

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041923\  
 Data File : BP014723.D  
 Acq On : 19 Apr 2023 11:07  
 Operator : CG/JU  
 Sample : PB152181BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK181

Quant Time: Apr 20 01:45:23 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP041823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 19 01:37:13 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.152  | 152  | 181000   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.981 | 136  | 741994   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.787 | 164  | 470204   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.534 | 188  | 1061158  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.610 | 240  | 996745   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.192 | 264  | 1137795  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.487  | 96   | 27152    | 5.962  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1 |          |
| 7) Phenol-d5                  | 7.287  | 99   | 403429   | 25.779 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.469  | 67   | 284028   | 29.419 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.675  | 132  | 338908   | 29.160 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.852  | 113  | 302146   | 24.766 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.322  | 128  | 143664   | 30.899 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.052 | 143  | 125759   | 29.947 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.587 | 165  | 307926   | 27.006 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.116 | 131  | 1266     | 0.073  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.187 | 166  | 1187160  | 31.886 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.481 | 160  | 1234248  | 28.606 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.945 | 143  | 133061   | 22.469 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.775 | 176  | 995342   | 31.156 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.875 | 200  | 80283    | 18.695 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.634 | 188  | 1587813  | 30.973 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.851 | 212  | 1883675  | 32.990 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.021 | 264  | 1757195  | 29.798 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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