

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN090722\  
 Data File : BN021621.D  
 Acq On : 07 Sep 2022 10:32  
 Operator : CG/JU  
 Sample : PB146666BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK666

Quant Time: Sep 07 12:18:14 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  | 229454   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.893 | 136  | 1042327  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.716 | 164  | 675179   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.469 | 188  | 1403439  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.657 | 240  | 1290165  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.180 | 264  | 1168317  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.369  | 96   | 36531    | 5.978  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.799  | 84   | 463694   | 27.839 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.210  | 99   | 662892   | 31.853 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 416717   | 33.081 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.587  | 132  | 519672   | 34.323 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.775  | 113  | 533985   | 32.702 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.240  | 128  | 273372   | 38.495 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.969  | 143  | 296223   | 46.701 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.504 | 165  | 491690   | 35.001 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.034 | 131  | 921140   | 38.611 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.122 | 166  | 1686011  | 36.535 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.416 | 160  | 1979322  | 36.790 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.910 | 143  | 307787   | 33.516 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.710 | 176  | 1381493  | 34.404 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.822 | 200  | 246895   | 43.982 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.569 | 188  | 2149609  | 34.757 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.863 | 212  | 2436354  | 39.709 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.010 | 264  | 2218528  | 40.063 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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