

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013025\  
 Data File : BP023809.D  
 Acq On : 30 Jan 2025 22:29  
 Operator : RC/JU  
 Sample : Q1192-06  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A44W1

Quant Time: Jan 31 08:15:17 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 31 08:09:38 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.787  | 152  | 467715   | 20.000 | ng/u1 | 0.02     |
| 20) Naphthalene-d8            | 10.557 | 136  | 1950514  | 20.000 | ng/u1 | 0.01     |
| 38) Acenaphthene-d10          | 14.404 | 164  | 1287708  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 2656045  | 20.000 | ng/u1 | -0.02    |
| 79) Chrysene-d12              | 21.633 | 240  | 2570025  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 25.010 | 264  | 2807720  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 49651    | 4.411  | ng/uL | 0.01     |
| 4) Pyridine-d5                | 3.717  | 84   | 132637   | 4.083  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 6.963  | 99   | 294004   | 7.103  | ng/u1 | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 699601   | 31.932 | ng/u1 | 0.02     |
| 11) 2-Chlorophenol-d4         | 7.328  | 132  | 894613   | 27.523 | ng/u1 | 0.02     |
| 15) 4-Methylphenol-d8         | 8.487  | 113  | 579316   | 17.133 | ng/u1 | 0.02     |
| 21) Nitrobenzene-d5           | 8.928  | 128  | 502924   | 32.228 | ng/u1 | 0.02     |
| 24) 2-Nitrophenol-d4          | 9.646  | 143  | 517131   | 30.711 | ng/u1 | 0.02     |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165  | 911193   | 29.498 | ng/u1 | 0.01     |
| 31) 4-Chloroaniline-d4        | 10.693 | 131  | 625527   | 13.540 | ng/u1 | 0.01     |
| 46) Dimethylphthalate-d6      | 13.810 | 166  | 3455605  | 35.977 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.093 | 160  | 3614899  | 33.369 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.616 | 143  | 129545   | 6.791  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.404 | 176  | 2976332  | 36.797 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200  | 468163   | 28.571 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.292 | 188  | 5097020  | 39.084 | ng/u1 | -0.02    |
| 81) Pyrene-d10                | 19.663 | 212  | 5774074  | 41.930 | ng/u1 | -0.02    |
| 92) Benzo(a)pyrene-d12        | 24.780 | 264  | 5903094  | 40.296 | ng/u1 | 0.00     |

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

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

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