

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP091023\  
 Data File : BP017097.D  
 Acq On : 11 Sep 2023 10:48  
 Operator : MA/JU  
 Sample : 04262-03  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BBJD4

Quant Time: Sep 12 02:27:04 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP091023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 11 06:13:29 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.958  | 152  | 337706   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.781 | 136  | 1474643  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.616 | 164  | 885448   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.387 | 188  | 1882389  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 1468866  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.021 | 264  | 1401144  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.317  | 96   | 45334    | 5.124  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.752  | 84   | 278691   | 10.445 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 248754   | 8.103  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.299  | 67   | 680919   | 33.073 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 597982   | 26.952 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.663  | 113  | 427822   | 17.618 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.158  | 128  | 363568   | 34.630 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.869  | 143  | 374990   | 34.607 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.393 | 165  | 620065   | 30.179 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.946 | 131  | 810919   | 24.613 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.034 | 166  | 2311260  | 36.684 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.316 | 160  | 2491640  | 34.731 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.834 | 143  | 78826    | 6.064  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.616 | 176  | 1957792  | 36.618 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.740 | 200  | 307349   | 28.933 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.487 | 188  | 3148244  | 39.019 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.728 | 212  | 3552449  | 45.925 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.857 | 264  | 2733900  | 40.644 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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