



# CASE NARRATIVE

G Environmental Project Name: Nelson Project # N/A Order ID # Q2073

**Test Name: SVOCMS Group2** 

# A. Number of Samples and Date of Receipt:

2 Water samples were received on 05/16/2025.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: SVOC-SIMGroup1, SVOCMS Group1, SVOCMS Group2 and VOCMS Group1. This data package contains results for SVOCMS Group2.

### C. Analytical Techniques:

The samples were analyzed on instrument BNA\_P using GC Column ZB-SemiVolatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGAThe analysis of SVOCMS Group2 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 RPD met criteria.

The Blank Spike for {PB168098BS} with File ID: BP024755.D met requirements for all samples except for 3,3-Dichlorobenzidine[63%], 3-Nitroaniline[41%] and 4-Chloroaniline[35%],these compounds did not meet the NJDKQP criteria but met the inhouse criteria.

The Blank Spike Duplicate for {PB168098BSD} with File ID: BP024756.D met requirements for all samples except for 3,3-Dichlorobenzidine[61%], 3-Nitroaniline[42%] and 4-Chloroaniline[30%],these compounds did not meet the NJDKQP criteria but met the in-house criteria.

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

The Initial Calibration met the requirements.



The Continuous Calibration File ID BP024788.D met the requirements except for 2,2-oxybis(1-Chloropropane) is marginally biased low and Hexachlorocyclopentadiene is biased high and 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 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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