

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100124\  
 Data File : BP022130.D  
 Acq On : 01 Oct 2024 11:06  
 Operator : RC/JU  
 Sample : PB163565BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK565

Quant Time: Oct 02 07:10:47 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 24 15:23:13 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.699  | 152  | 128962   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.451 | 136  | 461536   | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10          | 14.304 | 164  | 348762   | 20.000 | ng/u1 | -0.02    |
| 64) Phenanthrene-d10          | 17.098 | 188  | 817572   | 20.000 | ng/u1 | -0.02    |
| 79) Chrysene-d12              | 21.533 | 240  | 983899   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.815 | 264  | 1142448  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.216  | 96   | 15538    | 4.719  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.628  | 84   | 209176   | 22.227 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.881  | 99   | 243942   | 23.124 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.034  | 67   | 152875   | 23.534 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.240  | 132  | 193681   | 25.376 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.393  | 113  | 196546   | 23.513 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.834  | 128  | 91496    | 26.487 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.551  | 143  | 108831   | 27.786 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 206246   | 24.247 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.587 | 131  | 248256   | 24.461 | ng/u1 | -0.02    |
| 46) Dimethylphthalate-d6      | 13.704 | 166  | 683233   | 25.343 | ng/u1 | -0.02    |
| 49) Acenaphthylene-d8         | 13.992 | 160  | 713496   | 24.997 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.510 | 143  | 98651    | 21.393 | ng/u1 | -0.02    |
| 60) Fluorene-d10              | 15.304 | 176  | 579810   | 24.313 | ng/u1 | -0.02    |
| 65) 4,6-Dinitro-2-methylph... | 15.428 | 200  | 113562   | 22.373 | ng/u1 | -0.02    |
| 73) Anthracene-d10            | 17.198 | 188  | 976474   | 25.584 | ng/u1 | -0.02    |
| 81) Pyrene-d10                | 19.580 | 212  | 1279351  | 25.098 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 1485289  | 24.718 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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