

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062122\  
 Data File : BG054006.D  
 Acq On : 22 Jun 2022 11:22  
 Operator : CG/JU  
 Sample : N3358-17  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 EW4T0

Quant Time: Jun 23 01:13:05 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 20 15:29:13 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.076  | 152  | 20212    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.891 | 136  | 83528    | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.704 | 164  | 64465    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.442 | 188  | 168934   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.719 | 240  | 193103   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 24.980 | 264  | 192004   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.499  | 96   | 2272     | 5.348  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.917  | 84   | 14609    | 12.644 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.236  | 99   | 16032    | 11.072 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.395  | 67   | 35359    | 40.481 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.612  | 132  | 38305    | 33.173 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.776  | 113  | 28022    | 22.696 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.240  | 128  | 24461    | 39.454 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.962  | 143  | 27342    | 41.506 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.503 | 165  | 52651    | 36.125 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.014 | 131  | 51022    | 27.711 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.105 | 166  | 207862   | 40.987 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 222410   | 40.331 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143  | 7388     | 10.639 | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.691 | 176  | 187145   | 41.082 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.803 | 200  | 38340    | 41.604 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.542 | 188  | 332766   | 43.617 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.827 | 212  | 427672   | 42.477 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.757 | 264  | 425938   | 43.948 | ng/u1 | 0.00     |

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

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

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