

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111723\  
 Data File : BP018250.D  
 Acq On : 18 Nov 2023 08:44  
 Operator : MA/JU  
 Sample : 05370-02  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :

Quant Time: Nov 19 00:49:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP111023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 15 18:13:37 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.781  | 152  | 66474    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.593 | 136  | 265888   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 178752   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.239 | 188  | 415090   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 481451   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.851 | 264  | 522794   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.158  | 96   | 9045     | 4.800  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.599  | 84   | 42891    | 8.730  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 48931    | 7.625  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 92306    | 24.475 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.299  | 132  | 103869   | 20.438 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 81672    | 15.531 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.987  | 128  | 60677    | 25.329 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.699  | 143  | 66809    | 24.617 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.228 | 165  | 106268   | 21.970 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.781 | 131  | 135304   | 20.398 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.881 | 166  | 465221   | 28.207 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.157 | 160  | 468377   | 26.073 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.845 | 143  | 104      | 0.042  | ng/u1 | 0.11     |
| 60) Fluorene-d10              | 15.463 | 176  | 396295   | 29.103 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 263      | 0.088  | ng/u1 | 0.12     |
| 73) Anthracene-d10            | 17.345 | 188  | 712675   | 31.745 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 952569   | 29.621 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.692 | 264  | 1012232  | 33.298 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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