

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120822\  
 Data File : BP013011.D  
 Acq On : 09 Dec 2022 09:15  
 Operator : CG/JU  
 Sample : PB149468BL  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK468

Quant Time: Dec 09 21:50:41 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 08 00:35:07 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.975  | 152  | 684806   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.799 | 136  | 2968691  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.622 | 164  | 1963574  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.381 | 188  | 4390768  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.469 | 240  | 3909140  | 20.000 | ng/u1 | 0.01     |
| 88) Perylene-d12              | 23.963 | 264  | 3717729  | 20.000 | ng/u1 | 0.01     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.358  | 96   | 78881    | 4.925  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 1118343  | 24.580 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.128  | 99   | 1418843  | 25.437 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.305  | 67   | 897482   | 26.673 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.505  | 132  | 1181432  | 26.906 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.681  | 113  | 1154471  | 25.260 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.146  | 128  | 591521   | 27.119 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.875  | 143  | 676461   | 27.548 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.410 | 165  | 1208859  | 25.345 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.934 | 131  | 1688715  | 25.079 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.028 | 166  | 3997965  | 26.492 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.316 | 160  | 4386012  | 26.450 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.810 | 143  | 620183   | 22.847 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.616 | 176  | 3386574  | 26.526 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.734 | 200  | 598163   | 21.170 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.481 | 188  | 5482654  | 27.678 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.710 | 212  | 6368446  | 28.382 | ng/u1 | 0.01     |
| 92) Benzo(a)pyrene-d12        | 23.804 | 264  | 5184874  | 27.980 | ng/u1 | 0.02     |

Target Compounds Qvalue

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

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