

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011023\  
 Data File : BP013421.D  
 Acq On : 10 Jan 2023 09:56  
 Operator : CG/JU  
 Sample : PB150112BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK112

Quant Time: Jan 10 23:19:29 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP010523.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 06 01:03:32 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 488270   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.728 | 136  | 1955555  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.563 | 164  | 1194652  | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.322 | 188  | 2622942  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.410 | 240  | 2504519  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.874 | 264  | 2766511  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.311  | 96   | 59526    | 4.739 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.734  | 84   | 784919   | 22.403 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.075  | 99   | 917023   | 22.634 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.246  | 67   | 644091   | 25.199 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.446  | 132  | 758774   | 25.113 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.628  | 113  | 717099   | 23.231 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.087  | 128  | 339116   | 24.355 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.810  | 143  | 385397   | 25.285 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.346 | 165  | 717759   | 23.629 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.869 | 131  | 977508   | 22.843 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.969 | 166  | 2311113  | 26.553 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.257 | 160  | 2415464  | 25.071 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 278212   | 17.463 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.557 | 176  | 1964472  | 25.552 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.681 | 200  | 321668   | 19.341 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.422 | 188  | 3140728  | 27.090 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.657 | 212  | 3743675  | 27.473 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.710 | 264  | 3714900  | 27.337 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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