

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081222\  
 Data File : BM036247.D  
 Acq On : 13 Aug 2022 02:00  
 Operator : CG/JU  
 Sample : N4165-01  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BGBX7

Quant Time: Aug 13 02:30:02 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 12 11:23:11 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.810  | 152  | 202577   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.616 | 136  | 845007   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 466425   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.198 | 188  | 898519   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 896149   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.721 | 264  | 975532   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.216  | 96   | 22126    | 4.110  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.652  | 84   | 92323    | 6.260  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.004  | 99   | 109792   | 6.463  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.151  | 67   | 324683   | 28.670 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.346  | 132  | 328804   | 23.318 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.545  | 113  | 198135   | 14.397 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.981  | 128  | 207033   | 30.919 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 204536   | 31.353 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165  | 322567   | 27.339 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 452312   | 22.645 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.875 | 166  | 1009540  | 31.775 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 1244650  | 30.704 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.710 | 143  | 29907    | 4.584  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.445 | 176  | 934484   | 32.549 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.574 | 200  | 130060   | 26.929 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.298 | 188  | 1495531  | 36.815 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.592 | 212  | 1676051  | 33.746 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.568 | 264  | 1817568  | 36.699 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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