

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092824\  
 Data File : BM047688.D  
 Acq On : 28 Sep 2024 12:44  
 Operator : RC/JU  
 Sample : PB163765BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK765

Quant Time: Sep 29 00:48:40 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Sep 28 02:30:36 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.816  | 152  | 193863   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 760972   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.445 | 164  | 443105   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.186 | 188  | 858061   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.409 | 240  | 756789   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.415 | 264  | 881547   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.257  | 96   | 34453    | 5.743  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.669  | 84   | 494296   | 28.141 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.975  | 99   | 511221   | 27.887 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.145  | 67   | 377596   | 29.028 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.345  | 132  | 394758   | 29.598 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.516  | 113  | 379487   | 27.843 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.969  | 128  | 187219   | 30.900 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.686  | 143  | 181315   | 31.228 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.228 | 165  | 336297   | 28.994 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.739 | 131  | 528390   | 29.467 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 1033403  | 32.168 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.139 | 160  | 1173184  | 30.801 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.633 | 143  | 160483   | 25.091 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.439 | 176  | 878995   | 31.589 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.551 | 200  | 119747   | 23.848 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.280 | 188  | 1325533  | 31.417 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.568 | 212  | 1524539  | 35.494 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.209 | 264  | 1432772  | 31.325 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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