

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051923\  
 Data File : BP015200.D  
 Acq On : 20 May 2023 07:00  
 Operator : CG/JU  
 Sample : 02689-17  
 Misc :  
 ALS Vial : 37 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JREN9

Quant Time: May 22 00:26:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri May 19 01:07:51 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.122  | 152  | 235281   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.951 | 136  | 986257   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.763 | 164  | 589237   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.510 | 188  | 1152507  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.586 | 240  | 921776   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.145 | 264  | 1106765  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 32400    | 5.350  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.869  | 84   | 381336   | 22.216 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.275  | 99   | 778129   | 35.620 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.446  | 67   | 472055   | 37.415 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.652  | 132  | 622284   | 38.743 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.834  | 113  | 651133   | 37.406 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.299  | 128  | 318928   | 41.152 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.028 | 143  | 374112   | 43.384 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.569 | 165  | 640601   | 41.168 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.087 | 131  | 808697   | 35.126 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.163 | 166  | 2072175  | 46.539 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.457 | 160  | 2251937  | 44.221 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.951 | 143  | 260771   | 30.161 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.751 | 176  | 1660095  | 44.185 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.869 | 200  | 222951   | 30.696 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.610 | 188  | 2440584  | 46.246 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.833 | 212  | 2709634  | 50.684 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.980 | 264  | 2814735  | 51.232 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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