

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082624\  
 Data File : BP021695.D  
 Acq On : 28 Aug 2024 08:51  
 Operator : RC/JU  
 Sample : P3675-16  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCXY9

Quant Time: Aug 28 09:20:42 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 27 05:11:38 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 400797   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.581 | 136  | 1780320  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.440 | 164  | 1172950  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.234 | 188  | 2521934  | 20.000 | ng/u1 | -0.02    |
| 79) Chrysene-d12              | 21.698 | 240  | 2284798  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.116 | 264  | 2561243  | 20.000 | ng/u1 | 0.02     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.276  | 96   | 37582    | 3.113  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 565735   | 16.202 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.964  | 99   | 1004085  | 23.475 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 505267   | 21.706 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.328  | 132  | 634019   | 22.959 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 755917   | 22.870 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 330294   | 23.060 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.663  | 143  | 356092   | 24.492 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.205 | 165  | 691468   | 24.134 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.705 | 131  | 624804   | 14.989 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.834 | 166  | 2193771  | 24.443 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.122 | 160  | 2496094  | 24.771 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.634 | 143  | 404943   | 23.715 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.445 | 176  | 1903656  | 25.229 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 244209   | 17.606 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.339 | 188  | 2817225  | 23.538 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.716 | 212  | 3418870  | 26.176 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.868 | 264  | 3436900  | 25.165 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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