

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3434</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>10/28/2025 13:17</u>      |
| Lab File ID:    | <u>BF144097.D</u> | Init. Calib. Date(s):  | <u>10/24/2025 10/24/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>10:13 16:24</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|-----------------------------|-------|--------|---------|------|-------|
| 2-Fluorophenol              | 1.264 | 1.303  |         | 3.1  |       |
| Benzaldehyde                | 1.084 | 1.105  |         | 1.9  |       |
| Phenol-d6                   | 1.396 | 1.466  |         | 5.0  |       |
| Phenol                      | 1.722 | 1.775  |         | 3.1  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.245 | 1.329  |         | 6.7  |       |
| 2-Chlorophenol              | 1.245 | 1.300  |         | 4.4  |       |
| 2-Methylphenol              | 0.957 | 1.028  |         | 7.4  |       |
| 2,2-oxybis(1-Chloropropane) | 2.339 | 2.513  |         | 7.4  |       |
| Acetophenone                | 0.432 | 0.450  |         | 4.2  |       |
| 3+4-Methylphenols           | 1.132 | 1.216  |         | 7.4  |       |
| n-Nitroso-di-n-propylamine  | 0.903 | 1.004  | 0.050   | 11.2 |       |
| Nitrobenzene-d5             | 0.365 | 0.371  |         | 1.6  |       |
| Hexachloroethane            | 0.541 | 0.558  |         | 3.1  |       |
| Nitrobenzene                | 0.386 | 0.408  |         | 5.7  |       |
| Isophorone                  | 0.638 | 0.714  |         | 11.9 |       |
| 2-Nitrophenol               | 0.184 | 0.203  |         | 10.3 | 20.0  |
| 2,4-Dimethylphenol          | 0.273 | 0.303  |         | 11.0 |       |
| bis(2-Chloroethoxy)methane  | 0.398 | 0.444  |         | 11.6 |       |
| 2,4-Dichlorophenol          | 0.317 | 0.338  |         | 6.6  | 20.0  |
| Naphthalene                 | 0.944 | 0.987  |         | 4.6  |       |
| 4-Chloroaniline             | 0.329 | 0.364  |         | 10.6 |       |
| Hexachlorobutadiene         | 0.267 | 0.269  |         | 0.7  | 20.0  |
| Caprolactam                 | 0.080 | 0.097  |         | 21.3 |       |
| 4-Chloro-3-methylphenol     | 0.280 | 0.322  |         | 15.0 | 20.0  |
| 2-Methylnaphthalene         | 0.622 | 0.666  |         | 7.1  |       |
| Hexachlorocyclopentadiene   | 0.212 | 0.238  | 0.050   | 12.3 |       |
| 2,4,6-Trichlorophenol       | 0.447 | 0.474  |         | 6.0  | 20.0  |
| 2-Fluorobiphenyl            | 1.421 | 1.379  |         | -3.0 |       |
| 2,4,5-Trichlorophenol       | 0.469 | 0.522  |         | 11.3 |       |
| 1,1-Biphenyl                | 1.432 | 1.391  |         | -2.9 |       |
| 2-Chloronaphthalene         | 1.199 | 1.218  |         | 1.6  |       |
| 2-Nitroaniline              | 0.356 | 0.388  |         | 9.0  |       |
| Dimethylphthalate           | 1.311 | 1.432  |         | 9.2  |       |
| Acenaphthylene              | 1.617 | 1.684  |         | 4.1  |       |
| 2,6-Dinitrotoluene          | 0.259 | 0.297  |         | 14.7 |       |
| 3-Nitroaniline              | 0.288 | 0.326  |         | 13.2 |       |
| Acenaphthene                | 1.080 | 1.111  |         | 2.9  | 20.0  |
| 2,4-Dinitrophenol           | 0.140 | 0.166  | 0.050   | 18.6 |       |
| 4-Nitrophenol               | 0.195 | 0.221  | 0.050   | 13.3 |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3434</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>10/28/2025 13:17</u>      |
| Lab File ID:    | <u>BF144097.D</u> | Init. Calib. Date(s):  | <u>10/24/2025 10/24/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>10:13 16:24</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.499 | 1.565  |            | 4.4  |       |
| 2,4-Dinitrotoluene         | 0.350 | 0.407  |            | 16.3 |       |
| Diethylphthalate           | 1.233 | 1.355  |            | 9.9  |       |
| 4-Chlorophenyl-phenylether | 0.643 | 0.693  |            | 7.8  |       |
| Fluorene                   | 1.143 | 1.207  |            | 5.6  |       |
| 4-Nitroaniline             | 0.273 | 0.304  |            | 11.4 |       |
| 4,6-Dinitro-2-methylphenol | 0.128 | 0.136  |            | 6.3  |       |
| n-Nitrosodiphenylamine     | 0.639 | 0.653  |            | 2.2  | 20.0  |
| 2,4,6-Tribromophenol       | 0.255 | 0.287  |            | 12.5 |       |
| 4-Bromophenyl-phenylether  | 0.250 | 0.266  |            | 6.4  |       |
| Hexachlorobenzene          | 0.307 | 0.322  |            | 4.9  |       |
| Atrazine                   | 0.210 | 0.208  |            | -1.0 |       |
| Pentachlorophenol          | 0.146 | 0.170  |            | 16.4 | 20.0  |
| Phenanthrene               | 0.994 | 1.013  |            | 1.9  |       |
| Anthracene                 | 1.026 | 1.035  |            | 0.9  |       |
| Carbazole                  | 0.874 | 0.869  |            | -0.6 |       |
| Di-n-butylphthalate        | 1.071 | 1.082  |            | 1.0  |       |
| Fluoranthene               | 1.098 | 1.000  |            | -8.9 | 20.0  |
| Pyrene                     | 1.438 | 2.068  |            | 43.8 |       |
| Terphenyl-d14              | 1.144 | 1.573  |            | 37.5 |       |
| Butylbenzylphthalate       | 0.544 | 0.651  |            | 19.7 |       |
| 3,3-Dichlorobenzidine      | 0.470 | 0.517  |            | 10.0 |       |
| Benzo (a) anthracene       | 1.352 | 1.480  |            | 9.5  |       |
| Chrysene                   | 1.268 | 1.334  |            | 5.2  |       |
| Bis(2-ethylhexyl)phthalate | 0.743 | 0.775  |            | 4.3  |       |
| Di-n-octyl phthalate       | 1.277 | 1.323  |            | 3.6  | 20.0  |
| Benzo (b) fluoranthene     | 1.131 | 1.126  |            | -0.4 |       |
| Benzo (k) fluoranthene     | 1.043 | 1.047  |            | 0.4  |       |
| Benzo (a) pyrene           | 1.014 | 1.094  |            | 7.9  | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.373 | 1.546  |            | 12.6 |       |
| Dibenzo (a,h) anthracene   | 1.087 | 1.218  |            | 12.1 |       |
| Benzo (g,h,i) perylene     | 1.093 | 1.265  |            | 15.7 |       |
| 1,2,4,5-Tetrachlorobenzene | 0.687 | 0.668  |            | -2.8 |       |
| 1,4-Dioxane                | 0.546 | 0.520  |            | -4.8 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.330 | 0.372  |            | 12.7 |       |

All other compounds must meet a minimum RRF of 0.010.