

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100122\  
 Data File : BP011853.D  
 Acq On : 01 Oct 2022 10:10  
 Operator : CG/JU  
 Sample : PB147968BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK968

Quant Time: Oct 03 00:57:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP093022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Oct 01 00:53:59 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.893  | 152  | 215867   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.705 | 136  | 990802   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.540 | 164  | 676107   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.298 | 188  | 1489005  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 1526937  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.821 | 264  | 1694883  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 30542    | 6.240  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 371426   | 25.907 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.052  | 99   | 558947   | 29.993 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.222  | 67   | 338416   | 32.761 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.422  | 132  | 472762   | 32.136 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.599  | 113  | 487096   | 31.619 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.058  | 128  | 248111   | 32.876 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.781  | 143  | 273579   | 33.883 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.322 | 165  | 466147   | 31.212 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.840 | 131  | 731217   | 32.116 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.951 | 166  | 1903094  | 39.362 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.234 | 160  | 2012584  | 36.514 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.734 | 143  | 330661   | 33.026 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.534 | 176  | 1578090  | 37.584 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200  | 293519   | 32.568 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.398 | 188  | 2633854  | 38.832 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.628 | 212  | 3090808  | 39.969 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.663 | 264  | 3420987  | 40.321 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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