

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100423\  
 Data File : BP017558.D  
 Acq On : 05 Oct 2023 10:50  
 Operator : MA/JU  
 Sample : 04679-22  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :

Quant Time: Oct 06 01:58:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 04 16:33:47 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 102515   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.781 | 136  | 411221   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.640 | 164  | 259220   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.428 | 188  | 587689   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.563 | 240  | 611393   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.145 | 264  | 694467   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.293  | 96   | 9037     | 3.524  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.734  | 84   | 36042    | 5.370  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 46863    | 5.433  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.287  | 67   | 149281   | 28.364 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 144914   | 21.338 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.669  | 113  | 89494    | 12.961 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.152  | 128  | 91671    | 28.977 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.869  | 143  | 95867    | 26.867 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.399 | 165  | 163923   | 24.780 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.951 | 131  | 235170   | 24.956 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 599789   | 29.663 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.334 | 160  | 644317   | 28.898 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143  | 14698    | 4.151  | ng/ul | 0.02     |
| 60) Fluorene-d10              | 15.639 | 176  | 523482   | 30.523 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.781 | 200  | 75352    | 20.567 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.528 | 188  | 894733   | 32.890 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.780 | 212  | 1150766  | 33.873 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.974 | 264  | 1175327  | 33.536 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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