

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052224\  
 Data File : BG061259.D  
 Acq On : 22 May 2024 13:57  
 Operator : MA/JU  
 Sample : P2409-01  
 Misc : SFAM-MDL-WATER-01  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-WATER-QT2-2024-03

Quant Time: May 22 23:48:02 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG052124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 21 15:23:39 2024  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh  
 Patel

05/24/2024  
 Supervised By :mohammad  
 ahmed

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.881  | 152  | 27710    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.684 | 136  | 107520   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.521 | 164  | 90392    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.270 | 188  | 249345m  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.518 | 240  | 265045   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.609 | 264  | 282772   | 20.000 | ng/u1 | 0.00     |

05/24/2024

|                               |        |     |        |        |       |      |
|-------------------------------|--------|-----|--------|--------|-------|------|
| System Monitoring Compounds   |        |     |        |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.334  | 96  | 2411   | 4.869  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.745  | 84  | 40995  | 23.272 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.059  | 99  | 50892  | 22.365 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.212  | 67  | 33715  | 23.572 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.417  | 132 | 38841  | 23.063 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.598  | 113 | 37143  | 19.142 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.039  | 128 | 19590  | 23.664 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.762  | 143 | 22272  | 22.988 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.314 | 165 | 40691  | 21.544 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.819 | 131 | 53539  | 21.667 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.927 | 166 | 160856 | 22.945 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.215 | 160 | 170109 | 23.339 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.750 | 143 | 20510  | 23.663 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.514 | 176 | 144064 | 23.802 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.649 | 200 | 30628  | 20.309 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.370 | 188 | 231536 | 21.062 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.662 | 212 | 303946 | 22.334 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.397 | 264 | 320652 | 23.587 | ng/u1 | 0.00 |

| Target Compounds              | Qvalue |     |       |       |        |    |
|-------------------------------|--------|-----|-------|-------|--------|----|
| 2) 1,4-Dioxane                | 3.375  | 88  | 634m  | 1.067 | ng/uL  |    |
| 5) Pyridine                   | 3.763  | 79  | 9183  | 5.271 | ng/u1  | 93 |
| 6) Benzaldehyde               | 7.030  | 77  | 7011  | 5.892 | ng/u1  | 91 |
| 8) Phenol                     | 7.088  | 94  | 10520 | 4.537 | ng/u1  | 94 |
| 10) Bis(2-Chloroethyl)ether   | 7.306  | 93  | 8808  | 5.194 | ng/u1  | 98 |
| 12) 2-Chlorophenol            | 7.458  | 128 | 8945  | 5.026 | ng/u1  | 95 |
| 13) 2-Methylphenol            | 8.334  | 108 | 7275  | 4.049 | ng/u1  | 97 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.416  | 45  | 20790 | 4.539 | ng/u1  | 99 |
| 16) Acetophenone              | 8.704  | 105 | 14337 | 4.329 | ng/u1  | 91 |
| 17) N-Nitroso-di-n-propyla... | 8.686  | 70  | 7798  | 4.283 | ng/u1# | 97 |
| 18) 4-Methylphenol            | 8.669  | 108 | 7817  | 3.926 | ng/u1  | 97 |
| 19) Hexachloroethane          | 8.963  | 117 | 3564  | 4.501 | ng/u1  | 94 |
| 22) Nitrobenzene              | 9.086  | 77  | 11369 | 5.007 | ng/u1  | 93 |
| 23) Isophorone                | 9.609  | 82  | 22419 | 4.854 | ng/u1  | 98 |
| 25) 2-Nitrophenol             | 9.797  | 139 | 5260  | 5.002 | ng/u1# | 90 |
| 26) 2,4-Dimethylphenol        | 9.861  | 107 | 5802  | 2.784 | ng/u1  | 89 |
| 27) Bis(2-Chloroethoxy)met... | 10.091 | 93  | 11332 | 4.895 | ng/u1# | 88 |
| 29) 2,4-Dichlorophenol        | 10.337 | 162 | 8639  | 4.966 | ng/u1  | 98 |
| 30) Naphthalene               | 10.731 | 128 | 27763 | 4.892 | ng/u1  | 98 |
| 32) 4-Chloroaniline           | 10.843 | 127 | 11464 | 4.668 | ng/u1  | 94 |
| 33) Hexachlorobutadiene       | 11.019 | 225 | 8427  | 4.956 | ng/u1# | 84 |
| 34) Caprolactam               | 11.612 | 113 | 3782  | 5.682 | ng/u1  | 90 |
| 35) 4-Chloro-3-methylphenol   | 11.983 | 107 | 10647 | 4.734 | ng/u1  | 91 |
| 36) 2-Methylnaphthalene       | 12.347 | 142 | 19523 | 4.744 | ng/u1  | 95 |

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 Sample : P2409-01  
 Misc : SFAM-MDL-WATER-01  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-WATER-QT2-2024-03

Manual Integrations  
 APPROVED

Reviewed By :Yogesh  
 Patel

05/24/2024  
 Supervised By :mohammad  
 ahmed

Quant Time: May 22 23:48:02 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG052124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 21 15:23:39 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.564 | 142  | 20822    | 4.930 | ng/ul# | 85       |
| 39) 1,2,4,5-Tetrachloroben... | 12.717 | 216  | 14899    | 4.992 | ng/ul  | 93       |
| 40) Hexachlorocyclopentadiene | 12.693 | 237  | 5930     | 4.402 | ng/ul# | 94       |
| 41) 2,4,6-Trichlorophenol     | 12.964 | 196  | 7839     | 4.663 | ng/ul  | 91       |
| 42) 2,4,5-Trichlorophenol     | 13.040 | 196  | 8701     | 4.915 | ng/ul  | 95       |
| 43) 1,1'-Biphenyl             | 13.357 | 154  | 29701    | 5.081 | ng/ul  | 98       |
| 44) 2-Chloronaphthalene       | 13.393 | 162  | 22962    | 4.992 | ng/ul  | 94       |
| 45) 2-Nitroaniline            | 13.598 | 65   | 8962     | 4.721 | ng/ul  | 98       |
| 47) Dimethylphthalate         | 13.974 | 163  | 36045    | 4.933 | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 14.098 | 165  | 6904     | 4.818 | ng/ul  | 95       |
| 50) Acenaphthylene            | 14.245 | 152  | 41604    | 5.148 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.433 | 138  | 6661     | 5.356 | ng/ul  | 99       |
| 52) Acenaphthene              | 14.585 | 153  | 27275    | 5.144 | ng/ul  | 94       |
| 53) 2,4-Dinitrophenol         | 14.656 | 184  | 3294     | 4.335 | ng/ul  | 89       |
| 55) 4-Nitrophenol             | 14.762 | 109  | 5794     | 5.686 | ng/ul  | 91       |
| 56) Dibenzofuran              | 14.920 | 168  | 38407    | 5.067 | ng/ul  | 94       |
| 57) 2,4-Dinitrotoluene        | 14.891 | 165  | 10890    | 4.995 | ng/ul# | 96       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.161 | 232  | 6952     | 4.306 | ng/ul# | 97       |
| 59) Diethylphthalate          | 15.343 | 149  | 37312    | 5.171 | ng/ul  | 97       |
| 61) Fluorene                  | 15.572 | 166  | 34110    | 5.201 | ng/ul  | 93       |
| 62) 4-Chlorophenyl-phenyle... | 15.567 | 204  | 20092    | 5.223 | ng/ul  | 96       |
| 63) 4-Nitroaniline            | 15.590 | 138  | 6949     | 5.365 | ng/ul# | 75       |
| 66) 4,6-Dinitro-2-methylph... | 15.666 | 198  | 7756m    | 4.907 | ng/ul  |          |
| 67) N-Nitrosodiphenylamine    | 15.778 | 169  | 31050    | 4.954 | ng/ul  | 96       |
| 68) 4-Bromophenyl-phenylether | 16.460 | 248  | 13872    | 4.695 | ng/ul  | 95       |
| 69) Hexachlorobenzene         | 16.583 | 284  | 15101    | 4.598 | ng/ul  | 96       |
| 70) Atrazine                  | 16.730 | 200  | 15441    | 6.226 | ng/ul# | 96       |
| 71) Pentachlorophenol         | 16.941 | 266  | 5763m    | 5.042 | ng/ul  |          |
| 72) Phenanthrene              | 17.312 | 178  | 59890m   | 4.949 | ng/ul  |          |
| 74) Anthracene                | 17.406 | 178  | 56668    | 4.529 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.322 | 216  | 15147    | 4.596 | ng/ul  | 99       |
| 76) Pentachlorobenzene        | 14.850 | 250  | 17341    | 4.759 | ng/ul  | 92       |
| 77) Carbazole                 | 17.676 | 167  | 51422    | 5.052 | ng/ul  | 98       |
| 78) Di-n-butylphthalate       | 18.234 | 149  | 62958    | 4.937 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.327 | 202  | 78030    | 4.780 | ng/ul  | 99       |
| 82) Pyrene                    | 19.691 | 202  | 82658    | 4.878 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 20.578 | 149  | 28893    | 4.963 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.419 | 252  | 25502    | 4.533 | ng/ul# | 98       |
| 85) Benzo(a)anthracene        | 21.501 | 228  | 88933    | 4.863 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.419 | 149  | 42458    | 4.927 | ng/ul# | 99       |
| 87) Chrysene                  | 21.565 | 228  | 83118    | 4.936 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.582 | 149  | 71950    | 5.086 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.634 | 252  | 86528m   | 5.139 | ng/ul  |          |
| 91) Benzo(k)fluoranthene      | 23.692 | 252  | 85747    | 5.017 | ng/ul  | 95       |
| 93) Benzo(a)pyrene            | 24.462 | 252  | 80627    | 5.043 | ng/ul  | 95       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.105 | 276  | 95592    | 4.944 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 28.152 | 278  | 80169    | 5.029 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.198 | 276  | 77013    | 5.007 | ng/ul# | 96       |

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

