

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP113022\  
 Data File : BP012900.D  
 Acq On : 01 Dec 2022 03:29  
 Operator : CG/JU  
 Sample : N5770-02  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH1F8

Quant Time: Dec 01 04:00:49 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 29 22:53:34 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.010  | 152  | 670697   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.822 | 136  | 3078379  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 2036485  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.398 | 188  | 3882932  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.480 | 240  | 2484002  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 2502454  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 75586    | 4.829  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 297873   | 6.359  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.157  | 99   | 490951   | 8.908  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.328  | 67   | 1045249  | 30.778 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 1111205  | 26.460 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 843639   | 18.999 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.175  | 128  | 667507   | 35.428 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.898  | 143  | 717086   | 35.844 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.440 | 165  | 1363352  | 29.206 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 425190   | 6.652  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 4451489  | 31.757 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.339 | 160  | 4992712  | 30.526 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 111896   | 4.477  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.639 | 176  | 3826779  | 31.173 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200  | 501868   | 27.043 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.498 | 188  | 5424373  | 31.736 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.727 | 212  | 5181302  | 35.898 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.833 | 264  | 3971564  | 32.311 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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