08/12/24 | 08/13/24 | |
| | | | Metals Group4 | 6020B | | 08/23/24 | 08/25/24 | |



SAMPLE DATA

Report of Analysis

| | | | |
|-------------------|---------------------------------------|-----------------|----------------|
| Client: | JACOBS Engineering Group, Inc. | Date Collected: | 08/01/24 14:20 |
| Project: | Former Schlumberger Site Princeton NJ | Date Received: | 08/01/24 |
| Client Sample ID: | 921-J-WS-080124 | SDG No.: | P3451 |
| Lab Sample ID: | P3451-01 | 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 | | 08/02/24 14:12 | 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

LAB CHRONICLE

| | | | |
|----------|--------------------------------|------------|---------------------------------------|
| OrderID: | P3451 | OrderDate: | 8/1/2024 4:38:00 PM |
| Client: | JACOBS Engineering Group, Inc. | Project: | Former Schlumberger Site Princeton NJ |
| Contact: | Mary I. Murphy | Location: | D31,VOA Ref. #3 Water |

| LabID | ClientID | Matrix | Test | Method | Sample Date | Prep Date | Anal Date | Received |
|----------|-----------------|--------|---------------------|--------|-------------------|-----------|-------------------|----------|
| P3451-01 | 921-J-WS-080124 | WATER | Hexavalent Chromium | 7196A | 08/01/24 14:20 | | 08/02/24 14:12 | 08/01/24 |



SHIPPING DOCUMENTS

CHEMTECH

CHAIN OF CUSTODY RECORD

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

CHEMTECH PROJECT NO.

QUOTE NO.

COC Number

2041308

P3451

10

10.1

CLIENT INFORMATION

REPORT TO BE SENT TO:

COMPANY: Jacobs
ADDRESS: 412 Mt Kramb Ave Suite #100
CITY: Morrisstown STATE: NJ ZIP: 07960
ATTENTION: John Yphantis
PHONE: (281) 414-1719 FAX:

CLIENT PROJECT INFORMATION

PROJECT NAME: STC PTC
PROJECT NO.: D3779972 LOCATION: Princeton Junction
PROJECT MANAGER: Mary Murphy
e-mail: Mary.Murphy@Jacobs.com
PHONE: (761) 936-0586 FAX:

CLIENT BILLING INFORMATION

BILL TO: Mary Murphy PO#:
ADDRESS:
CITY STATE ZIP:
ATTENTION: PHONE:

ANALYSIS

DATA TURNAROUND INFORMATION

FAX (RUSH) Standard TAT DAYS*
HARDCOPY (DATA PACKAGE): DAYS*
EDD: DAYS*

*TO BE APPROVED BY CHEMTECH

STANDARD HARDCOPY TURNAROUND TIME IS 10 BUSINESS DAYS

DATA DELIVERABLE INFORMATION

☐ Level 1 (Results Only) ☐ Level 4 (QC + Full Raw Data)
☐ Level 2 (Results + QC) ☐ NJ Reduced ☐ US EPA CLP
☒ Level 3 (Results + QC + Raw Data) ☐ NYS ASP A ☐ NYS ASP B
☐ EDD FORMAT ☐ Other

VOCs 8260D
SVOCs 8260E
Metals 6020B, Hg
7172B
(PCV) 7196A

| CHEMTECH SAMPLE ID | PROJECT SAMPLE IDENTIFICATION | SAMPLE MATRIX | SAMPLE TYPE | | SAMPLE COLLECTION | | # OF BOTTLES | PRESERVATIVES | | | | | | | | | COMMENTS | |
|--------------------------|----------------------------------|------------------|----------------|------|----------------------|------|--------------|---------------|---|-----|---|--|--|--|--|--|----------|--|
| | | | COMP | GRAB | DATE | TIME | | A/E | E | B/E | E | | | | | | | ← Specify Preservatives A-HCl D-NaOH B-HNO3 E-ICE C-H2SO4 F-OTHER |
| 1. | 921-J-WS-650124 | WS | | X | 8/1/24 | 1420 | 6 | 2 | 2 | 1 | 1 | | | | | | | |
| 2. | TB-03-080124 | DI | | X | 8/1/24 | 1550 | 1 | 1 | | | | | | | | | | |
| 3. | | | | | | | | | | | | | | | | | | |
| 4. | | | | | | | | | | | | | | | | | | |
| 5. | | | | | | | | | | | | | | | | | | |
| 6. | | | | | | | | | | | | | | | | | | |
| 7. | | | | | | | | | | | | | | | | | | |
| 8. | | | | | | | | | | | | | | | | | | |
| 9. | | | | | | | | | | | | | | | | | | |
| 10. | | | | | | | | | | | | | | | | | | |

SAMPLE CUSTODY MUST BE DOCUMENTED BELOW EACH TIME SAMPLES CHANGE POSSESSION INCLUDING COURIER DELIVERY

| | | | |
|---|----------------------------------|--|---|
| RELINQUISHED BY SAMPLER: 1. <u>[Signature]</u> | DATE/TIME: <u>8/1/24 1630</u> | RECEIVED BY: <u>[Signature]</u> <u>1630</u> | Conditions of bottles or coolers at receipt: <input type="checkbox"/> COMPLIANT <input type="checkbox"/> NON COMPLIANT <input type="checkbox"/> COOLER TEMP <u>3.0 °C</u> |
| RELINQUISHED BY SAMPLER: 2. <u>[Signature]</u> | DATE/TIME: | RECEIVED BY: | Comments: <u>See attached table for required analytes list of ECO-VOCs, ECO-SVOCs, and ECO metals</u> |
| RELINQUISHED BY SAMPLER: 3. <u>[Signature]</u> | DATE/TIME: <u>8-1-24</u> | RECEIVED BY: | |

Page 1 of 1

CLIENT: ☐ Hand Delivered ☐ Other
CHEMTECH: ☐ Picked Up ☐ Field Sampling

Shipment Complete
☐ YES ☐ NO

P3451

WHITE - CHEMTECH COPY FOR RETURN TO CLIENT

YELLOW - CHEMTECH COPY

PINK - SAMPLER COPY

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 |

LOGIN REPORT/SAMPLE TRANSFER

Order ID : P3451 JACO05

Order Date : 8/1/2024 4:38:00 PM

Project Mgr :

Client Name : JACOBS Engineering Grou

Project Name : Former Schlumberger Site I

Report Type : Level 4

Client Contact : Mary I. Murphy

Receive DateTime : 8/1/2024 ~~12:00:00 AM~~

EDD Type : CH2MHILL

Invoice Name : JACOBS Engineering Grou

Purchase Order :

17:30

Hard Copy Date :

Invoice Contact : Mary I. Murphy

Date Signoff :

| LAB ID | CLIENT ID | MATRIX | SAMPLE DATE | SAMPLE TIME | TEST | TEST GROUP | METHOD | FAX DATE | DUE DATES |
|----------|-----------------|--------|-------------|-------------|--------------|------------|----------|--------------|-----------|
| P3451-01 | 921-J-WS-080124 | Water | 08/01/2024 | 14:20 | | | | | |
| | | | | | VOCMS Group6 | | 8260-Low | 10 Bus. Days | |
| P3451-02 | TB-03-080124 | Water | 08/01/2024 | 15:50 | | | | | |
| | | | | | VOCMS Group6 | | 8260-Low | 10 Bus. Days | |

Relinquished By :



Date / Time :

8/2/24 0730

Received By :



Date / Time :

8/2/24 7:10 Ref # 4

Storage Area : VOA Refridgerator Room