

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110123\  
 Data File : BP017883.D  
 Acq On : 02 Nov 2023 13:17  
 Operator : MA/JU  
 Sample : PB156594BL  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK594

Quant Time: Nov 02 23:53:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101823.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 19 01:32:15 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.875  | 152  | 145632   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.704 | 136  | 548483   | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10          | 14.575 | 164  | 316557   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.369 | 188  | 635488   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.516 | 240  | 529198   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.068 | 264  | 542447   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 22854    | 5.838  | ng/uL | 0.01     |
| 4) Pyridine-d5                | 3.699  | 84   | 270101   | 26.353 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.040  | 99   | 396177   | 32.691 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.222  | 67   | 245906   | 32.856 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 332224   | 33.611 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.599  | 113  | 317086   | 33.319 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.081  | 128  | 157846   | 35.516 | ng/u1 | -0.02    |
| 24) 2-Nitrophenol-d4          | 9.799  | 143  | 178964   | 36.444 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.328 | 165  | 308517   | 32.028 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.881 | 131  | 397295   | 31.515 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.992 | 166  | 894908   | 33.676 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.269 | 160  | 1012584  | 34.168 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 106087   | 28.017 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 737070   | 32.087 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.728 | 200  | 117367   | 28.039 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.469 | 188  | 1144716  | 35.249 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.733 | 212  | 1280391  | 38.591 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.898 | 264  | 1069123  | 35.387 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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