

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112423\  
 Data File : BP018332.D  
 Acq On : 24 Nov 2023 17:52  
 Operator : MA/JU  
 Sample : PB157095BL  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK095

Quant Time: Nov 24 22:28:54 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 24 22:02:00 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.869  | 152  | 297773   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.669 | 136  | 1192249  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.498 | 164  | 681966   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.251 | 188  | 1215517  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.351 | 240  | 896193   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.780 | 264  | 1100570  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 57072    | 7.179  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 670600   | 30.693 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 920183   | 33.813 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.205  | 67   | 558579   | 34.325 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 765426   | 33.500 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 732313   | 33.345 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.028  | 128  | 354349   | 36.592 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.752  | 143  | 319901   | 35.770 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165  | 708281   | 32.826 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.804 | 131  | 830061   | 28.523 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.916 | 166  | 2064148  | 34.016 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.192 | 160  | 2406984  | 34.812 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 222155   | 26.519 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.492 | 176  | 1706328  | 33.619 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.604 | 200  | 135461   | 26.179 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.351 | 188  | 2344710  | 36.090 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 2337157  | 32.767 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.621 | 264  | 2361845  | 37.016 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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