

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110222\  
 Data File : BP012514.D  
 Acq On : 04 Nov 2022 05:52  
 Operator : CG/JU  
 Sample : N5320-03  
 Misc :  
 ALS Vial : 71 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHAB5

Quant Time: Nov 04 06:33:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 04 02:53:36 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 342673   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.575 | 136  | 1639003  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.428 | 164  | 1127470  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 2517944  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.286 | 240  | 2237319  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.669 | 264  | 1995855  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.228  | 96   | 32664    | 3.838  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.658  | 84   | 235473   | 9.632  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 196246   | 6.385  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.116  | 67   | 508785   | 27.730 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.311  | 132  | 530228   | 23.624 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 381902   | 15.711 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.946  | 128  | 351148   | 33.038 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.663  | 143  | 380279   | 34.471 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.205 | 165  | 685336   | 28.666 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.728 | 131  | 904860   | 25.419 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.845 | 166  | 2405703  | 31.169 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.122 | 160  | 2570213  | 29.140 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 63170    | 4.325  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.428 | 176  | 2039317  | 30.861 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 339430   | 25.892 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.292 | 188  | 3477333  | 32.051 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.533 | 212  | 3995215  | 35.567 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.516 | 264  | 3224488  | 33.065 | ng/u1 | 0.00     |

| Target Compounds           |        |     |       |             | Qvalue |
|----------------------------|--------|-----|-------|-------------|--------|
| 67) N-Nitrosodiphenylamine | 15.698 | 169 | 89661 | 1.264 ng/u1 | 97     |

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

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