

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082823\  
 Data File : BP016869.D  
 Acq On : 30 Aug 2023 13:13  
 Operator : MA/JU  
 Sample : PB155122BL  
 Misc :  
 ALS Vial : 89 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK122

Quant Time: Aug 31 00:49:49 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 30 02:25:00 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 349131   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.852 | 136  | 1532429  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.675 | 164  | 951037   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.445 | 188  | 2068779  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 1528265  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.110 | 264  | 1396685  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.376  | 96   | 61562    | 7.288  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.811  | 84   | 846952   | 32.809 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.164  | 99   | 1097853  | 35.363 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.358  | 67   | 697298   | 35.241 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.540  | 132  | 828076   | 35.476 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.728  | 113  | 867973   | 34.249 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.222  | 128  | 417068   | 35.996 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.934  | 143  | 449490   | 36.804 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.458 | 165  | 806712   | 35.430 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.010 | 131  | 1150100  | 32.938 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.087 | 166  | 2576979  | 35.930 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 2935900  | 36.211 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.881 | 143  | 375440   | 25.326 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.669 | 176  | 2251558  | 37.279 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.793 | 200  | 276329   | 22.429 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.545 | 188  | 3536724  | 38.569 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.781 | 212  | 3844057  | 45.798 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.939 | 264  | 2688592  | 38.430 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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