

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062422\  
 Data File : BP010783.D  
 Acq On : 24 Jun 2022 09:22  
 Operator : CG/JU  
 Sample : PB145677BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK677

Quant Time: Jun 24 22:44:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 24 22:42:26 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 370690   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.522 | 136  | 1579122  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.381 | 164  | 931362   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.145 | 188  | 1846630  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.257 | 240  | 1545826  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.627 | 264  | 1351943  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.211  | 96   | 64636    | 6.596  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.605  | 84   | 856902   | 32.850 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.893  | 99   | 990665   | 33.760 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.063  | 67   | 658292   | 34.303 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.252  | 132  | 824091   | 33.850 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.434  | 113  | 800542   | 34.511 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.881  | 128  | 367665   | 36.574 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.604  | 143  | 311504   | 37.455 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.140 | 165  | 668811   | 32.546 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.657 | 131  | 1165519  | 34.746 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.798 | 166  | 2274056  | 35.644 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.069 | 160  | 2736523  | 33.608 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.581 | 143  | 292996   | 29.403 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.381 | 176  | 1905305  | 34.800 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.492 | 200  | 155028   | 25.339 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.245 | 188  | 2914485  | 35.474 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.492 | 212  | 3128495  | 37.148 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.474 | 264  | 2430277  | 36.377 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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