

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092424\  
 Data File : BP022032.D  
 Acq On : 25 Sep 2024 17:26  
 Operator : RC/JU  
 Sample : P4107-07  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0B66

Quant Time: Sep 25 18:21:34 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 24 15:23:13 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 80166    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.463 | 136  | 299638   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.310 | 164  | 227117   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.110 | 188  | 569006   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.533 | 240  | 680598   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.822 | 264  | 847083   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.223  | 96   | 12274    | 5.997  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 54838    | 9.374  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 99606    | 15.189 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.040  | 67   | 155063   | 38.401 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 182699   | 38.507 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.405  | 113  | 164157   | 31.591 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.840  | 128  | 94656    | 42.207 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.552  | 143  | 111672   | 43.916 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 242984   | 44.001 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.599 | 131  | 205166   | 31.138 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.722 | 166  | 793345   | 45.188 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.004 | 160  | 804110   | 43.260 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.516 | 143  | 35378    | 11.781 | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.316 | 176  | 684819   | 44.097 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.440 | 200  | 135427   | 38.336 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.204 | 188  | 1218114  | 45.857 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.586 | 212  | 1623292  | 46.037 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 2069158  | 46.442 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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