

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG032822\  
 Data File : BG052890.D  
 Acq On : 28 Mar 2022 14:09  
 Operator : CG/JU  
 Sample : N2082-14  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 EW606

Quant Time: Mar 28 15:48:56 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG031822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 23 13:16:32 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.140  | 152  | 28271    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.948 | 136  | 125190   | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.761 | 164  | 104656   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.498 | 188  | 254946   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.787 | 240  | 269538   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.100 | 264  | 268854   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.552  | 96   | 3619     | 4.522  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.975  | 84   | 17231    | 8.624  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.277  | 99   | 22884    | 9.038  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.453  | 67   | 55377    | 35.808 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.664  | 132  | 47045    | 27.363 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.828  | 113  | 36524    | 19.912 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.292  | 128  | 33437    | 37.104 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.020 | 143  | 36925    | 38.848 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.561 | 165  | 66444    | 32.132 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.072 | 131  | 82865    | 30.320 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.156 | 166  | 289426   | 35.998 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.449 | 160  | 326624   | 33.432 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.925 | 143  | 11418    | 8.447  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.748 | 176  | 254516   | 36.052 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.853 | 200  | 49028    | 37.789 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.598 | 188  | 461909   | 38.908 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.883 | 212  | 571148   | 37.878 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.871 | 264  | 559430   | 40.271 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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