

## LAB CHRONICLE

|                 |                      |                   |                                        |
|-----------------|----------------------|-------------------|----------------------------------------|
| <b>OrderID:</b> | Q3324                | <b>OrderDate:</b> | 10/10/2025 10:16:00 AM                 |
| <b>Client:</b>  | Tetra Tech NUS, Inc. | <b>Project:</b>   | NWIRP Bethpage 112G08005-WE13          |
| <b>Contact:</b> | Ernie Wu             | <b>Location:</b>  | D41,VOA Ref. #2 Soil,VOA Ref. #3 Water |

| LabID           | ClientID                           | Matrix       | Test           | Method        | Sample Date     | Prep Date | Anal Date | Received        |
|-----------------|------------------------------------|--------------|----------------|---------------|-----------------|-----------|-----------|-----------------|
| <b>Q3324-03</b> | <b>BP-VPB-184-EB-2025<br/>1009</b> | <b>Water</b> |                |               | <b>10/09/25</b> |           |           | <b>10/09/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 10/10/25  | 10/13/25  |                 |
| <b>Q3324-05</b> | <b>BP-VPB-184-GW-780-<br/>782</b>  | <b>Water</b> |                |               | <b>10/07/25</b> |           |           | <b>10/09/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 10/10/25  | 10/13/25  |                 |
| <b>Q3324-06</b> | <b>BP-VPB-184-GW-840-<br/>842</b>  | <b>Water</b> |                |               | <b>10/09/25</b> |           |           | <b>10/09/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 10/10/25  | 10/13/25  |                 |



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

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q3324  
**Client:** Tetra Tech NUS, Inc.

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



# SAMPLE DATA

## Report of Analysis

|                    |                               |                    |                 |                |
|--------------------|-------------------------------|--------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          |                    | Date Collected: | 10/09/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                    | Date Received:  | 10/09/25       |
| Client Sample ID:  | BP-VPB-184-EB-20251009        |                    | SDG No.:        | Q3324          |
| Lab Sample ID:     | Q3324-03                      |                    | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | Level : LOW        | % Solid:        | 0              |
| Sample Wt/Vol:     | 760 mL                        | Final Vol: 1000 uL | Test:           | SVOC-SIMGroup1 |
| Prep Method :      | Prep Date: 10/10/25           |                    |                 |                |

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.26              | U    | 1  | 0.090    | 0.26 | 0.26       | ug/L     | 10/13/25 19:48 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.34              |      |    | 30 - 150 |      | 86%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.43              |      |    | 30 - 150 |      | 106%       | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.35              |      |    | 55 - 111 |      | 87%        | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.42              |      |    | 53 - 106 |      | 104%       | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.53              | *    |    | 58 - 132 |      | 133%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         |                   |      |    |          |      |            |          |                |              |
|                           |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 5140              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 13300             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 6680              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 12100             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 10400             |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 9950              |      |    |          |      |            |          |                |              |

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:            | Tetra Tech NUS, Inc.          |                     | Date Collected: | 10/07/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                     | Date Received:  | 10/09/25       |
| Client Sample ID:  | BP-VPB-184-GW-780-782         |                     | SDG No.:        | Q3324          |
| Lab Sample ID:     | Q3324-05                      |                     | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | Level : LOW         | % Solid:        | 0              |
| Sample Wt/Vol:     | 650 mL                        | Final Vol: 1000 uL  | Test:           | SVOC-SIMGroup1 |
| Prep Method :      |                               | Prep Date: 10/10/25 |                 |                |

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.31              | U    | 1  | 0.10     | 0.31 | 0.31       | ug/L     | 10/13/25 20:24 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.39              |      |    | 30 - 150 |      | 96%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.39              |      |    | 30 - 150 |      | 98%        | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.40              |      |    | 55 - 111 |      | 100%       | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.49              | *    |    | 53 - 106 |      | 123%       | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.61              | *    |    | 58 - 132 |      | 152%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 5800              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 14900             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 7290              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 12200             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 8880              |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 8840              |      |    |          |      |            |          |                |              |

U = Not Detected

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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: Tetra Tech NUS, Inc.  
Project: NWIRP Bethpage 112G08005-WE13  
Client Sample ID: BP-VPB-184-GW-840-842  
Lab Sample ID: Q3324-06  
Analytical Method: SW8270ESIM Level : LOW  
Sample Wt/Vol: 710 mL Final Vol: 1000 uL  
Prep Method : Prep Date: 10/10/25

Date Collected: 10/09/25  
Date Received: 10/09/25  
SDG No.: Q3324  
Matrix: Water  
% Solid: 0  
Test: SVOC-SIMGroup1

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.28              | U    | 1  | 0.090    | 0.28 | 0.28       | ug/L     | 10/13/25 21:00 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.32              |      |    | 30 - 150 |      | 80%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.38              |      |    | 30 - 150 |      | 95%        | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.32              |      |    | 55 - 111 |      | 79%        | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.40              |      |    | 53 - 106 |      | 99%        | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.44              |      |    | 58 - 132 |      | 111%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         |                   |      |    |          |      |            |          |                |              |
|                           |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 5890              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 15900             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 7550              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 13300             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 12100             |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 7290              |      |    |          |      |            |          |                |              |

U = Not Detected

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# QC SUMMARY

## Surrogate Summary

SW-846

SDG No.: **Q3324**

Client: **Tetra Tech NUS, Inc.**

Analytical Method: **8270-Modified**

| Lab Sample ID | Client ID              | Parameter               | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|------------------------|-------------------------|-------------|--------------|--------------|------|------------|------|
|               |                        |                         |             |              |              |      | Low        | High |
| PB170061BL    | PB170061BL             | 2-Methylnaphthalene-d10 | 0.4         | 0.33         | 83           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.35         | 88           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.37         | 91           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.41         | 101          |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.49         | 122          |      | 58         | 132  |
| PB170061BS    | PB170061BS             | 2-Methylnaphthalene-d10 | 0.4         | 0.39         | 98           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.32         | 81           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.36         | 90           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.43         | 108          | *    | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.49         | 121          |      | 58         | 132  |
| PB170061BSD   | PB170061BSD            | 2-Methylnaphthalene-d10 | 0.4         | 0.39         | 98           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.33         | 82           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.38         | 95           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.43         | 108          | *    | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.49         | 123          |      | 58         | 132  |
| Q3324-03      | BP-VPB-184-EB-20251009 | 2-Methylnaphthalene-d10 | 0.4         | 0.34         | 86           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.43         | 106          |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.35         | 87           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.42         | 104          |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.53         | 133          | *    | 58         | 132  |
| Q3324-05      | BP-VPB-184-GW-780-782  | 2-Methylnaphthalene-d10 | 0.4         | 0.39         | 96           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.39         | 98           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.40         | 100          |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.49         | 123          | *    | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.61         | 152          | *    | 58         | 132  |
| Q3324-06      | BP-VPB-184-GW-840-842  | 2-Methylnaphthalene-d10 | 0.4         | 0.32         | 80           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.38         | 95           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.32         | 79           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.40         | 99           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.44         | 111          |      | 58         | 132  |

**Laboratory Control Sample/Laboratory Control Sample Duplicate Summary**

SW-846

SDG No.: Q3324

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN038079.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB170061BS    | 1,4-Dioxane | 0.4   | 0.37   | ug/L | 93  |     |      |      | 70     | 130  |     |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3324

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN038080.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB170061BSD   | 1,4-Dioxane | 0.4   | 0.39   | ug/L | 98  | 5   |      |      | 70     | 130  | 20  |

4B

SEMIVOLATILE METHOD BLANK SUMMARY

Client ID

PB170061BL

Lab Name: Alliance

Contract: TETR06

Lab Code: ACE

SDG NO.: Q3324

Lab File ID: BN038084.D

Lab Sample ID: PB170061BL

Instrument ID: BNA\_N

Date Extracted: 10/10/2025

Matrix: (soil/water) Water

Date Analyzed: 10/14/2025

Level: (low/med) LOW

Time Analyzed: 02:31

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| EPA<br>SAMPLE NO.      | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|------------------------|------------------|----------------|------------------|
| PB170061BS             | PB170061BS       | BN038079.D     | 10/13/2025       |
| BP-VPB-184-EB-20251009 | Q3324-03         | BN038076.D     | 10/13/2025       |
| BP-VPB-184-GW-780-782  | Q3324-05         | BN038077.D     | 10/13/2025       |
| BP-VPB-184-GW-840-842  | Q3324-06         | BN038078.D     | 10/13/2025       |
| PB170061BSD            | PB170061BSD      | BN038080.D     | 10/13/2025       |

COMMENTS: \_\_\_\_\_  
\_\_\_\_\_

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037860.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3324  
DFTPP Injection Date: 09/23/2025  
DFTPP Injection Time: 11:33

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.7 ( 1.8 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.9                  |
| 365 | Greater than 1% of mass 198        | 4.2                  |
| 441 | Present, but less than mass 443    | 72.6                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 17.7 (20.8) 2        |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC0.1        | SSTDICC0.1       | BN037861.D     | 09/23/2025       | 12:13            |
| SSTDICC0.2        | SSTDICC0.2       | BN037862.D     | 09/23/2025       | 12:49            |
| SSTDICCC0.4       | SSTDICCC0.4      | BN037863.D     | 09/23/2025       | 13:25            |
| SSTDICC0.8        | SSTDICC0.8       | BN037864.D     | 09/23/2025       | 14:02            |
| SSTDICC1.6        | SSTDICC1.6       | BN037865.D     | 09/23/2025       | 14:38            |
| SSTDICC3.2        | SSTDICC3.2       | BN037866.D     | 09/23/2025       | 15:15            |
| SSTDICC5.0        | SSTDICC5.0       | BN037867.D     | 09/23/2025       | 15:51            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN038065.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3324  
DFTPP Injection Date: 10/13/2025  
DFTPP Injection Time: 13:08

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.2 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.5 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.2                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.7                  |
| 365 | Greater than 1% of mass 198        | 3.4                  |
| 441 | Present, but less than mass 443    | 70.3                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.2 ( 21 ) 2        |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO.      | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|------------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4             | SSTDCCC0.4       | BN038066.D     | 10/13/2025       | 13:47            |
| BP-VPB-184-EB-20251009 | Q3324-03         | BN038076.D     | 10/13/2025       | 19:48            |
| BP-VPB-184-GW-780-782  | Q3324-05         | BN038077.D     | 10/13/2025       | 20:24            |
| BP-VPB-184-GW-840-842  | Q3324-06         | BN038078.D     | 10/13/2025       | 21:00            |
| PB170061BS             | PB170061BS       | BN038079.D     | 10/13/2025       | 21:36            |
| PB170061BSD            | PB170061BSD      | BN038080.D     | 10/13/2025       | 22:12            |
| SSTDCCC0.4EC           | SSTDCCC0.4       | BN038081.D     | 10/13/2025       | 23:24            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN038082.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3324  
DFTPP Injection Date: 10/14/2025  
DFTPP Injection Time: 00:40

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.7 ( 1.7 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.5 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.4                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.8                  |
| 365 | Greater than 1% of mass 198        | 3.5                  |
| 441 | Present, but less than mass 443    | 83.1                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 17 ( 18 ) 2          |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4        | SSTDCCC0.4       | BN038083.D     | 10/14/2025       | 01:19            |
| PB170061BL        | PB170061BL       | BN038084.D     | 10/14/2025       | 02:31            |
| SSTDCCC0.4EC      | SSTDCCC0.4       | BN038088.D     | 10/14/2025       | 04:56            |

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3324

Client ID : SSTDCCC0.4

Date Analyzed: 10/13/2025

Lab File ID: BN038066.D

Time Analyzed: 13:47

Instrument ID: BNA\_N

GC Column: ZB-GR ID: 0.25 (mm)

|                           | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|---------------------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD               | 7166                | 7.797 | 18651               | 10.59  | 10244               | 14.45  |
| UPPER LIMIT               | 14332               | 8.297 | 37302               | 11.094 | 20488               | 14.952 |
| LOWER LIMIT               | 3583                | 7.297 | 9325.5              | 10.094 | 5122                | 13.952 |
| EPA SAMPLE NO.            |                     |       |                     |        |                     |        |
| 01 PB170061BS             | 6519                | 7.80  | 17676               | 10.59  | 7952                | 14.45  |
| 02 PB170061BSD            | 6262                | 7.80  | 16679               | 10.59  | 7730                | 14.45  |
| 03 BP-VPB-184-EB-20251009 | 5139                | 7.80  | 13291               | 10.61  | 6680                | 14.45  |
| 04 BP-VPB-184-GW-780-782  | 5800                | 7.80  | 14884               | 10.59  | 7288                | 14.45  |
| 05 BP-VPB-184-GW-840-842  | 5887                | 7.80  | 15917               | 10.59  | 7552                | 14.45  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3324

Client ID: SSTDCCC0.4

Date Analyzed: 10/13/2025

Lab File ID: BN038066.D

Time Analyzed: 13:47

Instrument ID: BNA\_N

GC Column: ZB-GR ID: 0.25 (mm)

|                           | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|---------------------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD               | 18901               | 17.198 | 16827               | 21.393 | 17326               | 23.712 |
| UPPER LIMIT               | 37802               | 17.698 | 33654               | 21.893 | 34652               | 24.212 |
| LOWER LIMIT               | 9450.5              | 16.698 | 8413.5              | 20.893 | 8663                | 23.212 |
| EPA SAMPLE NO.            |                     |        |                     |        |                     |        |
| 01 PB170061BS             | 12853               | 17.20  | 8233 *              | 21.39  | 7491 *              | 23.71  |
| 02 PB170061BSD            | 12472               | 17.20  | 7875 *              | 21.39  | 7156 *              | 23.71  |
| 03 BP-VPB-184-EB-20251009 | 12083               | 17.20  | 10438               | 21.39  | 9949                | 23.71  |
| 04 BP-VPB-184-GW-780-782  | 12210               | 17.20  | 8879                | 21.39  | 8844                | 23.71  |
| 05 BP-VPB-184-GW-840-842  | 13324               | 17.20  | 12137               | 21.39  | 7289 *              | 23.71  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3324

Client ID : SSTDCCC0.4

Date Analyzed: 10/14/2025

Lab File ID: BN038083.D

Time Analyzed: 01:19

Instrument ID: BNA\_N

GC Column: ZB-GR ID: 0.25 (mm)

|                | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|----------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 7061                | 7.797 | 18151               | 10.59  | 9016                | 14.45  |
| UPPER LIMIT    | 14122               | 8.297 | 36302               | 11.094 | 18032               | 14.952 |
| LOWER LIMIT    | 3530.5              | 7.297 | 9075.5              | 10.094 | 4508                | 13.952 |
| EPA SAMPLE NO. |                     |       |                     |        |                     |        |
| 01 PB170061BL  | 6692                | 7.80  | 16789               | 10.61  | 7676                | 14.45  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3324

Client ID: SSTDCCC0.4

Date Analyzed: 10/14/2025

Lab File ID: BN038083.D

Time Analyzed: 01:19

Instrument ID: BNA\_N

GC Column: ZB-GR ID: 0.25 (mm)

|                | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|----------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 16225               | 17.198 | 13579               | 21.384 | 13727               | 23.709 |
| UPPER LIMIT    | 32450               | 17.698 | 27158               | 21.884 | 27454               | 24.209 |
| LOWER LIMIT    | 8112.5              | 16.698 | 6789.5              | 20.884 | 6863.5              | 23.209 |
| EPA SAMPLE NO. |                     |        |                     |        |                     |        |
| 01 PB170061BL  | 11455               | 17.21  | 7795                | 21.39  | 6815 *              | 23.71  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.



# QC SAMPLE DATA

## Report of Analysis

Client: Tetra Tech NUS, Inc.  
Project: NWIRP Bethpage 112G08005-WE13  
Client Sample ID: PB170061BL  
Lab Sample ID: PB170061BL  
Analytical Method: SW8270ESIM Level : LOW  
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL  
Prep Method : Prep Date: 10/10/25

Date Collected:  
Date Received:  
SDG No.: Q3324  
Matrix: Water  
% Solid: 0  
Test: SVOC-SIMGroup1

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.20              | U    | 1  | 0.070    | 0.20 | 0.20       | ug/L     | 10/14/25 02:31 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.33              |      |    | 30 - 150 |      | 83%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.35              |      |    | 30 - 150 |      | 88%        | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.37              |      |    | 55 - 111 |      | 91%        | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.41              |      |    | 53 - 106 |      | 101%       | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.49              |      |    | 58 - 132 |      | 122%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         |                   |      |    |          |      |            |          |                |              |
|                           |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6690              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 16800             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 7680              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 11500             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 7800              |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 6820              |      |    |          |      |            |          |                |              |

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: Tetra Tech NUS, Inc.  
Project: NWIRP Bethpage 112G08005-WE13  
Client Sample ID: PB170061BS  
Lab Sample ID: PB170061BS  
Analytical Method: SW8270ESIM Level : LOW  
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL  
Prep Method : Prep Date: 10/10/25

Date Collected:  
Date Received:  
SDG No.: Q3324  
Matrix: Water  
% Solid: 0  
Test: SVOC-SIMGroup1

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.37              |      | 1  | 0.070    | 0.20 | 0.20       | ug/L     | 10/13/25 21:36 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.39              |      |    | 30 - 150 |      | 98%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.32              |      |    | 30 - 150 |      | 81%        | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.36              |      |    | 55 - 111 |      | 90%        | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.43              | *    |    | 53 - 106 |      | 108%       | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.49              |      |    | 58 - 132 |      | 121%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6520              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 17700             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 7950              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 12900             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 8230              |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 7490              |      |    |          |      |            |          |                |              |

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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: Tetra Tech NUS, Inc.  
Project: NWIRP Bethpage 112G08005-WE13  
Client Sample ID: PB170061BSD  
Lab Sample ID: PB170061BSD  
Analytical Method: SW8270ESIM Level : LOW  
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL  
Prep Method : Prep Date: 10/10/25

Date Collected:  
Date Received:  
SDG No.: Q3324  
Matrix: Water  
% Solid: 0  
Test: SVOC-SIMGroup1

| CAS Number                | Parameter               | Conc.             | Qua. | DF | MDL      | LOD  | LOQ / CRQL | Units    | Date Ana.      | Prep BatchID |
|---------------------------|-------------------------|-------------------|------|----|----------|------|------------|----------|----------------|--------------|
| <b>TARGETS</b>            |                         |                   |      |    |          |      |            |          |                |              |
| 123-91-1                  | 1,4-Dioxane             | 0.39              |      | 1  | 0.070    | 0.20 | 0.20       | ug/L     | 10/13/25 22:12 | PB170061     |
| <b>SURROGATES</b>         |                         |                   |      |    |          |      |            |          |                |              |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.39              |      |    | 30 - 150 |      | 98%        | SPK: 0.4 |                |              |
| 93951-69-0                | Fluoranthene-d10        | 0.33              |      |    | 30 - 150 |      | 82%        | SPK: 0.4 |                |              |
| 4165-60-0                 | Nitrobenzene-d5         | 0.38              |      |    | 55 - 111 |      | 95%        | SPK: 0.4 |                |              |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.43              | *    |    | 53 - 106 |      | 108%       | SPK: 0.4 |                |              |
| 1718-51-0                 | Terphenyl-d14           | 0.49              |      |    | 58 - 132 |      | 123%       | SPK: 0.4 |                |              |
| <b>INTERNAL STANDARDS</b> |                         |                   |      |    |          |      |            |          |                |              |
|                           |                         | <b>Area Count</b> |      |    |          |      |            |          |                |              |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6260              |      |    |          |      |            |          |                |              |
| 1146-65-2                 | Naphthalene-d8          | 16700             |      |    |          |      |            |          |                |              |
| 15067-26-2                | Acenaphthene-d10        | 7730              |      |    |          |      |            |          |                |              |
| 1517-22-2                 | Phenanthrene-d10        | 12500             |      |    |          |      |            |          |                |              |
| 1719-03-5                 | Chrysene-d12            | 7880              |      |    |          |      |            |          |                |              |
| 1520-96-3                 | Perylene-d12            | 7160              |      |    |          |      |            |          |                |              |

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\* = Values outside of QC limits

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() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products



# CALIBRATION SUMMARY

Method Path : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\  
 Method File : 8270-SIM-BN092325.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Wed Sep 24 11:00:01 2025  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BN037861.D 0.2 =BN037862.D 0.4 =BN037863.D 0.8 =BN037864.D 1.6 =BN037865.D 3.2 =BN037866.D 5 =BN037867.D

|           | Compound              | 0.1            | 0.2   | 0.4   | 0.8   | 1.6   | 3.2   | 5     | Avg   | %RSD  |
|-----------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| -----     |                       |                |       |       |       |       |       |       |       |       |
| 1) I      | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |
| 2)        | 1,4-Dioxane           | 0.413          | 0.410 | 0.403 | 0.425 | 0.410 | 0.376 | 0.406 | 4.04  |       |
| 3)        | n-Nitrosodimet...     | 0.541          | 0.517 | 0.526 | 0.561 | 0.549 | 0.547 | 0.540 | 3.03  |       |
| 4) S      | 2-Fluorophenol        | 0.944          | 0.921 | 0.856 | 0.839 | 0.908 | 0.923 | 0.998 | 0.913 | 5.84  |
| 5) S      | Phenol-d6             | 1.190          | 1.200 | 1.118 | 1.088 | 1.211 | 1.236 | 1.353 | 1.199 | 7.15  |
| 6)        | bis(2-Chloroet...     | 1.133          | 1.096 | 1.082 | 1.066 | 1.150 | 1.124 | 1.143 | 1.113 | 2.89  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 7) I      | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |
| 8) S      | Nitrobenzene-d5       | 0.299          | 0.289 | 0.289 | 0.290 | 0.310 | 0.320 | 0.338 | 0.305 | 6.14  |
| 9)        | Naphthalene           | 1.105          | 1.089 | 1.048 | 1.051 | 1.123 | 1.137 | 1.160 | 1.102 | 3.84  |
| 10)       | Hexachlorobuta...     | 0.198          | 0.184 | 0.182 | 0.183 | 0.192 | 0.188 | 0.188 | 0.188 | 2.94  |
| 11) SURR2 | Methylnaphth...       | 0.515          | 0.518 | 0.499 | 0.504 | 0.553 | 0.587 | 0.714 | 0.556 | 13.74 |
| 12)       | 2-Methylnaphth...     | 0.678          | 0.681 | 0.668 | 0.682 | 0.754 | 0.775 | 0.802 | 0.720 | 7.68  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 13) I     | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 14) S     | 2,4,6-Tribromo...     | 0.148          | 0.142 | 0.135 | 0.139 | 0.163 | 0.184 | 0.152 | 12.15 |       |
| 15) S     | 2-Fluorobiphenyl      | 1.660          | 1.631 | 1.559 | 1.547 | 1.654 | 1.650 | 1.762 | 1.638 | 4.36  |
| 16)       | Acenaphthylene        | 1.698          | 1.721 | 1.663 | 1.680 | 1.907 | 1.951 | 2.094 | 1.816 | 9.23  |
| 17)       | Acenaphthene          | 1.262          | 1.230 | 1.198 | 1.208 | 1.351 | 1.371 | 1.437 | 1.294 | 7.16  |
| 18)       | Fluorene              | 1.455          | 1.455 | 1.452 | 1.466 | 1.664 | 1.717 | 1.748 | 1.565 | 8.77  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 19) I     | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 20)       | 4,6-Dinitro-2-...     | 0.041          | 0.044 | 0.046 | 0.058 | 0.071 | 0.084 | 0.057 | 29.75 |       |
| 21)       | 4-Bromophenyl-...     | 0.252          | 0.247 | 0.244 | 0.244 | 0.272 | 0.275 | 0.291 | 0.261 | 7.07  |
| 22)       | Hexachlorobenzene     | 0.317          | 0.312 | 0.305 | 0.300 | 0.324 | 0.323 | 0.332 | 0.316 | 3.53  |
| 23)       | Atrazine              | 0.180          | 0.181 | 0.172 | 0.178 | 0.205 | 0.224 | 0.249 | 0.199 | 14.70 |
| 24)       | Pentachlorophenol     | 0.101          | 0.092 | 0.097 | 0.116 | 0.136 | 0.108 | 16.50 |       |       |
| 25)       | Phenanthrene          | 1.301          | 1.271 | 1.230 | 1.224 | 1.358 | 1.379 | 1.452 | 1.316 | 6.37  |
| 26)       | Anthracene            | 1.034          | 1.035 | 1.023 | 1.057 | 1.209 | 1.275 | 1.361 | 1.142 | 12.10 |
| 27) SURR  | Fluoranthene-d10      | 0.880          | 0.918 | 0.887 | 0.900 | 1.004 | 1.085 | 1.370 | 1.006 | 17.59 |
| 28)       | Fluoranthene          | 1.278          | 1.315 | 1.311 | 1.358 | 1.555 | 1.617 | 1.713 | 1.450 | 12.07 |
|           |                       |                |       |       |       |       |       |       |       |       |
| 29) I     | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |
| 30)       | Pyrene                | 1.576          | 1.600 | 1.521 | 1.476 | 1.607 | 1.595 | 1.679 | 1.579 | 4.12  |
| 31) S     | Terphenyl-d14         | 0.799          | 0.791 | 0.753 | 0.735 | 0.798 | 0.809 | 0.893 | 0.797 | 6.34  |
| 32)       | Benzo(a)anthra...     | 1.376          | 1.359 | 1.294 | 1.285 | 1.423 | 1.483 | 1.567 | 1.398 | 7.27  |
| 33)       | Chrysene              | 1.508          | 1.526 | 1.462 | 1.446 | 1.559 | 1.512 | 1.569 | 1.512 | 3.03  |
| 34)       | Bis(2-ethylhex...     | 0.592          | 0.487 | 0.476 | 0.518 | 0.569 | 0.676 | 0.553 | 13.57 |       |
|           |                       |                |       |       |       |       |       |       |       |       |
| 35) I     | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\

Method File : 8270-SIM-BN092325.M

|       |                   |       |       |       |       |       |       |       |       |       |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 36)   | Indeno(1,2,3-c... | 1.715 | 1.672 | 1.658 | 1.707 | 1.887 | 1.996 | 2.159 | 1.828 | 10.52 |
| 37)   | Benzo(b)fluora... | 1.599 | 1.606 | 1.641 | 1.590 | 1.810 | 1.848 | 1.981 | 1.725 | 8.95  |
| 38)   | Benzo(k)fluora... | 1.620 | 1.586 | 1.597 | 1.628 | 1.875 | 1.917 | 1.926 | 1.736 | 9.26  |
| 39) C | Benzo(a)pyrene    | 1.254 | 1.296 | 1.231 | 1.254 | 1.413 | 1.492 | 1.610 | 1.364 | 10.61 |
| 40)   | Dibenzo(a,h)an... | 1.285 | 1.283 | 1.303 | 1.333 | 1.495 | 1.587 | 1.705 | 1.427 | 11.87 |
| 41)   | Benzo(g,h,i)pe... | 1.407 | 1.381 | 1.404 | 1.397 | 1.521 | 1.613 | 1.773 | 1.499 | 9.85  |

-----  
(#) = Out of Range

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>TETR06</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3324</u>                 |
| Instrument ID:  | <u>BNA_N</u>      | Calibration Date/Time: | <u>10/13/2025 13:47</u>      |
| Lab File ID:    | <u>BN038066.D</u> | Init. Calib. Date(s):  | <u>09/23/2025 09/23/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4</u> | Init. Calib. Time(s):  | <u>12:13 15:51</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.556 | 0.571  |            | 2.7   | 20.0  |
| Fluoranthene-d10        | 1.006 | 1.065  |            | 5.9   | 20.0  |
| 2-Fluorophenol          | 0.913 | 0.893  |            | -2.2  | 20.0  |
| Phenol-d6               | 1.199 | 0.939  |            | -21.7 | 20.0  |
| Nitrobenzene-d5         | 0.305 | 0.238  |            | -22.0 | 20.0  |
| 2-Fluorobiphenyl        | 1.638 | 1.428  |            | -12.8 | 20.0  |
| 2,4,6-Tribromophenol    | 0.152 | 0.137  |            | -9.9  | 20.0  |
| Terphenyl-d14           | 0.797 | 0.724  |            | -9.2  | 20.0  |
| 1,4-Dioxane             | 0.406 | 0.465  |            | 14.5  | 20.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q3324  
Instrument ID: BNA\_N Calibration Date/Time: 10/13/2025 23:24  
Lab File ID: BN038081.D Init. Calib. Date(s): 09/23/2025 09/23/2025  
EPA Sample No.: SSTDCCC0.4EC Init. Calib. Time(s): 12:13 15:51  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.556 | 0.494  |            | -11.2 | 50.0  |
| Fluoranthene-d10        | 1.006 | 0.861  |            | -14.4 | 50.0  |
| 2-Fluorophenol          | 0.913 | 0.864  |            | -5.4  | 50.0  |
| Phenol-d6               | 1.199 | 0.927  |            | -22.7 | 50.0  |
| Nitrobenzene-d5         | 0.305 | 0.270  |            | -11.5 | 50.0  |
| 2-Fluorobiphenyl        | 1.638 | 1.596  |            | -2.6  | 50.0  |
| 2,4,6-Tribromophenol    | 0.152 | 0.107  |            | -29.6 | 50.0  |
| Terphenyl-d14           | 0.797 | 0.757  |            | -5.0  | 50.0  |
| 1,4-Dioxane             | 0.406 | 0.485  |            | 19.5  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q3324  
Instrument ID: BNA\_N Calibration Date/Time: 10/14/2025 01:19  
Lab File ID: BN038083.D Init. Calib. Date(s): 09/23/2025 09/23/2025  
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 12:13 15:51  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.556 | 0.546  |            | -1.8  | 20.0  |
| Fluoranthene-d10        | 1.006 | 1.016  |            | 1.0   | 20.0  |
| 2-Fluorophenol          | 0.913 | 0.823  |            | -9.9  | 20.0  |
| Phenol-d6               | 1.199 | 0.940  |            | -21.6 | 20.0  |
| Nitrobenzene-d5         | 0.305 | 0.248  |            | -18.7 | 20.0  |
| 2-Fluorobiphenyl        | 1.638 | 1.500  |            | -8.4  | 20.0  |
| 2,4,6-Tribromophenol    | 0.152 | 0.116  |            | -23.7 | 20.0  |
| Terphenyl-d14           | 0.797 | 0.709  |            | -11.0 | 20.0  |
| 1,4-Dioxane             | 0.406 | 0.473  |            | 16.5  | 20.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                     |                        |                              |
|-----------------|---------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>     | Contract:              | <u>TETR06</u>                |
| Lab Code:       | <u>ACE</u>          | SDG No.:               | <u>Q3324</u>                 |
| Instrument ID:  | <u>BNA_N</u>        | Calibration Date/Time: | <u>10/14/2025 04:56</u>      |
| Lab File ID:    | <u>BN038088.D</u>   | Init. Calib. Date(s):  | <u>09/23/2025 09/23/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4EC</u> | Init. Calib. Time(s):  | <u>12:13 15:51</u>           |
| GC Column:      | <u>ZB-GR</u>        | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.556 | 0.537  |            | -3.4  | 50.0  |
| Fluoranthene-d10        | 1.006 | 0.979  |            | -2.7  | 50.0  |
| 2-Fluorophenol          | 0.913 | 0.880  |            | -3.6  | 50.0  |
| Phenol-d6               | 1.199 | 0.975  |            | -18.7 | 50.0  |
| Nitrobenzene-d5         | 0.305 | 0.255  |            | -16.4 | 50.0  |
| 2-Fluorobiphenyl        | 1.638 | 1.577  |            | -3.7  | 50.0  |
| 2,4,6-Tribromophenol    | 0.152 | 0.119  |            | -21.7 | 50.0  |
| Terphenyl-d14           | 0.797 | 0.748  |            | -6.1  | 50.0  |
| 1,4-Dioxane             | 0.406 | 0.494  |            | 21.7  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.