

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN031522\  
 Data File : BN018971.D  
 Acq On : 16 Mar 2022 05:44  
 Operator : CG/JU  
 Sample : N1818-12  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DBSF6

Quant Time: Mar 16 06:30:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN031522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 15 15:17:15 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.893  | 152  | 218951   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.699 | 136  | 959800   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.528 | 164  | 636081   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.263 | 188  | 1293921  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.445 | 240  | 1042595  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.833 | 264  | 833050   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 37743    | 7.187  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 497567   | 35.637 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.046  | 99   | 710930   | 37.876 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.216  | 67   | 453039   | 38.854 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.416  | 132  | 547802   | 38.278 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.593  | 113  | 581640   | 38.639 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.052  | 128  | 281368   | 38.352 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.775  | 143  | 329647   | 40.452 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.316 | 165  | 558702   | 37.720 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.828 | 131  | 794636   | 38.921 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.934 | 166  | 1862324  | 39.684 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.222 | 160  | 2259992  | 39.021 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.710 | 143  | 296362   | 34.402 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.516 | 176  | 1549302  | 39.460 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.628 | 200  | 272453   | 34.308 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.363 | 188  | 2354895  | 39.271 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.657 | 212  | 2506524  | 38.562 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.680 | 264  | 1730441  | 40.728 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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