

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
 Data File : BP022604.D  
 Acq On : 25 Oct 2024 09:53  
 Operator : RC/JU  
 Sample : PB164352BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK352

Quant Time: Oct 25 23:43:13 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 22 05:59:59 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.652  | 152  | 108323   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.399 | 136  | 388352   | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10      | 14.263 | 164  | 306053   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10      | 17.069 | 188  | 746273   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12          | 21.504 | 240  | 862777   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.763 | 264  | 958798   | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.187  | 96  | 10879   | 3.903  | ng/uL | -0.01 |
| 4) Pyridine-d5                | 3.599  | 84  | 159319  | 20.819 | ng/u1 | -0.01 |
| 7) Phenol-d5                  | 6.846  | 99  | 196725  | 21.575 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 6.993  | 67  | 117638  | 21.951 | ng/u1 | -0.01 |
| 11) 2-Chlorophenol-d4         | 7.193  | 132 | 149173  | 23.056 | ng/u1 | -0.01 |
| 15) 4-Methylphenol-d8         | 8.358  | 113 | 156112  | 21.453 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.793  | 128 | 75902   | 25.419 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.505  | 143 | 87645   | 25.352 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.040 | 165 | 166937  | 22.720 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.546 | 131 | 201431  | 23.201 | ng/u1 | -0.01 |
| 46) Dimethylphthalate-d6      | 13.669 | 166 | 590498  | 24.284 | ng/u1 | -0.01 |
| 49) Acenaphthylene-d8         | 13.951 | 160 | 610728  | 23.647 | ng/u1 | -0.01 |
| 54) 4-Nitrophenol-d4          | 14.498 | 143 | 84117   | 20.620 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.269 | 176 | 493502  | 23.398 | ng/u1 | -0.02 |
| 65) 4,6-Dinitro-2-methylph... | 15.422 | 200 | 105369  | 21.468 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.175 | 188 | 875736  | 24.945 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.557 | 212 | 1122630 | 24.605 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.533 | 264 | 1299243 | 25.373 | ng/u1 | 0.00  |

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

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

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