

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN090622\  
 Data File : BN021609.D  
 Acq On : 06 Sep 2022 20:15  
 Operator : CG/JU  
 Sample : N4482-12  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 EW349

Quant Time: Sep 07 01:07:17 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN081922.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 07 01:01:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.063  | 152  | 195763   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.899 | 136  | 897964   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.722 | 164  | 568685   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.469 | 188  | 1215411  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.657 | 240  | 1114067  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.174 | 264  | 959250   | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.364  | 96   | 22168    | 4.252 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.799  | 84   | 109631   | 7.715 ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.211  | 99   | 118105   | 6.652 ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 308030   | 28.661 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.587  | 132  | 297522   | 23.032 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.775  | 113  | 215940   | 15.500 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.240  | 128  | 194967   | 31.868 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.969  | 143  | 195786   | 35.829 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.510 | 165  | 331676   | 27.406 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.034 | 131  | 443470   | 21.577 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.128 | 166  | 1302605  | 33.512 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.416 | 160  | 1393417  | 30.750 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.910 | 143  | 46632    | 6.029 ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.710 | 176  | 1117710  | 33.047 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.822 | 200  | 176404   | 36.287 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.569 | 188  | 1850708  | 34.553 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.857 | 212  | 2069670  | 39.064 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.010 | 264  | 1715675  | 37.734 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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