

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110822\  
 Data File : BP012581.D  
 Acq On : 09 Nov 2022 09:39  
 Operator : CG/JU  
 Sample : PB148835BL  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK835

Quant Time: Nov 09 10:13:17 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 08 22:40:23 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.763  | 152  | 212456   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.563 | 136  | 1079690  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.422 | 164  | 781737   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.180 | 188  | 1801172  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.280 | 240  | 1617031  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.662 | 264  | 1482983  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.222  | 96   | 32190    | 6.101  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.652  | 84   | 205292   | 13.544 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.957  | 99   | 666101   | 34.956 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.105  | 67   | 368264   | 32.373 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.305  | 132  | 490261   | 35.231 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 551866   | 36.617 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.934  | 128  | 258418   | 36.909 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 287057   | 39.501 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.204 | 165  | 534271   | 33.924 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.722 | 131  | 504698   | 21.522 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.839 | 166  | 1791116  | 33.470 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.110 | 160  | 2016113  | 32.967 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 277822   | 27.431 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.422 | 176  | 1595004  | 34.812 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 275833   | 29.414 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.286 | 188  | 2695553  | 34.733 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.527 | 212  | 3080945  | 37.949 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.504 | 264  | 2589379  | 35.735 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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