

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061623\  
 Data File : BP015587.D  
 Acq On : 16 Jun 2023 18:37  
 Operator : MA/JU  
 Sample : PB153428BS  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK428

Quant Time: Jun 16 22:08:25 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 15 05:36:46 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.087  | 152  | 101312   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.910 | 136  | 451809   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.727 | 164  | 333203   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.486 | 188  | 812047   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.568 | 240  | 789174   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.109 | 264  | 841099   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.404  | 96   | 19349    | 7.072  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.846  | 84   | 238021   | 33.493 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.240  | 99   | 288349   | 30.894 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.404  | 67   | 198193   | 35.556 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.610  | 132  | 237919   | 35.159 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.798  | 113  | 248717   | 32.469 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.263  | 128  | 122785   | 34.615 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.986  | 143  | 141380   | 34.902 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.533 | 165  | 238522   | 32.279 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.051 | 131  | 371840   | 35.057 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.133 | 166  | 1026665  | 39.074 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.422 | 160  | 1053046  | 36.652 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.945 | 143  | 113486   | 24.119 | ng/ul | 0.01     |
| 60) Fluorene-d10              | 15.721 | 176  | 833777   | 37.630 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.845 | 200  | 152353   | 29.615 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.580 | 188  | 1390165  | 37.626 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.804 | 212  | 1733193  | 41.034 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.945 | 264  | 1634335  | 38.726 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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