

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP050522\  
 Data File : BP010254.D  
 Acq On : 05 May 2022 19:06  
 Operator : CG/JU  
 Sample : N2678-03  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH0B8

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/10/2022  
 Supervised By :mohammad ahmed 05/10/2022

Quant Time: May 06 02:07:52 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri May 06 01:05:07 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 184301   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.716 | 136  | 814426   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.546 | 164  | 527829   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.298 | 188  | 1026004  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.380 | 240  | 649502   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.822 | 264  | 612196   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.358  | 96   | 20077m   | 4.750  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 88945    | 7.481  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.081  | 99   | 103137   | 6.152  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.240  | 67   | 318436   | 31.657 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.446  | 132  | 294382   | 23.711 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.622  | 113  | 201175   | 14.484 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.075  | 128  | 201638   | 31.129 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.799  | 143  | 207576   | 29.342 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.340 | 165  | 338372   | 24.570 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.852 | 131  | 323207   | 16.368 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.957 | 166  | 1341894  | 32.682 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.240 | 160  | 1492396m | 30.032 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.751 | 143  | 28737m   | 3.802  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.540 | 176  | 1143238  | 32.369 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.657 | 200  | 139074   | 21.437 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.398 | 188  | 1763655  | 35.531 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.628 | 212  | 1710012  | 46.295 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.669 | 264  | 1190602  | 37.096 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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