

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100924\  
 Data File : BP022317.D  
 Acq On : 10 Oct 2024 10:46  
 Operator : RC/JU  
 Sample : PB163718BL  
 Misc :  
 ALS Vial : 37 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK718

Quant Time: Oct 11 02:23:20 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 08 02:14:47 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 117158   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.463 | 136  | 448301   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.316 | 164  | 342334   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.116 | 188  | 848326   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 932797   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.821 | 264  | 1034622  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.222  | 96   | 13269    | 4.401  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 181350   | 21.911 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.893  | 99   | 234104   | 23.738 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.040  | 67   | 138103   | 23.826 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 177860   | 25.417 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.405  | 113  | 187245   | 23.791 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.846  | 128  | 86626    | 25.131 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 106138   | 26.596 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 205613   | 24.242 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.593 | 131  | 236092   | 23.557 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.722 | 166  | 656500   | 24.137 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.004 | 160  | 713431   | 24.696 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.534 | 143  | 99025    | 21.702 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.322 | 176  | 586322   | 24.853 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 104593   | 18.747 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.216 | 188  | 1014831  | 25.430 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 1297690  | 26.307 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.592 | 264  | 1411985  | 25.554 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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