

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100322\  
 Data File : BP011869.D  
 Acq On : 03 Oct 2022 13:24  
 Operator : CG/JU  
 Sample : PB147928BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK928

Quant Time: Oct 03 23:30:38 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP093022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Oct 01 00:53:59 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.893  | 152  | 279525   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.698 | 136  | 1163951  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.539 | 164  | 709373   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.292 | 188  | 1518867  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.374 | 240  | 1589593  | 20.000 ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.810 | 264  | 1700707  | 20.000 ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 36990    | 5.836 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 453598   | 24.433 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.058  | 99   | 664700   | 27.545 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.222  | 67   | 371955   | 27.808 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.422  | 132  | 553116   | 29.036 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.604  | 113  | 527196   | 26.428 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.057  | 128  | 266329   | 30.040 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.781  | 143  | 280886   | 29.613 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.322 | 165  | 494623   | 28.192 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.840 | 131  | 761617   | 28.475 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.951 | 166  | 1551072  | 30.576 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.228 | 160  | 1783343  | 30.838 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.739 | 143  | 279545   | 26.612 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.534 | 176  | 1384377  | 31.424 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200  | 221204   | 24.062 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.392 | 188  | 2240939  | 32.389 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.627 | 212  | 2714331  | 33.717 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.657 | 264  | 2838521  | 33.341 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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