

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062122\  
 Data File : BG053977.D  
 Acq On : 21 Jun 2022 14:09  
 Operator : CG/JU  
 Sample : N3285-23  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :

Quant Time: Jun 21 23:07:09 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     | 8.077  | 152  | 20658    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.885 | 136  | 80454    | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.704 | 164  | 63518    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.442 | 188  | 168040   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.719 | 240  | 184989   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 24.980 | 264  | 186464   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.505  | 96   | 1902     | 4.381  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.923  | 84   | 9811     | 8.308  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.230  | 99   | 11589    | 7.831  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.395  | 67   | 27647    | 30.968 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.606  | 132  | 30319    | 25.690 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.776  | 113  | 21279    | 16.863 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.240  | 128  | 19655    | 32.913 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.957  | 143  | 21327    | 33.612 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.503 | 165  | 41655    | 29.673 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.014 | 131  | 37830    | 21.331 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.099 | 166  | 171488   | 34.319 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.393 | 160  | 175856   | 32.365 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.886 | 143  | 4443     | 6.494  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.691 | 176  | 153011   | 34.090 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.803 | 200  | 28881    | 31.506 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.542 | 188  | 274905   | 36.224 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.827 | 212  | 357311   | 37.045 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.757 | 264  | 352093   | 37.408 | ng/u1 | 0.00     |

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

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

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