

Cover Page

- Order ID : Q2231
- Project ID : Con Ed Non MGP - Atlantic Ave 453957.600024.05
 - Client : PARSONS Engineering of New York, Inc.

Lab Sample Number

Client Sample Number Q2231-01 MW-10D-20250604 Q2231-02 MW-14-20250604 Q2231-03 MW-15-20250604 Q2231-04 MW-16D-20250604 Q2231-05 MW-17-20250604 Q2231-06 TB-20250604

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: 6/17/2025

NYDOH CERTIFICATION NO - 11376

NJDEP CERTIFICATION NO - 20012



CASE NARRATIVE

PARSONS Engineering of New York, Inc. Project Name: Con Ed Non MGP – Atlantic Ave 453957.600024.05 Project # N/A Order ID # Q2231 Test Name: VOCMS Group1

A. Number of Samples and Date of Receipt:

6 Water samples were received on 06/04/2025.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Sulfate, SVOCMS Group1, TDS and VOCMS Group1. This data package contains results for VOCMS Group1.

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 Group1 was based on method 8260D.

D. QA/ QC Samples:

The Holding Times were met for all analysis. The Surrogate recoveries met the acceptable criteria. The Internal Standards Areas met the acceptable requirements. The Retention Times were acceptable for all samples.

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

The Blank Spike Duplicate for {VN0611WBSD02} with File ID: VN086945.D met requirements for all samples except for Bromoform[112%] is failing high but no positive hit in associate sample therefore no corrective action taken.

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

The Initial Calibration met the requirements.

The Continuous Calibration File ID VN086940.D met the requirements except for Bromoform is failing high but no positive hit in associate sample therefore no corrective action taken.

The Tuning criteria met requirements.



E. Additional Comments:

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

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

F. Manual Integration Comments:

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

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CASE NARRATIVE

PARSONS Engineering of New York, Inc. Project Name: Con Ed Non MGP – Atlantic Ave 453957.600024.05 Project # N/A Order ID # Q2231 Test Name: SVOCMS Group1

A. Number of Samples and Date of Receipt:

6 Water samples were received on 06/04/2025.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Sulfate, SVOCMS Group1, TDS and VOCMS Group1. This data package contains results for SVOCMS Group1.

C. Analytical Techniques:

The samples were analyzed on instrument BNA_M using GC Column ZB-SemiVolatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGAThe samples were analyzed on instrument BNA_P using GC Column ZB-SemiVolatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGAThe analysis of SVOCMS Group1 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 {Q2230-03MS} with File ID: BP024879.D recoveries met the requirements for all compounds except for 1,4-Dioxane[35%], Recovery failed due to matrix interference, Therefor no further corrective action was taken.

The MSD {Q2230-04MSD} with File ID: BP024880.D recoveries met the acceptable requirements except for 1,4-Dioxane[36%], Recovery failed due to matrix interference, Therefor no further corrective action was taken.

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 8270-BM060525.M) for 2,4-Dinitrophenol this Compound is passing on Linear regression. And Method is failed for Di-n-octyl phthalate Under this Initial Calibration only diluted samples were analyzed and failed compounds in Initial Calibration are not associated , therefor no further corrective action was required.

The %RSD is greater than 20% in the Initial Calibration (Method 8270-BP060625.M) for 2,4-Dinitrophenol, 4-Nitrophenol these compound are passing on Linear Regression.

The Continuous Calibration File ID BM050262.D met the requirements except for 2,4-Dinitrophenol,4,6-Dinitro-2-methylphenol and Pentachlorophenol . Under this Continuous Calibration only diluted samples were analyzed and failed compounds in Continuous Calibration are not associated , therefor no further corrective action was required.

The Tuning criteria met requirements.

Samples MW-14-20250604, MW-14-20250604DL were diluted due to high concentrations.

E. Additional Comments:

Alliance has analyzed samples for SVOCMS Group1 by Method 8270 E for Project "Con Ed Non MGP – Atlantic Ave". Alliance certification was in applied status for compound "2,4-Dimethylphenol" with NJDEP for Method 8270E for SVOC group 1 at the time when samples for Project "Con Ed Non MGP – Atlantic Ave "were analyzed.

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

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

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_____



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CASE NARRATIVE

PARSONS Engineering of New York, Inc. Project Name: Con Ed Non MGP – Atlantic Ave 453957.600024.05 Project # N/A Order ID # Q2231 Test Name: Sulfate,TDS

A. Number of Samples and Date of Receipt:

6 Water samples were received on 06/04/2025.

B. Parameters:

According to the Chain of Custody document, the following analyses were requested: Sulfate, SVOCMS Group1, TDS and VOCMS Group1. This data package contains results for Sulfate, TDS.

C. Analytical Techniques:

The analysis of Sulfate was based on method 300.0 and The analysis of TDS was based on method SM2540 C.

D. QA/ QC Samples:

The Holding Times were met for all analysis. Sample MW-16D-20250604 was diluted due to high concentrations for Sulfate. The Blank Spike met requirements for all samples. The Duplicate analysis met criteria for all samples. The Matrix Spike (MW-16D-20250604MS) analysis met criteria for all analysts except for Sulfate due to matrix interference. The Matrix Spike Duplicate (MW-16D-20250604MSD) analysis met criteria for all analysts except for Sulfate due to matrix interference. The Blank analysis did not indicate the presence of lab contamination. The Calibration met the requirements.

E. Additional Comments:

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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 "AV" for automated Cold Vapor AA "CA" for MIDI-Distillation Spectrophotometric "AS" for Semi – Automated Spectrophotometric "AS" for Semi – Automated Spectrophotometric "C" for Manual Spectrophotometric "T" for Titrimetric "NR" for analyte not required to be analyzed 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
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 #: Q2231

Completed

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<u>✓</u>
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QA Review Signature: SOHIL JODHANI