

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
 Data File : BP013319.D  
 Acq On : 25 Dec 2022 02:31  
 Operator : CG/JU  
 Sample : N6018-10  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBYJ0

Quant Time: Dec 25 04:26:08 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 24 00:19:17 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.940  | 152  | 561970   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.751 | 136  | 2099166  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 1097680  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.339 | 188  | 2281890  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2511163  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 2907017  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.328  | 96   | 48398    | 3.682  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.758  | 84   | 643903   | 17.246 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 873471   | 19.082 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.263  | 67   | 605599   | 21.932 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.469  | 132  | 759807   | 21.086 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 585212   | 15.603 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.104  | 128  | 364918   | 23.660 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.828  | 143  | 405172   | 23.335 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 717817   | 21.284 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.893 | 131  | 806979   | 16.949 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.987 | 166  | 1903023  | 22.558 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.275 | 160  | 2182824  | 23.547 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.787 | 143  | 244959   | 16.142 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.575 | 176  | 1660453  | 23.265 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.698 | 200  | 148791   | 10.133 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.439 | 188  | 2381319  | 23.132 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.675 | 212  | 3210272  | 22.272 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264  | 3430955  | 23.678 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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