

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110722\  
 Data File : BP012534.D  
 Acq On : 07 Nov 2022 16:46  
 Operator : CG/JU  
 Sample : N5353-19  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBXB2

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 11/08/2022  
 Supervised By :mohammad ahmed 11/14/2022

Quant Time: Nov 07 21:41:40 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 04 02:53:36 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 278956   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.575 | 136  | 1366827  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.428 | 164  | 941189   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 2157410  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.286 | 240  | 1955536  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.668 | 264  | 1788723  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.228  | 96   | 25875    | 3.735  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.663  | 84   | 101617   | 5.106  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.957  | 99   | 122064   | 4.879  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.116  | 67   | 370613   | 24.813 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.310  | 132  | 354687   | 19.412 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.499  | 113  | 227138   | 11.478 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.946  | 128  | 256836   | 28.977 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.669  | 143  | 270284   | 29.379 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.204 | 165  | 465004   | 23.323 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.734 | 131  | 518374   | 17.462 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.845 | 166  | 1845590  | 28.645 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.122 | 160  | 1957235  | 26.582 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.686 | 143  | 30577m   | 2.508  | ng/u1 | 0.03     |
| 60) Fluorene-d10              | 15.428 | 176  | 1636171  | 29.661 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 250980   | 22.344 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.292 | 188  | 2578372  | 27.737 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.533 | 212  | 3112943  | 31.706 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.515 | 264  | 2600625  | 29.756 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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