

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080922\  
 Data File : BP011363.D  
 Acq On : 10 Aug 2022 11:09  
 Operator : CG/JU  
 Sample : N4069-14  
 Misc :  
 ALS Vial : 45 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GBJL8

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 08/11/2022  
 Supervised By :Sohil Jodhani 08/12/2022

Quant Time: Aug 10 23:56:32 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 08 18:47:19 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.587  | 152  | 105824   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.375 | 136  | 453513   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.257 | 164  | 258510   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.028 | 188  | 508829   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.157 | 240  | 465010   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.468 | 264  | 461073   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.087  | 96   | 16239    | 4.704  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.505  | 84   | 55867    | 6.416  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.769  | 99   | 67705    | 7.229  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.934  | 67   | 205850   | 31.000 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.116  | 132  | 186886   | 25.433 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.305  | 113  | 121081   | 16.466 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.752  | 128  | 117378   | 38.030 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.469  | 143  | 112808   | 39.920 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.004 | 165  | 170574   | 29.890 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.528 | 131  | 246135   | 24.671 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.681 | 166  | 685410   | 38.213 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.945 | 160  | 791546   | 36.660 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.492 | 143  | 13637m   | 4.163  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.263 | 176  | 570558   | 37.195 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.392 | 200  | 86504    | 36.601 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.128 | 188  | 982131   | 42.943 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.392 | 212  | 1101911  | 43.011 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.321 | 264  | 1033396  | 44.443 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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