

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041223\  
 Data File : BP014707.D  
 Acq On : 13 Apr 2023 11:36  
 Operator : CG/JU  
 Sample : 02255-01 10X  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AN5

Quant Time: Apr 14 00:55:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 13 02:01:50 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.975  | 152  | 99309    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.793 | 136  | 402549   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.622 | 164  | 245329   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.392 | 188  | 467194   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.480 | 240  | 436767   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.974 | 264  | 522584   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.364  | 96   | 77       | 0.030  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.799  | 84   | 337      | 0.051  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.152  | 99   | 11       | 0.001  | ng/u1 | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.293  | 67   | 64       | 0.011  | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.511  | 132  | 13       | 0.002  | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.711  | 113  | 11       | 0.001  | ng/u1 | 0.02     |
| 21) Nitrobenzene-d5           | 9.134  | 128  | 11       | 0.003  | ng/u1 | -0.02    |
| 24) 2-Nitrophenol-d4          | 9.875  | 143  | 12       | 0.003  | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.369 | 165  | 6        | 0.001  | ng/u1 | -0.05    |
| 31) 4-Chloroaniline-d4        | 10.946 | 131  | 12       | 0.001  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.022 | 166  | 6        | 0.000  | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.328 | 160  | 191      | 0.009  | ng/u1 | 0.01     |
| 54) 4-Nitrophenol-d4          | 14.851 | 143  | 13       | 0.004  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.616 | 176  | 73       | 0.004  | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 2        | 0.001  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.392 | 188  | 467194   | 20.663 | ng/u1 | -0.09    |
| 81) Pyrene-d10                | 19.745 | 212  | 1268     | 0.043  | ng/u1 | 0.02     |
| 92) Benzo(a)pyrene-d12        | 23.810 | 264  | 463      | 0.017  | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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