

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103024\  
 Data File : BP022741.D  
 Acq On : 30 Oct 2024 18:29  
 Operator : RC/JU  
 Sample : P4492-17  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4EZ8

Quant Time: Oct 31 22:24:17 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 30 10:56:39 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.640  | 152  | 88288    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.393 | 136  | 339873   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.269 | 164  | 258506   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.069 | 188  | 607251   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.504 | 240  | 644956   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.762 | 264  | 777359   | 20.000 | ng/u1  | 0.01     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.181  | 96   | 5621     | 2.474  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.593  | 84   | 84814    | 13.598 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.834  | 99   | 124150   | 16.705 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.981  | 67   | 72971    | 16.706 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.181  | 132  | 93722    | 17.773 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.346  | 113  | 103847   | 17.509 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.775  | 128  | 45950    | 17.583 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.493  | 143  | 57870    | 19.127 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.034 | 165  | 116812   | 18.166 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.534 | 131  | 126329   | 16.626 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.669 | 166  | 385519   | 18.770 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 13.957 | 160  | 412513   | 18.910 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.510 | 143  | 47624    | 13.821 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.269 | 176  | 353909   | 19.866 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.416 | 200  | 39161    | 9.805  | ng/u1  | 0.01     |
| 73) Anthracene-d10            | 17.169 | 188  | 562338   | 19.685 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 726413   | 21.298 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.539 | 264  | 850147   | 20.478 | ng/u1  | 0.02     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.116 | 178  | 49252    | 1.516  | ng/u1  | 98       |
| 80) Fluoranthene              | 19.204 | 202  | 93739    | 2.391  | ng/u1  | 99       |
| 82) Pyrene                    | 19.580 | 202  | 85612    | 2.037  | ng/u1  | 99       |
| 85) Benzo(a)anthracene        | 21.486 | 228  | 46806    | 1.035  | ng/u1  | 97       |
| 87) Chrysene                  | 21.545 | 228  | 46483    | 1.109  | ng/u1  | 98       |
| 90) Benzo(b)fluoranthene      | 23.745 | 252  | 65884    | 1.346  | ng/u1  | 98       |
| -----                         |        |      |          |        |        |          |

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

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