

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: JPCL01  
Lab Code: CHEM Case No.: P4768 SAS No.: P4768 SDG No.: P4768  
Instrument ID: BNA\_F Calibration Date(s): 11/05/2024 11/05/2024  
Calibration Time(s): 12:26 15:30

| LAB FILE ID:                |        | RRF2.5 = BF140231.D |        |        | RRF005 = BF140232.D |        | RRF010 = BF140233.D |       |
|-----------------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
|                             |        | RRF020 = BF140234.D |        |        | RRF040 = BF140235.D |        | RRF050 = BF140236.D |       |
| COMPOUND                    | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| 2-Fluorophenol              |        | 1.359               | 1.298  | 1.261  | 1.159               | 1.138  | 1.207               | 8.3   |
| Benzaldehyde                |        | 1.159               | 1.137  | 1.085  | 0.949               | 0.839  | 0.956               | 18.5  |
| Phenol-d6                   |        | 1.765               | 1.679  | 1.604  | 1.495               | 1.462  | 1.552               | 8.7   |
| Phenol                      |        | 1.863               | 1.779  | 1.752  | 1.593               | 1.575  | 1.649               | 9.2   |
| bis(2-Chloroethyl) ether    |        | 1.417               | 1.354  | 1.285  | 1.209               | 1.201  | 1.252               | 8.6   |
| 2-Chlorophenol              |        | 1.441               | 1.385  | 1.311  | 1.214               | 1.187  | 1.263               | 9.4   |
| 2-Methylphenol              |        | 1.155               | 1.105  | 1.086  | 1.028               | 1.005  | 1.053               | 6.1   |
| 2,2-oxybis(1-Chloropropane) |        | 2.305               | 2.213  | 2.143  | 1.958               | 1.910  | 2.027               | 9.7   |
| Acetophenone                |        | 0.564               | 0.535  | 0.512  | 0.474               | 0.461  | 0.496               | 8.4   |
| 3+4-Methylphenols           |        | 1.503               | 1.446  | 1.403  | 1.277               | 1.240  | 1.323               | 9.7   |
| n-Nitroso-di-n-propylamine  | 1.035  | 1.117               | 1.062  | 1.015  | 0.949               | 0.923  | 0.990               | 8.0   |
| Nitrobenzene-d5             |        | 0.435               | 0.420  | 0.415  | 0.392               | 0.387  | 0.402               | 5.2   |
| Hexachloroethane            |        | 0.612               | 0.603  | 0.576  | 0.533               | 0.524  | 0.556               | 7.4   |
| Nitrobenzene                |        | 0.452               | 0.442  | 0.436  | 0.414               | 0.407  | 0.422               | 5.0   |
| Isophorone                  |        | 0.754               | 0.722  | 0.699  | 0.675               | 0.659  | 0.693               | 5.1   |
| 2-Nitrophenol               |        | 0.160               | 0.167  | 0.170  | 0.171               | 0.167  | 0.168               | 2.5   |
| 2,4-Dimethylphenol          |        | 0.240               | 0.237  | 0.234  | 0.224               | 0.221  | 0.228               | 3.9   |
| bis(2-Chloroethoxy) methane |        | 0.459               | 0.441  | 0.428  | 0.400               | 0.390  | 0.415               | 6.9   |
| 2,4-Dichlorophenol          |        | 0.297               | 0.299  | 0.291  | 0.278               | 0.270  | 0.283               | 4.7   |
| Naphthalene                 |        | 1.206               | 1.119  | 1.071  | 0.992               | 0.956  | 1.033               | 9.9   |
| 4-Chloroaniline             |        | 0.372               | 0.367  | 0.354  | 0.330               | 0.322  | 0.340               | 7.1   |
| Hexachlorobutadiene         |        | 0.251               | 0.241  | 0.238  | 0.222               | 0.218  | 0.229               | 6.2   |
| Caprolactam                 |        | 0.090               | 0.087  | 0.088  | 0.085               | 0.084  | 0.086               | 2.9   |
| 4-Chloro-3-methylphenol     |        | 0.337               | 0.326  | 0.319  | 0.312               | 0.303  | 0.315               | 4.3   |
| 2-Methylnaphthalene         |        | 0.772               | 0.743  | 0.702  | 0.644               | 0.626  | 0.674               | 9.6   |
| Hexachlorocyclopentadiene   |        | 0.124               | 0.145  | 0.169  | 0.183               | 0.177  | 0.164               | 13.3  |
| 2,4,6-Trichlorophenol       |        | 0.377               | 0.369  | 0.363  | 0.361               | 0.354  | 0.363               | 2.2   |
| 2-Fluorobiphenyl            |        | 1.481               | 1.381  | 1.312  | 1.183               | 1.148  | 1.253               | 11.3  |
| 2,4,5-Trichlorophenol       |        | 0.391               | 0.398  | 0.396  | 0.373               | 0.374  | 0.381               | 4.0   |
| 1,1-Biphenyl                |        | 1.663               | 1.590  | 1.521  | 1.376               | 1.342  | 1.446               | 10.0  |
| 2-Chloronaphthalene         |        | 1.195               | 1.171  | 1.135  | 1.050               | 1.036  | 1.088               | 7.2   |
| 2-Nitroaniline              |        | 0.357               | 0.372  | 0.371  | 0.369               | 0.363  | 0.364               | 2.0   |
| Dimethylphthalate           |        | 1.363               | 1.310  | 1.262  | 1.191               | 1.182  | 1.232               | 6.7   |
| Acenaphthylene              |        | 1.730               | 1.668  | 1.618  | 1.486               | 1.465  | 1.540               | 8.6   |

All other compounds must meet a minimum RRF of 0.010.

Form VI SV-1

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: JPCL01  
Lab Code: CHEM Case No.: P4768 SAS No.: P4768 SDG No.: P4768  
Instrument ID: BNA\_F Calibration Date(s): 11/05/2024 11/05/2024  
Calibration Time(s): 12:26 15:30

|                              |        |                     |        |        |                     |        |                     |       |
|------------------------------|--------|---------------------|--------|--------|---------------------|--------|---------------------|-------|
| LAB FILE ID:                 |        | RRF2.5 = BF140231.D |        |        | RRF005 = BF140232.D |        | RRF010 = BF140233.D |       |
|                              |        | RRF020 = BF140234.D |        |        | RRF040 = BF140235.D |        | RRF050 = BF140236.D |       |
| COMPOUND                     | RRF2.5 | RRF005              | RRF010 | RRF020 | RRF040              | RRF050 | RRF                 | % RSD |
| 2,6-Dinitrotoluene           |        | 0.302               | 0.297  | 0.299  | 0.286               | 0.286  | 0.290               | 3.6   |
| 3-Nitroaniline               |        | 0.294               | 0.286  | 0.282  | 0.269               | 0.267  | 0.273               | 5.8   |
| Acenaphthene                 |        | 1.204               | 1.153  | 1.134  | 1.059               | 1.055  | 1.093               | 6.6   |
| 2,4-Dinitrophenol            |        |                     | 0.071  | 0.101  | 0.128               | 0.135  | 0.120               | 23.2  |
| 4-Nitrophenol                |        | 0.160               | 0.178  | 0.196  | 0.203               | 0.206  | 0.192               | 8.8   |
| Dibenzofuran                 |        | 1.773               | 1.683  | 1.603  | 1.467               | 1.424  | 1.528               | 10.6  |
| 2,4-Dinitrotoluene           |        | 0.384               | 0.385  | 0.388  | 0.374               | 0.373  | 0.376               | 3.2   |
| Diethylphthalate             |        | 1.360               | 1.319  | 1.271  | 1.185               | 1.170  | 1.224               | 7.7   |
| 4-Chlorophenyl-phenylether   |        | 0.715               | 0.673  | 0.638  | 0.581               | 0.573  | 0.613               | 10.3  |
| Fluorene                     |        | 1.446               | 1.345  | 1.267  | 1.154               | 1.127  | 1.216               | 11.6  |
| 4-Nitroaniline               |        | 0.278               | 0.276  | 0.282  | 0.270               | 0.267  | 0.271               | 3.5   |
| 4,6-Dinitro-2-methylphenol   |        |                     | 0.080  | 0.100  | 0.110               | 0.112  | 0.105               | 12.6  |
| n-Nitrosodiphenylamine       |        | 0.646               | 0.608  | 0.600  | 0.551               | 0.543  | 0.575               | 7.5   |
| 2,4,6-Tribromophenol         |        | 0.199               | 0.200  | 0.197  | 0.194               | 0.197  | 0.198               | 1.4   |
| 4-Bromophenyl-phenylether    |        | 0.230               | 0.219  | 0.223  | 0.209               | 0.208  | 0.215               | 4.1   |
| Hexachlorobenzene            |        | 0.261               | 0.255  | 0.251  | 0.237               | 0.235  | 0.245               | 4.3   |
| Atrazine                     |        | 0.204               | 0.189  | 0.155  | 0.156               | 0.127  | 0.157               | 18.5  |
| Pentachlorophenol            |        | 0.105               | 0.121  | 0.133  | 0.138               | 0.139  | 0.132               | 10.6  |
| Phenanthrene                 |        | 1.082               | 1.044  | 1.004  | 0.915               | 0.885  | 0.954               | 9.3   |
| Anthracene                   |        | 1.071               | 1.014  | 0.984  | 0.894               | 0.881  | 0.936               | 9.3   |
| Carbazole                    |        | 0.973               | 0.919  | 0.901  | 0.822               | 0.805  | 0.855               | 8.9   |
| Di-n-butylphthalate          |        | 1.119               | 1.096  | 1.062  | 0.975               | 0.961  | 1.012               | 7.9   |
| Fluoranthene                 |        | 1.157               | 1.101  | 1.044  | 0.950               | 0.928  | 0.998               | 10.5  |
| Pyrene                       |        | 1.851               | 1.733  | 1.728  | 1.590               | 1.572  | 1.651               | 7.5   |
| Terphenyl-d14                |        | 1.383               | 1.309  | 1.258  | 1.159               | 1.164  | 1.221               | 8.2   |
| Butylbenzylphthalate         |        | 0.580               | 0.596  | 0.613  | 0.597               | 0.601  | 0.595               | 2.9   |
| 3,3-Dichlorobenzidine        |        | 0.387               | 0.409  | 0.410  | 0.417               | 0.406  | 0.403               | 2.6   |
| Benzo (a) anthracene         |        | 1.386               | 1.344  | 1.345  | 1.232               | 1.232  | 1.284               | 5.7   |
| Chrysene                     |        | 1.320               | 1.264  | 1.206  | 1.154               | 1.177  | 1.202               | 5.8   |
| Bis (2-ethylhexyl) phthalate |        | 0.868               | 0.858  | 0.858  | 0.819               | 0.825  | 0.833               | 3.8   |
| Di-n-octyl phthalate         |        | 1.201               | 1.234  | 1.283  | 1.267               | 1.265  | 1.254               | 2.5   |
| Benzo (b) fluoranthene       |        | 1.300               | 1.184  | 1.214  | 1.245               | 1.125  | 1.210               | 6.0   |
| Benzo (k) fluoranthene       |        | 1.214               | 1.230  | 1.206  | 1.027               | 1.085  | 1.110               | 9.5   |
| Benzo (a) pyrene             |        | 1.023               | 0.999  | 1.022  | 0.993               | 0.975  | 0.997               | 2.2   |
| Indeno (1,2,3-cd) pyrene     |        | 1.107               | 1.110  | 1.213  | 1.195               | 1.172  | 1.174               | 4.1   |

All other compounds must meet a minimum RRF of 0.010.

6C

SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: JPCL01  
Lab Code: CHEM Case No.: P4768 SAS No.: P4768 SDG No.: P4768  
Instrument ID: BNA\_F Calibration Date(s): 11/05/2024 11/05/2024  
Calibration Time(s): 12:26 15:30

|                            |        |                     |        |                     |        |                     |       |       |
|----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:               |        | RRF2.5 = BF140231.D |        | RRF005 = BF140232.D |        | RRF010 = BF140233.D |       |       |
|                            |        | RRF020 = BF140234.D |        | RRF040 = BF140235.D |        | RRF050 = BF140236.D |       |       |
| COMPOUND                   | RRF2.5 | RRF005              | RRF010 | RRF020              | RRF040 | RRF050              | RRF   | % RSD |
| Dibenzo (a,h) anthracene   |        | 0.917               | 0.923  | 0.998               | 0.979  | 0.968               | 0.966 | 3.4   |
| Benzo (g,h,i) perylene     |        | 0.949               | 0.941  | 1.017               | 0.994  | 0.977               | 0.985 | 3.1   |
| 1,2,4,5-Tetrachlorobenzene |        | 0.657               | 0.619  | 0.600               | 0.562  | 0.559               | 0.586 | 7.2   |
| 1,4-Dioxane                |        | 0.613               | 0.582  | 0.576               | 0.550  | 0.526               | 0.555 | 6.6   |
| 2,3,4,6-Tetrachlorophenol  |        | 0.323               | 0.324  | 0.321               | 0.305  | 0.309               | 0.313 | 3.3   |