

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070823\  
 Data File : BG058112.D  
 Acq On : 9 Jul 2023 4:19  
 Operator : CG/JU  
 Sample : 03349-05  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 YBW00

Quant Time: Jul 09 05:01:37 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jul 08 10:01:01 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.917  | 152  | 56843    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.720 | 136  | 252013   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.551 | 164  | 189317   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.300 | 188  | 485969   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.548 | 240  | 451101   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.704 | 264  | 552208   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.358  | 96   | 3849     | 2.371  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.763  | 84   | 56152    | 11.394 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.077  | 99   | 75720    | 13.044 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.242  | 67   | 48477    | 13.857 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.447  | 132  | 54366    | 13.579 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.622  | 113  | 60927    | 13.909 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.075  | 128  | 29310    | 14.270 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.797  | 143  | 32527    | 15.068 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.338 | 165  | 61618    | 14.879 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.855 | 131  | 87096    | 13.872 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.957 | 166  | 233583   | 15.637 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.245 | 160  | 259122   | 15.897 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.750 | 143  | 28552    | 10.819 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.544 | 176  | 215701   | 17.003 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.661 | 200  | 27023    | 10.241 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.394 | 188  | 373253   | 16.552 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.686 | 212  | 451555   | 15.894 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.486 | 264  | 469680   | 16.695 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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