

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG103124\  
 Data File : BG063131.D  
 Acq On : 31 Oct 2024 13:30  
 Operator : RC/JU  
 Sample : SSTDICC010  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA\_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 11/04/2024

Supervised By :mohammad ahmed 11/04/2024

Quant Time: Nov 01 04:53:35 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG103124.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 01 04:41:50 2024

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.849  | 152  | 89789    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.634 | 136  | 360846   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 14.477 | 164  | 290744   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.226 | 188  | 754489   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.480 | 240  | 841662   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 24.518 | 264  | 963367   | 20.000 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.440  | 112  | 112853   | 19.489 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.015  | 99   | 159830   | 21.063 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 8.995  | 82   | 166029   | 20.011 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.969 | 330  | 97569    | 19.756 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.096 | 172  | 397885   | 20.269 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.858 | 244  | 946066   | 23.131 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.360  | 88   | 23519    | 10.339 | ng    | 92       |
| 3) Pyridine                   | 3.748  | 79   | 58738    | 10.694 | ng    | 87       |
| 4) n-Nitrosodimethylamine     | 3.666  | 42   | 35618    | 10.440 | ng    | # 93     |
| 6) Aniline                    | 7.173  | 93   | 74056    | 10.852 | ng    | # 92     |
| 8) 2-Chlorophenol             | 7.408  | 128  | 55622    | 10.343 | ng    | 96       |
| 9) Benzaldehyde               | 6.985  | 77   | 53377    | 11.320 | ng    | # 77     |
| 10) Phenol                    | 7.044  | 94   | 84650    | 10.413 | ng    | 97       |
| 11) bis(2-Chloroethyl)ether   | 7.267  | 93   | 56896    | 10.378 | ng    | 88       |
| 12) 1,3-Dichlorobenzene       | 7.737  | 146  | 68637    | 10.508 | ng    | 100      |
| 13) 1,4-Dichlorobenzene       | 7.878  | 146  | 69035    | 10.380 | ng    | 92       |
| 14) 1,2-Dichlorobenzene       | 8.196  | 146  | 67809    | 10.577 | ng    | 96       |
| 15) Benzyl Alcohol            | 8.078  | 79   | 56615    | 9.285  | ng    | 95       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.372  | 45   | 106362   | 10.695 | ng    | 96       |
| 17) 2-Methylphenol            | 8.290  | 107  | 50374    | 9.741  | ng    | 95       |
| 18) Hexachloroethane          | 8.918  | 117  | 25327    | 10.641 | ng    | 91       |
| 19) n-Nitroso-di-n-propyla... | 8.642  | 70   | 57317    | 10.846 | ng    | 98       |
| 20) 3+4-Methylphenols         | 8.613  | 107  | 70965    | 10.262 | ng    | 95       |
| 22) Acetophenone              | 8.660  | 105  | 100160   | 10.359 | ng    | # 98     |
| 24) Nitrobenzene              | 9.042  | 77   | 90687    | 10.109 | ng    | # 93     |
| 25) Isophorone                | 9.559  | 82   | 145228   | 10.120 | ng    | 96       |
| 26) 2-Nitrophenol             | 9.758  | 139  | 32029    | 9.707  | ng    | # 85     |
| 27) 2,4-Dimethylphenol        | 9.811  | 122  | 41019    | 10.113 | ng    | 95       |
| 28) bis(2-Chloroethoxy)met... | 10.046 | 93   | 75326    | 10.010 | ng    | 94       |
| 29) 2,4-Dichlorophenol        | 10.287 | 162  | 60370    | 9.854  | ng    | 94       |
| 30) 1,2,4-Trichlorobenzene    | 10.499 | 180  | 72800    | 9.661  | ng    | 95       |
| 31) Naphthalene               | 10.687 | 128  | 192167   | 10.095 | ng    | 97       |
| 32) Benzoic acid              | 9.917  | 122  | 25334m   | 7.150  | ng    |          |
| 33) 4-Chloroaniline           | 10.793 | 127  | 62750    | 9.768  | ng    | # 89     |
| 34) Hexachlorobutadiene       | 10.969 | 225  | 60224    | 9.711  | ng    | 94       |
| 35) Caprolactam               | 11.539 | 113  | 20699m   | 10.579 | ng    |          |
| 36) 4-Chloro-3-methylphenol   | 11.921 | 107  | 68858    | 9.756  | ng    | # 89     |
| 37) 2-Methylnaphthalene       | 12.297 | 142  | 140518   | 10.407 | ng    | 91       |
| 38) 1-Methylnaphthalene       | 12.514 | 142  | 138478   | 10.419 | ng    | 97       |
| 40) 1,2,4,5-Tetrachloroben... | 12.667 | 216  | 103206   | 9.878  | ng    | 97       |
| 41) Hexachlorocyclopentadiene | 12.643 | 237  | 26682    | 7.860  | ng    | 90       |
| 43) 2,4,6-Trichlorophenol     | 12.908 | 196  | 61097m   | 9.766  | ng    |          |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG103124\  
 Data File : BG063131.D  
 Acq On : 31 Oct 2024 13:30  
 Operator : RC/JU  
 Sample : SSTDICC010  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTDICC010

**Manual Integrations****APPROVED**

Reviewed By :Yogesh Patel 11/04/2024

Supervised By :mohammad ahmed 11/04/2024

Quant Time: Nov 01 04:53:35 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG103124.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 01 04:41:50 2024

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 12.984 | 196  | 67525    | 9.818  | ng    | 91       |
| 46) 1,1'-Biphenyl             | 13.313 | 154  | 197252   | 9.996  | ng    | 97       |
| 47) 2-Chloronaphthalene       | 13.354 | 162  | 162021   | 9.804  | ng    | 98       |
| 48) 2-Nitroaniline            | 13.554 | 65   | 56607    | 9.759  | ng    | 94       |
| 49) Acenaphthylene            | 14.200 | 152  | 241251   | 10.340 | ng    | 98       |
| 50) Dimethylphthalate         | 13.930 | 163  | 212675   | 10.263 | ng    | 98       |
| 51) 2,6-Dinitrotoluene        | 14.053 | 165  | 45583    | 10.303 | ng    | # 90     |
| 52) Acenaphthene              | 14.541 | 154  | 146141   | 10.118 | ng    | 99       |
| 53) 3-Nitroaniline            | 14.383 | 138  | 45206    | 10.762 | ng    | # 94     |
| 54) 2,4-Dinitrophenol         | 14.600 | 184  | 18392    | 6.816  | ng    | 85       |
| 55) Dibenzofuran              | 14.876 | 168  | 256061   | 10.054 | ng    | 95       |
| 56) 4-Nitrophenol             | 14.706 | 139  | 31283    | 9.837  | ng    | # 87     |
| 57) 2,4-Dinitrotoluene        | 14.847 | 165  | 63276    | 10.000 | ng    | # 93     |
| 58) Fluorene                  | 15.528 | 166  | 206200   | 10.217 | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.111 | 232  | 67516    | 10.230 | ng    | 98       |
| 60) Diethylphthalate          | 15.299 | 149  | 223359   | 10.498 | ng    | 98       |
| 61) 4-Chlorophenyl-phenyle... | 15.522 | 204  | 127781   | 10.186 | ng    | 96       |
| 62) 4-Nitroaniline            | 15.546 | 138  | 46318    | 9.995  | ng    | # 81     |
| 63) Azobenzene                | 15.816 | 77   | 222034   | 10.089 | ng    | 98       |
| 65) 4,6-Dinitro-2-methylph... | 15.610 | 198  | 43947    | 9.060  | ng    | 94       |
| 66) n-Nitrosodiphenylamine    | 15.734 | 169  | 182337   | 9.687  | ng    | 94       |
| 67) 4-Bromophenyl-phenylether | 16.415 | 248  | 89382    | 9.667  | ng    | 94       |
| 68) Hexachlorobenzene         | 16.533 | 284  | 102895   | 9.606  | ng    | # 92     |
| 69) Atrazine                  | 16.686 | 200  | 77354    | 8.940  | ng    | 97       |
| 70) Pentachlorophenol         | 16.885 | 266  | 44895    | 8.424  | ng    | 98       |
| 71) Phenanthrene              | 17.267 | 178  | 369189   | 10.094 | ng    | 97       |
| 72) Anthracene                | 17.361 | 178  | 358597   | 9.989  | ng    | 99       |
| 73) Carbazole                 | 17.632 | 167  | 338372   | 10.013 | ng    | 98       |
| 74) Di-n-butylphthalate       | 18.196 | 149  | 384698   | 10.026 | ng    | 99       |
| 75) Fluoranthene              | 19.289 | 202  | 520487   | 10.053 | ng    | 98       |
| 77) Benzidine                 | 19.471 | 184  | 187631   | 14.469 | ng    | 99       |
| 78) Pyrene                    | 19.653 | 202  | 542266   | 10.869 | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.552 | 149  | 158932   | 9.887  | ng    | 94       |
| 81) Benzo(a)anthracene        | 21.462 | 228  | 537433   | 10.140 | ng    | 96       |
| 82) 3,3'-Dichlorobenzidine    | 21.386 | 252  | 204564   | 10.410 | ng    | 97       |
| 83) Chrysene                  | 21.527 | 228  | 515379   | 10.351 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.380 | 149  | 227572   | 9.759  | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.526 | 149  | 381622   | 9.781  | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene    | 27.931 | 276  | 631090   | 9.760  | ng    | # 97     |
| 88) Benzo(b)fluoranthene      | 23.560 | 252  | 559454   | 9.961  | ng    | 95       |
| 89) Benzo(k)fluoranthene      | 23.613 | 252  | 522236   | 9.746  | ng    | 99       |
| 90) Benzo(a)pyrene            | 24.371 | 252  | 495998   | 10.117 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene    | 27.984 | 278  | 536750   | 9.945  | ng    | # 97     |
| 92) Benzo(g,h,i)perylene      | 29.001 | 276  | 532054   | 9.973  | ng    | 93       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG103124\  
Data File : BG063131.D  
Acq On : 31 Oct 2024 13:30  
Operator : RC/JU  
Sample : SSTDICC010  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTDICC010

Quant Time: Nov 01 04:53:35 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG103124.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Fri Nov 01 04:41:50 2024  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Reviewed By :Yogesh Patel 11/04/2024  
Supervised By :mohammad ahmed 11/04/2024

