

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080922\  
 Data File : BP011373.D  
 Acq On : 10 Aug 2022 16:55  
 Operator : CG/JU  
 Sample : N4128-06  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Aug 11 00:02:10 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     | 7.581  | 152  | 84826    | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.375 | 136  | 354426   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.257 | 164  | 198928   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.128 | 188  | 694768   | 20.000 | ng/u1 | 0.09     |
| 79) Chrysene-d12              | 21.157 | 240  | 396030   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.468 | 264  | 418849   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.140  | 96   | 130      | 0.047  | ng/uL | 0.05     |
| 4) Pyridine-d5                | 3.511  | 84   | 32346    | 4.634  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.769  | 99   | 55660    | 7.415  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.934  | 67   | 172074   | 32.328 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.116  | 132  | 157643   | 26.764 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.305  | 113  | 103022   | 17.478 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.752  | 128  | 94351    | 39.115 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.469  | 143  | 88886    | 40.249 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.004 | 165  | 144787   | 32.465 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.528 | 131  | 187443   | 24.041 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.681 | 166  | 536894   | 38.898 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.945 | 160  | 635438   | 38.245 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.563 | 143  | 461      | 0.183  | ng/u1 | 0.07     |
| 60) Fluorene-d10              | 15.263 | 176  | 440185   | 37.291 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.469 | 200  | 323      | 0.100  | ng/u1 | 0.07     |
| 73) Anthracene-d10            | 17.128 | 188  | 699367   | 22.395 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.392 | 212  | 829671   | 38.026 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.321 | 264  | 835502   | 39.554 | ng/u1 | 0.00     |

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

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

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