

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032824\  
 Data File : BP020018.D  
 Acq On : 28 Mar 2024 20:13  
 Operator : MA/JU  
 Sample : P1856-01 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0LZ0

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 03/29/2024  
 Supervised By :Jagrut Upadhyay 03/29/2024

Quant Time: Mar 28 23:20:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 28 14:08:48 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.605  | 152  | 110659   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.363 | 136  | 453091   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.234 | 164  | 305327   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.075 | 188  | 650352   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.533 | 240  | 586441   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.816 | 264  | 635290   | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.240  | 96   | 7839     | 3.104  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.664  | 84   | 17995    | 2.360  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 6.799  | 99   | 67824    | 7.267  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.969  | 67   | 234582   | 38.580 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.152  | 132  | 182700   | 27.259 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.305  | 113  | 127266   | 17.625 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.758  | 128  | 128707   | 37.061 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.463  | 143  | 135361   | 34.836 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.999  | 165  | 218000   | 30.720 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.557 | 131  | 5387m    | 0.582  | ng/u1 | 0.04     |
| 46) Dimethylphthalate-d6      | 13.669 | 166  | 888451   | 41.502 | ng/u1 | 0.02     |
| 49) Acenaphthylene-d8         | 13.928 | 160  | 942457   | 37.688 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.534 | 143  | 22628m   | 6.724  | ng/u1 | 0.05     |
| 60) Fluorene-d10              | 15.263 | 176  | 746890   | 40.957 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.404 | 200  | 136255   | 35.001 | ng/u1 | -0.02    |
| 73) Anthracene-d10            | 17.181 | 188  | 1203119  | 41.792 | ng/u1 | 0.01     |
| 81) Pyrene-d10                | 19.569 | 212  | 1490453  | 43.318 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 1387356  | 45.456 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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