

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070524\  
 Data File : BG061908.D  
 Acq On : 5 Jul 2024 23:11  
 Operator : MA/JU  
 Sample : P2982-11  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A4CF7

Quant Time: Jul 06 00:24:34 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 03 02:38:46 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.094  | 152  | 56116    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.909 | 136  | 273825   | 20.000 | ng/ul  | # 0.00   |
| 38) Acenaphthene-d10          | 14.716 | 164  | 197110   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.460 | 188  | 458519   | 20.000 | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.714 | 240  | 411585   | 20.000 | ng/ul  | -0.01    |
| 88) Perylene-d12              | 24.957 | 264  | 496231   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.482  | 96   | 3967     | 2.657  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.905  | 84   | 50516    | 11.002 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.248  | 99   | 94791    | 15.435 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.419  | 67   | 58287    | 14.988 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.624  | 132  | 69225    | 16.433 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.793  | 113  | 75513    | 15.617 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.258  | 128  | 36122    | 17.345 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.980  | 143  | 45825    | 19.269 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.521 | 165  | 80347    | 17.516 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.038 | 131  | 107313   | 16.114 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.117 | 166  | 299395   | 18.475 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.416 | 160  | 326021   | 19.240 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.904 | 143  | 42935    | 16.591 | ng/ul  | -0.01    |
| 60) Fluorene-d10              | 15.703 | 176  | 261741   | 19.186 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.815 | 200  | 40929    | 16.997 | ng/ul  | -0.01    |
| 73) Anthracene-d10            | 17.554 | 188  | 398874   | 18.632 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.833 | 212  | 452124   | 18.676 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.739 | 264  | 486079   | 19.751 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.501 | 178  | 30837    | 1.275  | ng/ul# | 93       |
| 80) Fluoranthene              | 19.499 | 202  | 54325    | 1.896  | ng/ul  | 96       |
| 82) Pyrene                    | 19.863 | 202  | 59165    | 1.999  | ng/ul  | 97       |
| 87) Chrysene                  | 21.766 | 228  | 28084    | 1.007  | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.923 | 252  | 32362    | 1.035  | ng/ul# | 95       |
| -----                         |        |      |          |        |        |          |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070524\  
Data File : BG061908.D  
Acq On : 5 Jul 2024 23:11  
Operator : MA/JU  
Sample : P2982-11  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CF7

Quant Time: Jul 06 00:24:34 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070224.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Jul 03 02:38:46 2024  
Response via : Initial Calibration

