

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100424\  
 Data File : BP022241.D  
 Acq On : 04 Oct 2024 23:54  
 Operator : RC/JU  
 Sample : PB163635BL  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK635

Quant Time: Oct 05 05:15:36 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 04 05:25:28 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.692  | 152  | 97831    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.445 | 136  | 366440   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.304 | 164  | 284245   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.098 | 188  | 680565   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.527 | 240  | 699504   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.803 | 264  | 772422   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.216  | 96   | 13534    | 5.419  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.628  | 84   | 191688   | 26.850 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.875  | 99   | 218374   | 27.288 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.028  | 67   | 131354   | 26.656 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.234  | 132  | 171512   | 29.622 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.387  | 113  | 177823   | 28.042 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.828  | 128  | 81920    | 29.869 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.539  | 143  | 96451    | 31.016 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 188585   | 27.925 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.586 | 131  | 228702   | 28.382 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.704 | 166  | 636970   | 28.989 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.992 | 160  | 682176   | 29.324 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.527 | 143  | 75129    | 19.990 | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.310 | 176  | 554897   | 28.550 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.433 | 200  | 75351    | 17.833 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.204 | 188  | 925725   | 29.137 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.580 | 212  | 1153407  | 31.827 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 1205516  | 29.673 | ng/u1 | 0.01     |

Target Compounds Qvalue

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

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