

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM122022\  
 Data File : BM038155.D  
 Acq On : 21 Dec 2022 04:19  
 Operator : CG/JU  
 Sample : N6008-05  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBYB2

Quant Time: Dec 21 05:26:31 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  | 172944   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.857 | 136  | 751139   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.669 | 164  | 467068   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 1013494  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.574 | 240  | 858295   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.021 | 264  | 779097   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 16896    | 4.436  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 187720   | 16.614 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.198  | 99   | 258101   | 18.366 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67   | 192353   | 22.478 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.569  | 132  | 255022   | 22.121 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.751  | 113  | 135270   | 12.331 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.216  | 128  | 150343   | 25.280 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.934  | 143  | 167853   | 26.932 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.475 | 165  | 288456   | 24.367 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.998 | 131  | 333780   | 20.426 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.074 | 166  | 965555   | 28.997 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.363 | 160  | 1107141  | 27.204 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.863 | 143  | 134019   | 23.616 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.657 | 176  | 863842   | 29.118 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.774 | 200  | 113443   | 20.936 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.504 | 188  | 1338294  | 28.293 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 1581110  | 30.667 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.868 | 264  | 1153452  | 29.641 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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