

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080922\  
 Data File : BP011339.D  
 Acq On : 09 Aug 2022 20:44  
 Operator : CG/JU  
 Sample : N3900-01  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Aug 09 23:16:02 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 08 18:47:19 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.587  | 152  | 135572   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.381 | 136  | 574839   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.257 | 164  | 332805   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.034 | 188  | 668076   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.157 | 240  | 625126   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.463 | 264  | 651766   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.099  | 96  | 18525   | 4.189  | ng/uL | 0.01  |
| 4) Pyridine-d5                | 3.517  | 84  | 59820   | 5.362  | ng/u1 | 0.02  |
| 7) Phenol-d5                  | 6.775  | 99  | 78526   | 6.545  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 6.934  | 67  | 255470  | 30.031 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.122  | 132 | 232682  | 24.717 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.311  | 113 | 144665  | 15.357 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.752  | 128 | 140594  | 35.937 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.469  | 143 | 133317  | 37.221 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.005 | 165 | 213573  | 29.526 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.528 | 131 | 294418  | 23.282 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.687 | 166 | 808998  | 35.034 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 13.951 | 160 | 938874  | 33.776 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.493 | 143 | 15457   | 3.665  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.263 | 176 | 655149  | 33.176 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.398 | 200 | 92255   | 29.730 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.134 | 188 | 1082059 | 36.034 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.392 | 212 | 1235908 | 35.885 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.316 | 264 | 1189183 | 36.179 | ng/u1 | -0.01 |

Target Compounds Qvalue

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

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