

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100522\  
 Data File : BP011963.D  
 Acq On : 06 Oct 2022 11:03  
 Operator : CG/JU  
 Sample : N4956-11  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CB8T1

Quant Time: Oct 06 23:04:39 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP093022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 05 23:18:55 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.875  | 152  | 261606   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.687 | 136  | 1107075  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.522 | 164  | 672261   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.281 | 188  | 1401448  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.369 | 240  | 1331257  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.798 | 264  | 1333286  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 22634    | 3.816  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 91400    | 5.260  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.040  | 99   | 108157   | 4.789  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.211  | 67   | 303855   | 24.273 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.405  | 132  | 335062   | 18.794 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.587  | 113  | 207339   | 11.106 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.046  | 128  | 223188   | 26.468 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.769  | 143  | 214488   | 23.774 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.305 | 165  | 368240   | 22.067 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.828 | 131  | 323687   | 12.724 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.934 | 166  | 1335299  | 27.776 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.216 | 160  | 1484096  | 27.080 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.740 | 143  | 33573    | 3.372  | ng/ul | 0.01     |
| 60) Fluorene-d10              | 15.522 | 176  | 1191308  | 28.535 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.640 | 200  | 174875   | 20.616 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.381 | 188  | 1997491  | 31.290 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.616 | 212  | 2330667  | 34.569 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.639 | 264  | 2205464  | 33.044 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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