

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082123\  
 Data File : BP016685.D  
 Acq On : 21 Aug 2023 21:22  
 Operator : MA/JU  
 Sample : 04034-01  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYN5

Quant Time: Aug 22 00:34:31 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 17 01:38:34 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.046  | 152  | 308906   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.875 | 136  | 1404305  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.693 | 164  | 869617   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.463 | 188  | 1937079  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.557 | 240  | 1603379  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.133 | 264  | 1473488  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 19753    | 2.430  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 135502   | 5.479  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.187  | 99   | 201098   | 7.061  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 611613   | 32.607 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.564  | 132  | 521215   | 25.481 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.746  | 113  | 373750   | 16.505 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.240  | 128  | 358792   | 36.311 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.957  | 143  | 354940   | 37.820 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 608123   | 32.927 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.028 | 131  | 617008   | 19.471 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 14.104 | 166  | 2222147  | 35.329 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.393 | 160  | 2541229  | 35.795 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.898 | 143  | 68830    | 5.372  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.693 | 176  | 1973215  | 36.799 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.810 | 200  | 279911   | 30.375 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.563 | 188  | 3423861  | 40.863 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 3857923  | 45.920 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.963 | 264  | 2953713  | 41.828 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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