

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112522\  
 Data File : BP012798.D  
 Acq On : 27 Nov 2022 09:29  
 Operator : CG/JU  
 Sample : PB149206BL  
 Misc :  
 ALS Vial : 73 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK206

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 11/28/2022  
 Supervised By :mohammad ahmed 11/30/2022

Quant Time: Nov 28 00:02:16 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Nov 27 22:55:08 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.016  | 152  | 500721   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.834 | 136  | 2270895  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.651 | 164  | 1573404  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 3508153  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 3439566  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.010 | 264  | 3405327  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 61477    | 5.261  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1 |          |
| 7) Phenol-d5                  | 7.164  | 99   | 1092950  | 26.563 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67   | 717623   | 28.304 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.540  | 132  | 884419   | 28.209 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 900427   | 27.161 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.181  | 128  | 430724   | 30.989 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.905  | 143  | 479793   | 32.510 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.446 | 165  | 920628   | 26.735 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.981 | 131  | 3185m    | 0.068  | ng/u1 | 0.01     |
| 46) Dimethylphthalate-d6      | 14.057 | 166  | 3355883  | 30.987 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.345 | 160  | 3575009  | 28.291 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 489538   | 25.354 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.645 | 176  | 2875670  | 30.319 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 386188   | 23.032 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.510 | 188  | 4682274  | 30.321 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.733 | 212  | 5468313  | 27.361 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.845 | 264  | 5048286  | 30.181 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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