

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032824\  
 Data File : BP020021.D  
 Acq On : 28 Mar 2024 22:23  
 Operator : MA/JU  
 Sample : P1856-04 10X  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0LZ1

Manual Integrations  
 APPROVED

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

Quant Time: Mar 28 23:27:33 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     | 7.611  | 152  | 89980    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.369 | 136  | 364288   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.257 | 164  | 252722   | 20.000 | ng/u1 | 0.01     |
| 64) Phenanthrene-d10          | 17.063 | 188  | 553378   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.545 | 240  | 510609   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.827 | 264  | 537034   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.240  | 96   | 6665     | 3.245  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.664  | 84   | 14693m   | 2.369  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 6.811  | 99   | 60153    | 7.927  | ng/u1 | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.969  | 67   | 227848   | 46.084 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.152  | 132  | 171027   | 31.381 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.310  | 113  | 116046   | 19.765 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.758  | 128  | 125296   | 44.874 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.469  | 143  | 126110   | 40.367 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.999  | 165  | 200110   | 35.073 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.540 | 131  | 31520m   | 4.238  | ng/u1 | 0.02     |
| 46) Dimethylphthalate-d6      | 13.657 | 166  | 849743   | 47.957 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.940 | 160  | 935013   | 45.173 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.534 | 143  | 20060m   | 7.201  | ng/u1 | 0.05     |
| 60) Fluorene-d10              | 15.269 | 176  | 726935   | 48.161 | ng/u1 | 0.01     |
| 65) 4,6-Dinitro-2-methylph... | 15.416 | 200  | 126177   | 38.092 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.175 | 188  | 1215829  | 49.634 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.569 | 212  | 1445732  | 48.258 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.604 | 264  | 1342282  | 52.026 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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