

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081622\  
 Data File : BP011450.D  
 Acq On : 16 Aug 2022 18:35  
 Operator : CG/JU  
 Sample : SSTD16048  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160601

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 08/17/2022  
 Supervised By :Sohil Jodhani 08/19/2022

Quant Time: Aug 17 01:10:55 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 17 01:01:09 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.558  | 152  | 85073    | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.352 | 136  | 390355   | 20.000  | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.234 | 164  | 236773   | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.010 | 188  | 477200   | 20.000  | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.139 | 240  | 460069   | 20.000  | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.433 | 264  | 438396   | 20.000  | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 3.475  | 84   | 1166471  | 166.635 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.758  | 99   | 1327602  | 176.338 | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.917  | 67   | 807732   | 151.311 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/ul  |          |
| 15) 4-Methylphenol-d8         | 8.299  | 113  | 1046774  | 177.076 | ng/ul  | 0.02     |
| 21) Nitrobenzene-d5           | 0.000  | 128  | 0d       | 0.000   | ng/ul  |          |
| 24) 2-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000   | ng/ul  |          |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165  | 0d       | 0.000   | ng/ul  |          |
| 31) 4-Chloroaniline-d4        | 10.510 | 131  | 1430635  | 166.598 | ng/ul  | 0.01     |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.000   | ng/ul  |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.000   | ng/ul  |          |
| 54) 4-Nitrophenol-d4          | 14.493 | 143  | 561207   | 187.032 | ng/ul  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/ul  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.393 | 200  | 480236   | 216.661 | ng/ul  | 0.02     |
| 73) Anthracene-d10            | 0.000  | 188  | 0d       | 0.000   | ng/ul  |          |
| 81) Pyrene-d10                | 0.000  | 212  | 0d       | 0.000   | ng/ul  |          |
| 92) Benzo(a)pyrene-d12        | 0.000  | 264  | 0d       | 0.000   | ng/ul  |          |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 5) Pyridine                   | 3.493  | 79   | 1165150  | 159.795 | ng/ul  | 99       |
| 6) Benzaldehyde               | 6.716  | 77   | 491356   | 149.755 | ng/ul  | 98       |
| 8) Phenol                     | 6.787  | 94   | 1349548  | 165.602 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.011  | 93   | 1003510  | 152.901 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.022  | 108  | 995985   | 169.700 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.105  | 45   | 1426223  | 144.886 | ng/ul  | 100      |
| 16) Acetophenone              | 8.405  | 105  | 1684140  | 176.087 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.375  | 108  | 1082871  | 172.134 | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.540 | 127  | 1474625  | 171.318 | ng/ul  | 97       |
| 34) Caprolactam               | 11.381 | 113  | 341269m  | 198.676 | ng/ul  |          |
| 40) Hexachlorocyclopentadiene | 12.375 | 237  | 693017   | 207.475 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.175 | 138  | 607361   | 212.215 | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 14.381 | 184  | 385507   | 249.737 | ng/ul  | 97       |
| 55) 4-Nitrophenol             | 14.510 | 109  | 652903   | 234.555 | ng/ul  | 93       |
| 63) 4-Nitroaniline            | 15.363 | 138  | 556477m  | 199.819 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.410 | 198  | 484553   | 212.301 | ng/ul# | 97       |
| 70) Atrazine                  | 16.504 | 200  | 858355   | 183.129 | ng/ul  | 97       |
| 71) Pentachlorophenol         | 16.657 | 266  | 555497   | 193.959 | ng/ul  | 99       |
| 77) Carbazole                 | 17.434 | 167  | 3940749  | 158.313 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.069 | 252  | 1493046  | 178.319 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 21.963 | 149  | 6223930  | 204.618 | ng/ul  | 100      |
| -----                         |        |      |          |         |        |          |

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

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SSTD160601

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