

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG030323\  
 Data File : BG056820.D  
 Acq On : 4 Mar 2023 2:05  
 Operator : CG/JU  
 Sample : 01711-11  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DBAC7

Quant Time: Mar 04 03:08:02 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022323.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 02 00:12:28 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.421  | 152  | 14230    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 11.253 | 136  | 66096    | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 15.019 | 164  | 54227    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.751 | 188  | 135877   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 22.064 | 240  | 129497   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.589 | 264  | 141530   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.691  | 96   | 955      | 3.723  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1 |          |
| 7) Phenol-d5                  | 7.540  | 99   | 21961    | 15.422 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.722  | 67   | 13589    | 15.645 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.939  | 132  | 14592    | 15.431 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 9.102  | 113  | 16915    | 15.546 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.596  | 128  | 8057     | 14.546 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.319 | 143  | 9927     | 15.056 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.853 | 165  | 18003    | 14.493 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.376 | 131  | 12966    | 8.003  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.402 | 166  | 65830    | 15.046 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.719 | 160  | 74118    | 14.905 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.172 | 143  | 9492     | 12.583 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 16.000 | 176  | 61802    | 15.271 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.100 | 200  | 12610    | 13.871 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.851 | 188  | 99586    | 15.107 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 20.113 | 212  | 120854   | 13.792 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.342 | 264  | 114378   | 14.972 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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