

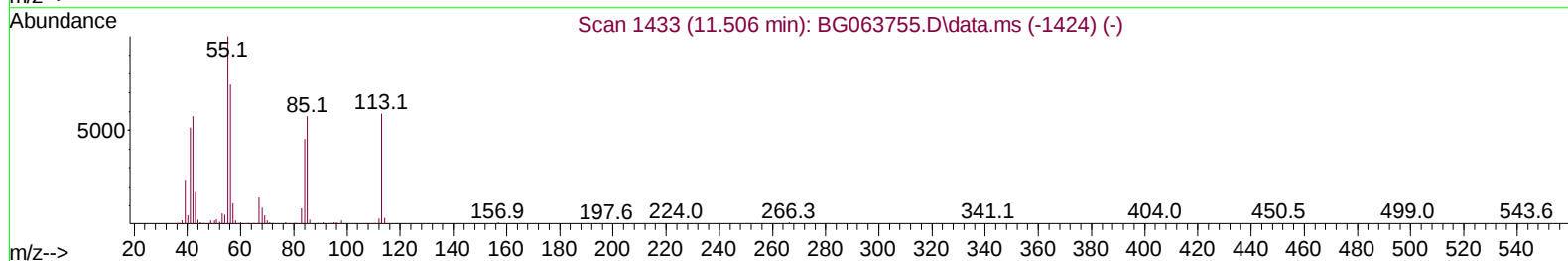
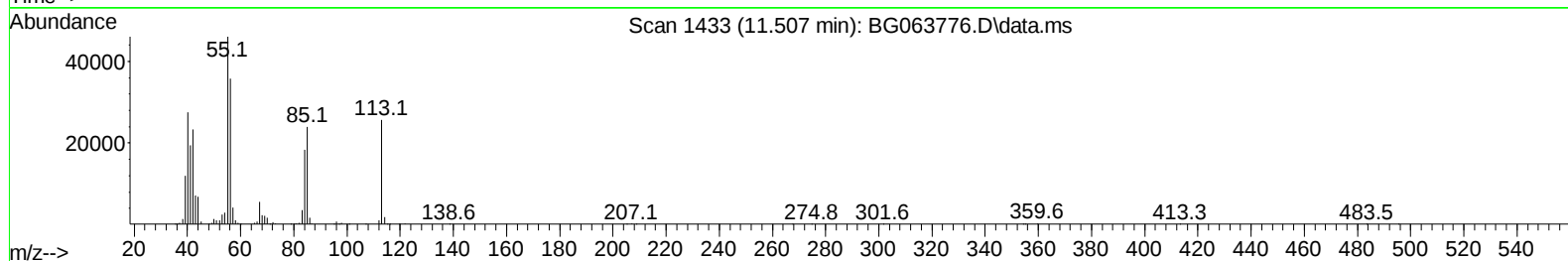
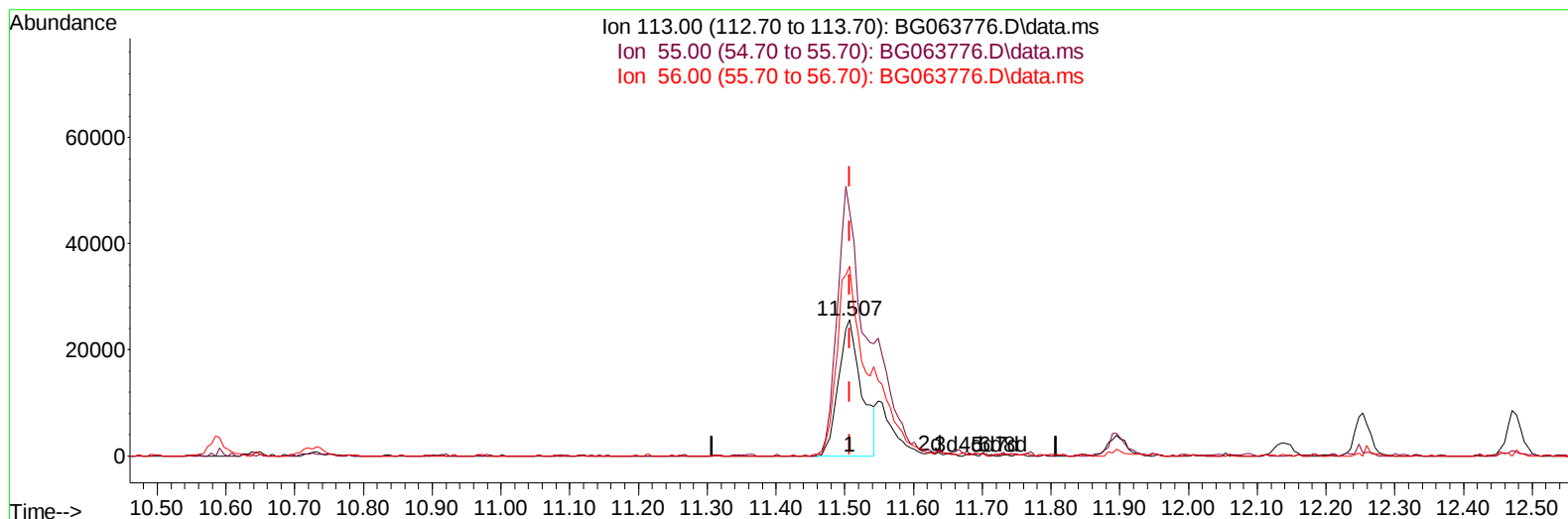
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
Data File : BG063776.D  
Acq On : 20 Dec 2024 17:51  
Operator : RC/JU  
Sample : PB165768BS  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:12:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Dec 11 23:59:23 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
Supervised By :mohammad ahmed 12/24/2024



TIC: BG063776.D\data.ms

(34) Caprolactam

11.507min (-0.000) 24.89 ng/ul

response 60382

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 202.90 | 179.56 |
| 56.00  | 133.90 | 139.34 |
| 0.00   | 0.00   | 0.00   |

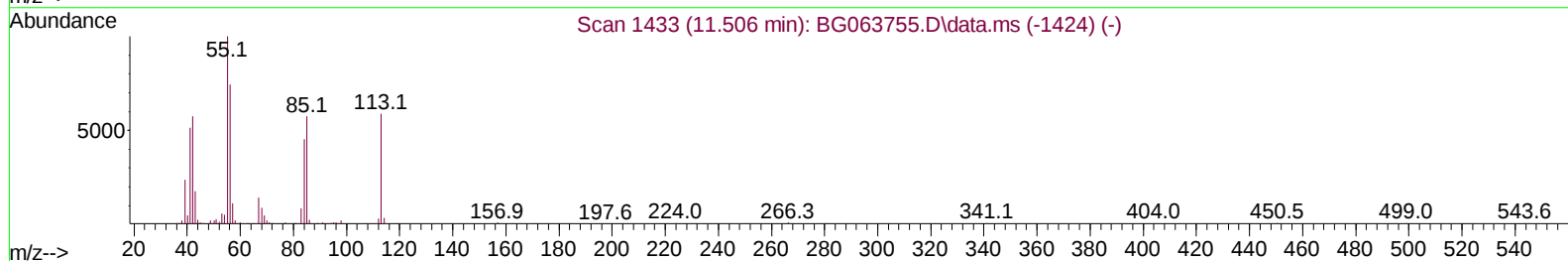
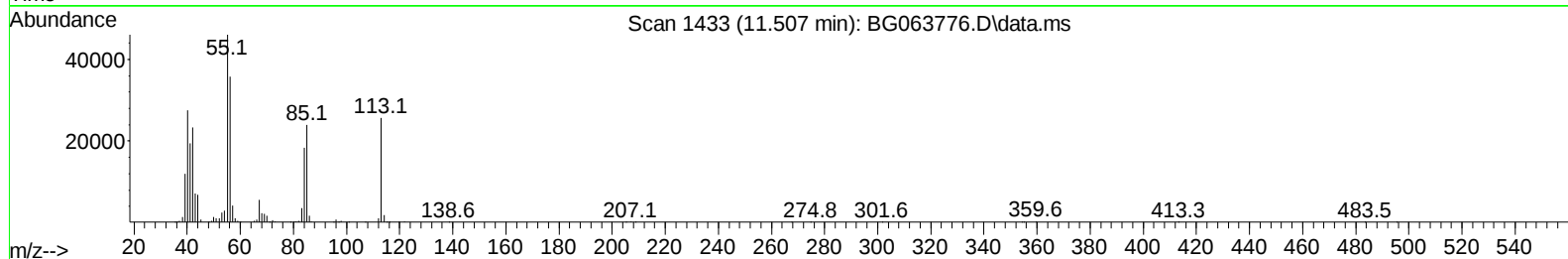
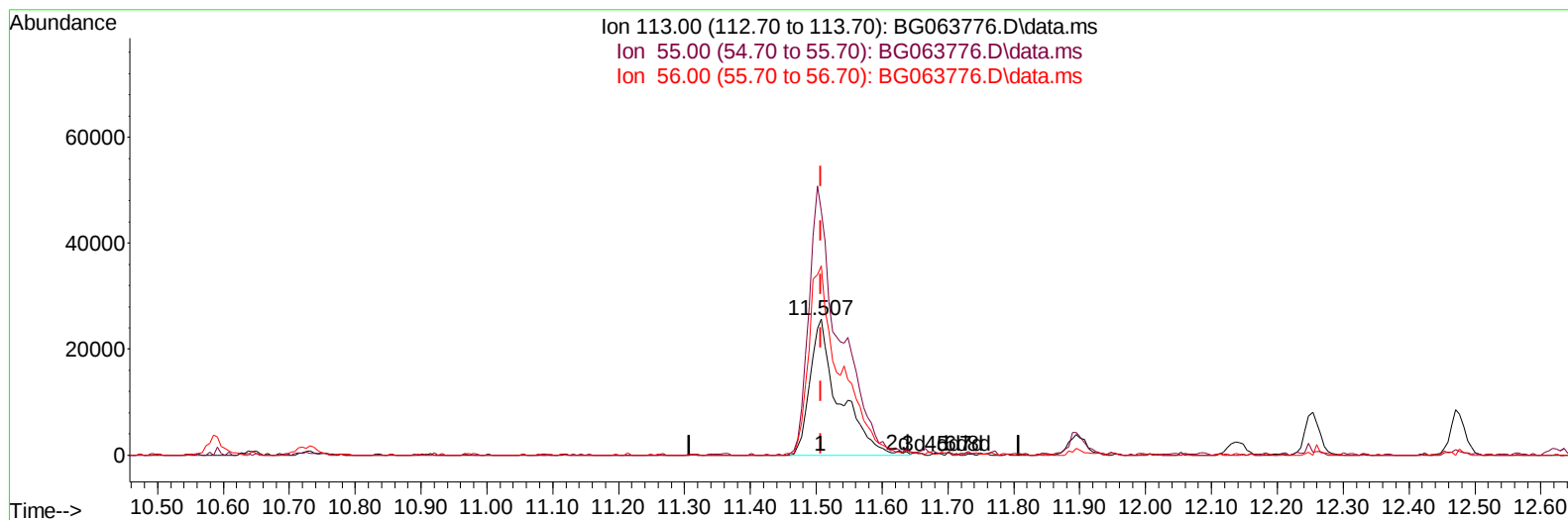
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
Data File : BG063776.D  
Acq On : 20 Dec 2024 17:51  
Operator : RC/JU  
Sample : PB165768BS  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:12:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Dec 11 23:59:23 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
Supervised By :mohammad ahmed 12/24/2024



TIC: BG063776.D\data.ms

(34) Caprolactam

11.507min (-0.000) 32.51 ng/ul m

response 78861

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 202.90 | 179.56 |
| 56.00  | 133.90 | 139.34 |
| 0.00   | 0.00   | 0.00   |

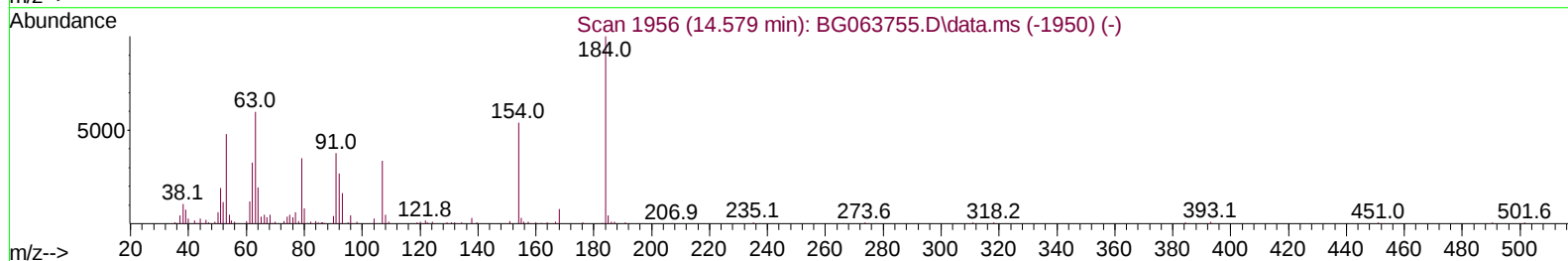
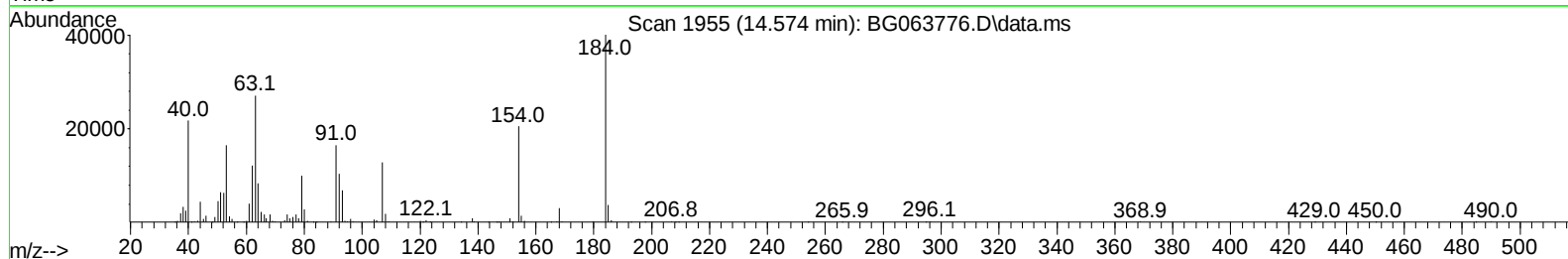
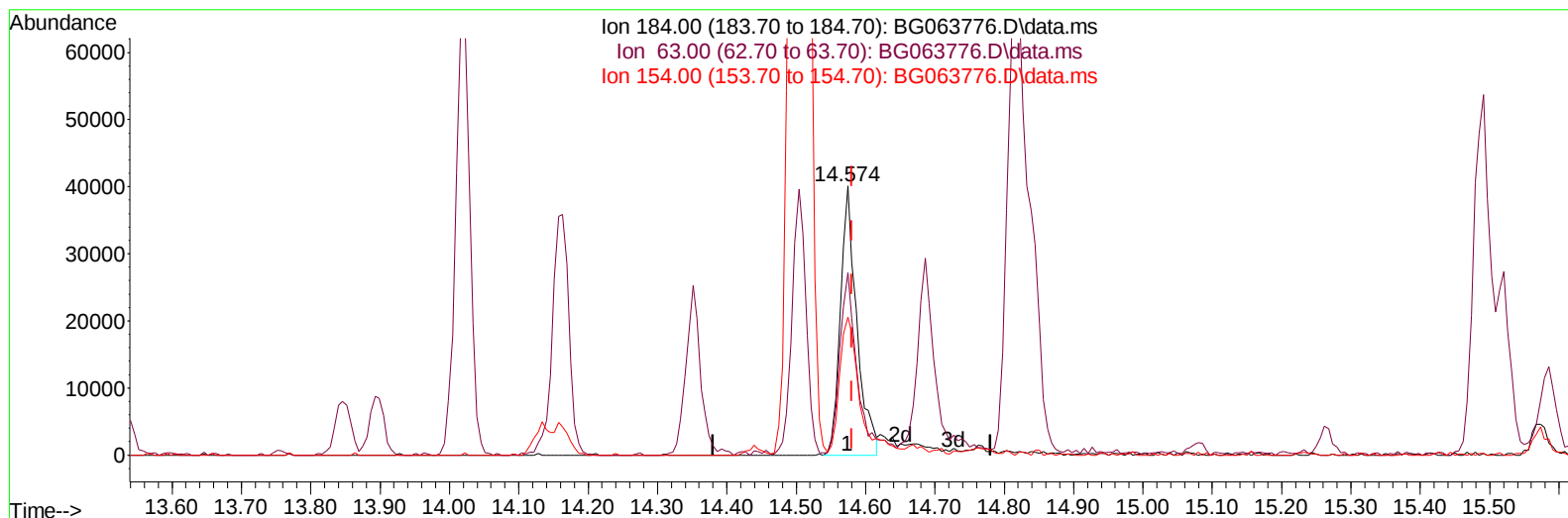
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
Data File : BG063776.D  
Acq On : 20 Dec 2024 17:51  
Operator : RC/JU  
Sample : PB165768BS  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:12:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Dec 11 23:59:23 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
Supervised By :mohammad ahmed 12/24/2024



TIC: BG063776.D\data.ms

(53) 2,4-Dinitrophenol

14.574min (-0.006) 30.24 ng/ul

response 65813

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 65.10  | 67.72  |
| 154.00 | 59.70  | 51.32  |
| 0.00   | 0.00   | 0.00   |

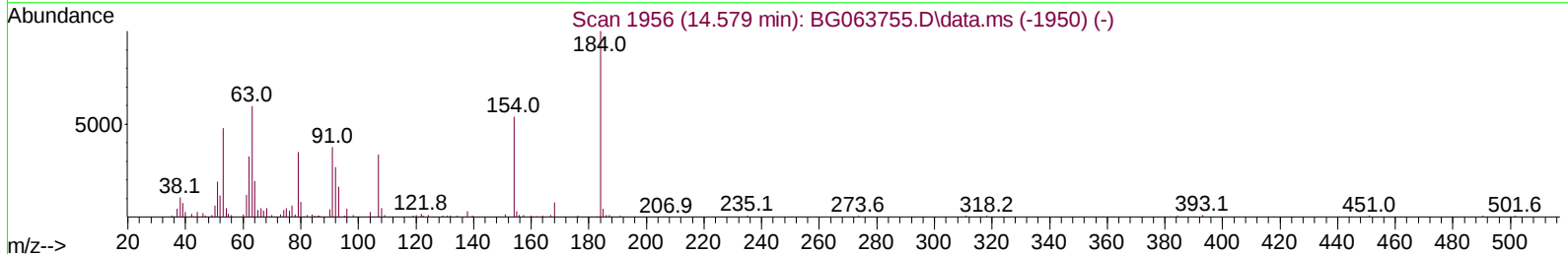
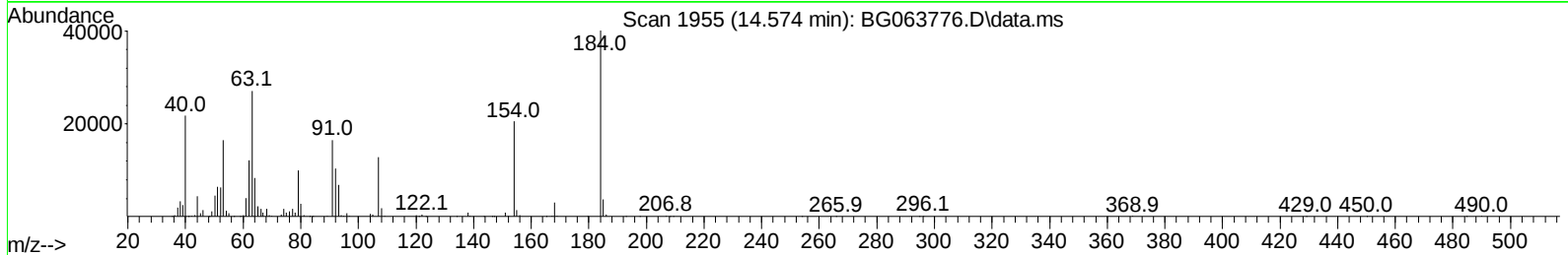
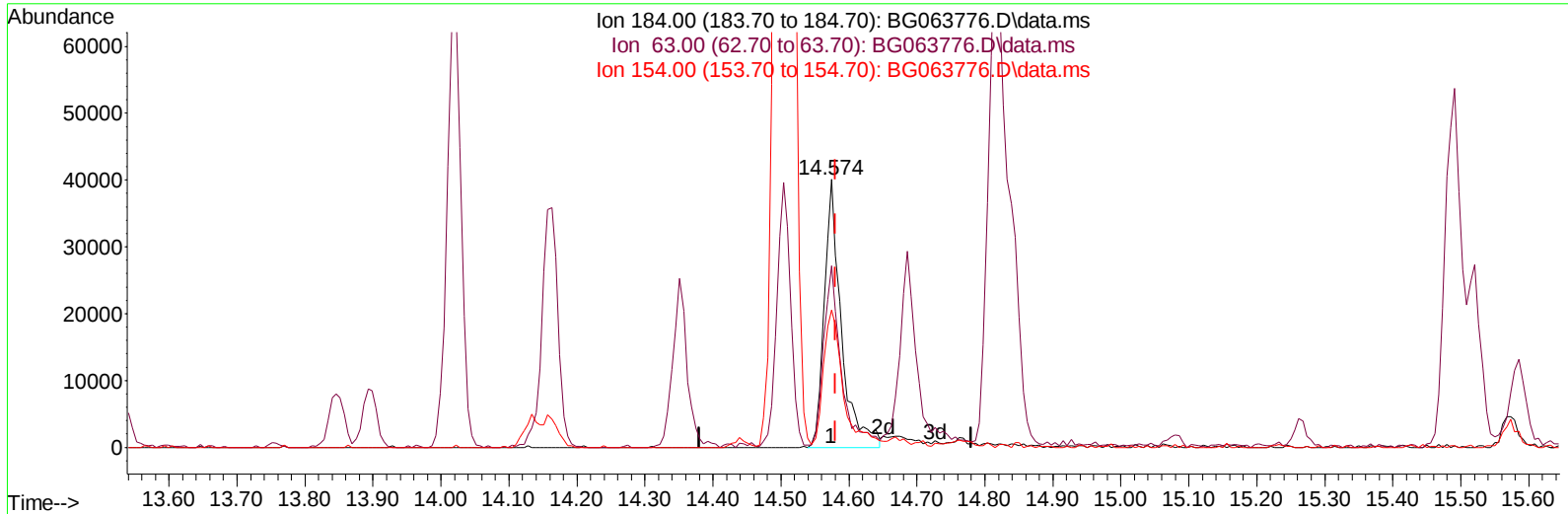
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
Data File : BG063776.D  
Acq On : 20 Dec 2024 17:51  
Operator : RC/JU  
Sample : PB165768BS  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:12:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Dec 11 23:59:23 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
Supervised By :mohammad ahmed 12/24/2024



TIC: BG063776.D\data.ms

(53) 2,4-Dinitrophenol

14.574min (-0.006) 32.16 ng/ul m

response 70002

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 65.10  | 67.72  |
| 154.00 | 59.70  | 51.32  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
 Data File : BG063776.D  
 Acq On : 20 Dec 2024 17:51  
 Operator : RC/JU  
 Sample : PB165768BS  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:14:02 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 11 23:59:23 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
 Supervised By :mohammad ahmed 12/24/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.800  | 152  | 113313   | 20.000 | ng/ul  | -0.01    |
| 20) Naphthalene-d8            | 10.591 | 136  | 455719   | 20.000 | ng/ul  | -0.01    |
| 38) Acenaphthene-d10          | 14.439 | 164  | 336245   | 20.000 | ng/ul  | -0.01    |
| 64) Phenanthrene-d10          | 17.195 | 188  | 895319   | 20.000 | ng/ul  | -0.01    |
| 79) Chrysene-d12              | 21.443 | 240  | 1021658  | 20.000 | ng/ul  | -0.02    |
| 88) Perylene-d12              | 24.457 | 264  | 1093163  | 20.000 | ng/ul  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.258  | 96   | 15722    | 5.437  | ng/uL  | -0.02    |
| 4) Pyridine-d5                | 3.675  | 84   | 254581   | 27.910 | ng/ul  | -0.01    |
| 7) Phenol-d5                  | 6.989  | 99   | 307354   | 29.451 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.130  | 67   | 191665   | 26.591 | ng/ul  | -0.02    |
| 11) 2-Chlorophenol-d4         | 7.342  | 132  | 204309   | 30.240 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.523  | 113  | 231446   | 28.565 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.957  | 128  | 97291    | 29.900 | ng/ul  | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.674  | 143  | 112899   | 30.953 | ng/ul  | -0.02    |
| 28) 2,4-Dichlorophenol-d3     | 10.221 | 165  | 222319   | 31.482 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.726 | 131  | 241904   | 26.979 | ng/ul  | -0.02    |
| 46) Dimethylphthalate-d6      | 13.846 | 166  | 686223   | 29.992 | ng/ul  | -0.01    |
| 49) Acenaphthylene-d8         | 14.134 | 160  | 793865   | 30.437 | ng/ul  | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.674 | 143  | 110227   | 34.002 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.438 | 176  | 647040   | 31.985 | ng/ul  | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.573 | 200  | 137327   | 29.542 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.295 | 188  | 1195863  | 31.067 | ng/ul  | -0.01    |
| 81) Pyrene-d10                | 19.592 | 212  | 1535234  | 28.877 | ng/ul  | -0.01    |
| 92) Benzo(a)pyrene-d12        | 24.245 | 264  | 1525839  | 29.723 | ng/ul  | -0.03    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.299  | 88   | 35842    | 10.474 | ng/ul# | 94       |
| 5) Pyridine                   | 3.693  | 79   | 280041   | 28.663 | ng/ul  | 89       |
| 6) Benzaldehyde               | 6.942  | 77   | 22459    | 3.508  | ng/ul  | 91       |
| 8) Phenol                     | 7.018  | 94   | 346695   | 31.171 | ng/ul  | 92       |
| 10) Bis(2-Chloroethyl)ether   | 7.224  | 93   | 238062   | 27.975 | ng/ul  | 96       |
| 12) 2-Chlorophenol            | 7.371  | 128  | 216977   | 31.726 | ng/ul  | 95       |
| 13) 2-Methylphenol            | 8.252  | 108  | 231062   | 28.744 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.317  | 45   | 341541   | 27.035 | ng/ul  | 98       |
| 16) Acetophenone              | 8.617  | 105  | 365818   | 26.619 | ng/ul  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.605  | 70   | 211893   | 25.840 | ng/ul  | 91       |
| 18) 4-Methylphenol            | 8.581  | 108  | 262541   | 29.841 | ng/ul  | 100      |
| 19) Hexachloroethane          | 8.875  | 117  | 89526    | 26.641 | ng/ul  | 95       |
| 22) Nitrobenzene              | 8.999  | 77   | 309971   | 28.527 | ng/ul  | 94       |
| 23) Isophorone                | 9.516  | 82   | 568459   | 27.689 | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.709  | 139  | 117868   | 31.444 | ng/ul  | 91       |
| 26) 2,4-Dimethylphenol        | 9.780  | 107  | 261481   | 30.741 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.003 | 93   | 316821   | 28.531 | ng/ul# | 95       |
| 29) 2,4-Dichlorophenol        | 10.250 | 162  | 232276   | 33.103 | ng/ul  | 96       |
| 30) Naphthalene               | 10.638 | 128  | 665096   | 29.191 | ng/ul  | 97       |
| 32) 4-Chloroaniline           | 10.755 | 127  | 160024   | 18.758 | ng/ul  | 95       |
| 33) Hexachlorobutadiene       | 10.926 | 225  | 152673   | 26.782 | ng/ul  | 97       |
| 34) Caprolactam               | 11.507 | 113  | 78861m   | 32.513 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.895 | 107  | 253322   | 30.751 | ng/ul  | 95       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122024\  
 Data File : BG063776.D  
 Acq On : 20 Dec 2024 17:51  
 Operator : RC/JU  
 Sample : PB165768BS  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLCS768

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:14:02 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 11 23:59:23 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/23/2024  
 Supervised By :mohammad ahmed 12/24/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.254 | 142  | 477049   | 29.132 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene       | 12.477 | 142  | 481408   | 28.861 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.630 | 216  | 299924   | 27.863 | ng/ul  | 96       |
| 40) Hexachlorocyclopentadiene | 12.606 | 237  | 72057    | 16.934 | ng/ul  | 93       |
| 41) 2,4,6-Trichlorophenol     | 12.876 | 196  | 197199   | 32.414 | ng/ul  | 95       |
| 42) 2,4,5-Trichlorophenol     | 12.953 | 196  | 215668   | 33.968 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.270 | 154  | 663815   | 29.936 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.311 | 162  | 542354   | 30.458 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.523 | 65   | 200191   | 33.563 | ng/ul  | 93       |
| 47) Dimethylphthalate         | 13.899 | 163  | 706575   | 30.814 | ng/ul  | 96       |
| 48) 2,6-Dinitrotoluene        | 14.022 | 165  | 141398   | 33.202 | ng/ul# | 92       |
| 50) Acenaphthylene            | 14.163 | 152  | 870877   | 31.140 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.351 | 138  | 148134   | 37.622 | ng/ul  | 93       |
| 52) Acenaphthene              | 14.504 | 153  | 607662   | 30.656 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol         | 14.574 | 184  | 70002m   | 32.165 | ng/ul  |          |
| 55) 4-Nitrophenol             | 14.686 | 109  | 101776   | 30.235 | ng/ul  | 90       |
| 56) Dibenzofuran              | 14.839 | 168  | 888521   | 31.794 | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 14.815 | 165  | 221870   | 35.062 | ng/ul# | 92       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.080 | 232  | 186317   | 34.043 | ng/ul  | 95       |
| 59) Diethylphthalate          | 15.262 | 149  | 722199   | 32.579 | ng/ul  | 98       |
| 61) Fluorene                  | 15.491 | 166  | 736932   | 33.025 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.485 | 204  | 391487   | 31.491 | ng/ul  | 94       |
| 63) 4-Nitroaniline            | 15.520 | 138  | 176932   | 44.886 | ng/ul  | 91       |
| 66) 4,6-Dinitro-2-methylph... | 15.591 | 198  | 147537   | 30.808 | ng/ul# | 87       |
| 67) N-Nitrosodiphenylamine    | 15.702 | 169  | 648871   | 29.929 | ng/ul  | 96       |
| 68) 4-Bromophenyl-phenylether | 16.384 | 248  | 270261   | 29.619 | ng/ul  | 97       |
| 69) Hexachlorobenzene         | 16.501 | 284  | 302655   | 29.605 | ng/ul  | 97       |
| 70) Atrazine                  | 16.654 | 200  | 283092   | 30.889 | ng/ul  | 96       |
| 71) Pentachlorophenol         | 16.854 | 266  | 123899   | 28.745 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.236 | 178  | 1393662  | 32.302 | ng/ul  | 99       |
| 74) Anthracene                | 17.324 | 178  | 1414192  | 32.384 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.235 | 216  | 304887   | 25.523 | ng/ul  | 98       |
| 76) Pentachlorobenzene        | 14.768 | 250  | 316573   | 26.751 | ng/ul  | 99       |
| 77) Carbazole                 | 17.600 | 167  | 1307554  | 36.990 | ng/ul  | 96       |
| 78) Di-n-butylphthalate       | 18.158 | 149  | 1489083  | 34.881 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.257 | 202  | 1798057  | 29.756 | ng/ul  | 97       |
| 82) Pyrene                    | 19.621 | 202  | 1936446  | 30.121 | ng/ul  | 95       |
| 83) Butylbenzylphthalate      | 20.514 | 149  | 677378   | 34.324 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.343 | 252  | 558333   | 30.273 | ng/ul  | 95       |
| 85) Benzo(a)anthracene        | 21.425 | 228  | 1910135  | 29.833 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.337 | 149  | 986518   | 34.364 | ng/ul  | 100      |
| 87) Chrysene                  | 21.484 | 228  | 1845194  | 30.237 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.465 | 149  | 1684169  | 35.694 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.505 | 252  | 1894448  | 30.491 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.564 | 252  | 1897868  | 30.623 | ng/ul  | 100      |
| 93) Benzo(a)pyrene            | 24.316 | 252  | 1751024  | 30.026 | ng/ul  | 96       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.865 | 276  | 2166917  | 30.426 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 27.906 | 278  | 1757729  | 30.844 | ng/ul  | 94       |
| 96) Benzo(g,h,i)perylene      | 28.916 | 276  | 1697569  | 30.038 | ng/ul  | 98       |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

SLCS768

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/23/2024  
Supervised By :mohammad ahmed 12/24/2024

