

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022423\  
 Data File : BG056719.D  
 Acq On : 24 Feb 2023 12:34  
 Operator : CG/JU  
 Sample : 01378-03  
 Misc : SFAM-MDL-SOIL-01  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-SOIL-QT1-2023-01

Quant Time: Feb 24 13:09:32 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022323.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Feb 23 15:40:42 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 02/27/2023  
 Supervised By :Jagrut Upadhyay 02/27/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.415  | 152  | 17115    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.258 | 136  | 77312    | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 15.025 | 164  | 61239    | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.757 | 188  | 156833   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 22.069 | 240  | 141588   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 25.600 | 264  | 142341   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.703  | 96   | 1236     | 4.006  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 4.137  | 84   | 16639m   | 13.361 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.539  | 99   | 30178    | 17.620 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.715  | 67   | 18639    | 17.841 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.945  | 132  | 20364    | 17.905 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 9.108  | 113  | 23543    | 17.990 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.596  | 128  | 11402    | 17.598 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.324 | 143  | 14010    | 18.166 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.865 | 165  | 25304    | 17.416 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.382 | 131  | 33565    | 17.712 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.414 | 166  | 89800    | 18.175 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.725 | 160  | 99658    | 17.747 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.177 | 143  | 14818    | 17.395 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 16.006 | 176  | 81464    | 17.824 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.106 | 200  | 16793    | 16.004 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.857 | 188  | 130413   | 17.140 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 20.119 | 212  | 159370   | 16.634 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.354 | 264  | 135948   | 17.695 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.738  | 88   | 395m     | 1.038  | ng/uL  |          |
| 5) Pyridine                   | 4.161  | 79   | 5057     | 3.834  | ng/u1  | 91       |
| 6) Benzaldehyde               | 7.539  | 77   | 4271     | 4.760  | ng/u1  | 92       |
| 8) Phenol                     | 7.563  | 94   | 7236     | 4.142  | ng/u1  | 94       |
| 10) Bis(2-Chloroethyl)ether   | 7.815  | 93   | 5259     | 4.237  | ng/u1# | 87       |
| 12) 2-Chlorophenol            | 7.974  | 128  | 4792     | 4.254  | ng/u1  | 90       |
| 13) 2-Methylphenol            | 8.844  | 108  | 5464     | 4.147  | ng/u1  | 94       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.932  | 45   | 7507     | 4.501  | ng/u1# | 94       |
| 16) Acetophenone              | 9.243  | 105  | 9437     | 4.211  | ng/u1  | 88       |
| 17) N-Nitroso-di-n-propyla... | 9.214  | 70   | 5711     | 4.442  | ng/u1  | 90       |
| 18) 4-Methylphenol            | 9.167  | 108  | 5631     | 3.933  | ng/u1  | 83       |
| 19) Hexachloroethane          | 9.525  | 117  | 2013     | 4.203  | ng/u1  | 95       |
| 22) Nitrobenzene              | 9.631  | 77   | 8272     | 4.403  | ng/u1  | 98       |
| 23) Isophorone                | 10.154 | 82   | 15558    | 4.236  | ng/u1# | 92       |
| 25) 2-Nitrophenol             | 10.359 | 139  | 2997     | 3.937  | ng/u1  | 97       |
| 26) 2,4-Dimethylphenol        | 10.389 | 107  | 6001     | 3.530  | ng/u1  | 85       |
| 27) Bis(2-Chloroethoxy)met... | 10.630 | 93   | 7720     | 4.158  | ng/u1  | 96       |
| 29) 2,4-Dichlorophenol        | 10.882 | 162  | 5763     | 4.096  | ng/u1  | 92       |
| 30) Naphthalene               | 11.311 | 128  | 17575    | 4.130  | ng/u1# | 97       |
| 32) 4-Chloroaniline           | 11.405 | 127  | 7675     | 4.089  | ng/u1  | 98       |
| 33) Hexachlorobutadiene       | 11.587 | 225  | 4617     | 4.117  | ng/u1  | 90       |
| 34) Caprolactam               | 12.140 | 113  | 2313     | 4.836  | ng/u1  | 81       |
| 35) 4-Chloro-3-methylphenol   | 12.486 | 107  | 6377     | 3.945  | ng/u1# | 84       |
| 36) 2-Methylnaphthalene       | 12.880 | 142  | 12372    | 4.059  | ng/u1  | 96       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022423\  
 Data File : BG056719.D  
 Acq On : 24 Feb 2023 12:34  
 Operator : CG/JU  
 Sample : 01378-03  
 Misc : SFAM-MDL-SOIL-01  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-SOIL-QT1-2023-01

Quant Time: Feb 24 13:09:32 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022323.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Feb 23 15:40:42 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 02/27/2023  
 Supervised By :Jagrut Upadhyay 02/27/2023

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 37) 1-Methylnaphthalene       | 13.097 | 142  | 12510    | 4.163 | ng/ul  | 94       |
| 39) 1,2,4,5-Tetrachloroben... | 13.238 | 216  | 9003     | 4.147 | ng/ul# | 95       |
| 40) Hexachlorocyclopentadiene | 13.221 | 237  | 6341     | 4.076 | ng/ul  | 94       |
| 41) 2,4,6-Trichlorophenol     | 13.468 | 196  | 5788     | 3.990 | ng/ul  | 98       |
| 42) 2,4,5-Trichlorophenol     | 13.532 | 196  | 6799     | 4.316 | ng/ul  | 83       |
| 43) 1,1'-Biphenyl             | 13.867 | 154  | 19409    | 4.310 | ng/ul  | 95       |
| 44) 2-Chloronaphthalene       | 13.914 | 162  | 16000    | 4.294 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 14.096 | 65   | 5177     | 4.003 | ng/ul  | 89       |
| 47) Dimethylphthalate         | 14.455 | 163  | 20886    | 4.157 | ng/ul# | 92       |
| 48) 2,6-Dinitrotoluene        | 14.584 | 165  | 3870     | 3.757 | ng/ul  | 94       |
| 50) Acenaphthylene            | 14.754 | 152  | 23901    | 4.213 | ng/ul  | 97       |
| 51) 3-Nitroaniline            | 14.913 | 138  | 4523     | 4.906 | ng/ul# | 78       |
| 52) Acenaphthene              | 15.089 | 153  | 16218    | 4.161 | ng/ul  | 95       |
| 53) 2,4-Dinitrophenol         | 15.113 | 184  | 2896     | 4.389 | ng/ul# | 86       |
| 55) 4-Nitrophenol             | 15.195 | 109  | 3909     | 3.964 | ng/ul  | 88       |
| 56) Dibenzofuran              | 15.418 | 168  | 24520    | 4.145 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 15.359 | 165  | 6381     | 4.321 | ng/ul# | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.636 | 232  | 5573     | 3.802 | ng/ul# | 91       |
| 59) Diethylphthalate          | 15.806 | 149  | 21360    | 4.247 | ng/ul  | 94       |
| 61) Fluorene                  | 16.065 | 166  | 21110    | 4.456 | ng/ul  | 91       |
| 62) 4-Chlorophenyl-phenyle... | 16.047 | 204  | 11823    | 4.214 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 16.065 | 138  | 3836     | 4.259 | ng/ul  | 89       |
| 66) 4,6-Dinitro-2-methylph... | 16.117 | 198  | 4515     | 4.470 | ng/ul# | 98       |
| 67) N-Nitrosodiphenylamine    | 16.253 | 169  | 18231    | 4.180 | ng/ul  | 95       |
| 68) 4-Bromophenyl-phenylether | 16.940 | 248  | 8303     | 4.178 | ng/ul  | 89       |
| 69) Hexachlorobenzene         | 17.063 | 284  | 8807     | 4.125 | ng/ul  | 95       |
| 70) Atrazine                  | 17.181 | 200  | 8641     | 4.562 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.404 | 266  | 4605     | 3.798 | ng/ul# | 87       |
| 72) Phenanthrene              | 17.798 | 178  | 35861    | 4.221 | ng/ul  | 97       |
| 74) Anthracene                | 17.892 | 178  | 34595    | 4.030 | ng/ul  | 96       |
| 75) 1,2,3,4-Tetrachloroben... | 13.838 | 216  | 8704     | 3.628 | ng/uL  | 93       |
| 76) Pentachlorobenzene        | 15.336 | 250  | 9004     | 3.700 | ng/uL  | 98       |
| 77) Carbazole                 | 18.150 | 167  | 31680    | 4.410 | ng/ul  | 95       |
| 78) Di-n-butylphthalate       | 18.679 | 149  | 35016    | 4.230 | ng/ul  | 98       |
| 80) Fluoranthene              | 19.784 | 202  | 44640    | 3.853 | ng/ul# | 97       |
| 82) Pyrene                    | 20.142 | 202  | 45491    | 3.986 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 21.006 | 149  | 14754    | 4.075 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.946 | 252  | 15910    | 4.482 | ng/ul  | 100      |
| 85) Benzo(a)anthracene        | 22.046 | 228  | 41408    | 4.018 | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.911 | 149  | 21432    | 4.214 | ng/ul  | 99       |
| 87) Chrysene                  | 22.116 | 228  | 39597    | 4.126 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 23.227 | 149  | 36730    | 4.746 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 24.461 | 252  | 43185    | 4.532 | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene      | 24.537 | 252  | 41133    | 4.432 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 25.430 | 252  | 37197    | 4.441 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.649 | 276  | 49014m   | 4.308 | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 29.725 | 278  | 40474m   | 4.254 | ng/ul  |          |
| 96) Benzo(g,h,i)perylene      | 30.929 | 276  | 38702m   | 4.258 | ng/ul  |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022423\  
 Data File : BG056719.D  
 Acq On : 24 Feb 2023 12:34  
 Operator : CG/JU  
 Sample : 01378-03  
 Misc : SFAM-MDL-SOIL-01  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-SOIL-QT1-2023-01

Quant Time: Feb 24 13:09:32 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022323.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Feb 23 15:40:42 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 02/27/2023  
 Supervised By :Jagrut Upadhyay 02/27/2023

