

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070722\  
 Data File : BP010934.D  
 Acq On : 07 Jul 2022 11:42  
 Operator : CG/JU  
 Sample : PB146060BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK060

Quant Time: Jul 08 00:58:07 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 06 23:13:53 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 336669   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.498 | 136  | 1300517  | 20.000 | ng/ul | # 0.00   |
| 38) Acenaphthene-d10          | 14.363 | 164  | 670467   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.128 | 188  | 1306184  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12              | 21.239 | 240  | 1294968  | 20.000 | ng/ul | # 0.00   |
| 88) Perylene-d12              | 23.598 | 264  | 1324432  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.199  | 96   | 67834    | 7.622  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.611  | 84   | 852013   | 35.964 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.881  | 99   | 942744   | 35.374 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.046  | 67   | 617130   | 35.407 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.240  | 132  | 796622   | 36.028 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.422  | 113  | 728160   | 34.563 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 340602   | 41.140 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 316218   | 46.167 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.122 | 165  | 592635   | 35.017 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.646 | 131  | 979199   | 35.445 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.781 | 166  | 1666200  | 36.279 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.051 | 160  | 2017691  | 34.422 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 241671   | 33.689 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.363 | 176  | 1459071  | 37.020 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.481 | 200  | 141359   | 32.665 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.228 | 188  | 2218328  | 38.172 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.474 | 212  | 2485200  | 35.226 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.445 | 264  | 2515891  | 38.441 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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