

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: AECO02

Lab Code: CHEM Case No.: P4861 SAS No.: P4861 SDG No.: P4861

Instrument ID: BNA\_F Calibration Date/Time: 11/25/2024 15:49

Lab File ID: BF140604.D Init. Calib. Date(s): 11/21/2024 11/21/2024

EPA Sample No.: SSTDCCC040 Init. Calib. Time(s): 11:13 14:18

GC Column: DB-UI ID: 0.18 (mm)

| COMPOUND              | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|-----------------------|-------|--------|------------|------|-------|
| Pyridine              | 1.084 | 1.248  |            | 15.1 |       |
| 2-Fluorophenol        | 1.172 | 1.153  |            | -1.6 |       |
| Phenol-d6             | 1.550 | 1.521  |            | -1.9 |       |
| 1,4-Dichlorobenzene   | 1.435 | 1.401  |            | -2.4 | 20.0  |
| 2-Methylphenol        | 1.010 | 1.019  |            | 0.9  |       |
| 3+4-Methylphenols     | 1.298 | 1.259  |            | -3.0 |       |
| Nitrobenzene-d5       | 0.391 | 0.378  |            | -3.3 |       |
| Hexachloroethane      | 0.536 | 0.521  |            | -2.8 |       |
| Nitrobenzene          | 0.404 | 0.392  |            | -3.0 |       |
| Hexachlorobutadiene   | 0.215 | 0.210  |            | -2.3 | 20.0  |
| 2,4,6-Trichlorophenol | 0.367 | 0.358  |            | -2.5 | 20.0  |
| 2-Fluorobiphenyl      | 1.342 | 1.282  |            | -4.5 |       |
| 2,4,5-Trichlorophenol | 0.399 | 0.395  |            | -1.0 |       |
| 2,4-Dinitrotoluene    | 0.397 | 0.389  |            | -2.0 |       |
| 2,4,6-Tribromophenol  | 0.214 | 0.206  |            | -3.7 |       |
| Hexachlorobenzene     | 0.238 | 0.236  |            | -0.8 |       |
| Pentachlorophenol     | 0.105 | 0.106  |            | 1.0  | 20.0  |
| Terphenyl-d14         | 1.284 | 1.258  |            | -2.0 |       |

All other compounds must meet a minimum RRF of 0.010.