

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP053123\  
 Data File : BP015353.D  
 Acq On : 31 May 2023 11:30  
 Operator : CG/JU  
 Sample : PB153064BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK064

Quant Time: May 31 14:16:16 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 30 15:23:22 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.105  | 152  | 135654   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.934 | 136  | 585407   | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.751 | 164  | 427792   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.504 | 188  | 1062458  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.586 | 240  | 1162190  | 20.000 | ng/u1 | #-0.01   |
| 88) Perylene-d12              | 24.145 | 264  | 1267756  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.417  | 96   | 20645    | 5.913  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.852  | 84   | 293323   | 29.639 | ng/u1 | -0.01    |
| 7) Phenol-d5                  | 7.258  | 99   | 350101   | 27.797 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.428  | 67   | 217035   | 29.836 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.634  | 132  | 292166   | 31.549 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.822  | 113  | 303587   | 30.249 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.287  | 128  | 153299   | 33.325 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.010 | 143  | 181325   | 35.426 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.557 | 165  | 294279   | 31.862 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.075 | 131  | 526425   | 38.522 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.151 | 166  | 1200945  | 37.151 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.446 | 160  | 1259881  | 34.077 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.951 | 143  | 196246   | 31.264 | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.740 | 176  | 992284   | 36.378 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.863 | 200  | 226682   | 33.855 | ng/u1 | -0.02    |
| 73) Anthracene-d10            | 17.604 | 188  | 1713145  | 35.213 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.828 | 212  | 2165202  | 32.122 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.980 | 264  | 2318525  | 36.841 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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