

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN102621\  
 Data File : BN017118.D  
 Acq On : 27 Oct 2021 08:48  
 Operator : CG/JU  
 Sample : PB140099BL  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK099

Quant Time: Oct 27 11:01:05 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN102621.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 26 15:29:15 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.534  | 152  | 258733   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.310 | 136  | 1223347  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.186 | 164  | 854644   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.945 | 188  | 1874090  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.157 | 240  | 1984869  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.368 | 264  | 2085360  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.069  | 96   | 31396    | 4.396  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.458  | 84   | 429906   | 22.471 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.722  | 99   | 561101   | 23.338 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.881  | 67   | 354279   | 24.826 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.069  | 132  | 471586   | 24.942 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.252  | 113  | 483700   | 24.645 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.687  | 128  | 239918   | 25.323 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.399  | 143  | 274642   | 26.185 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.934  | 165  | 453295   | 23.590 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.451 | 131  | 740303   | 25.867 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.610 | 166  | 1837887  | 28.726 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.875 | 160  | 2088984  | 25.900 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.410 | 143  | 326069   | 25.829 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.186 | 176  | 1517645  | 27.772 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.316 | 200  | 298411   | 24.590 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.039 | 188  | 2591580  | 28.642 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.351 | 212  | 2983942  | 25.634 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.227 | 264  | 3261571  | 29.428 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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