

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052323\  
 Data File : BP015257.D  
 Acq On : 24 May 2023 00:55  
 Operator : CG/JU  
 Sample : 02861-12  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 H0FL5

Quant Time: May 24 01:42:00 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 23 01:25:41 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.116  | 152  | 228303   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.946 | 136  | 1029834  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.757 | 164  | 648663   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.510 | 188  | 1376269  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.580 | 240  | 983465   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.139 | 264  | 1036583  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.428  | 96   | 24858    | 4.230  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.875  | 84   | 111590   | 6.700  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.269  | 99   | 143735   | 6.781  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.434  | 67   | 423674   | 34.607 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.640  | 132  | 420558   | 26.984 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.828  | 113  | 274366   | 16.243 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.293  | 128  | 284682   | 35.179 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.022 | 143  | 310627   | 34.498 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.557 | 165  | 513189   | 31.585 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.081 | 131  | 390379   | 16.239 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.157 | 166  | 1879883  | 38.352 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.451 | 160  | 2019984  | 36.032 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.951 | 143  | 45834    | 4.816  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.745 | 176  | 1540918  | 37.256 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.863 | 200  | 263554   | 30.387 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.610 | 188  | 2320384  | 36.820 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.828 | 212  | 2566605  | 44.997 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.974 | 264  | 2115206  | 41.106 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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