

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM051822\  
 Data File : BM035148.D  
 Acq On : 18 May 2022 17:34  
 Operator : CG/JU  
 Sample : N2766-08  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EW4F7

Quant Time: May 19 00:50:11 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM051722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 17 16:13:52 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.910  | 152  | 209007   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.710 | 136  | 802446   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.545 | 164  | 475852   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.286 | 188  | 948017   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.468 | 240  | 878924   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.856 | 264  | 918167   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.375  | 96   | 13467    | 2.662  | ng/uL | 0.01     |
| 4) Pyridine-d5                | 3.787  | 84   | 177981   | 13.640 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.075  | 99   | 248647   | 16.142 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.240  | 67   | 163204   | 16.793 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.440  | 132  | 223826   | 16.391 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.616  | 113  | 205727   | 15.899 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.069  | 128  | 112508   | 17.883 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.792  | 143  | 128399   | 18.227 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.328 | 165  | 221323   | 17.664 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.845 | 131  | 287360   | 16.632 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.951 | 166  | 764749   | 22.059 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.233 | 160  | 873827   | 21.089 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.733 | 143  | 99257    | 16.149 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.533 | 176  | 649155   | 22.133 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200  | 82196    | 13.427 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.386 | 188  | 970991   | 22.004 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.680 | 212  | 1096760  | 23.429 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.703 | 264  | 1055827  | 22.690 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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