

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
 Data File : BP013327.D  
 Acq On : 25 Dec 2022 07:35  
 Operator : CG/JU  
 Sample : PB149784BL  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK784

Quant Time: Dec 25 21:39:32 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 24 00:19:17 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.940  | 152  | 766437   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.752 | 136  | 3024865  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 1576176  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 2798458  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2846627  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.916 | 264  | 3362441  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.328  | 96   | 106802   | 5.958  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.752  | 84   | 1559563  | 30.627 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 2055140  | 32.920 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.263  | 67   | 1272497  | 33.790 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.469  | 132  | 1652782  | 33.631 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 1604928  | 31.376 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 814371   | 36.643 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 917727   | 36.679 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 1575282  | 32.414 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.893 | 131  | 2070761  | 30.182 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.993 | 166  | 4072219  | 33.617 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.275 | 160  | 4515782  | 33.926 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.787 | 143  | 526617   | 24.168 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 3118669  | 30.431 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.698 | 200  | 367343   | 20.398 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.439 | 188  | 4435929  | 35.136 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.675 | 212  | 5426664  | 33.212 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264  | 5996343  | 35.778 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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