

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP022223\  
 Data File : BP013869.D  
 Acq On : 22 Feb 2023 10:07  
 Operator : CG/JU  
 Sample : PB150979BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK979

Quant Time: Feb 23 03:14:24 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP020723.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Feb 22 01:41:33 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.104  | 152  | 144689   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.934 | 136  | 602757   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.739 | 164  | 425142   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.498 | 188  | 1007047  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.568 | 240  | 985413   | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.121 | 264  | 1031752  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.452  | 96   | 22126    | 6.338 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.887  | 84   | 299733   | 28.961 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.252  | 99   | 375469   | 29.125 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.422  | 67   | 221420   | 29.310 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.628  | 132  | 298505   | 31.455 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.810  | 113  | 316065   | 29.996 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.281  | 128  | 155069   | 36.976 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.004 | 143  | 180248   | 37.224 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.546 | 165  | 326689   | 33.107 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.063 | 131  | 448609   | 32.073 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.145 | 166  | 1204669  | 35.976 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.439 | 160  | 1267086  | 34.957 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.928 | 143  | 176532   | 29.592 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.733 | 176  | 1004436  | 35.217 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.845 | 200  | 196250   | 31.765 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.598 | 188  | 1645205  | 35.728 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.816 | 212  | 1922388  | 34.401 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.957 | 264  | 1824746  | 35.297 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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