

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM111124\  
 Data File : BM048603.D  
 Acq On : 12 Nov 2024 03:35  
 Operator : RC/JU  
 Sample : PB164732BL  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK732

Quant Time: Nov 12 04:12:05 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Nov 07 22:50:20 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.657  | 152  | 170143   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.445 | 136  | 691828   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.310 | 164  | 425432   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.051 | 188  | 886612   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.280 | 240  | 854379   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.185 | 264  | 989871   | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.122  | 96   | 28987    | 6.792  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.540  | 84   | 395799   | 33.651 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.840  | 99   | 453560   | 35.018 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.992  | 67   | 264145   | 33.954 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.192  | 132  | 386085   | 35.986 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.375  | 113  | 359525   | 35.303 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.816  | 128  | 180995   | 36.096 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.533  | 143  | 198053   | 37.134 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.075 | 165  | 329145   | 33.220 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.592 | 131  | 492250   | 34.978 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.721 | 166  | 991833   | 36.522 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.998 | 160  | 1168711  | 36.501 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.539 | 143  | 139045   | 27.566 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.304 | 176  | 877063   | 36.998 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.439 | 200  | 119930   | 26.889 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.151 | 188  | 1406999  | 37.501 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.450 | 212  | 1660127  | 38.718 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.991 | 264  | 1748835  | 37.813 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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