

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100423\  
 Data File : BP017602.D  
 Acq On : 06 Oct 2023 15:21  
 Operator : MA/JU  
 Sample : PB156034BL  
 Misc :  
 ALS Vial : 82 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK034

Quant Time: Oct 06 23:32:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 06 05:55:05 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 78438    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.769 | 136  | 322539   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.628 | 164  | 204615   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.416 | 188  | 479884   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.551 | 240  | 474935   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.127 | 264  | 539384   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 15528    | 7.915  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.722  | 84   | 234281   | 45.620 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.087  | 99   | 277625   | 42.067 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 172386   | 42.807 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.452  | 132  | 223352   | 42.982 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 214700   | 40.638 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.140  | 128  | 106080   | 42.750 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.857  | 143  | 115429   | 41.244 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.387 | 165  | 210486   | 40.568 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.940 | 131  | 300868   | 40.706 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 705789   | 44.221 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.322 | 160  | 742894   | 42.211 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.869 | 143  | 99893    | 35.740 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.628 | 176  | 600936   | 44.389 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.775 | 200  | 86955    | 29.065 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.516 | 188  | 996034   | 44.839 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.769 | 212  | 1220419  | 46.244 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.957 | 264  | 1241292  | 45.601 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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