

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG030623\  
 Data File : BG056839.D  
 Acq On : 6 Mar 2023 13:16  
 Operator : CG/JU  
 Sample : 01712-08MSD  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

DBZX0MSD

## Manual Integrations

## APPROVED

Reviewed By :Christian Giraldo 03/07/2023  
 Supervised By :Jagrut Upadhyay 03/07/2023

Quant Time: Mar 07 01:38:22 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022323.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Mar 02 00:12:28 2023

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.415  | 152  | 14436    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.253 | 136  | 61852    | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 15.019 | 164  | 49997    | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.751 | 188  | 125138   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 22.058 | 240  | 117064   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 25.583 | 264  | 128278   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.697  | 96   | 1954     | 7.508  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 4.132  | 84   | 36858    | 35.089 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.539  | 99   | 50016    | 34.623 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.716  | 67   | 29349    | 33.306 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.939  | 132  | 31101    | 32.419 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 9.108  | 113  | 36807    | 33.345 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.584  | 128  | 17720    | 34.186 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.318 | 143  | 20669    | 33.499 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.853 | 165  | 40943    | 35.223 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.376 | 131  | 47317    | 31.210 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.402 | 166  | 135842   | 33.675 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.719 | 160  | 156023   | 34.031 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.177 | 143  | 22182    | 31.894 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 16.000 | 176  | 131098   | 35.133 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.106 | 200  | 26771    | 31.976 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.851 | 188  | 200412   | 33.012 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 20.113 | 212  | 247836   | 31.287 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.342 | 264  | 238055   | 34.381 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        |        | Qvalue   |
| 2) 1,4-Dioxane                | 3.738  | 88   | 5186m    | 16.162 | ng/uL  |          |
| 5) Pyridine                   | 4.155  | 79   | 41013    | 36.868 | ng/u1  | 94       |
| 6) Benzaldehyde               | 7.539  | 77   | 36091    | 47.687 | ng/u1  | 95       |
| 8) Phenol                     | 7.569  | 94   | 53541    | 36.336 | ng/u1  | 95       |
| 10) Bis(2-Chloroethyl)ether   | 7.815  | 93   | 36218    | 34.594 | ng/u1  | 97       |
| 12) 2-Chlorophenol            | 7.974  | 128  | 32602    | 34.317 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.844  | 108  | 37602    | 33.832 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.932  | 45   | 49417    | 35.126 | ng/u1  | 98       |
| 16) Acetophenone              | 9.243  | 105  | 65351    | 34.573 | ng/u1  | 99       |
| 17) N-Nitroso-di-n-propyla... | 9.220  | 70   | 37377    | 34.463 | ng/u1  | 97       |
| 18) 4-Methylphenol            | 9.173  | 108  | 42337    | 35.056 | ng/u1  | 93       |
| 19) Hexachloroethane          | 9.519  | 117  | 14016    | 34.694 | ng/u1  | 91       |
| 22) Nitrobenzene              | 9.631  | 77   | 57936    | 38.544 | ng/u1  | 99       |
| 23) Isophorone                | 10.154 | 82   | 103698   | 35.288 | ng/u1  | 98       |
| 25) 2-Nitrophenol             | 10.348 | 139  | 20719    | 34.018 | ng/u1# | 92       |
| 26) 2,4-Dimethylphenol        | 10.389 | 107  | 47966    | 35.264 | ng/u1  | 96       |
| 27) Bis(2-Chloroethoxy)met... | 10.630 | 93   | 52959    | 35.654 | ng/u1  | 98       |
| 29) 2,4-Dichlorophenol        | 10.882 | 162  | 41630    | 36.984 | ng/u1  | 97       |
| 30) Naphthalene               | 11.300 | 128  | 120150   | 35.291 | ng/u1  | 97       |
| 32) 4-Chloroaniline           | 11.400 | 127  | 46476    | 30.947 | ng/u1# | 88       |
| 33) Hexachlorobutadiene       | 11.582 | 225  | 32500    | 36.227 | ng/u1  | 91       |
| 34) Caprolactam               | 12.140 | 113  | 14403m   | 37.638 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.481 | 107  | 46754    | 36.155 | ng/u1  | 95       |
| 36) 2-Methylnaphthalene       | 12.874 | 142  | 87641    | 35.944 | ng/u1  | 100      |

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 Supervised By : Jagrut Upadhyay 03/07/2023

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Quant Title : SVOA CALIBRATION

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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 13.092 | 142  | 89393    | 37.185 | ng/u1  | 95       |
| 39) 1,2,4,5-Tetrachloroben... | 13.233 | 216  | 65056    | 36.701 | ng/u1  | 99       |
| 40) Hexachlorocyclopentadiene | 13.209 | 237  | 36620    | 28.831 | ng/u1  | 99       |
| 41) 2,4,6-Trichlorophenol     | 13.462 | 196  | 41775    | 35.270 | ng/u1  | 93       |
| 42) 2,4,5-Trichlorophenol     | 13.532 | 196  | 46281    | 35.986 | ng/u1  | 99       |
| 43) 1,1'-Biphenyl             | 13.861 | 154  | 128812   | 35.037 | ng/u1  | 99       |
| 44) 2-Chloronaphthalene       | 13.908 | 162  | 107476   | 35.332 | ng/u1  | 94       |
| 45) 2-Nitroaniline            | 14.096 | 65   | 38003    | 35.993 | ng/u1  | 93       |
| 47) Dimethylphthalate         | 14.449 | 163  | 139717   | 34.059 | ng/u1# | 98       |
| 48) 2,6-Dinitrotoluene        | 14.578 | 165  | 29779    | 35.412 | ng/u1# | 82       |
| 50) Acenaphthylene            | 14.749 | 152  | 163858   | 35.378 | ng/u1  | 99       |
| 51) 3-Nitroaniline            | 14.907 | 138  | 25499    | 33.881 | ng/u1  | 88       |
| 52) Acenaphthene              | 15.083 | 153  | 112985   | 35.511 | ng/u1  | 96       |
| 53) 2,4-Dinitrophenol         | 15.113 | 184  | 18152    | 33.695 | ng/u1# | 82       |
| 55) 4-Nitrophenol             | 15.195 | 109  | 29503    | 36.648 | ng/u1  | 90       |
| 56) Dibenzofuran              | 15.412 | 168  | 170621   | 35.326 | ng/u1  | 95       |
| 57) 2,4-Dinitrotoluene        | 15.360 | 165  | 41718    | 34.604 | ng/u1# | 94       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.630 | 232  | 42826    | 35.783 | ng/u1  | 97       |
| 59) Diethylphthalate          | 15.800 | 149  | 140001   | 34.094 | ng/u1  | 98       |
| 61) Fluorene                  | 16.059 | 166  | 133577   | 34.539 | ng/u1  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 16.035 | 204  | 81461    | 35.561 | ng/u1  | 96       |
| 63) 4-Nitroaniline            | 16.065 | 138  | 26482    | 36.013 | ng/u1  | 92       |
| 66) 4,6-Dinitro-2-methylph... | 16.118 | 198  | 26675    | 33.095 | ng/u1  | 96       |
| 67) N-Nitrosodiphenylamine    | 16.247 | 169  | 116535   | 33.489 | ng/u1  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.928 | 248  | 55254    | 34.842 | ng/u1  | 97       |
| 69) Hexachlorobenzene         | 17.058 | 284  | 59664    | 35.020 | ng/u1  | 95       |
| 70) Atrazine                  | 17.181 | 200  | 50112    | 33.157 | ng/u1  | 97       |
| 71) Pentachlorophenol         | 17.398 | 266  | 32672m   | 33.773 | ng/u1  |          |
| 72) Phenanthrene              | 17.792 | 178  | 235761   | 34.781 | ng/u1  | 99       |
| 74) Anthracene                | 17.886 | 178  | 229922   | 33.565 | ng/u1  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.832 | 216  | 68540    | 35.802 | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 15.330 | 250  | 68257    | 35.155 | ng/uL  | 92       |
| 77) Carbazole                 | 18.145 | 167  | 200257   | 34.939 | ng/u1  | 98       |
| 78) Di-n-butylphthalate       | 18.673 | 149  | 223679   | 33.861 | ng/u1  | 99       |
| 80) Fluoranthene              | 19.778 | 202  | 294476   | 30.745 | ng/u1  | 99       |
| 82) Pyrene                    | 20.136 | 202  | 293917   | 31.148 | ng/u1  | 99       |
| 83) Butylbenzylphthalate      | 20.994 | 149  | 96273    | 32.161 | ng/u1  | 91       |
| 84) 3,3'-Dichlorobenzidine    | 21.934 | 252  | 85131    | 29.005 | ng/u1  | 97       |
| 85) Benzo(a)anthracene        | 22.040 | 228  | 298474   | 35.032 | ng/u1  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.899 | 149  | 141055   | 33.544 | ng/u1  | 98       |
| 87) Chrysene                  | 22.110 | 228  | 278201   | 35.058 | ng/u1  | 100      |
| 89) Di-n-octyl phthalate      | 23.209 | 149  | 235826   | 33.815 | ng/u1  | 100      |
| 90) Benzo(b)fluoranthene      | 24.455 | 252  | 304026m  | 35.403 | ng/u1  |          |
| 91) Benzo(k)fluoranthene      | 24.525 | 252  | 293444   | 35.082 | ng/u1  | 98       |
| 93) Benzo(a)pyrene            | 25.418 | 252  | 259315   | 34.353 | ng/u1  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.643 | 276  | 360651   | 35.171 | ng/u1# | 97       |
| 95) Dibenzo(a,h)anthracene    | 29.707 | 278  | 295122   | 34.421 | ng/u1# | 95       |
| 96) Benzo(g,h,i)perylene      | 30.930 | 276  | 288855   | 35.265 | ng/u1  | 98       |

03/07/2023

Supervised By : Jagrut Upadhyay

03/07/2023

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

