

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>ENVI60</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2594</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>07/18/2025 11:24</u>      |
| Lab File ID:    | <u>BF143159.D</u> | Init. Calib. Date(s):  | <u>07/17/2025 07/17/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>11:04 14:34</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|-----------------------------|-------|--------|------------|------|-------|
| 2-Fluorophenol              | 1.264 | 1.224  |            | -3.2 |       |
| Benzaldehyde                | 1.193 | 1.315  |            | 10.2 |       |
| Phenol-d6                   | 1.591 | 1.558  |            | -2.1 |       |
| Phenol                      | 1.733 | 1.713  |            | -1.2 | 20.0  |
| bis(2-Chloroethyl)ether     | 1.327 | 1.320  |            | -0.5 |       |
| 2-Chlorophenol              | 1.325 | 1.308  |            | -1.3 |       |
| 2-Methylphenol              | 1.114 | 1.090  |            | -2.2 |       |
| 2,2-oxybis(1-Chloropropane) | 2.461 | 2.451  |            | -0.4 |       |
| Acetophenone                | 0.474 | 0.456  |            | -3.8 |       |
| 3+4-Methylphenols           | 1.351 | 1.350  |            | -0.1 |       |
| n-Nitroso-di-n-propylamine  | 1.021 | 0.986  | 0.050      | -3.4 |       |
| Nitrobenzene-d5             | 0.401 | 0.411  |            | 2.5  |       |
| Hexachloroethane            | 0.502 | 0.499  |            | -0.6 |       |
| Nitrobenzene                | 0.374 | 0.376  |            | 0.5  |       |
| Isophorone                  | 0.708 | 0.694  |            | -2.0 |       |
| 2-Nitrophenol               | 0.162 | 0.167  |            | 3.1  | 20.0  |
| 2,4-Dimethylphenol          | 0.331 | 0.323  |            | -2.4 |       |
| bis(2-Chloroethoxy)methane  | 0.425 | 0.413  |            | -2.8 |       |
| 2,4-Dichlorophenol          | 0.279 | 0.273  |            | -2.2 | 20.0  |
| Naphthalene                 | 0.985 | 0.953  |            | -3.2 |       |
| 4-Chloroaniline             | 0.402 | 0.387  |            | -3.7 |       |
| Hexachlorobutadiene         | 0.184 | 0.182  |            | -1.1 | 20.0  |
| Caprolactam                 | 0.084 | 0.084  |            | 0.0  |       |
| 4-Chloro-3-methylphenol     | 0.303 | 0.296  |            | -2.3 | 20.0  |
| 2-Methylnaphthalene         | 0.599 | 0.575  |            | -4.0 |       |
| Hexachlorocyclopentadiene   | 0.376 | 0.376  | 0.050      | 0.0  |       |
| 2,4,6-Trichlorophenol       | 0.400 | 0.386  |            | -3.5 | 20.0  |
| 2-Fluorobiphenyl            | 1.462 | 1.373  |            | -6.1 |       |
| 2,4,5-Trichlorophenol       | 0.400 | 0.398  |            | -0.5 |       |
| 1,1-Biphenyl                | 1.560 | 1.495  |            | -4.2 |       |
| 2-Chloronaphthalene         | 1.155 | 1.124  |            | -2.7 |       |
| 2-Nitroaniline              | 0.362 | 0.371  |            | 2.5  |       |
| Dimethylphthalate           | 1.304 | 1.253  |            | -3.9 |       |
| Acenaphthylene              | 1.935 | 1.861  |            | -3.8 |       |
| 2,6-Dinitrotoluene          | 0.265 | 0.270  |            | 1.9  |       |
| 3-Nitroaniline              | 0.310 | 0.306  |            | -1.3 |       |
| Acenaphthene                | 1.144 | 1.092  |            | -4.5 | 20.0  |
| 2,4-Dinitrophenol           | 0.098 | 0.102  | 0.050      | 4.1  |       |
| 4-Nitrophenol               | 0.232 | 0.221  | 0.050      | -4.7 |       |

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| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2594</u>                 |
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| Lab File ID:    | <u>BF143159.D</u> | Init. Calib. Date(s):  | <u>07/17/2025 07/17/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>11:04 14:34</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.698 | 1.630  |            | -4.0 |       |
| 2,4-Dinitrotoluene         | 0.333 | 0.335  |            | 0.6  |       |
| Diethylphthalate           | 1.289 | 1.253  |            | -2.8 |       |
| 4-Chlorophenyl-phenylether | 0.635 | 0.598  |            | -5.8 |       |
| Fluorene                   | 1.274 | 1.203  |            | -5.6 |       |
| 4-Nitroaniline             | 0.266 | 0.261  |            | -1.9 |       |
| 4,6-Dinitro-2-methylphenol | 0.094 | 0.097  |            | 3.2  |       |
| n-Nitrosodiphenylamine     | 0.704 | 0.686  |            | -2.6 | 20.0  |
| 2,4,6-Tribromophenol       | 0.207 | 0.198  |            | -4.3 |       |
| 4-Bromophenyl-phenylether  | 0.238 | 0.236  |            | -0.8 |       |
| Hexachlorobenzene          | 0.249 | 0.243  |            | -2.4 |       |
| Atrazine                   | 0.194 | 0.194  |            | 0.0  |       |
| Pentachlorophenol          | 0.147 | 0.151  |            | 2.7  | 20.0  |
| Phenanthrene               | 1.062 | 1.008  |            | -5.1 |       |
| Anthracene                 | 1.083 | 1.044  |            | -3.6 |       |
| Carbazole                  | 0.946 | 0.911  |            | -3.7 |       |
| Di-n-butylphthalate        | 1.070 | 1.053  |            | -1.6 |       |
| Fluoranthene               | 0.975 | 0.946  |            | -3.0 | 20.0  |
| Pyrene                     | 1.707 | 1.562  |            | -8.5 |       |
| Terphenyl-d14              | 1.344 | 1.220  |            | -9.2 |       |
| Butylbenzylphthalate       | 0.523 | 0.515  |            | -1.5 |       |
| 3,3-Dichlorobenzidine      | 0.446 | 0.437  |            | -2.0 |       |
| Benzo (a) anthracene       | 1.339 | 1.282  |            | -4.3 |       |
| Chrysene                   | 1.204 | 1.187  |            | -1.4 |       |
| Bis(2-ethylhexyl)phthalate | 0.791 | 0.789  |            | -0.3 |       |
| Di-n-octyl phthalate       | 1.434 | 1.381  |            | -3.7 | 20.0  |
| Benzo (b) fluoranthene     | 1.222 | 1.169  |            | -4.3 |       |
| Benzo (k) fluoranthene     | 1.090 | 1.072  |            | -1.7 |       |
| Benzo (a) pyrene           | 1.126 | 1.108  |            | -1.6 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.410 | 1.396  |            | -1.0 |       |
| Dibenzo (a,h) anthracene   | 1.148 | 1.145  |            | -0.3 |       |
| Benzo (g,h,i) perylene     | 1.110 | 1.112  |            | 0.2  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.577 | 0.563  |            | -2.4 |       |
| 1,4-Dioxane                | 0.598 | 0.597  |            | -0.2 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.331 | 0.324  |            | -2.1 |       |

All other compounds must meet a minimum RRF of 0.010.