

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP033023\  
 Data File : BP014410.D  
 Acq On : 30 Mar 2023 19:48  
 Operator : CG/JU  
 Sample : PB151800BL  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK800

Quant Time: Mar 30 22:43:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 30 12:32:14 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.016  | 152  | 145862   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.840 | 136  | 647988   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.669 | 164  | 445979   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.434 | 188  | 989758   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.522 | 240  | 765864   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.039 | 264  | 726053   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 22292    | 5.927  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.817  | 84   | 310704   | 32.234 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.175  | 99   | 400529   | 29.774 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67   | 273873   | 31.732 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 300670   | 30.924 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.728  | 113  | 329641   | 30.379 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.193  | 128  | 157502   | 30.109 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.922  | 143  | 183548   | 30.642 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.463 | 165  | 302684   | 27.747 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.987 | 131  | 422186   | 29.381 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.075 | 166  | 1184155  | 32.747 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.363 | 160  | 1237335  | 30.838 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.887 | 143  | 133095   | 24.363 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.663 | 176  | 965476   | 32.017 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 196829   | 27.027 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.534 | 188  | 1531897  | 31.981 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 1683752  | 32.731 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.880 | 264  | 1280519  | 33.196 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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