

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080922\  
 Data File : BP011362.D  
 Acq On : 10 Aug 2022 10:35  
 Operator : CG/JU  
 Sample : N4069-12  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Aug 10 23:56:11 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  | 122138   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.375 | 136  | 513009   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.257 | 164  | 294049   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.028 | 188  | 604782   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.157 | 240  | 550871   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.468 | 264  | 553922   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.087  | 96   | 14186    | 3.561  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.511  | 84   | 54372    | 5.410  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.769  | 99   | 70214    | 6.496  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.934  | 67   | 206125   | 26.895 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.116  | 132  | 194882   | 22.979 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.305  | 113  | 106285   | 12.523 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.746  | 128  | 116991   | 33.508 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.469  | 143  | 115624   | 36.172 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.005 | 165  | 175300   | 27.156 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.528 | 131  | 143082   | 12.678 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.681 | 166  | 697340   | 34.179 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.945 | 160  | 818502   | 33.327 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.493 | 143  | 15650    | 4.200  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.263 | 176  | 598963   | 34.328 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.392 | 200  | 75058    | 26.719 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.128 | 188  | 939508   | 34.561 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.392 | 212  | 1094328  | 36.058 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.321 | 264  | 973159   | 34.837 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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