

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN011323\  
 Data File : BN023648.D  
 Acq On : 12 Jan 2023 17:24  
 Operator : CG/JU  
 Sample : 01050-04MEDL 10X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A4477MEDL

Quant Time: Jan 12 17:55:16 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN010523.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jan 05 03:34:47 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min)     |
|-------------------------------|--------|------|----------|--------|-------|--------------|
| -----                         |        |      |          |        |       |              |
| Internal Standards            |        |      |          |        |       |              |
| 1) 1,4-Dichlorobenzene-d4     | 7.940  | 152  | 186118   | 20.000 | ng/u1 | -0.01        |
| 20) Naphthalene-d8            | 10.757 | 136  | 859636   | 20.000 | ng/u1 | -0.01        |
| 38) Acenaphthene-d10          | 14.598 | 164  | 535601   | 20.000 | ng/u1 | 0.00         |
| 64) Phenanthrene-d10          | 17.345 | 188  | 1167920  | 20.000 | ng/u1 | 0.00         |
| 79) Chrysene-d12              | 21.533 | 240  | 1063422  | 20.000 | ng/u1 | -0.01        |
| 88) Perylene-d12              | 23.963 | 264  | 1086401  | 20.000 | ng/u1 | 0.00         |
| System Monitoring Compounds   |        |      |          |        |       |              |
| 3) 1,4-Dioxane-d8             | 3.293  | 96   | 3272     | 0.757  | ng/uL | 0.00         |
| 4) Pyridine-d5                | 3.740  | 84   | 33077    | 2.549  | ng/u1 | 0.01         |
| 7) Phenol-d5                  | 7.105  | 99   | 53357    | 3.235  | ng/u1 | 0.00         |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 37344    | 3.690  | ng/u1 | 0.00         |
| 11) 2-Chlorophenol-d4         | 7.469  | 132  | 43250    | 3.424  | ng/u1 | 0.00         |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 41835    | 3.178  | ng/u1 | -0.01        |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 19909    | 3.071  | ng/u1 | -0.01        |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 18372    | 2.677  | ng/u1 | 0.00         |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 38903    | 2.972  | ng/u1 | -0.01        |
| 31) 4-Chloroaniline-d4        | 10.904 | 131  | 61806    | 3.002  | ng/u1 | 0.00         |
| 46) Dimethylphthalate-d6      | 14.004 | 166  | 131998   | 3.313  | ng/u1 | -0.01        |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 132460   | 3.001  | ng/u1 | -0.01        |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1 |              |
| 60) Fluorene-d10              | 15.587 | 176  | 119582   | 3.512  | ng/u1 | -0.01        |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/u1 |              |
| 73) Anthracene-d10            | 17.445 | 188  | 181254   | 3.468  | ng/u1 | 0.00         |
| 81) Pyrene-d10                | 19.739 | 212  | 218105   | 3.586  | ng/u1 | 0.00         |
| 92) Benzo(a)pyrene-d12        | 23.798 | 264  | 189001   | 3.518  | ng/u1 | -0.01        |
| Target Compounds              |        |      |          |        |       |              |
| 86) Bis(2-ethylhexyl)phtha... | 21.445 | 149  | 1170630  | 27.104 | ng/u1 | Qvalue<br>98 |
| -----                         |        |      |          |        |       |              |

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

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