

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: ROYF02

Lab Code: ACE

SDG No.: Q2641

Instrument ID: BNA\_F

Calibration Date(s): 07/17/2025 07/17/2025

Calibration Time(s): 11:04 14:34

| LAB FILE ID:          |        | RRF2.5 = BF143140.D |        |        | RRF005 = BF143141.D |        | RRF010 = BF143142.D |       |
|-----------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
|                       |        | RRF020 = BF143143.D |        |        | RRF040 = BF143144.D |        | RRF050 = BF143145.D |       |
| COMPOUND              | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| Pyridine              |        | 1.631               | 1.557  | 1.551  | 1.530               | 1.618  | 1.552               | 3.9   |
| 2-Fluorophenol        |        | 1.385               | 1.313  | 1.319  | 1.231               | 1.284  | 1.264               | 7.0   |
| Phenol-d6             |        | 1.733               | 1.640  | 1.657  | 1.554               | 1.632  | 1.591               | 6.7   |
| 1,4-Dichlorobenzene   |        | 1.602               | 1.518  | 1.502  | 1.418               | 1.473  | 1.449               | 7.5   |
| 2-Methylphenol        |        | 1.178               | 1.111  | 1.141  | 1.097               | 1.160  | 1.114               | 4.6   |
| 3+4-Methylphenols     |        |                     | 1.467  | 1.469  | 1.341               | 1.402  | 1.351               | 8.9   |
| Nitrobenzene-d5       |        | 0.407               | 0.394  | 0.417  | 0.399               | 0.417  | 0.401               | 3.5   |
| Hexachloroethane      |        | 0.522               | 0.498  | 0.522  | 0.495               | 0.525  | 0.502               | 4.7   |
| Nitrobenzene          |        | 0.380               | 0.370  | 0.388  | 0.375               | 0.388  | 0.374               | 3.9   |
| Hexachlorobutadiene   |        | 0.193               | 0.190  | 0.191  | 0.181               | 0.188  | 0.184               | 4.6   |
| 2,4,6-Trichlorophenol |        | 0.391               | 0.406  | 0.412  | 0.384               | 0.429  | 0.400               | 4.6   |
| 2-Fluorobiphenyl      |        | 1.761               | 1.644  | 1.547  | 1.358               | 1.414  | 1.462               | 13.6  |
| 2,4,5-Trichlorophenol |        | 0.404               | 0.387  | 0.411  | 0.405               | 0.411  | 0.400               | 2.8   |
| 2,4-Dinitrotoluene    |        | 0.269               | 0.296  | 0.342  | 0.354               | 0.372  | 0.333               | 11.0  |
| 2,4,6-Tribromophenol  |        | 0.200               | 0.202  | 0.212  | 0.210               | 0.218  | 0.207               | 3.5   |
| Hexachlorobenzene     |        | 0.269               | 0.248  | 0.252  | 0.239               | 0.258  | 0.249               | 4.8   |
| Pentachlorophenol     |        |                     | 0.119  | 0.146  | 0.148               | 0.160  | 0.147               | 9.7   |
| Terphenyl-d14         |        | 1.567               | 1.456  | 1.419  | 1.311               | 1.342  | 1.344               | 11.4  |

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1