

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG031822\  
 Data File : BG052717.D  
 Acq On : 18 Mar 2022 19:39  
 Operator : CG/JU  
 Sample : N1882-22  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DBRZ2

Quant Time: Mar 19 00:01:14 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG031822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 18 00:55:42 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.146  | 152  | 34520    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.965 | 136  | 139330   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.766 | 164  | 100118   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.510 | 188  | 227587   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.798 | 240  | 261556   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.117 | 264  | 257659   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.570  | 96   | 6332     | 6.479  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.981  | 84   | 86230    | 35.344 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.294  | 99   | 116231   | 37.593 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.470  | 67   | 70630    | 37.403 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.682  | 132  | 74829    | 35.644 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.839  | 113  | 90707    | 40.500 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.309  | 128  | 43142    | 43.015 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.037 | 143  | 44860    | 42.406 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.572 | 165  | 88369    | 38.398 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.083 | 131  | 126004   | 41.426 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.167 | 166  | 344204   | 44.751 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.461 | 160  | 405955   | 43.435 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.937 | 143  | 54601    | 42.222 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.759 | 176  | 297661   | 44.075 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.859 | 200  | 47758    | 41.235 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.609 | 188  | 495645   | 46.769 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.889 | 212  | 627227   | 42.867 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.894 | 264  | 611842   | 45.957 | ng/u1 | 0.00     |

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

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

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