

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP031023\  
 Data File : BP014057.D  
 Acq On : 11 Mar 2023 07:30  
 Operator : CG/JU  
 Sample : PB151314BL  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK314

Quant Time: Mar 13 00:49:56 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030723.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 10 22:47:08 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.069  | 152  | 149501   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.893 | 136  | 626250   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.710 | 164  | 445580   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.469 | 188  | 1121447  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.545 | 240  | 1153564  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.080 | 264  | 1107852  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 24861    | 6.283  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.864  | 84   | 343100   | 31.317 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.222  | 99   | 414615   | 31.427 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.393  | 67   | 249013   | 32.364 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.599  | 132  | 311602   | 31.000 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.781  | 113  | 319764   | 29.672 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.246  | 128  | 155339   | 31.088 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.969  | 143  | 186275   | 31.142 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.510 | 165  | 318448   | 28.295 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.034 | 131  | 413926   | 27.791 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.116 | 166  | 1138083  | 30.096 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.404 | 160  | 1213579  | 30.295 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.904 | 143  | 173433   | 26.026 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.704 | 176  | 955414   | 29.812 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.822 | 200  | 201025   | 23.717 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.563 | 188  | 1655284  | 30.405 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.786 | 212  | 2074544  | 32.607 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.916 | 264  | 1852729  | 31.440 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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