

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120423\  
 Data File : BP018577.D  
 Acq On : 04 Dec 2023 22:50  
 Operator : MA/JU  
 Sample : 05439-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AA1

Quant Time: Dec 04 23:19:24 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 04 21:39:40 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.852  | 152  | 390995   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.652 | 136  | 1700899  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.481 | 164  | 1151938  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.228 | 188  | 2490946  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.316 | 240  | 1881895  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 1920308  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 42959    | 3.744  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 265470   | 8.638  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.022  | 99   | 238684   | 6.083  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.187  | 67   | 731675   | 31.462 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 669364   | 21.862 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.563  | 113  | 451809   | 14.228 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.016  | 128  | 442484   | 29.229 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.734  | 143  | 433585   | 27.625 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.269 | 165  | 815268   | 25.403 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.793 | 131  | 982601   | 23.428 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.898 | 166  | 3205308  | 31.177 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.175 | 160  | 3421518  | 30.223 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.681 | 143  | 81410    | 4.824  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.475 | 176  | 2796172  | 32.401 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 407637   | 27.293 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.328 | 188  | 4538744  | 34.659 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.563 | 212  | 5132816  | 35.039 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 4009943  | 35.813 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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