

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112522\  
 Data File : BP012731.D  
 Acq On : 25 Nov 2022 17:42  
 Operator : CG/JU  
 Sample : PB149034BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK034

Quant Time: Nov 26 06:52:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 25 15:59:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 477623   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.840 | 136  | 2138680  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.657 | 164  | 1468345  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.416 | 188  | 3153139  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.504 | 240  | 2862673  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.021 | 264  | 2505489  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 67765    | 6.080  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.817  | 84   | 940043   | 28.182 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.169  | 99   | 1171591  | 29.851 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.346  | 67   | 768618   | 31.782 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.552  | 132  | 944211   | 31.573 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.722  | 113  | 969688   | 30.665 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.193  | 128  | 448702   | 34.278 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.916  | 143  | 504417   | 36.292 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.452 | 165  | 991129   | 30.562 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 1429219  | 32.183 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 3515831  | 34.786 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.351 | 160  | 3866432  | 32.787 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.834 | 143  | 494776   | 27.458 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.651 | 176  | 2994038  | 33.826 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.763 | 200  | 397727   | 26.391 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.516 | 188  | 4882082  | 35.174 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.739 | 212  | 5517057  | 33.168 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.857 | 264  | 4326529  | 35.156 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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