

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092424\  
 Data File : BP022011.D  
 Acq On : 25 Sep 2024 01:49  
 Operator : RC/JU  
 Sample : P4092-05  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0B19

Quant Time: Sep 25 02:18:27 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     | 7.710  | 152  | 90332    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.469 | 136  | 332862   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.322 | 164  | 254654   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.110 | 188  | 624925   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 787742   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.821 | 264  | 927906   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.223  | 96   | 11443    | 4.962  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 62474    | 9.477  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.893  | 99   | 68241    | 9.235  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.046  | 67   | 149739   | 32.910 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.252  | 132  | 164446   | 30.759 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.405  | 113  | 120571   | 20.592 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.852  | 128  | 90286    | 36.240 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 104027   | 36.826 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 223949   | 36.506 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.599 | 131  | 202182   | 27.622 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.728 | 166  | 775043   | 39.372 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.010 | 160  | 778219   | 37.340 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.528 | 143  | 26871    | 7.980  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.328 | 176  | 675052   | 38.768 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.445 | 200  | 132691   | 34.200 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.204 | 188  | 1178461  | 40.394 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.586 | 212  | 1588219  | 38.916 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.598 | 264  | 1984000  | 40.652 | ng/u1 | -0.01    |

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

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

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