

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122624\  
 Data File : BG063912.D  
 Acq On : 28 Dec 2024 20:22  
 Operator : RC/JU  
 Sample : SP6699  
 Misc : SPIKE  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SP6699

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/30/2024  
 Supervised By :mohammad ahmed 12/30/2024

Quant Time: Dec 30 00:37:53 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 24 00:44:39 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.801  | 152  | 101467   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.586 | 136  | 408179   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.440 | 164  | 279233   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.201 | 188  | 637441   | 20.000 | ng/u1  | 0.01     |
| 79) Chrysene-d12              | 21.455 | 240  | 742222   | 20.000 | ng/u1  | 0.02     |
| 88) Perylene-d12              | 24.475 | 264  | 767517   | 20.000 | ng/u1  | 0.03     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000  | ng/uL  |          |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 0.000  | 99   | 0d       | 0.000  | ng/u1  |          |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000  | 67   | 0d       | 0.000  | ng/u1  |          |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000  | ng/u1  |          |
| 15) 4-Methylphenol-d8         | 0.000  | 113  | 0d       | 0.000  | ng/u1  |          |
| 21) Nitrobenzene-d5           | 0.000  | 128  | 0d       | 0.000  | ng/u1  |          |
| 24) 2-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165  | 0d       | 0.000  | ng/u1  |          |
| 31) 4-Chloroaniline-d4        | 0.000  | 131  | 0d       | 0.000  | ng/u1  |          |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.000  | ng/u1  |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.000  | ng/u1  |          |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000  | ng/u1  |          |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/u1  |          |
| 73) Anthracene-d10            | 0.000  | 188  | 0d       | 0.000  | ng/u1  |          |
| 81) Pyrene-d10                | 0.000  | 212  | 0d       | 0.000  | ng/u1  |          |
| 92) Benzo(a)pyrene-d12        | 0.000  | 264  | 0d       | 0.000  | ng/u1  |          |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.306  | 88   | 95956    | 31.315 | ng/uL# | 97       |
| 5) Pyridine                   | 3.705  | 79   | 728292   | 83.245 | ng/u1  | 87       |
| 6) Benzaldehyde               | 6.943  | 77   | 384680   | 67.109 | ng/u1  | 95       |
| 8) Phenol                     | 7.019  | 94   | 814340   | 81.765 | ng/u1  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.225  | 93   | 604811   | 79.371 | ng/u1  | 97       |
| 12) 2-Chlorophenol            | 7.372  | 128  | 545609   | 89.093 | ng/u1  | 100      |
| 13) 2-Methylphenol            | 8.259  | 108  | 589844   | 81.942 | ng/u1  | 96       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.318  | 45   | 882912   | 78.048 | ng/u1  | 98       |
| 16) Acetophenone              | 8.623  | 105  | 915698   | 74.410 | ng/u1  | 95       |
| 17) N-Nitroso-di-n-propyla... | 8.606  | 70   | 535547   | 72.935 | ng/u1  | 92       |
| 18) 4-Methylphenol            | 8.588  | 108  | 624603   | 79.282 | ng/u1  | 95       |
| 19) Hexachloroethane          | 8.870  | 117  | 230653   | 76.652 | ng/u1  | 97       |
| 22) Nitrobenzene              | 8.999  | 77   | 804608   | 82.674 | ng/u1  | 98       |
| 23) Isophorone                | 9.522  | 82   | 1456231  | 79.192 | ng/u1  | 99       |
| 25) 2-Nitrophenol             | 9.710  | 139  | 296791   | 88.398 | ng/u1  | 92       |
| 26) 2,4-Dimethylphenol        | 9.781  | 107  | 645788   | 84.764 | ng/u1  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.004 | 93   | 822591   | 82.706 | ng/u1  | 96       |
| 29) 2,4-Dichlorophenol        | 10.251 | 162  | 554230   | 88.186 | ng/u1  | 98       |
| 30) Naphthalene               | 10.638 | 128  | 1659765  | 81.330 | ng/u1  | 98       |
| 32) 4-Chloroaniline           | 10.750 | 127  | 495280   | 64.817 | ng/u1  | 96       |
| 33) Hexachlorobutadiene       | 10.920 | 225  | 395924   | 77.543 | ng/u1  | 97       |
| 34) Caprolactam               | 11.537 | 113  | 179597m  | 82.668 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 11.902 | 107  | 602065   | 81.597 | ng/u1  | 98       |
| 36) 2-Methylnaphthalene       | 12.254 | 142  | 1147122  | 78.210 | ng/u1  | 97       |

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 24 00:44:39 2024  
 Response via : Initial Calibration

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|-------------------------------|--------|------|----------|---------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.477 | 142  | 1157809  | 77.497  | ng/ul  | 95       |
| 39) 1,2,4,5-Tetrachloroben... | 12.630 | 216  | 722521   | 80.828  | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.607 | 237  | 218853   | 61.933  | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.877 | 196  | 468078   | 92.647  | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.959 | 196  | 482391   | 91.489  | ng/ul  | 96       |
| 43) 1,1'-Biphenyl             | 13.271 | 154  | 1539625  | 83.607  | ng/ul  | 97       |
| 44) 2-Chloronaphthalene       | 13.318 | 162  | 1257031  | 85.006  | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.529 | 65   | 464316   | 93.740  | ng/ul  | 93       |
| 47) Dimethylphthalate         | 13.899 | 163  | 1557165  | 81.772  | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 14.029 | 165  | 333446   | 94.285  | ng/ul# | 90       |
| 50) Acenaphthylene            | 14.164 | 152  | 1939770  | 83.523  | ng/ul  | 97       |
| 51) 3-Nitroaniline            | 14.358 | 138  | 289860   | 88.648  | ng/ul  | 94       |
| 52) Acenaphthene              | 14.510 | 153  | 1359798  | 82.607  | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 14.581 | 184  | 84989    | 47.024  | ng/ul  | 90       |
| 55) 4-Nitrophenol             | 14.698 | 109  | 203544   | 72.813  | ng/ul  | 91       |
| 56) Dibenzofuran              | 14.845 | 168  | 1908153  | 82.222  | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.822 | 165  | 478226   | 91.003  | ng/ul  | 93       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.080 | 232  | 398901   | 87.766  | ng/ul  | 96       |
| 59) Diethylphthalate          | 15.268 | 149  | 1496215  | 81.276  | ng/ul  | 98       |
| 61) Fluorene                  | 15.497 | 166  | 1451176  | 78.310  | ng/ul  | 97       |
| 62) 4-Chlorophenyl-phenyle... | 15.492 | 204  | 792292   | 76.743  | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 15.533 | 138  | 338173   | 103.307 | ng/ul  | 93       |
| 66) 4,6-Dinitro-2-methylph... | 15.597 | 198  | 197822   | 58.019  | ng/ul# | 89       |
| 67) N-Nitrosodiphenylamine    | 15.709 | 169  | 1306945  | 84.669  | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.385 | 248  | 549040   | 84.515  | ng/ul  | 100      |
| 69) Hexachlorobenzene         | 16.514 | 284  | 604209   | 83.013  | ng/ul  | 97       |
| 70) Atrazine                  | 16.667 | 200  | 566692   | 86.849  | ng/ul  | 96       |
| 71) Pentachlorophenol         | 16.861 | 266  | 255022   | 83.101  | ng/ul  | 98       |
| 72) Phenanthrene              | 17.242 | 178  | 2569599  | 83.651  | ng/ul  | 99       |
| 74) Anthracene                | 17.331 | 178  | 2596737  | 83.521  | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.241 | 216  | 715503   | 84.129  | ng/ul  | 100      |
| 76) Pentachlorobenzene        | 14.769 | 250  | 711584   | 84.455  | ng/ul  | 99       |
| 77) Carbazole                 | 17.607 | 167  | 2343581  | 93.121  | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.165 | 149  | 2630056  | 86.532  | ng/ul  | 99       |
| 80) Fluoranthene              | 19.264 | 202  | 3173896  | 72.300  | ng/ul  | 96       |
| 82) Pyrene                    | 19.628 | 202  | 3366801  | 72.086  | ng/ul# | 91       |
| 83) Butylbenzylphthalate      | 20.521 | 149  | 1275774  | 88.983  | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.355 | 252  | 990071   | 73.893  | ng/ul  | 93       |
| 85) Benzo(a)anthracene        | 21.438 | 228  | 3534771  | 75.991  | ng/ul  | 96       |
| 86) Bis(2-ethylhexyl)phtha... | 21.344 | 149  | 1847162  | 88.568  | ng/ul  | 99       |
| 87) Chrysene                  | 21.502 | 228  | 3379720  | 76.233  | ng/ul  | 93       |
| 89) Di-n-octyl phthalate      | 22.483 | 149  | 3294129  | 99.437  | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.529 | 252  | 3736267  | 85.651  | ng/ul  | 100      |
| 91) Benzo(k)fluoranthene      | 23.594 | 252  | 3617831  | 83.144  | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.340 | 252  | 3500264  | 85.487  | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.906 | 276  | 4446335  | 88.920  | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 27.959 | 278  | 3572405  | 89.286  | ng/ul  | 95       |
| 96) Benzo(g,h,i)perylene      | 28.976 | 276  | 3483116  | 87.782  | ng/ul  | 97       |

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

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 ALS Vial : 40 Sample Multiplier: 1

## Instrument :

BNA\_G

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Quant Time: Dec 30 00:37:53 2024

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Quant Title : SVOA CALIBRATION

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