

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010083.D  
 Acq On : 25 Apr 2022 15:25  
 Operator : CG/JU  
 Sample : N2471-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BFFG2

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022

Quant Time: Apr 26 01:02:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 170819   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.740 | 136  | 723336   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.569 | 164  | 501934   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.322 | 188  | 1082903  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.404 | 240  | 914926   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.857 | 264  | 771631   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 17962    | 4.585  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 41280    | 3.746  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.093  | 99   | 87117    | 5.607  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.258  | 67   | 287626   | 30.851 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 261780   | 22.750 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.640  | 113  | 170694   | 13.260 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.093  | 128  | 177339   | 30.825 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 184701   | 29.397 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.357 | 165  | 299835   | 24.513 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.875 | 131  | 365893m  | 20.863 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.975 | 166  | 1301269  | 33.328 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.257 | 160  | 1400626  | 29.640 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 24990m   | 3.477  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.563 | 176  | 1098366  | 32.702 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.675 | 200  | 187580   | 27.395 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.422 | 188  | 1818120  | 34.704 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.651 | 212  | 2065816  | 39.703 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.698 | 264  | 1457009  | 36.017 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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