

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062323\  
 Data File : BP015687.D  
 Acq On : 24 Jun 2023 00:31  
 Operator : MA/JU  
 Sample : PB153415BL  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK415

Quant Time: Jun 24 02:36:30 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 22 22:01:49 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.081  | 152  | 118940   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.905 | 136  | 430547   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.728 | 164  | 226780   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.492 | 188  | 463448   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.580 | 240  | 497961   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.139 | 264  | 589501   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 23516    | 7.321  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.840  | 84   | 282945   | 33.913 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.246  | 99   | 351758   | 32.102 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.399  | 67   | 235947   | 36.055 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.605  | 132  | 296878   | 37.369 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.805  | 113  | 280196   | 31.157 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.263  | 128  | 135465   | 40.076 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.987  | 143  | 147673   | 38.255 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.540 | 165  | 232274   | 32.986 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.057 | 131  | 327966   | 32.448 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.134 | 166  | 706294   | 39.495 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.422 | 160  | 791244   | 40.464 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.975 | 143  | 69455    | 21.688 | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.728 | 176  | 556409   | 36.897 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.875 | 200  | 79368    | 27.033 | ng/u1 | 0.01     |
| 73) Anthracene-d10            | 17.592 | 188  | 853189   | 40.462 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.816 | 212  | 1101646  | 41.335 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.974 | 264  | 1232788  | 41.678 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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