

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110824\  
 Data File : BM048571.D  
 Acq On : 09 Nov 2024 18:31  
 Operator : RC/JU  
 Sample : PB164674BL  
 Misc :  
 ALS Vial : 53 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK674

Quant Time: Nov 09 21:51:47 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     | 7.663  | 152  | 230778   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.445 | 136  | 949258   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.310 | 164  | 581156   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.057 | 188  | 1163425  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.280 | 240  | 1042778  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.191 | 264  | 1225199  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.128  | 96   | 31349    | 5.415  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.546  | 84   | 427379   | 26.789 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.846  | 99   | 511506   | 29.116 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.998  | 67   | 309925   | 29.371 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.198  | 132  | 441312   | 30.326 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.375  | 113  | 407531   | 29.502 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.816  | 128  | 209530   | 30.454 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.539  | 143  | 231847   | 31.681 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 382839   | 28.161 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.592 | 131  | 527453   | 27.315 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.727 | 166  | 1122186  | 30.249 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.004 | 160  | 1327446  | 30.349 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.533 | 143  | 167937   | 24.373 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.304 | 176  | 979525   | 30.248 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.439 | 200  | 138246   | 23.621 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.157 | 188  | 1519646  | 30.866 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.451 | 212  | 1727458  | 33.009 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.991 | 264  | 1825599  | 31.891 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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