

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081822\  
 Data File : BG054650.D  
 Acq On : 18 Aug 2022 10:56  
 Operator : CG/JU  
 Sample : PB146656BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK656

Quant Time: Aug 18 11:32:51 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 18 03:22:26 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.159  | 152  | 39463    | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.991 | 136  | 166016   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.798 | 164  | 110701   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.542 | 188  | 248389   | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.837 | 240  | 240552   | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.221 | 264  | 240038   | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.511  | 96   | 5651     | 5.664 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.934  | 84   | 71551    | 24.067 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.295  | 99   | 91553    | 27.287 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.477  | 67   | 60125    | 28.006 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.689  | 132  | 66949    | 28.213 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.852  | 113  | 71457    | 28.193 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.334  | 128  | 31142    | 29.884 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.062 | 143  | 25372    | 27.195 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.597 | 165  | 66100    | 27.025 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.114 | 131  | 110876   | 31.578 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.193 | 166  | 232565   | 30.575 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.492 | 160  | 280488   | 30.798 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.962 | 143  | 30005    | 24.961 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.785 | 176  | 205948   | 29.859 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.885 | 200  | 18516    | 21.890 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.642 | 188  | 336259   | 31.124 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.921 | 212  | 391570   | 31.381 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.986 | 264  | 371918   | 31.551 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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