

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121823\  
 Data File : BP018932.D  
 Acq On : 19 Dec 2023 06:03  
 Operator : MA/JU  
 Sample : PB157772TB 10X  
 Misc :  
 ALS Vial : 36 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLEB772

Quant Time: Dec 19 06:39:01 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 19 05:28:03 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.834  | 152  | 211167   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.628 | 136  | 755745   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.463 | 164  | 442611   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.216 | 188  | 1010412  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.310 | 240  | 1056362  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.680 | 264  | 1261874  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.258  | 96   | 36535    | 5.896  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 482775   | 29.086 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.011  | 99   | 582172   | 27.472 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67   | 331338   | 26.381 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.363  | 132  | 469662   | 28.403 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.552  | 113  | 445532   | 25.979 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.993  | 128  | 218272   | 32.450 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.716  | 143  | 236351   | 33.892 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 442264   | 31.015 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 575640   | 30.889 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.881 | 166  | 1201423  | 30.413 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.157 | 160  | 1420067  | 32.646 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 201465   | 31.069 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.457 | 176  | 1070756  | 32.291 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.587 | 200  | 57943    | 9.564  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.316 | 188  | 1796235  | 33.815 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.557 | 212  | 2280482  | 27.733 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.527 | 264  | 2405777  | 32.698 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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