

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112522\  
 Data File : BP012742.D  
 Acq On : 25 Nov 2022 23:58  
 Operator : CG/JU  
 Sample : N5592-19  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 H0AC4

Quant Time: Nov 26 07:16:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 25 15:59:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 493944   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.840 | 136  | 2181556  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.657 | 164  | 1440643  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.416 | 188  | 2954477  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.498 | 240  | 1986929  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.015 | 264  | 1540646  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 54895    | 4.762  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 345804   | 10.025 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.163  | 99   | 297909   | 7.340  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.346  | 67   | 923358   | 36.919 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 887531   | 28.697 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.722  | 113  | 587372   | 17.961 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.187  | 128  | 543059   | 40.671 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.910  | 143  | 568344   | 40.087 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.451 | 165  | 1083358  | 32.749 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 1365261  | 30.138 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 3894628  | 39.275 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.351 | 160  | 4434940  | 38.331 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 72044    | 4.075  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.651 | 176  | 3392869  | 39.069 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.757 | 200  | 411884   | 29.168 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.510 | 188  | 5435859  | 41.797 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.739 | 212  | 5712260  | 49.477 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.851 | 264  | 3257017  | 43.039 | ng/u1 | 0.00     |

|                  |       |    |       |       |        |    |
|------------------|-------|----|-------|-------|--------|----|
| Target Compounds |       |    |       |       | Qvalue |    |
| 2) 1,4-Dioxane   | 3.434 | 88 | 48057 | 3.882 | ng/uL  | 98 |

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

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