

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111723\  
 Data File : BP018249.D  
 Acq On : 18 Nov 2023 08:10  
 Operator : MA/JU  
 Sample : PB157217BL  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :

Quant Time: Nov 19 00:48:52 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP111023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 15 18:13:37 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 72054    | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.593 | 136  | 273824   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 171590   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.239 | 188  | 394759   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.375 | 240  | 408820   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.851 | 264  | 483624   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.158  | 96   | 12754    | 6.244  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.587  | 84   | 132088   | 24.802 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.946  | 99   | 170873   | 24.564 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 105845   | 25.891 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.299  | 132  | 145058   | 26.333 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.505  | 113  | 137715   | 24.160 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.987  | 128  | 69244    | 28.068 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.699  | 143  | 77545    | 27.745 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.228 | 165  | 126076   | 25.310 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.781 | 131  | 164738   | 24.115 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.881 | 166  | 453139   | 28.621 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.157 | 160  | 473630   | 27.466 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.740 | 143  | 39610    | 16.743 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.463 | 176  | 381690   | 29.201 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.628 | 200  | 65196    | 22.845 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.339 | 188  | 629916   | 29.504 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 799921   | 29.293 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.686 | 264  | 849080   | 30.193 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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