

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112723\  
 Data File : BP018442.D  
 Acq On : 28 Nov 2023 15:51  
 Operator : MA/JU  
 Sample : 05391-01RX  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4A06RX

Quant Time: Nov 29 01:22:08 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 29 01:22:40 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.857  | 152  | 492710   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.651 | 136  | 1906776  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.492 | 164  | 1052175  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 1994954  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.416 | 240  | 1901342  | 20.000 | ng/u1 | -0.02    |
| 88) Perylene-d12              | 24.051 | 264  | 2260402  | 20.000 | ng/u1 | -0.09    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 52460    | 3.988  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 457804   | 12.664 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.016  | 99   | 168004   | 3.731  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.187  | 67   | 714614   | 26.540 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.381  | 132  | 134384   | 3.555  | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.563  | 113  | 178426   | 4.910  | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.016  | 128  | 493247   | 31.848 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.734  | 143  | 49333    | 3.449  | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165  | 115835   | 3.357  | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.793 | 131  | 1296960  | 27.867 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.904 | 166  | 1943834  | 20.762 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 3141637  | 29.450 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 24034    | 1.860  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.487 | 176  | 2256573  | 28.817 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200  | 14703    | 1.731  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 3192824  | 29.943 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 3714907  | 24.549 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.874 | 264  | 4019852  | 30.675 | ng/u1 | -0.08    |

Target Compounds Qvalue

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

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