

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP050422\  
 Data File : BP010227.D  
 Acq On : 04 May 2022 17:26  
 Operator : CG/JU  
 Sample : N2668-04  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXL67

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/05/2022  
 Supervised By :mohammad ahmed 05/09/2022

Quant Time: May 05 02:40:11 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 03 00:58:32 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 164715   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.722 | 136  | 726666   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.551 | 164  | 490737   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.304 | 188  | 999510   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.386 | 240  | 681685   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.827 | 264  | 608276   | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.358  | 96   | 17615m   | 4.663  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.787  | 84   | 48870    | 4.599  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.081  | 99   | 65662    | 4.383  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.246  | 67   | 278931m  | 31.027 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.452  | 132  | 235334   | 21.209 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.622  | 113  | 143225   | 11.538 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.075  | 128  | 171832   | 29.731 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.799  | 143  | 174237   | 27.604 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.340 | 165  | 284144m  | 23.124 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.857 | 131  | 332835m  | 18.891 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.957 | 166  | 1164849  | 30.514 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.239 | 160  | 1311845  | 28.394 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.751 | 143  | 17440m   | 2.482  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.545 | 176  | 1009608  | 30.746 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.663 | 200  | 143201   | 22.658 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.398 | 188  | 1637466  | 33.863 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.633 | 212  | 1669687  | 43.069 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.668 | 264  | 1158499  | 36.329 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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