

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM122022\  
 Data File : BM038167.D  
 Acq On : 21 Dec 2022 12:10  
 Operator : CG/JU  
 Sample : N6014-13  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBXS5

Quant Time: Dec 21 23:30:43 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     | 8.040  | 152  | 202408   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.857 | 136  | 858688   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.669 | 164  | 515739   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 1067040  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.574 | 240  | 775186   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.027 | 264  | 758581   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 16628    | 3.730  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 206147   | 15.589 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.198  | 99   | 300979   | 18.299 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67   | 175237   | 17.497 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.569  | 132  | 269032   | 19.939 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.751  | 113  | 244999   | 19.082 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.216  | 128  | 132923   | 19.551 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.939  | 143  | 151739   | 21.297 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.475 | 165  | 279307   | 20.639 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.998 | 131  | 336345   | 18.005 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.080 | 166  | 798069   | 21.705 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.363 | 160  | 923693   | 20.555 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.869 | 143  | 112947   | 18.024 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.657 | 176  | 697910   | 21.305 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.774 | 200  | 62650    | 10.982 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.504 | 188  | 1044482  | 20.974 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 1091414  | 23.438 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.868 | 264  | 795781   | 21.003 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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