

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121922\  
 Data File : BP013219.D  
 Acq On : 21 Dec 2022 21:28  
 Operator : CG/JU  
 Sample : N6032-06  
 Misc :  
 ALS Vial : 101 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0L89

Quant Time: Dec 21 23:17:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 19 21:43:11 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.952  | 152  | 558997   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 2362941  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.593 | 164  | 1491822  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 2865545  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.433 | 240  | 1989669  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 2421559  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 53802    | 4.115  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 80366    | 2.164  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 7.111  | 99   | 283805   | 6.233  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 772473   | 28.124 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 649479   | 18.120 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 450900   | 12.086 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 507169   | 29.213 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.846  | 143  | 502593   | 25.714 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 842568   | 22.194 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.899 | 131  | 527184   | 9.836  | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.999 | 166  | 3465926  | 30.230 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 3886356  | 30.848 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.793 | 143  | 61866    | 3.000  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.587 | 176  | 3043975  | 31.382 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 354277   | 19.212 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 4373348  | 33.829 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.681 | 212  | 4164883  | 36.468 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264  | 4114494  | 34.088 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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