

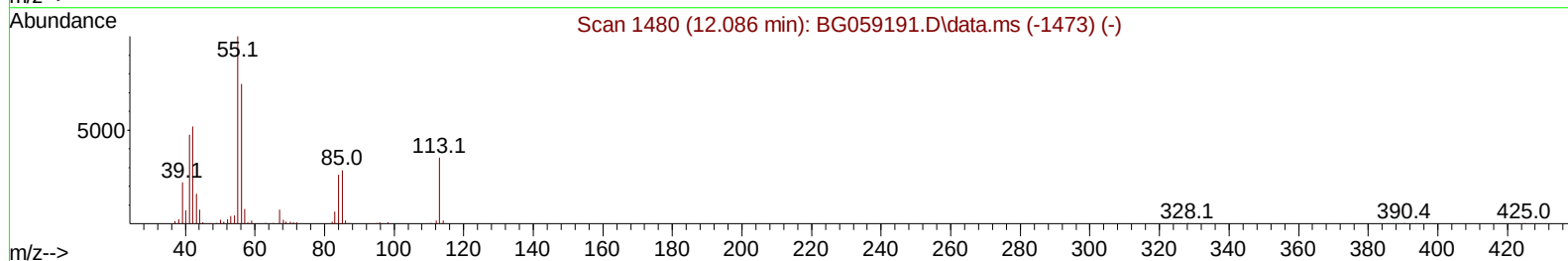
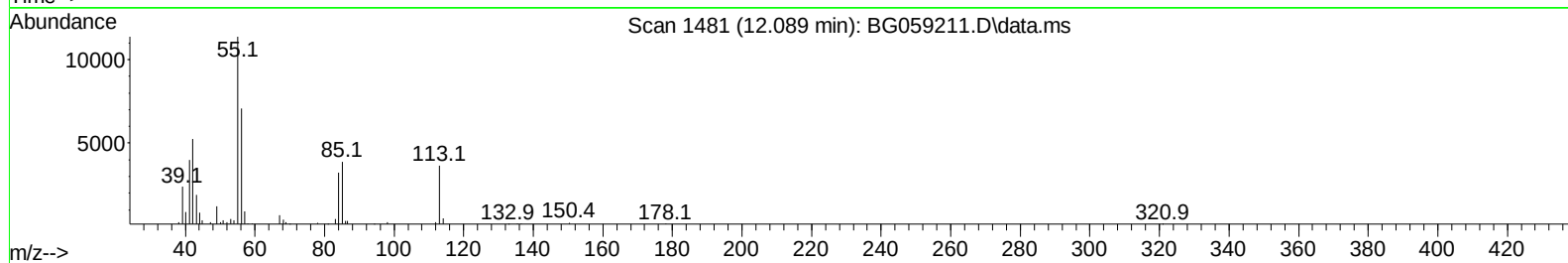
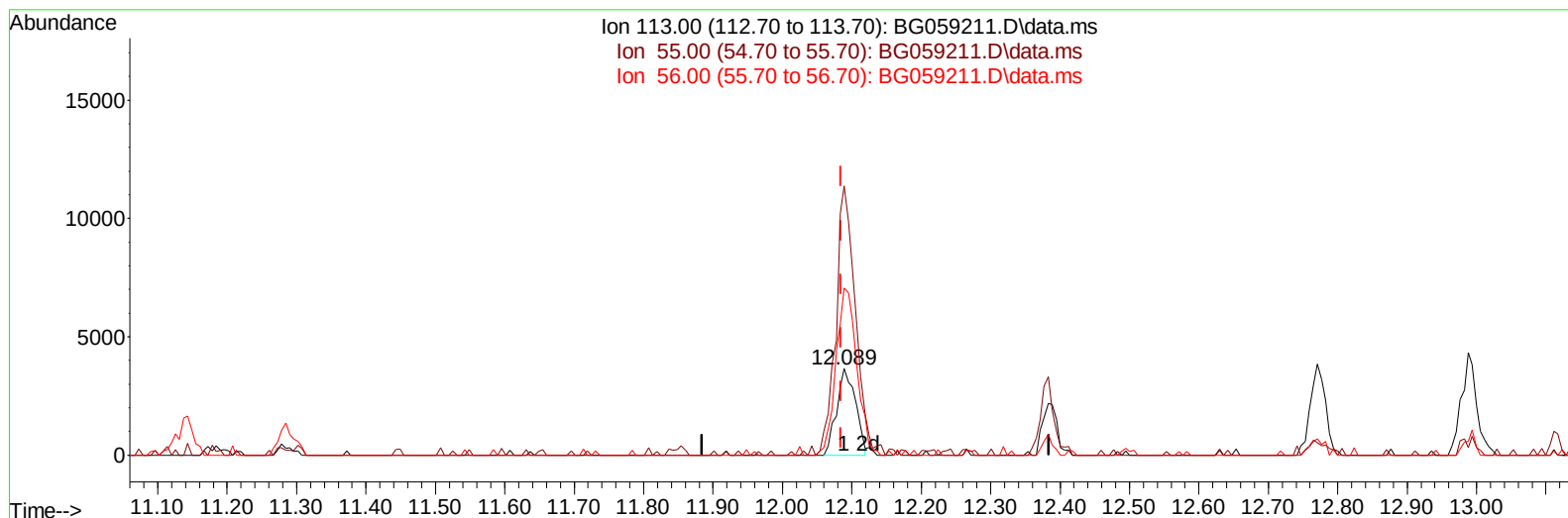
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(34) Caprolactam**

**12.089min (+ 0.005) 7.75 ng/ul**

**response 6861**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 197.00 | 311.34# |
| 56.00  | 140.40 | 193.59# |
| 0.00   | 0.00   | 0.00    |

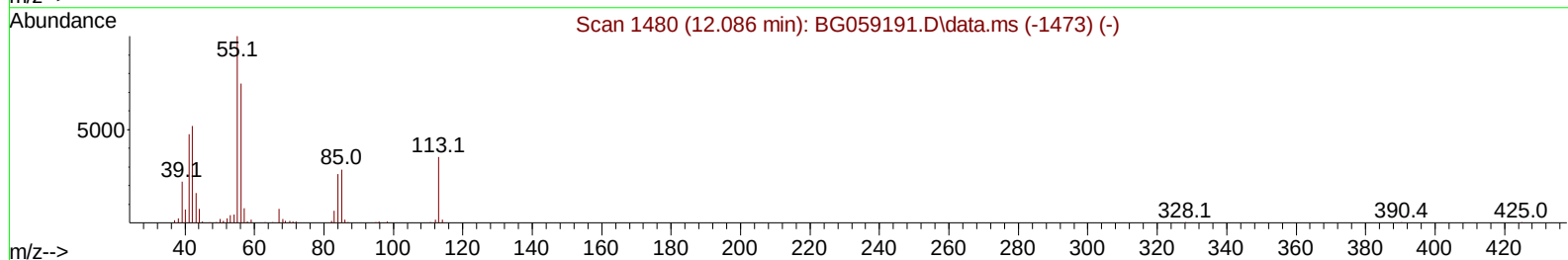
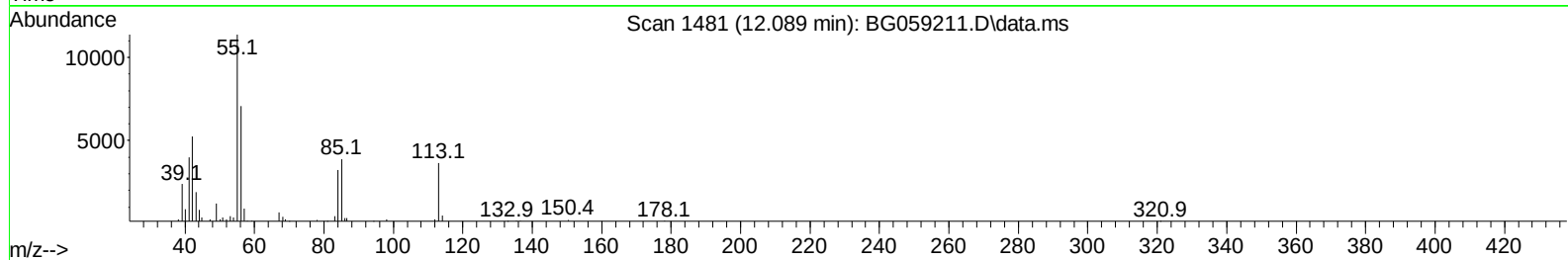
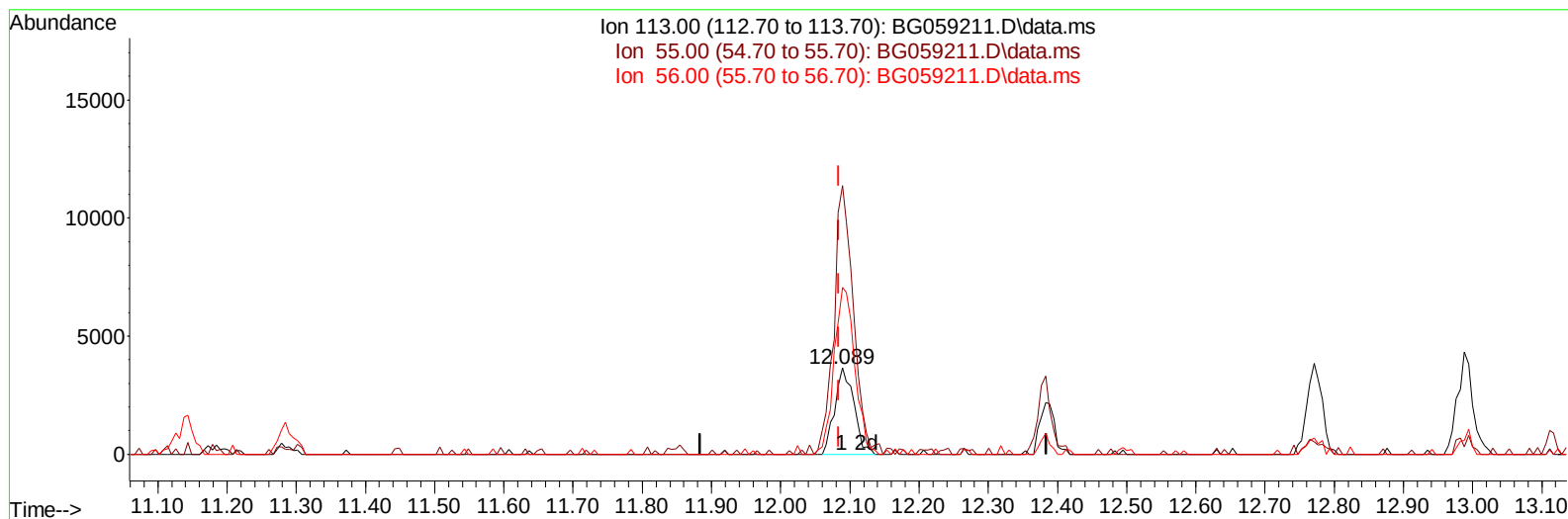
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(34) Caprolactam**

12.089min (+ 0.005) 7.95 ng/ul m

response 7041

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 197.00 | 311.34# |
| 56.00  | 140.40 | 193.59# |
| 0.00   | 0.00   | 0.00    |

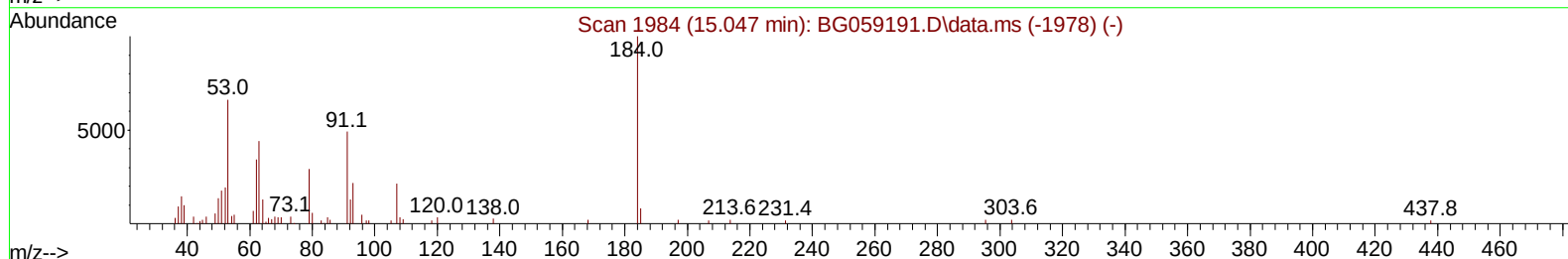
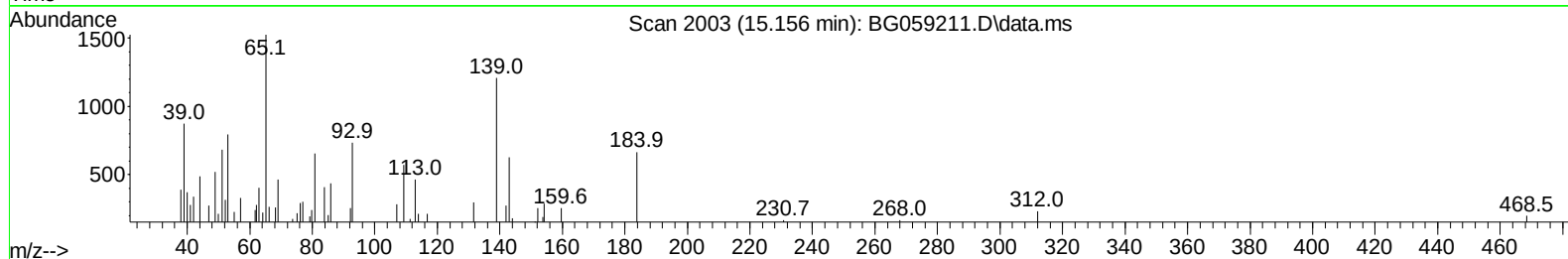
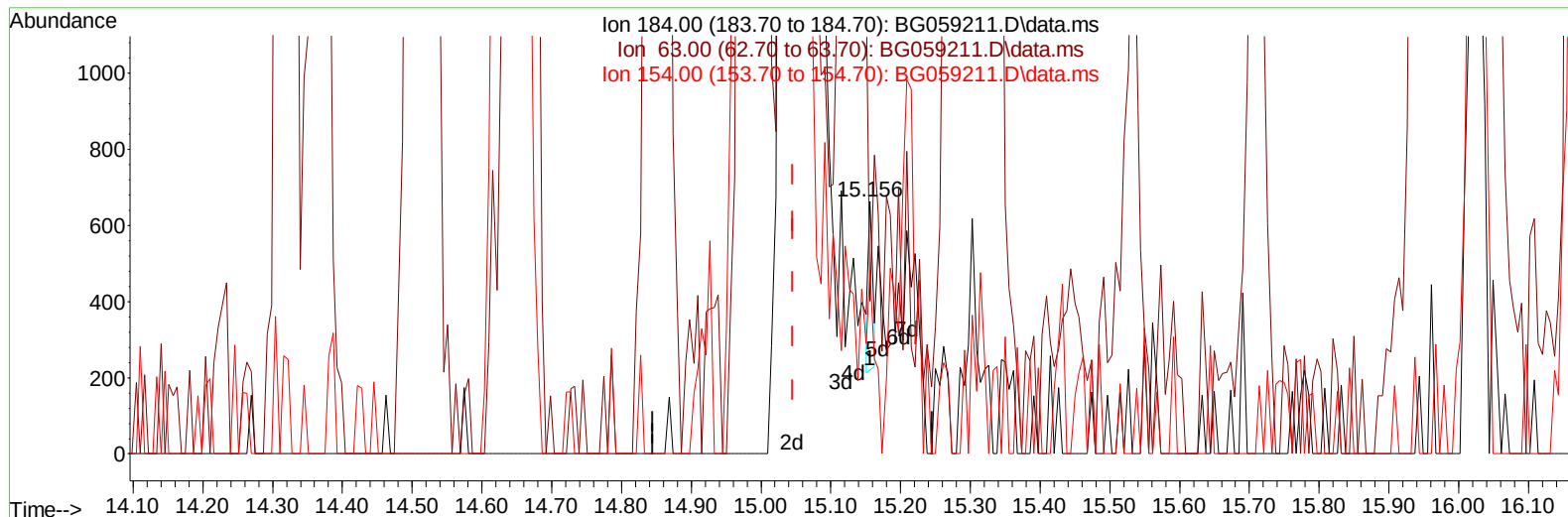
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(53) 2,4-Dinitrophenol**

**15.156min (+ 0.111) 0.19 ng/ul**

**response 199**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 66.90  | 60.39  |
| 154.00 | 68.20  | 71.99  |
| 0.00   | 0.00   | 0.00   |

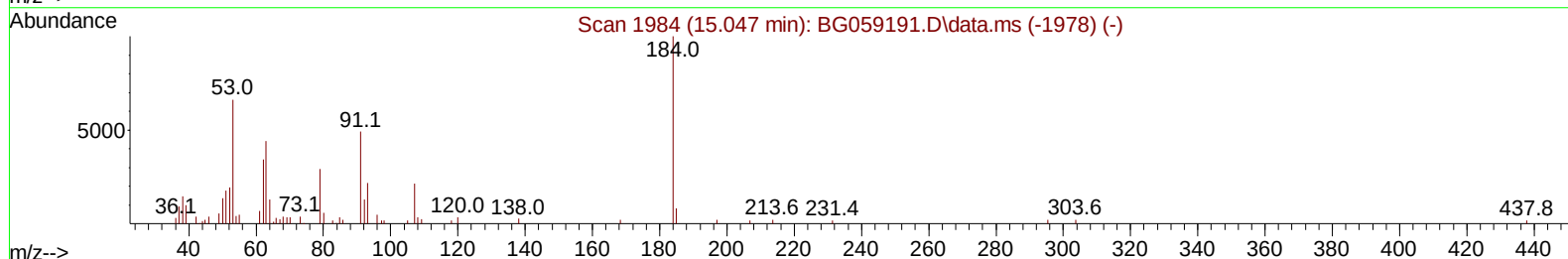
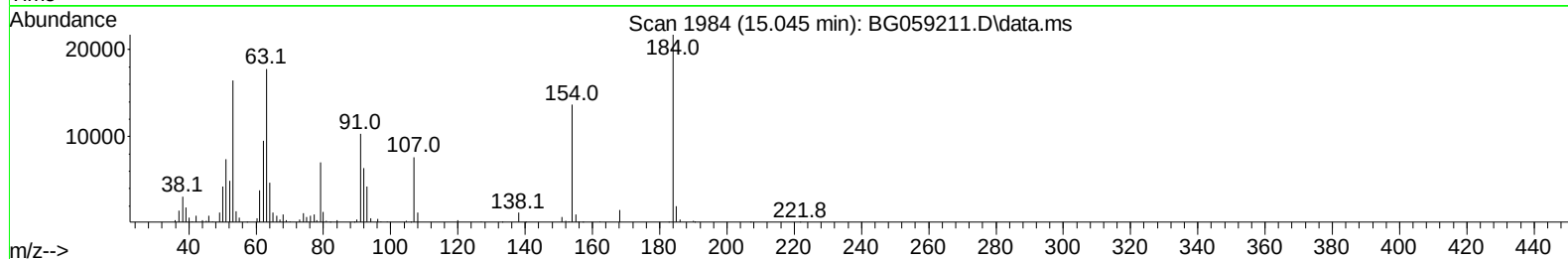
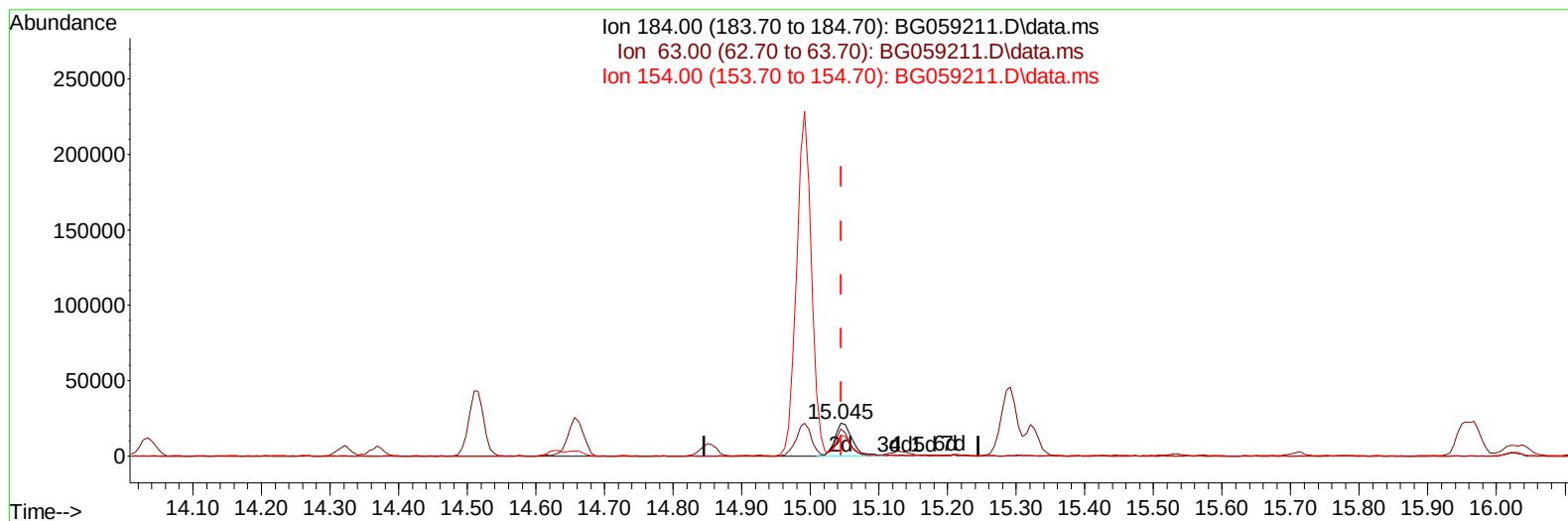
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(53) 2,4-Dinitrophenol**

**15.045min (-0.001) 37.03 ng/ul m**

**response 38161**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 66.90  | 81.98# |
| 154.00 | 68.20  | 62.93  |
| 0.00   | 0.00   | 0.00   |

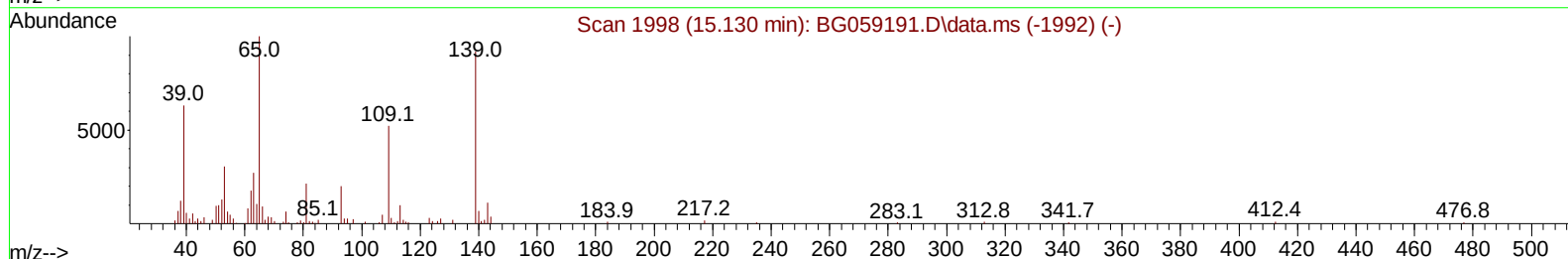
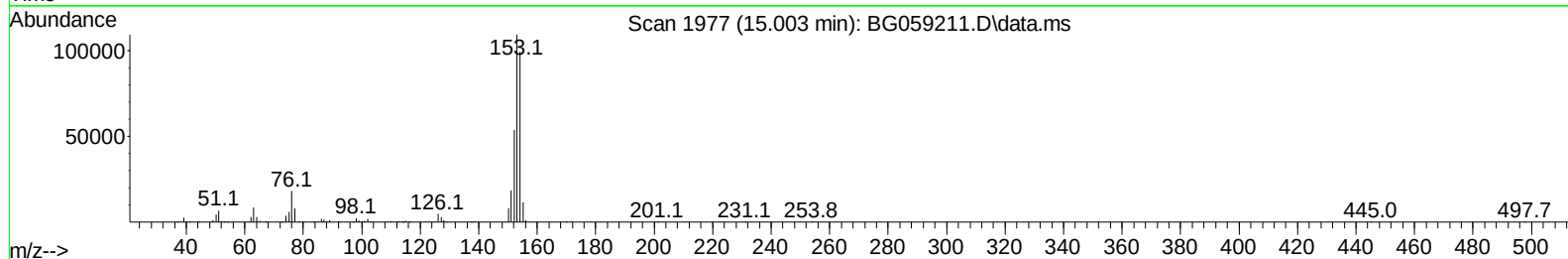
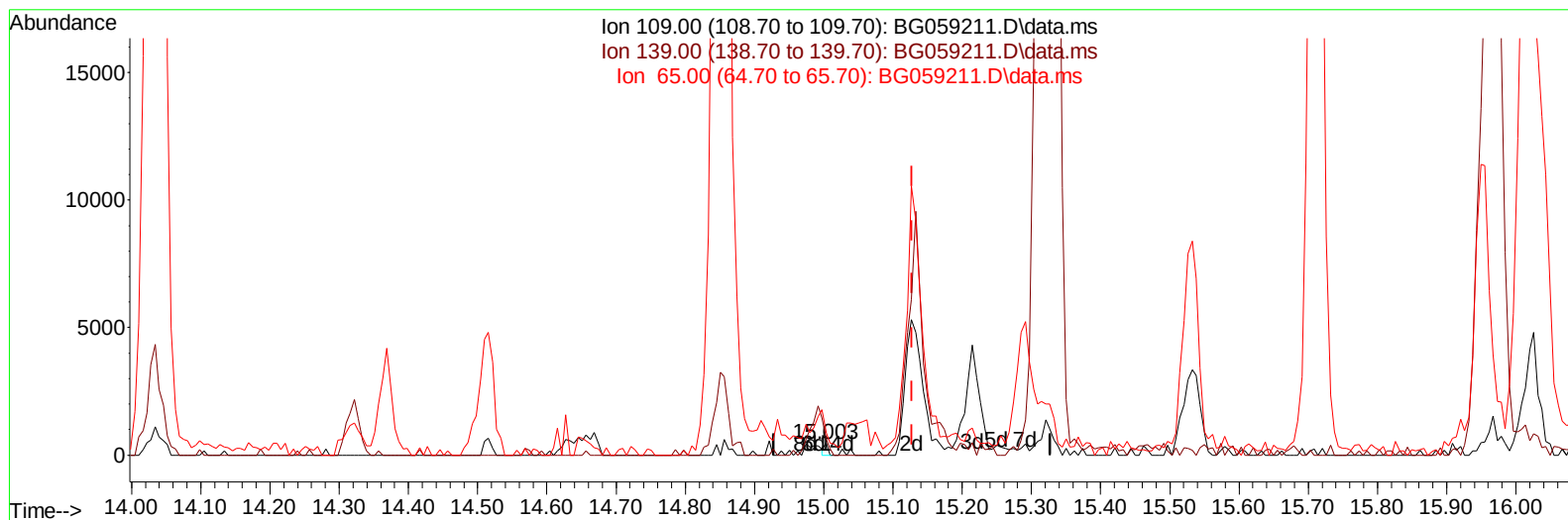
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
Data File : BG059211.D  
Acq On : 27 Sep 2023 1:12  
Operator : MA/JU  
Sample : 04450-19MSD  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Sep 26 23:48:55 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(55) 4-Nitrophenol**

**15.003min (-0.124) 0.15 ng/ul**

**response 163**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 109.00 | 100.00 | 100.00 |
| 139.00 | 143.10 | 133.41 |
| 65.00  | 165.10 | 185.47 |
| 0.00   | 0.00   | 0.00   |

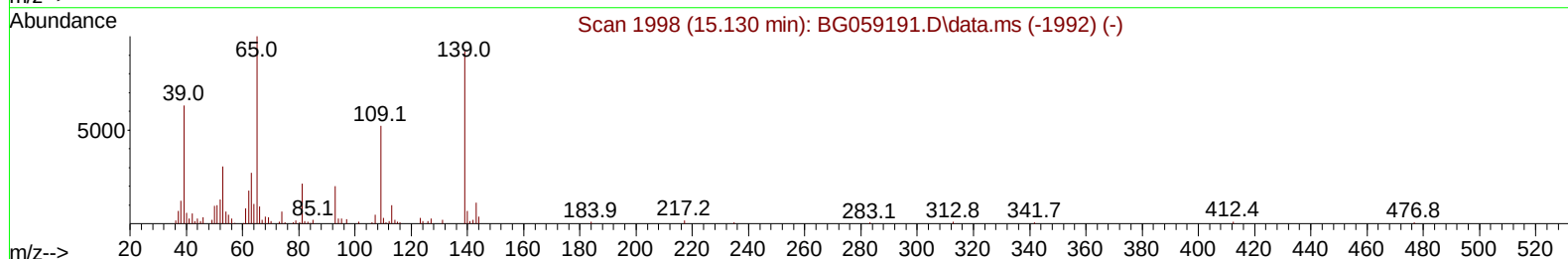
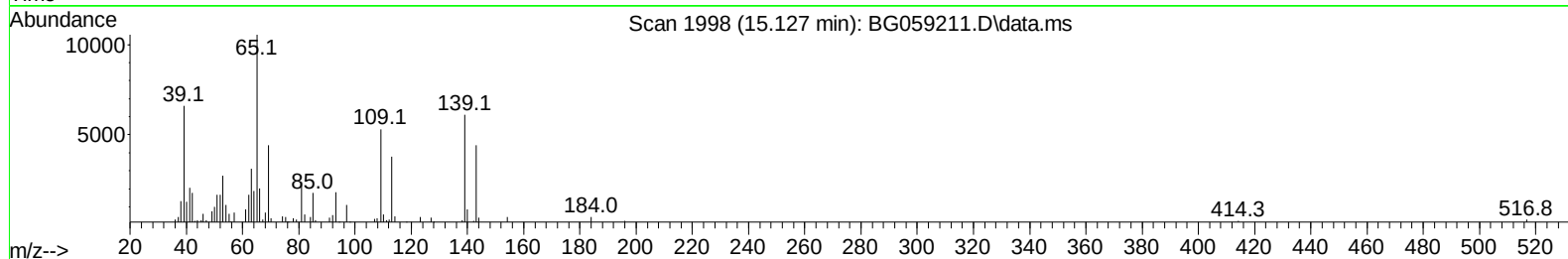
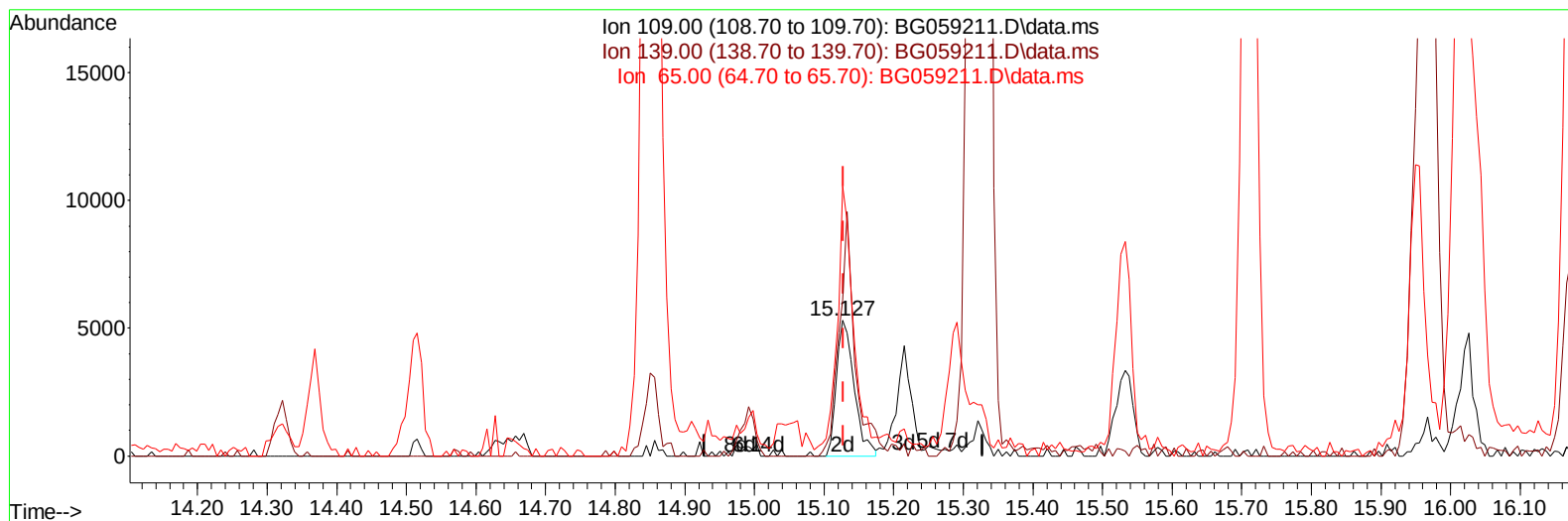
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(55) 4-Nitrophenol**

**15.127min (-0.001) 8.58 ng/ul m**

**response 9370**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 109.00 | 100.00 | 100.00  |
| 139.00 | 143.10 | 114.96  |
| 65.00  | 165.10 | 198.64# |
| 0.00   | 0.00   | 0.00    |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :

BNA\_G

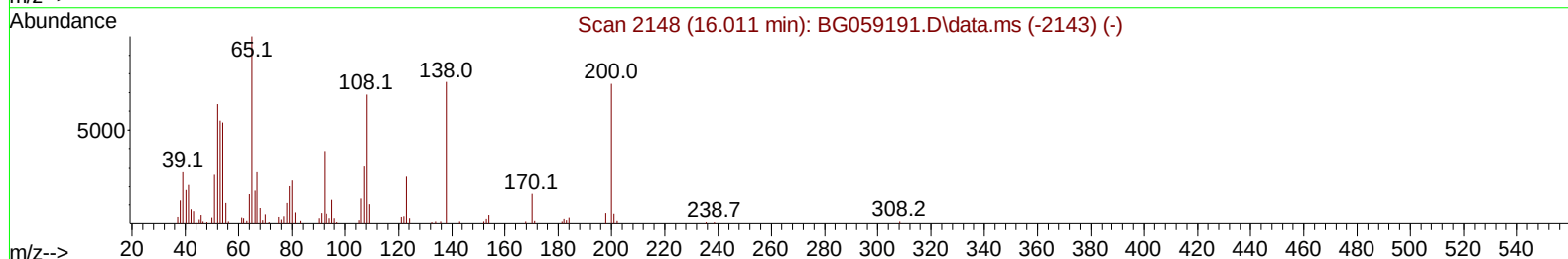
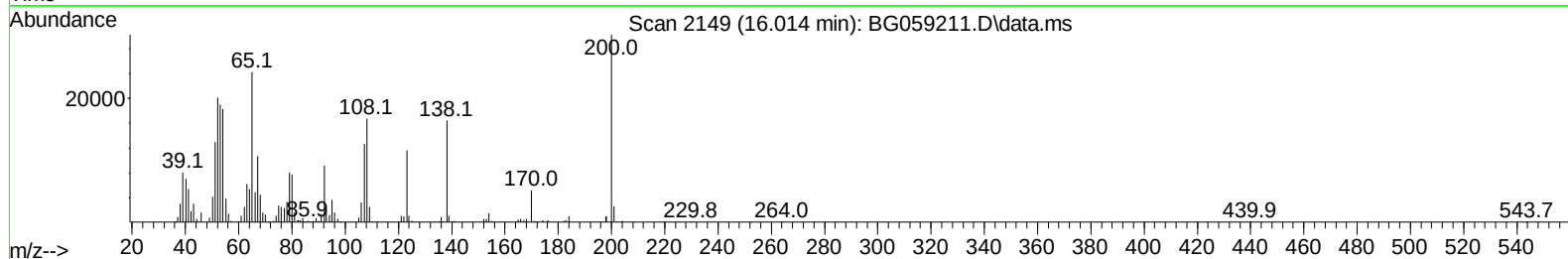
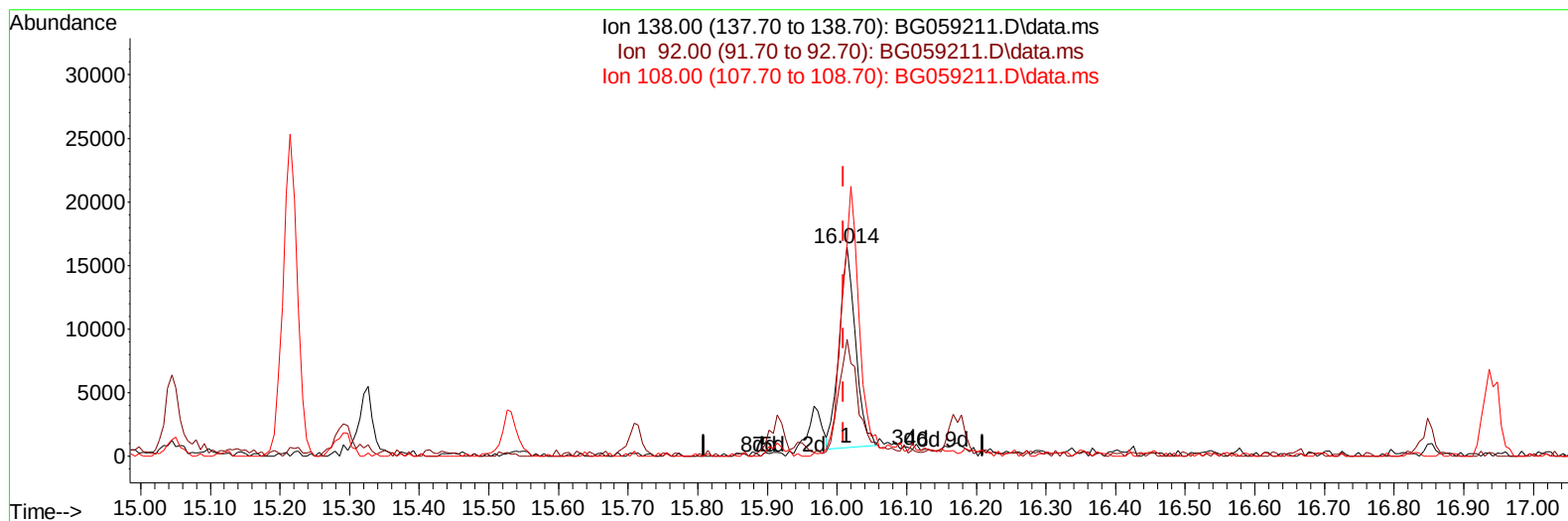
ClientSampleId :

BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(63) 4-Nitroaniline**

16.014min (+ 0.005) 14.75 ng/ul

response 25800

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 138.00 | 100.00 | 100.00  |
| 92.00  | 48.80  | 55.96   |
| 108.00 | 66.90  | 101.80# |
| 0.00   | 0.00   | 0.00    |

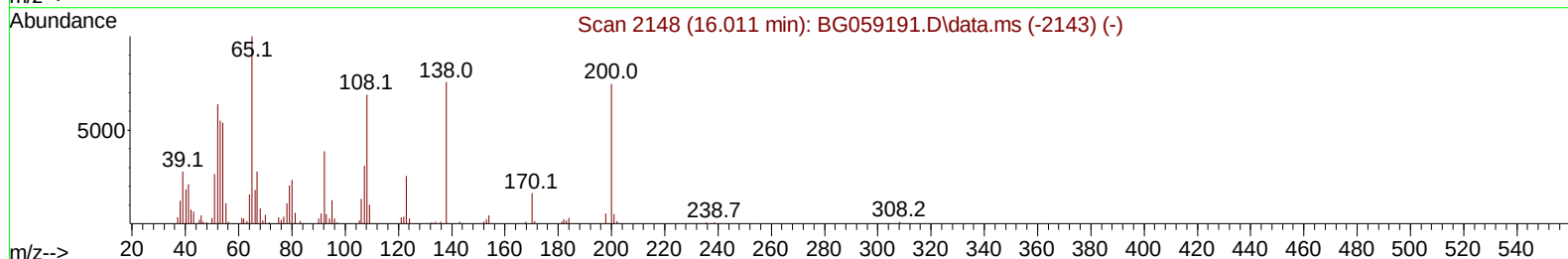
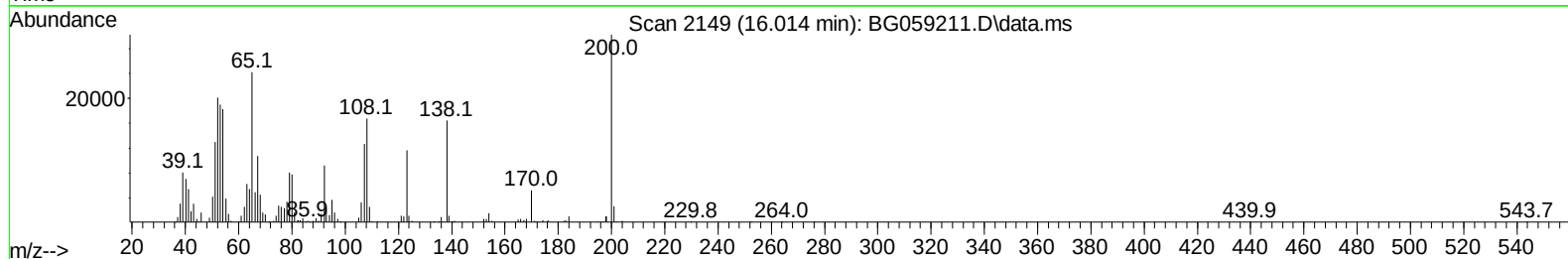
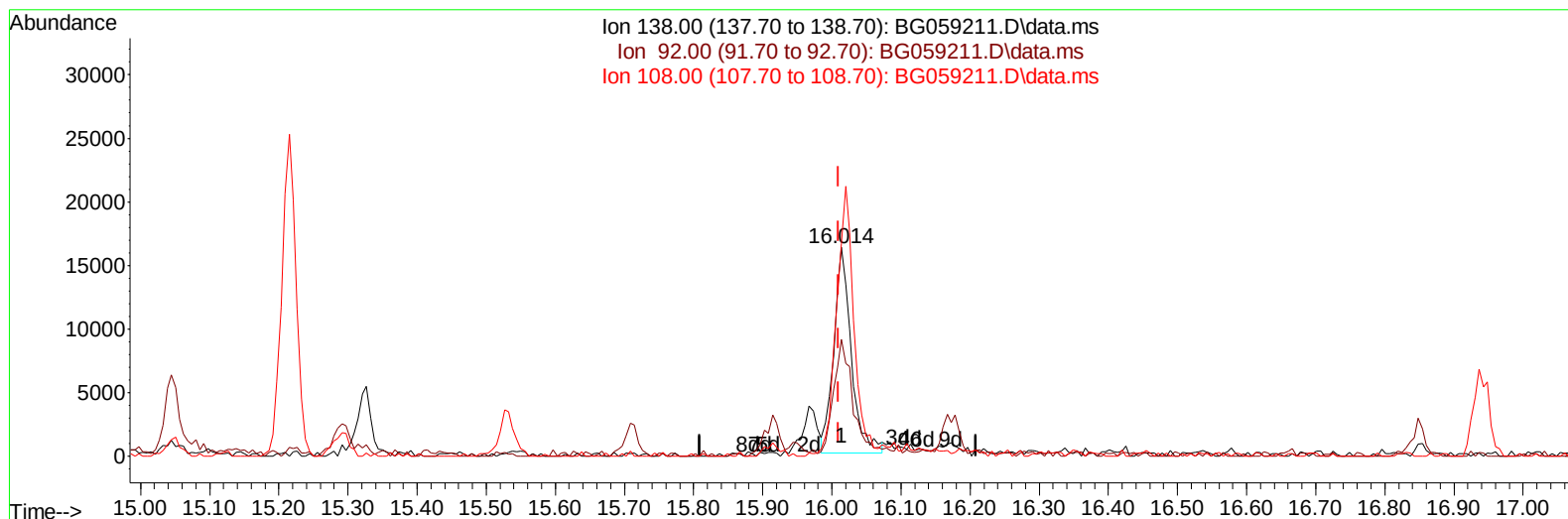
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:27:55 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059211.D\data.ms

**(63) 4-Nitroaniline**

**16.014min (+ 0.005) 16.31 ng/ul m**

**response 28530**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 138.00 | 100.00 | 100.00  |
| 92.00  | 48.80  | 55.96   |
| 108.00 | 66.90  | 101.80# |
| 0.00   | 0.00   | 0.00    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :

BNA\_G

ClientSampleId :

BBHR2MSD

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/27/2023

Supervised By :mohammad ahmed 09/29/2023

Quant Time: Sep 27 02:48:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.300  | 152  | 29736    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 11.143 | 136  | 161651   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.927 | 164  | 117830   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.677 | 188  | 284049   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 22.001 | 240  | 228903   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 25.550 | 264  | 266428   | 20.000 | ng/ul  | 0.01     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.605  | 96   | 4567     | 6.012  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 4.040  | 84   | 35551    | 15.951 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.424  | 99   | 38247    | 13.081 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.612  | 67   | 84711    | 47.211 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.812  | 132  | 86495    | 40.086 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.987  | 113  | 65572    | 28.159 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.492  | 128  | 56380    | 45.199 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.215 | 143  | 62293    | 49.435 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.744 | 165  | 115048   | 48.092 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.284 | 131  | 145478   | 38.963 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.322 | 166  | 483787   | 53.386 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.627 | 160  | 515609   | 46.729 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.121 | 143  | 13844    | 8.373  | ng/ul  | 0.01     |
| 60) Fluorene-d10              | 15.914 | 176  | 434530   | 52.385 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.026 | 200  | 79264    | 49.083 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.777 | 188  | 667086   | 51.614 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 20.050 | 212  | 773177   | 57.789 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.303 | 264  | 701387   | 53.914 | ng/ul  | 0.01     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.646  | 88   | 5799     | 6.751  | ng/ul# | 87       |
| 5) Pyridine                   | 4.063  | 79   | 36940    | 16.266 | ng/ul# | 89       |
| 6) Benzaldehyde               | 7.442  | 77   | 75598    | 49.863 | ng/ul  | 91       |
| 8) Phenol                     | 7.448  | 94   | 40058    | 13.329 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.712  | 93   | 112912   | 46.796 | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.847  | 128  | 84123    | 38.212 | ng/ul  | 96       |
| 13) 2-Methylphenol            | 8.723  | 108  | 70673    | 31.687 | ng/ul  | 94       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.799  | 45   | 231372   | 46.186 | ng/ul  | 97       |
| 16) Acetophenone              | 9.140  | 105  | 176674   | 49.281 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 9.099  | 70   | 86559    | 47.761 | ng/ul  | 95       |
| 18) 4-Methylphenol            | 9.052  | 108  | 67697    | 28.324 | ng/ul  | 82       |
| 19) Hexachloroethane          | 9.381  | 117  | 34535    | 39.900 | ng/ul  | 99       |
| 22) Nitrobenzene              | 9.539  | 77   | 133757   | 47.353 | ng/ul  | 92       |
| 23) Isophorone                | 10.044 | 82   | 286094   | 47.232 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 10.244 | 139  | 63834    | 45.331 | ng/ul  | 95       |
| 26) 2,4-Dimethylphenol        | 10.268 | 107  | 73473    | 27.907 | ng/ul  | 90       |
| 27) Bis(2-Chloroethoxy)met... | 10.526 | 93   | 167121   | 44.348 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.767 | 162  | 108333   | 44.798 | ng/ul# | 94       |
| 30) Naphthalene               | 11.196 | 128  | 385859   | 44.608 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 11.308 | 127  | 132702   | 36.105 | ng/ul  | 97       |
| 33) Hexachlorobutadiene       | 11.419 | 225  | 64224    | 41.262 | ng/ul# | 90       |
| 34) Caprolactam               | 12.089 | 113  | 7041m    | 7.950  | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.383 | 107  | 121071   | 47.530 | ng/ul  | 96       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092623\  
 Data File : BG059211.D  
 Acq On : 27 Sep 2023 1:12  
 Operator : MA/JU  
 Sample : 04450-19MSD  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BBHR2MSD

Manual IntegrationsAPPROVED

Quant Time: Sep 27 02:48:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 23:48:55 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023  
 Supervised By :mohammad ahmed 09/29/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.777 | 142  | 265482   | 45.953 | ng/ul  | 96       |
| 37) 1-Methylnaphthalene       | 12.988 | 142  | 274098   | 46.967 | ng/ul  | 94       |
| 39) 1,2,4,5-Tetrachloroben... | 13.111 | 216  | 140726   | 40.739 | ng/ul  | 98       |
| 40) Hexachlorocyclopentadiene | 13.064 | 237  | 57797    | 26.244 | ng/ul# | 94       |
| 41) 2,4,6-Trichlorophenol     | 13.358 | 196  | 107397   | 47.611 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 13.429 | 196  | 118562   | 48.591 | ng/ul  | 93       |
| 43) 1,1'-Biphenyl             | 13.764 | 154  | 410847   | 43.683 | ng/ul  | 97       |
| 44) 2-Chloronaphthalene       | 13.817 | 162  | 309421   | 43.087 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 14.034 | 65   | 94713    | 41.607 | ng/ul  | 96       |
| 47) Dimethylphthalate         | 14.369 | 163  | 482657   | 52.147 | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 14.516 | 165  | 98803    | 57.767 | ng/ul  | 89       |
| 50) Acenaphthylene            | 14.657 | 152  | 541364   | 45.934 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.851 | 138  | 58420    | 30.956 | ng/ul  | 97       |
| 52) Acenaphthene              | 14.992 | 153  | 378245   | 47.326 | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 15.045 | 184  | 38161m   | 37.028 | ng/ul  |          |
| 55) 4-Nitrophenol             | 15.127 | 109  | 9370m    | 8.575  | ng/ul  |          |
| 56) Dibenzofuran              | 15.321 | 168  | 511477   | 48.248 | ng/ul  | 95       |
| 57) 2,4-Dinitrotoluene        | 15.291 | 165  | 144269   | 58.507 | ng/ul  | 97       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.532 | 232  | 118969   | 52.675 | ng/ul# | 90       |
| 59) Diethylphthalate          | 15.714 | 149  | 487290   | 53.428 | ng/ul  | 99       |
| 61) Fluorene                  | 15.973 | 166  | 465921   | 51.473 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.949 | 204  | 225816   | 50.265 | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 16.014 | 138  | 28530m   | 16.309 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 16.037 | 198  | 86579    | 49.742 | ng/ul  | 99       |
| 67) N-Nitrosodiphenylamine    | 16.173 | 169  | 367074   | 48.539 | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.848 | 248  | 143261   | 51.839 | ng/ul  | 97       |
| 69) Hexachlorobenzene         | 16.942 | 284  | 168975   | 53.380 | ng/ul  | 94       |
| 70) Atrazine                  | 17.107 | 200  | 148317   | 49.290 | ng/ul  | 95       |
| 71) Pentachlorophenol         | 17.301 | 266  | 89002    | 43.286 | ng/ul  | 95       |
| 72) Phenanthrene              | 17.724 | 178  | 821198   | 51.662 | ng/ul  | 98       |
| 74) Anthracene                | 17.812 | 178  | 819071   | 50.429 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.723 | 216  | 146334   | 38.642 | ng/uL  | 93       |
| 76) Pentachlorobenzene        | 15.215 | 250  | 160882   | 45.930 | ng/uL  | 94       |
| 77) Carbazole                 | 18.088 | 167  | 641729   | 47.523 | ng/ul  | 97       |
| 78) Di-n-butylphthalate       | 18.593 | 149  | 872114   | 52.426 | ng/ul  | 100      |
| 80) Fluoranthene              | 19.716 | 202  | 933090   | 55.966 | ng/ul  | 99       |
| 82) Pyrene                    | 20.080 | 202  | 971906   | 55.885 | ng/ul  | 96       |
| 83) Butylbenzylphthalate      | 20.932 | 149  | 374240   | 59.575 | ng/ul  | 95       |
| 84) 3,3'-Dichlorobenzidine    | 21.889 | 252  | 114096   | 20.627 | ng/ul  | 97       |
| 85) Benzo(a)anthracene        | 21.978 | 228  | 892773   | 53.720 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.801 | 149  | 571077   | 60.761 | ng/ul  | 98       |
| 87) Chrysene                  | 22.048 | 228  | 825742   | 53.279 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 23.123 | 149  | 964664   | 59.223 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 24.398 | 252  | 911572   | 56.149 | ng/ul  | 96       |
| 91) Benzo(k)fluoranthene      | 24.475 | 252  | 930070   | 55.255 | ng/ul  | 100      |
| 93) Benzo(a)pyrene            | 25.379 | 252  | 819282   | 52.509 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.657 | 276  | 877453   | 49.448 | ng/ul  | 96       |
| 95) Dibenzo(a,h)anthracene    | 29.745 | 278  | 649773   | 45.565 | ng/ul  | 97       |
| 96) Benzo(g,h,i)perylene      | 30.967 | 276  | 759736   | 53.149 | ng/ul  | 96       |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

BBHR2MSD

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 09/27/2023

Supervised By :mohammad ahmed 09/29/2023

