

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112723\  
 Data File : BP018409.D  
 Acq On : 27 Nov 2023 20:03  
 Operator : MA/JU  
 Sample : 05332-13  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Nov 27 21:58:34 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 27 21:50:01 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.858  | 152  | 444161   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.657 | 136  | 1778035  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.493 | 164  | 1014629  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 1778183  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 1466687  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.892 | 264  | 1726272  | 20.000 | ng/ul | 0.02     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 40632    | 3.426  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 248773   | 7.634  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.022  | 99   | 215683   | 5.313  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67   | 656747   | 27.057 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 696830   | 20.446 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.569  | 113  | 420887   | 12.848 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.022  | 128  | 442843   | 30.664 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.740  | 143  | 396762   | 29.748 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165  | 779590   | 24.227 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.799 | 131  | 936294   | 21.574 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 2573610  | 28.506 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 2909978  | 28.288 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 33324    | 2.674  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.487 | 176  | 2129106  | 28.195 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200  | 201906   | 26.673 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 3259161  | 34.291 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.592 | 212  | 3413748  | 29.245 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.727 | 264  | 3707967  | 37.049 | ng/ul | 0.03     |

Target Compounds Qvalue

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

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