

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: CORE02

Lab Code: ACE

SDG No.: Q3741

Instrument ID: BNA\_F

Calibration Date(s): 11/05/2025 11/05/2025

Calibration Time(s): 12:50 16:49

| LAB FILE ID:                |        | RRF2.5 = BF144169.D |        |        | RRF005 = BF144170.D |        | RRF010 = BF144171.D |       |
|-----------------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
|                             |        | RRF020 = BF144172.D |        |        | RRF040 = BF144173.D |        | RRF050 = BF144174.D |       |
| COMPOUND                    | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| 2-Fluorophenol              |        | 1.332               | 1.303  | 1.316  | 1.315               | 1.157  | 1.278               | 5.6   |
| Benzaldehyde                |        |                     | 1.033  | 0.917  | 0.816               | 0.672  | 0.818               | 17.0  |
| Phenol-d6                   |        | 1.566               | 1.518  | 1.511  | 1.460               | 1.281  | 1.448               | 7.1   |
| Phenol                      |        | 1.845               | 1.896  | 1.840  | 1.858               | 1.562  | 1.769               | 7.6   |
| bis(2-Chloroethyl)ether     |        | 1.278               | 1.333  | 1.342  | 1.316               | 1.155  | 1.289               | 5.2   |
| 2-Chlorophenol              |        | 1.242               | 1.309  | 1.292  | 1.281               | 1.152  | 1.253               | 4.6   |
| 2-Methylphenol              |        | 1.031               | 1.054  | 1.007  | 1.033               | 0.895  | 0.997               | 5.4   |
| 2,2-oxybis(1-Chloropropane) |        | 2.408               | 2.470  | 2.496  | 2.418               | 2.147  | 2.374               | 5.5   |
| Acetophenone                |        | 0.479               | 0.464  | 0.434  | 0.434               | 0.380  | 0.429               | 8.3   |
| 3+4-Methylphenols           |        |                     | 1.330  | 1.223  | 1.215               | 1.030  | 1.175               | 9.2   |
| n-Nitroso-di-n-propylamine  | 0.998  | 1.010               | 1.023  | 0.996  | 0.950               | 0.813  | 0.953               | 7.4   |
| Nitrobenzene-d5             |        | 0.343               | 0.352  | 0.349  | 0.360               | 0.321  | 0.346               | 4.1   |
| Hexachloroethane            |        | 0.512               | 0.513  | 0.532  | 0.532               | 0.467  | 0.515               | 4.9   |
| Nitrobenzene                |        | 0.370               | 0.390  | 0.376  | 0.388               | 0.342  | 0.376               | 4.9   |
| Isophorone                  |        | 0.659               | 0.668  | 0.669  | 0.673               | 0.581  | 0.655               | 5.1   |
| 2-Nitrophenol               |        | 0.167               | 0.187  | 0.191  | 0.197               | 0.174  | 0.186               | 6.5   |
| 2,4-Dimethylphenol          |        | 0.281               | 0.278  | 0.265  | 0.298               | 0.260  | 0.279               | 5.1   |
| bis(2-Chloroethoxy)methane  |        | 0.426               | 0.417  | 0.412  | 0.423               | 0.372  | 0.409               | 5.0   |
| 2,4-Dichlorophenol          |        | 0.334               | 0.342  | 0.330  | 0.331               | 0.289  | 0.322               | 6.2   |
| Naphthalene                 |        | 1.019               | 1.008  | 0.982  | 0.969               | 0.850  | 0.951               | 6.7   |
| 4-Chloroaniline             |        | 0.358               | 0.364  | 0.355  | 0.362               | 0.309  | 0.347               | 6.1   |
| Hexachlorobutadiene         |        | 0.256               | 0.250  | 0.252  | 0.259               | 0.232  | 0.249               | 4.0   |
| Caprolactam                 |        |                     | 0.094  | 0.093  | 0.092               | 0.075  | 0.088               | 8.1   |
| 4-Chloro-3-methylphenol     |        | 0.283               | 0.307  | 0.304  | 0.297               | 0.259  | 0.288               | 5.8   |
| 2-Methylnaphthalene         |        | 0.702               | 0.672  | 0.665  | 0.647               | 0.559  | 0.636               | 8.0   |
| Hexachlorocyclopentadiene   |        |                     | 0.121  | 0.142  | 0.200               | 0.204  | 0.195               | 27.8  |
| 2,4,6-Trichlorophenol       |        | 0.440               | 0.449  | 0.449  | 0.456               | 0.409  | 0.446               | 5.5   |
| 2-Fluorobiphenyl            |        | 1.598               | 1.428  | 1.361  | 1.350               | 1.177  | 1.354               | 10.7  |
| 2,4,5-Trichlorophenol       |        | 0.497               | 0.489  | 0.497  | 0.489               | 0.438  | 0.481               | 5.1   |
| 1,1-Biphenyl                |        | 1.535               | 1.446  | 1.425  | 1.408               | 1.239  | 1.396               | 7.5   |
| 2-Chloronaphthalene         |        | 1.303               | 1.216  | 1.206  | 1.204               | 1.068  | 1.191               | 7.1   |
| 2-Nitroaniline              |        | 0.302               | 0.347  | 0.349  | 0.361               | 0.326  | 0.344               | 6.8   |
| Dimethylphthalate           |        | 1.437               | 1.343  | 1.359  | 1.345               | 1.168  | 1.325               | 6.7   |
| Acenaphthylene              |        | 1.727               | 1.686  | 1.677  | 1.645               | 1.440  | 1.623               | 6.6   |

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: CORE02

Lab Code: ACE

SDG No.: Q3741

Instrument ID: BNA\_F

Calibration Date(s): 11/05/2025 11/05/2025

Calibration Time(s): 12:50 16:49

| LAB FILE ID:                 |        | RRF2.5 = BF144169.D |        |        | RRF005 = BF144170.D |        | RRF010 = BF144171.D |       |
|------------------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
|                              |        | RRF020 = BF144172.D |        |        | RRF040 = BF144173.D |        | RRF050 = BF144174.D |       |
| COMPOUND                     | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| 2,6-Dinitrotoluene           |        | 0.251               | 0.271  | 0.271  | 0.276               | 0.252  | 0.268               | 5.2   |
| 3-Nitroaniline               |        | 0.271               | 0.301  | 0.302  | 0.315               | 0.264  | 0.293               | 7.1   |
| Acenaphthene                 |        | 1.138               | 1.162  | 1.120  | 1.096               | 0.956  | 1.085               | 6.9   |
| 2,4-Dinitrophenol            |        |                     | 0.202  | 0.166  | 0.140               | 0.133  | 0.162               | 15.0  |
| 4-Nitrophenol                |        |                     | 0.273  | 0.202  | 0.204               | 0.185  | 0.212               | 14.6  |
| Dibenzofuran                 |        | 1.644               | 1.600  | 1.586  | 1.551               | 1.337  | 1.520               | 8.0   |
| 2,4-Dinitrotoluene           |        | 0.322               | 0.377  | 0.372  | 0.385               | 0.328  | 0.359               | 7.1   |
| Diethylphthalate             |        | 1.294               | 1.273  | 1.254  | 1.254               | 1.075  | 1.221               | 6.3   |
| 4-Chlorophenyl-phenylether   |        | 0.711               | 0.701  | 0.678  | 0.685               | 0.570  | 0.658               | 7.9   |
| Fluorene                     |        | 1.295               | 1.223  | 1.197  | 1.180               | 0.997  | 1.156               | 8.9   |
| 4-Nitroaniline               |        | 0.273               | 0.290  | 0.293  | 0.288               | 0.254  | 0.279               | 5.1   |
| 4,6-Dinitro-2-methylphenol   |        |                     | 0.147  | 0.138  | 0.130               | 0.120  | 0.136               | 7.0   |
| n-Nitrosodiphenylamine       |        | 0.695               | 0.676  | 0.635  | 0.646               | 0.577  | 0.643               | 6.1   |
| 2,4,6-Tribromophenol         |        | 0.242               | 0.257  | 0.257  | 0.265               | 0.227  | 0.251               | 5.3   |
| 4-Bromophenyl-phenylether    |        | 0.264               | 0.260  | 0.255  | 0.263               | 0.228  | 0.255               | 5.2   |
| Hexachlorobenzene            |        | 0.306               | 0.306  | 0.305  | 0.300               | 0.273  | 0.301               | 4.7   |
| Atrazine                     |        | 0.215               | 0.211  | 0.215  | 0.206               | 0.180  | 0.205               | 6.3   |
| Pentachlorophenol            |        |                     | 0.135  | 0.144  | 0.156               | 0.137  | 0.150               | 8.8   |
| Phenanthrene                 |        | 1.103               | 1.042  | 0.986  | 1.009               | 0.892  | 0.996               | 6.9   |
| Anthracene                   |        | 1.120               | 1.059  | 1.027  | 1.022               | 0.897  | 1.013               | 7.4   |
| Carbazole                    |        | 0.938               | 0.917  | 0.882  | 0.878               | 0.762  | 0.861               | 7.4   |
| Di-n-butylphthalate          |        | 1.075               | 1.101  | 1.077  | 1.070               | 0.917  | 1.041               | 6.3   |
| Fluoranthene                 |        | 1.117               | 1.066  | 1.053  | 1.054               | 0.910  | 1.029               | 6.8   |
| Pyrene                       |        | 1.763               | 1.745  | 1.631  | 1.704               | 1.372  | 1.613               | 8.8   |
| Terphenyl-d14                |        | 1.403               | 1.415  | 1.274  | 1.317               | 1.042  | 1.259               | 10.7  |
| Butylbenzylphthalate         |        | 0.565               | 0.590  | 0.550  | 0.590               | 0.499  | 0.559               | 5.6   |
| 3,3-Dichlorobenzidine        |        |                     | 0.453  | 0.461  | 0.482               | 0.441  | 0.475               | 6.8   |
| Benzo (a) anthracene         |        | 1.455               | 1.340  | 1.330  | 1.430               | 1.290  | 1.386               | 5.4   |
| Chrysene                     |        | 1.276               | 1.245  | 1.279  | 1.403               | 1.164  | 1.280               | 5.7   |
| Bis (2-ethylhexyl) phthalate |        | 0.773               | 0.739  | 0.755  | 0.816               | 0.705  | 0.765               | 4.9   |
| Di-n-octyl phthalate         |        |                     | 1.163  | 1.272  | 1.410               | 1.269  | 1.320               | 7.9   |
| Benzo (b) fluoranthene       |        | 1.199               | 1.133  | 1.114  | 1.233               | 1.002  | 1.137               | 6.4   |
| Benzo (k) fluoranthene       |        | 1.077               | 1.094  | 1.083  | 1.041               | 0.999  | 1.064               | 4.4   |
| Benzo (a) pyrene             |        | 1.064               | 1.050  | 1.058  | 1.077               | 0.963  | 1.049               | 4.2   |
| Indeno (1,2,3-cd) pyrene     |        | 1.450               | 1.411  | 1.447  | 1.485               | 1.311  | 1.436               | 4.6   |

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: CORE02

Lab Code: ACE

SDG No.: Q3741

Instrument ID: BNA\_F

Calibration Date(s): 11/05/2025 11/05/2025

Calibration Time(s): 12:50 16:49

|                            |        |                     |        |                     |        |                     |       |       |
|----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:               |        | RRF2.5 = BF144169.D |        | RRF005 = BF144170.D |        | RRF010 = BF144171.D |       |       |
|                            |        | RRF020 = BF144172.D |        | RRF040 = BF144173.D |        | RRF050 = BF144174.D |       |       |
| COMPOUND                   | RRF2.5 | RRF005              | RRF010 | RRF020              | RRF040 | RRF050              | RRF   | % RSD |
| Dibenzo(a,h)anthracene     |        | 1.155               | 1.155  | 1.175               | 1.198  | 1.040               | 1.153 | 4.8   |
| Benzo(g,h,i)perylene       |        | 1.095               | 1.125  | 1.153               | 1.191  | 1.028               | 1.139 | 5.7   |
| 1,2,4,5-Tetrachlorobenzene |        | 0.677               | 0.662  | 0.659               | 0.675  | 0.588               | 0.655 | 6.6   |
| 2,3,4,6-Tetrachlorophenol  |        | 0.308               | 0.347  | 0.336               | 0.358  | 0.316               | 0.340 | 6.6   |