

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121323\  
 Data File : BP018855.D  
 Acq On : 14 Dec 2023 22:06  
 Operator : MA/JU  
 Sample : 05627-21DL 5X  
 Misc :  
 ALS Vial : 49 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCLD3DL

Quant Time: Dec 14 22:47:59 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 13 23:34:47 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.828  | 152  | 276204   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.622 | 136  | 988872   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.463 | 164  | 533595   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.210 | 188  | 1185500  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 1373173  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.657 | 264  | 1625790  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.258  | 96   | 8933     | 1.102  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 33071    | 1.523  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.999  | 99   | 38275    | 1.381  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67   | 103378   | 6.293  | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.364  | 132  | 101627   | 4.699  | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.546  | 113  | 50135    | 2.235  | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.993  | 128  | 51812    | 5.887  | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.710  | 143  | 46504    | 5.096  | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.246 | 165  | 96915    | 5.194  | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 125516   | 5.147  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.875 | 166  | 334366   | 7.021  | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 343138   | 6.543  | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.681 | 143  | 7632     | 0.976  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.451 | 176  | 265739   | 6.648  | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.569 | 200  | 25784    | 3.627  | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.304 | 188  | 497651   | 7.985  | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 725476   | 6.787  | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.498 | 264  | 784619   | 8.277  | ng/u1 | -0.02    |
| Target Compounds              |        |      |          |        |       |          |
| 78) Di-n-butylphthalate       | 18.175 | 149  | 1695074  | 22.256 | ng/u1 | 99       |
| -----                         |        |      |          |        |       |          |

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

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