

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121323\  
 Data File : BP018865.D  
 Acq On : 15 Dec 2023 04:18  
 Operator : MA/JU  
 Sample : PB157726BL  
 Misc :  
 ALS Vial : 60 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK726

Quant Time: Dec 15 04:54:39 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 13 23:34:47 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.828  | 152  | 264594   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.622 | 136  | 971373   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 579493   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.210 | 188  | 1231000  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 1007371  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.656 | 264  | 1222235  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.257  | 96   | 40300    | 5.190  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.675  | 84   | 525971   | 25.290 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.004  | 99   | 635975   | 23.951 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67   | 386230   | 24.542 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.363  | 132  | 517740   | 24.988 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.540  | 113  | 491219   | 22.859 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.992  | 128  | 234656   | 27.142 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.710  | 143  | 248469   | 27.720 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.245 | 165  | 438578   | 23.929 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 639011   | 26.678 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.880 | 166  | 1451427  | 28.063 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 1562419  | 27.434 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 209498   | 24.677 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.451 | 176  | 1220841  | 28.121 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.574 | 200  | 169851   | 23.012 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.310 | 188  | 1960501  | 30.294 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 1841699  | 23.486 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.503 | 264  | 2217517  | 31.116 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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