

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112423\  
 Data File : BP018371.D  
 Acq On : 25 Nov 2023 17:01  
 Operator : MA/JU  
 Sample : PB157191BL  
 Misc :  
 ALS Vial : 54 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK191

Quant Time: Nov 27 01:50:39 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  | 530457   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.663 | 136  | 2060875  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.498 | 164  | 1147656  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 2011480  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.351 | 240  | 1832785  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.792 | 264  | 2216543  | 20.000 | ng/u1 | 0.01     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 66937    | 4.726  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 893453   | 22.956 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 1133016  | 23.371 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.199  | 67   | 655799   | 22.622 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.393  | 132  | 966055   | 23.734 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 896169   | 22.907 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.028  | 128  | 453705   | 27.104 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.746  | 143  | 458804   | 29.678 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.281 | 165  | 894527   | 23.984 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.804 | 131  | 1245806  | 24.766 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.916 | 166  | 2422055  | 23.718 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.192 | 160  | 2856719  | 24.551 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.686 | 143  | 274206   | 19.451 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.492 | 176  | 1965194  | 23.008 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200  | 180761   | 21.110 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 2682177  | 24.948 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 2857621  | 19.591 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.633 | 264  | 3238196  | 25.199 | ng/u1 | 0.01     |

Target Compounds Qvalue

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

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