

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121123\  
 Data File : BP018773.D  
 Acq On : 12 Dec 2023 18:35  
 Operator : MA/JU  
 Sample : PB157629BL  
 Misc :  
 ALS Vial : 52 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK629

Quant Time: Dec 12 22:18:09 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 11 22:02:13 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.840  | 152  | 158856   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.634 | 136  | 599143   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.469 | 164  | 392398   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.216 | 188  | 964092   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.304 | 240  | 1080161  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.668 | 264  | 1247999  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 28108    | 6.030  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 358572   | 28.717 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.010  | 99   | 417164   | 26.168 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.175  | 67   | 248709   | 26.323 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.369  | 132  | 341793   | 27.477 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.552  | 113  | 331132   | 25.666 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.999  | 128  | 158603   | 29.742 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.716  | 143  | 176003   | 31.835 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 339950   | 30.071 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.775 | 131  | 441643   | 29.893 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.887 | 166  | 1065355  | 30.420 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.163 | 160  | 1129930  | 29.300 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 172120   | 29.941 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.457 | 176  | 912445   | 31.038 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.581 | 200  | 152628   | 26.403 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.316 | 188  | 1534321  | 30.272 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 1995012  | 23.727 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.515 | 264  | 2219709  | 30.504 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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