

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122023\  
 Data File : BP018984.D  
 Acq On : 21 Dec 2023 05:03  
 Operator : MA/JU  
 Sample : PB157897BL  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK897

Quant Time: Dec 21 05:55:43 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 21 04:46:12 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.822  | 152  | 156569   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.616 | 136  | 566277   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 351160   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 851387   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.298 | 240  | 1074317  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.663 | 264  | 1307439  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.252  | 96   | 20161    | 4.388  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.670  | 84   | 266324   | 21.640 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.999  | 99   | 312235   | 19.872 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.158  | 67   | 184482   | 19.810 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.352  | 132  | 255702   | 20.856 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.540  | 113  | 243875   | 19.179 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.987  | 128  | 117139   | 23.242 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.705  | 143  | 132560   | 25.369 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.240 | 165  | 241871   | 22.637 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.763 | 131  | 324070   | 23.208 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.875 | 166  | 735835   | 23.478 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 812251   | 23.536 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 122869   | 23.883 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.451 | 176  | 646960   | 24.592 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200  | 119485   | 23.406 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 1093518  | 24.431 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 1477343  | 17.666 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.510 | 264  | 1841681  | 24.159 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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