

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092823\  
 Data File : BG059230.D  
 Acq On : 28 Sep 2023 12:14  
 Operator : MA/JU  
 Sample : PB155856BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK856

Quant Time: Sep 29 00:41:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG092723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 28 04:33:25 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.231  | 152  | 39201    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 11.075 | 136  | 208329   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.865 | 164  | 139948   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.615 | 188  | 313356   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.915 | 240  | 289670   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 25.400 | 264  | 334938   | 20.000 | ng/ul | -0.01    |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.531  | 96  | 7670   | 7.385  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.960  | 84  | 113568 | 37.474 | ng/ul | 0.00  |
| 7) Phenol-d5                  | 7.356  | 99  | 139334 | 35.509 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.550  | 67  | 92331  | 36.834 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.750  | 132 | 103919 | 39.397 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8         | 8.925  | 113 | 105343 | 34.730 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5           | 9.430  | 128 | 52123  | 35.456 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4          | 10.147 | 143 | 55398  | 38.386 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.676 | 165 | 98156  | 33.108 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4        | 11.216 | 131 | 161849 | 33.742 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6      | 14.260 | 166 | 358653 | 35.446 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8         | 14.571 | 160 | 423645 | 33.357 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 15.053 | 143 | 53193  | 30.105 | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.852 | 176 | 318566 | 33.080 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.964 | 200 | 48710  | 30.241 | ng/ul | 0.00  |
| 73) Anthracene-d10            | 17.714 | 188 | 488332 | 34.202 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.988 | 212 | 569446 | 35.104 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 25.153 | 264 | 566429 | 35.192 | ng/ul | -0.01 |

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

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

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