

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP091023\  
 Data File : BP017104.D  
 Acq On : 11 Sep 2023 14:48  
 Operator : MA/JU  
 Sample : 04268-06  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 H0GH4

Quant Time: Sep 12 02:47:23 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP091023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 11 06:13:29 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.957  | 152  | 254362   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.787 | 136  | 1128864  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.616 | 164  | 687285   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.386 | 188  | 1494952  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 1240059  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.015 | 264  | 1280427  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.316  | 96   | 47058    | 7.061  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.752  | 84   | 191704   | 9.539  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.110  | 99   | 156117   | 6.752  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.299  | 67   | 498525   | 32.148 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 412918   | 24.709 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.669  | 113  | 287696   | 15.729 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.157  | 128  | 259170   | 32.248 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.869  | 143  | 263894   | 31.814 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.393 | 165  | 445618   | 28.332 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.945 | 131  | 590300   | 23.405 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.034 | 166  | 1810749  | 37.027 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.316 | 160  | 1875458  | 33.679 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.833 | 143  | 50559    | 5.011  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.616 | 176  | 1531054  | 36.893 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.739 | 200  | 228173   | 27.046 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.486 | 188  | 2476556  | 38.649 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.727 | 212  | 2867439  | 43.909 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.851 | 264  | 2494459  | 40.581 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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