

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM122022\  
 Data File : BM038146.D  
 Acq On : 20 Dec 2022 22:16  
 Operator : CG/JU  
 Sample : N6008-15  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBXX7

Quant Time: Dec 21 00:17:17 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 21 00:11:56 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.045  | 152  | 189723   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.863 | 136  | 814202   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.674 | 164  | 500250   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.409 | 188  | 1042250  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.574 | 240  | 826232   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.027 | 264  | 750813   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 18508    | 4.429  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 242995   | 19.604 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.204  | 99   | 375522   | 24.358 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67   | 216030   | 23.012 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.575  | 132  | 336245   | 26.587 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.757  | 113  | 294357   | 24.460 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.216  | 128  | 167004   | 25.906 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.939  | 143  | 191960   | 28.415 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 354412   | 27.620 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.998 | 131  | 390475   | 22.045 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.080 | 166  | 1038544  | 29.120 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.368 | 160  | 1208629  | 27.728 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.868 | 143  | 155899   | 25.649 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.657 | 176  | 907901   | 28.573 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.780 | 200  | 129669   | 23.270 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.509 | 188  | 1348350  | 27.719 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 1530065  | 30.828 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.862 | 264  | 1074741  | 28.659 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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