

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100522\  
 Data File : BP011926.D  
 Acq On : 05 Oct 2022 12:47  
 Operator : CG/JU  
 Sample : N4890-05  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EY0E4

Quant Time: Oct 05 23:21:19 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     | 7.881  | 152  | 261783   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.692 | 136  | 1041314  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.528 | 164  | 578379   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.286 | 188  | 1153782  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.368 | 240  | 1249265  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.804 | 264  | 1296064  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 30199    | 5.088  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.734  | 84   | 48815    | 2.808  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.046  | 99   | 162829   | 7.205  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.210  | 67   | 384464   | 30.691 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.410  | 132  | 454132   | 25.455 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.593  | 113  | 280357   | 15.007 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.051  | 128  | 261876   | 33.017 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.775  | 143  | 266559   | 31.412 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.310 | 165  | 446498   | 28.447 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.834 | 131  | 313997   | 13.122 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.939 | 166  | 1422198  | 34.386 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.222 | 160  | 1628133  | 34.530 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.739 | 143  | 44363    | 5.180  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.522 | 176  | 1283766  | 35.741 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.639 | 200  | 188308   | 26.965 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.386 | 188  | 2067126  | 39.331 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.621 | 212  | 2486203  | 39.296 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.645 | 264  | 2631508  | 40.560 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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