

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112522\  
 Data File : BP012762.D  
 Acq On : 26 Nov 2022 11:54  
 Operator : CG/JU  
 Sample : PB149078BL  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK078

Quant Time: Nov 27 21:50:30 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 25 15:59:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 484988   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.834 | 136  | 2101785  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.657 | 164  | 1391418  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 2819534  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 1739238  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.010 | 264  | 1452283  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 79638    | 7.036  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.823  | 84   | 1109214  | 32.749 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.164  | 99   | 1339164  | 33.603 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67   | 857699   | 34.927 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 1077944  | 35.497 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 1076980  | 33.541 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.187  | 128  | 497418   | 38.667 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.911  | 143  | 552213   | 40.428 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.446 | 165  | 1093035  | 34.296 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 1553985  | 35.606 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 3577882  | 37.358 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.351 | 160  | 4039846  | 36.151 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.834 | 143  | 450690   | 26.394 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.646 | 176  | 3081180  | 36.735 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.757 | 200  | 350824   | 26.033 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.510 | 188  | 4695558  | 37.833 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.734 | 212  | 4748189  | 46.984 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.845 | 264  | 2774143  | 38.889 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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