

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092523\  
 Data File : BP017340.D  
 Acq On : 26 Sep 2023 02:06  
 Operator : MA/JU  
 Sample : 04472-13  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYW9

Quant Time: Sep 26 03:10:54 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 21 23:51:07 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.928  | 152  | 187296   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.746 | 136  | 803482   | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10          | 14.587 | 164  | 484877   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.357 | 188  | 1040095  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.463 | 240  | 923107   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.974 | 264  | 1027213  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 22877    | 5.017  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 314590   | 24.678 | ng/u1 | -0.01    |
| 7) Phenol-d5                  | 7.081  | 99   | 416842   | 27.098 | ng/u1 | -0.02    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 267097   | 28.772 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.446  | 132  | 326533   | 26.764 | ng/u1 | -0.02    |
| 15) 4-Methylphenol-d8         | 8.640  | 113  | 330041   | 27.597 | ng/u1 | -0.02    |
| 21) Nitrobenzene-d5           | 9.122  | 128  | 164240   | 28.275 | ng/u1 | -0.02    |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 143261   | 22.714 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.363 | 165  | 312661   | 26.239 | ng/u1 | -0.02    |
| 31) 4-Chloroaniline-d4        | 10.916 | 131  | 481778   | 27.911 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 14.004 | 166  | 1102051  | 31.727 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 1224797  | 30.659 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.822 | 143  | 18021    | 2.978  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 968848   | 32.399 | ng/u1 | -0.02    |
| 65) 4,6-Dinitro-2-methylph... | 15.722 | 200  | 3850     | 0.657  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.457 | 188  | 1462051  | 31.695 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.698 | 212  | 1669943  | 33.453 | ng/u1 | -0.02    |
| 92) Benzo(a)pyrene-d12        | 23.810 | 264  | 1513119  | 30.594 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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