

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122422\  
 Data File : BG056118.D  
 Acq On : 25 Dec 2022 11:23  
 Operator : CG/JU  
 Sample : N6017-17  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DBYG3

Quant Time: Dec 25 22:37:32 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG122422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 24 22:54:38 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.343  | 152  | 39415    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 11.175 | 136  | 163334   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.953 | 164  | 145087   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.691 | 188  | 350792   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.986 | 240  | 339341   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.459 | 264  | 358198   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.649  | 96   | 4563     | 4.846  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 4.084  | 84   | 64232    | 20.922 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.474  | 99   | 94824    | 26.522 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.650  | 67   | 37928    | 26.915 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.867  | 132  | 61074    | 28.903 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 9.037  | 113  | 73778    | 27.419 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.518  | 128  | 33038    | 27.772 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.247 | 143  | 44869    | 28.868 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.782 | 165  | 91112    | 26.380 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.299 | 131  | 81043    | 20.745 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.342 | 166  | 300211   | 26.735 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.654 | 160  | 338610   | 26.701 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.112 | 143  | 40600    | 24.019 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.934 | 176  | 284688   | 25.685 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.040 | 200  | 59522    | 26.395 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.785 | 188  | 442019   | 27.267 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 20.053 | 212  | 526614   | 28.100 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.218 | 264  | 512571   | 28.153 | ng/u1 | 0.00     |

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

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

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