

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN012323\  
 Data File : BN023748.D  
 Acq On : 23 Jan 2023 12:19  
 Operator : CG/JU  
 Sample : PB150349BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK349

Quant Time: Jan 24 00:11:21 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN010523.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 20 23:02:18 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.887  | 152  | 233238   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.698 | 136  | 943183   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.539 | 164  | 529674   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.286 | 188  | 1052364  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.480 | 240  | 844287   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.909 | 264  | 822294   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 39754    | 7.335  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 542828   | 33.380 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.051  | 99   | 663122   | 32.079 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.216  | 67   | 431297   | 34.011 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.416  | 132  | 539721   | 34.092 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.598  | 113  | 506309   | 30.687 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.057  | 128  | 253048   | 35.579 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.781  | 143  | 269359   | 35.776 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.316 | 165  | 466739   | 32.501 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.839 | 131  | 702564   | 31.102 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.951 | 166  | 1379511  | 35.011 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.233 | 160  | 1547513  | 35.447 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.733 | 143  | 212828   | 26.533 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.533 | 176  | 1152937  | 34.244 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200  | 144961   | 22.834 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.386 | 188  | 1722389  | 36.575 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.680 | 212  | 1853967  | 38.396 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.750 | 264  | 1519208  | 37.359 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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