

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092424\  
 Data File : BP022058.D  
 Acq On : 26 Sep 2024 11:47  
 Operator : RC/JU  
 Sample : P4084-06  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCYM3

Quant Time: Sep 27 01:16:30 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 24 15:23:13 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 84394    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.463 | 136  | 313933   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.304 | 164  | 235972   | 20.000 | ng/ul | -0.02    |
| 64) Phenanthrene-d10          | 17.098 | 188  | 580735   | 20.000 | ng/ul | -0.02    |
| 79) Chrysene-d12              | 21.528 | 240  | 732828   | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12              | 24.804 | 264  | 885375   | 20.000 | ng/ul | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.217  | 96   | 12792    | 5.937  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 39383    | 6.395  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 71879    | 10.412 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.046  | 67   | 139885   | 32.907 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 160150   | 32.063 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.405  | 113  | 122109   | 22.322 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.846  | 128  | 83829    | 35.677 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.552  | 143  | 100997   | 37.909 | ng/ul | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 198532   | 34.314 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.599 | 131  | 124805   | 18.079 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.722 | 166  | 713458   | 39.113 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.004 | 160  | 681058   | 35.265 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.516 | 143  | 24951    | 7.997  | ng/ul | -0.01    |
| 60) Fluorene-d10              | 15.316 | 176  | 606857   | 37.611 | ng/ul | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.434 | 200  | 119947   | 33.268 | ng/ul | -0.01    |
| 73) Anthracene-d10            | 17.198 | 188  | 1081778  | 39.902 | ng/ul | -0.02    |
| 81) Pyrene-d10                | 19.575 | 212  | 1486900  | 39.164 | ng/ul | -0.02    |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 1844899  | 39.618 | ng/ul | -0.02    |

Target Compounds Qvalue

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

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