

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071124\  
 Data File : BP020922.D  
 Acq On : 12 Jul 2024 05:01  
 Operator : MA/JU  
 Sample : P3087-13  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4BP1

Quant Time: Jul 12 05:29:56 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 10 18:38:19 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 245233   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.493 | 136  | 1012069  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.351 | 164  | 685287   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.163 | 188  | 1507611  | 20.000 | ng/u1 | -0.02    |
| 79) Chrysene-d12              | 21.622 | 240  | 1471056  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.968 | 264  | 1681910  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.258  | 96   | 14633    | 2.715  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.675  | 84   | 214097   | 13.723 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.922  | 99   | 340599   | 17.036 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.063  | 67   | 212220   | 17.515 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.269  | 132  | 290040   | 17.607 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.434  | 113  | 295011   | 17.766 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 143951   | 18.312 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.581  | 143  | 164277   | 18.259 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.134 | 165  | 297214   | 18.269 | ng/u1 | 0.01     |
| 31) 4-Chloroaniline-d4        | 10.634 | 131  | 398862   | 18.506 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.763 | 166  | 1029724  | 20.670 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.045 | 160  | 1140005  | 20.380 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.610 | 143  | 156423   | 22.069 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.363 | 176  | 899125   | 22.000 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.510 | 200  | 135841   | 14.892 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.263 | 188  | 1467216  | 21.912 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.651 | 212  | 1801703  | 19.811 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 24.745 | 264  | 1839606  | 22.177 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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