

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>HDRI08</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3466</u>                 |
| Instrument ID:  | <u>BNA_G</u>      | Calibration Date/Time: | <u>10/29/2025 11:22</u>      |
| Lab File ID:    | <u>BG064618.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>09:00 15:32</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D    | MAX%D |
|-----------------------------|-------|--------|---------|-------|-------|
| 2-Fluorophenol              | 1.192 | 1.266  |         | 6.2   |       |
| Benzaldehyde                | 0.893 | 0.838  |         | -6.2  |       |
| Phenol-d6                   | 1.635 | 1.656  |         | 1.3   |       |
| Phenol                      | 1.808 | 1.838  |         | 1.7   | 20.0  |
| bis(2-Chloroethyl)ether     | 1.329 | 1.382  |         | 4.0   |       |
| 2-Chlorophenol              | 1.247 | 1.377  |         | 10.4  |       |
| 2-Methylphenol              | 1.201 | 1.261  |         | 5.0   |       |
| 2,2-oxybis(1-Chloropropane) | 3.487 | 3.674  |         | 5.4   |       |
| Acetophenone                | 0.547 | 0.581  |         | 6.2   |       |
| 3+4-Methylphenols           | 1.688 | 1.742  |         | 3.2   |       |
| n-Nitroso-di-n-propylamine  | 1.130 | 1.223  | 0.050   | 8.2   |       |
| Nitrobenzene-d5             | 0.333 | 0.401  |         | 20.4  |       |
| Hexachloroethane            | 0.512 | 0.567  |         | 10.7  |       |
| Nitrobenzene                | 0.362 | 0.429  |         | 18.5  |       |
| Isophorone                  | 0.799 | 0.833  |         | 4.3   |       |
| 2-Nitrophenol               | 0.121 | 0.168  |         | 38.8  | 20.0  |
| 2,4-Dimethylphenol          | 0.297 | 0.311  |         | 4.7   |       |
| bis(2-Chloroethoxy)methane  | 0.455 | 0.480  |         | 5.5   |       |
| 2,4-Dichlorophenol          | 0.320 | 0.352  |         | 10.0  | 20.0  |
| Naphthalene                 | 1.109 | 1.159  |         | 4.5   |       |
| 4-Chloroaniline             | 0.431 | 0.435  |         | 0.9   |       |
| Hexachlorobutadiene         | 0.284 | 0.322  |         | 13.4  | 20.0  |
| Caprolactam                 | 0.123 | 0.108  |         | -12.2 |       |
| 4-Chloro-3-methylphenol     | 0.384 | 0.389  |         | 1.3   | 20.0  |
| 2-Methylnaphthalene         | 0.814 | 0.836  |         | 2.7   |       |
| Hexachlorocyclopentadiene   | 0.302 | 0.400  | 0.050   | 32.5  |       |
| 2,4,6-Trichlorophenol       | 0.384 | 0.455  |         | 18.5  | 20.0  |
| 2-Fluorobiphenyl            | 1.319 | 1.473  |         | 11.7  |       |
| 2,4,5-Trichlorophenol       | 0.443 | 0.492  |         | 11.1  |       |
| 1,1-Biphenyl                | 1.422 | 1.535  |         | 7.9   |       |
| 2-Chloronaphthalene         | 1.073 | 1.175  |         | 9.5   |       |
| 2-Nitroaniline              | 0.311 | 0.378  |         | 21.5  |       |
| Dimethylphthalate           | 1.530 | 1.552  |         | 1.4   |       |
| Acenaphthylene              | 1.693 | 1.770  |         | 4.5   |       |
| 2,6-Dinitrotoluene          | 0.218 | 0.272  |         | 24.8  |       |
| 3-Nitroaniline              | 0.263 | 0.288  |         | 9.5   |       |
| Acenaphthene                | 1.157 | 1.225  |         | 5.9   | 20.0  |
| 2,4-Dinitrophenol           | 0.132 | 0.164  | 0.050   | 24.2  |       |
| 4-Nitrophenol               | 0.244 | 0.241  | 0.050   | -1.2  |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>HDRI08</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q3466</u>                 |
| Instrument ID:  | <u>BNA_G</u>      | Calibration Date/Time: | <u>10/29/2025 11:22</u>      |
| Lab File ID:    | <u>BG064618.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>09:00 15:32</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.870 | 1.930  |            | 3.2  |       |
| 2,4-Dinitrotoluene         | 0.340 | 0.396  |            | 16.5 |       |
| Diethylphthalate           | 1.563 | 1.507  |            | -3.6 |       |
| 4-Chlorophenyl-phenylether | 0.820 | 0.840  |            | 2.4  |       |
| Fluorene                   | 1.460 | 1.476  |            | 1.1  |       |
| 4-Nitroaniline             | 0.293 | 0.306  |            | 4.4  |       |
| 4,6-Dinitro-2-methylphenol | 0.079 | 0.104  |            | 31.6 |       |
| n-Nitrosodiphenylamine     | 0.540 | 0.582  |            | 7.8  | 20.0  |
| 2,4,6-Tribromophenol       | 0.289 | 0.312  |            | 8.0  |       |
| 4-Bromophenyl-phenylether  | 0.237 | 0.262  |            | 10.5 |       |
| Hexachlorobenzene          | 0.270 | 0.300  |            | 11.1 |       |
| Atrazine                   | 0.232 | 0.233  |            | 0.4  |       |
| Pentachlorophenol          | 0.169 | 0.189  |            | 11.8 | 20.0  |
| Phenanthrene               | 1.094 | 1.132  |            | 3.5  |       |
| Anthracene                 | 1.113 | 1.125  |            | 1.1  |       |
| Carbazole                  | 0.981 | 0.970  |            | -1.1 |       |
| Di-n-butylphthalate        | 1.135 | 1.140  |            | 0.4  |       |
| Fluoranthene               | 1.370 | 1.343  |            | -2.0 | 20.0  |
| Pyrene                     | 1.307 | 1.322  |            | 1.1  |       |
| Terphenyl-d14              | 1.060 | 1.074  |            | 1.3  |       |
| Butylbenzylphthalate       | 0.395 | 0.462  |            | 17.0 |       |
| 3,3-Dichlorobenzidine      | 0.443 | 0.484  |            | 9.3  |       |
| Benzo (a) anthracene       | 1.326 | 1.395  |            | 5.2  |       |
| Chrysene                   | 1.262 | 1.298  |            | 2.9  |       |
| Bis(2-ethylhexyl)phthalate | 0.590 | 0.671  |            | 13.7 |       |
| Di-n-octyl phthalate       | 0.995 | 1.203  |            | 20.9 | 20.0  |
| Benzo (b) fluoranthene     | 1.226 | 1.236  |            | 0.8  |       |
| Benzo (k) fluoranthene     | 1.227 | 1.235  |            | 0.7  |       |
| Benzo (a) pyrene           | 1.127 | 1.157  |            | 2.7  | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.471 | 1.519  |            | 3.3  |       |
| Dibenzo (a,h) anthracene   | 1.195 | 1.241  |            | 3.8  |       |
| Benzo (g,h,i) perylene     | 1.142 | 1.182  |            | 3.5  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.671 | 0.754  |            | 12.4 |       |
| 1,4-Dioxane                | 0.528 | 0.560  |            | 6.1  | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.424 | 0.469  |            | 10.6 |       |

All other compounds must meet a minimum RRF of 0.010.