

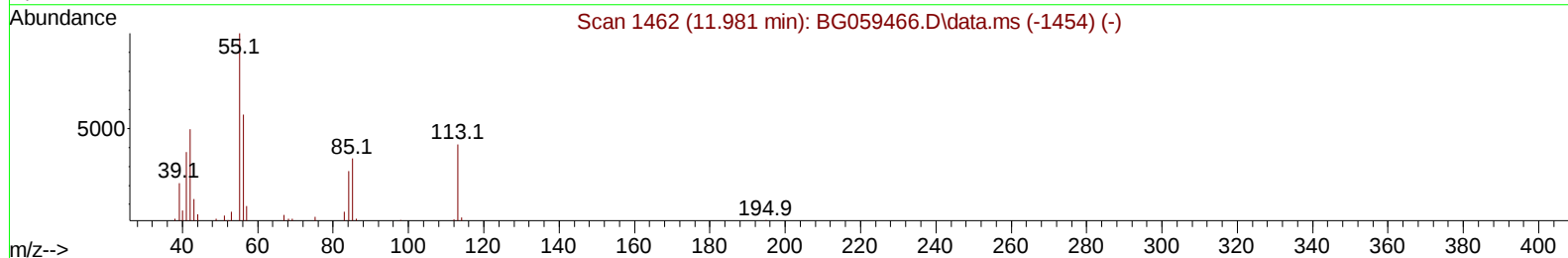
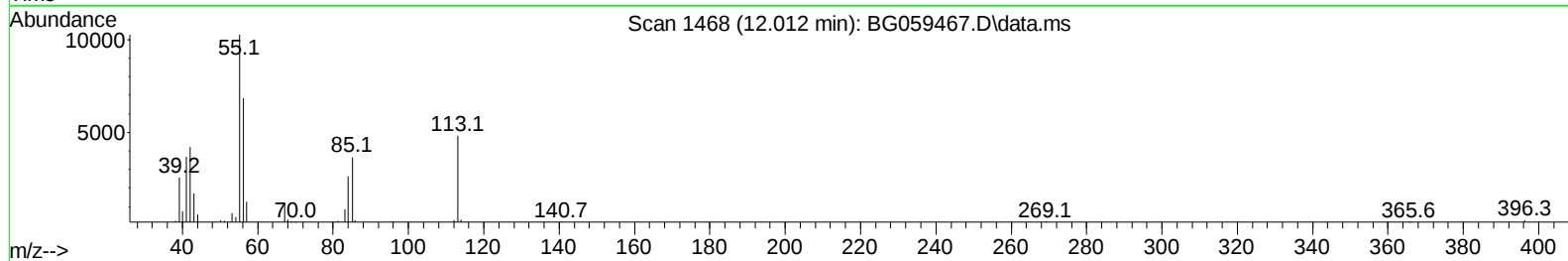
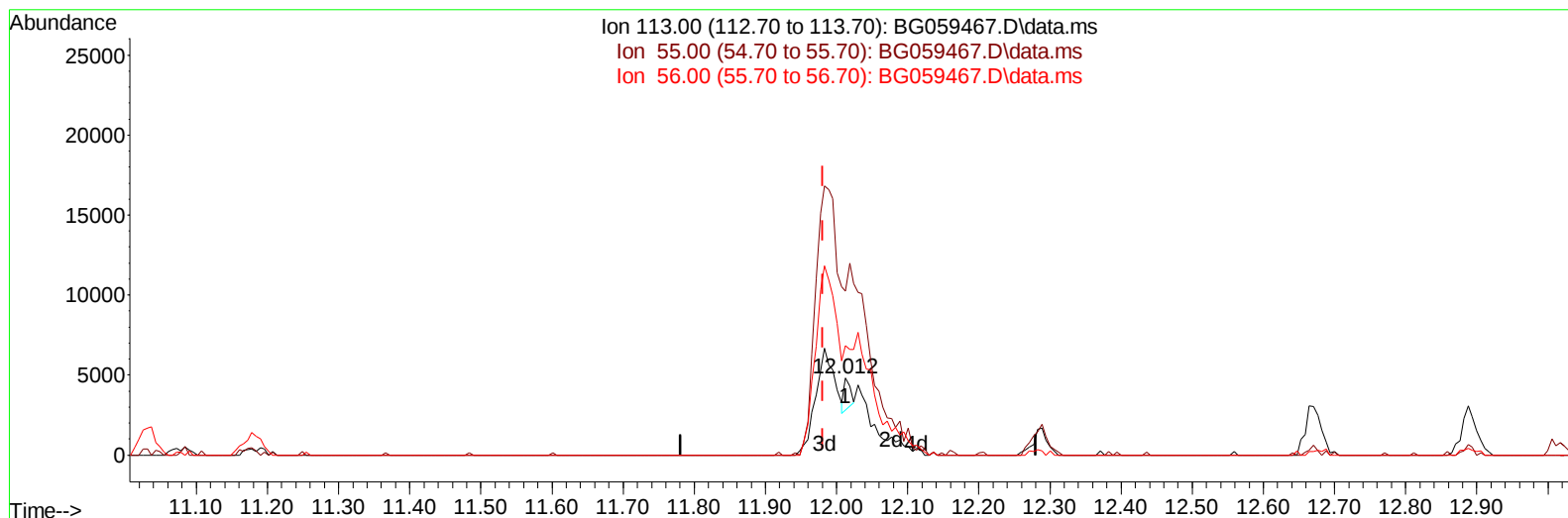
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
 Data File : BG059467.D  
 Acq On : 17 Oct 2023 21:02  
 Operator : MA/JU  
 Sample : SST04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 18 03:51:17 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
 Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

**(34) Caprolactam**

**12.012min (+ 0.032) 1.24 ng/ul**

**response 1270**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 239.70 | 212.91 |
| 56.00  | 137.00 | 141.81 |
| 0.00   | 0.00   | 0.00   |

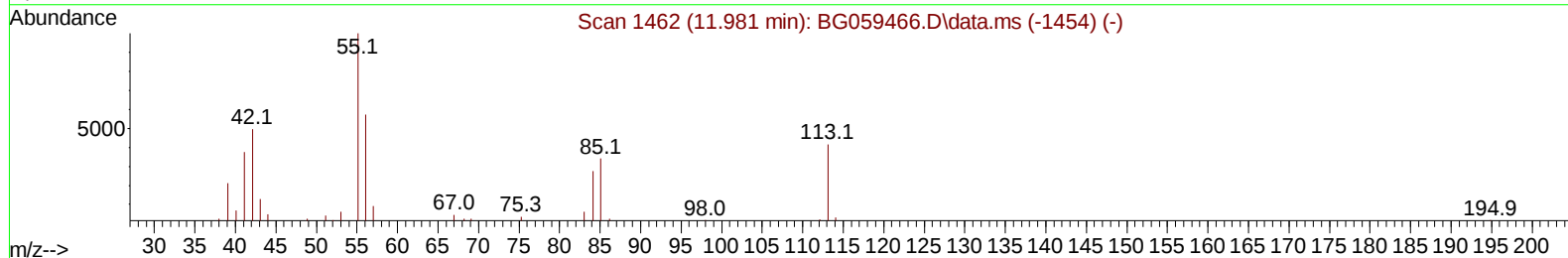
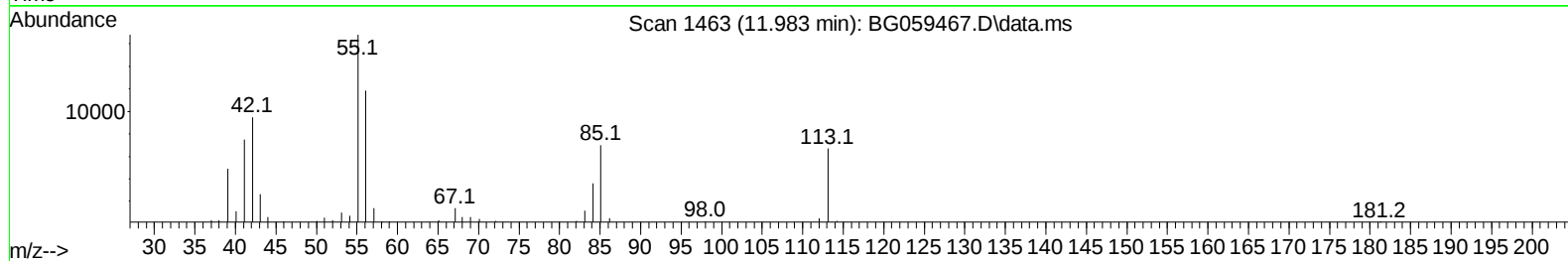
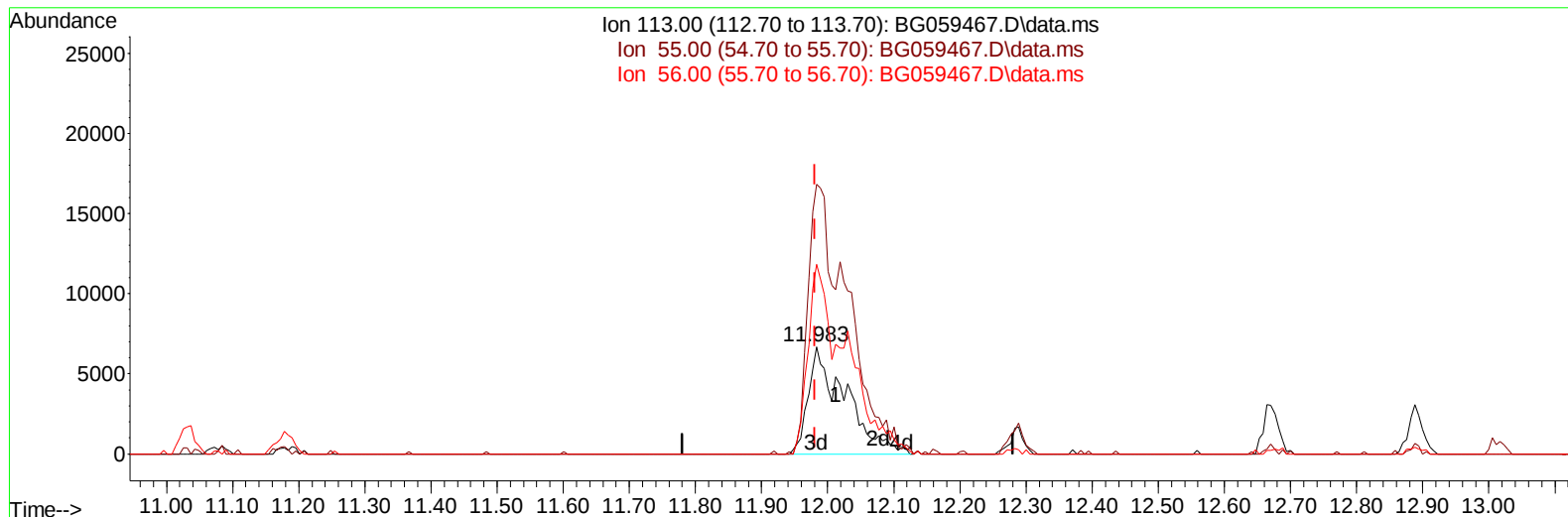
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
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Manual IntegrationsAPPROVED

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TIC: BG059467.D\data.ms

(34) Caprolactam

11.983min (+ 0.002) 25.53 ng/ul m

response 26124

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 239.70 | 251.43  |
| 56.00  | 137.00 | 177.13# |
| 0.00   | 0.00   | 0.00    |

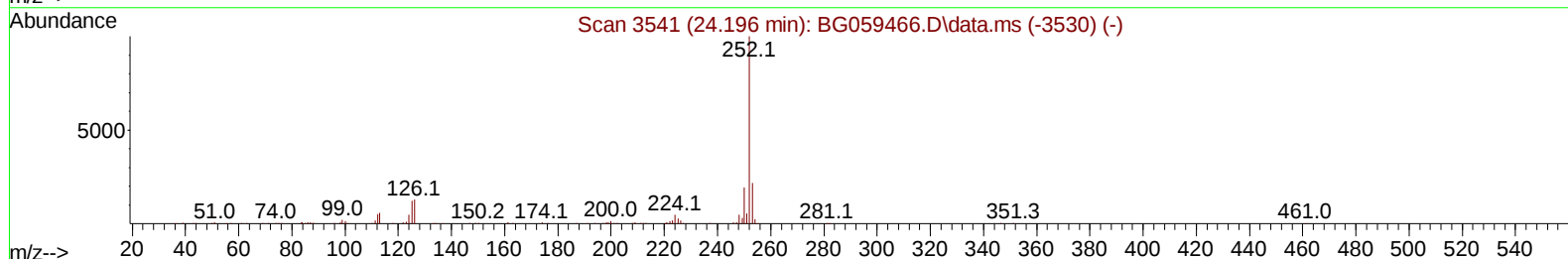
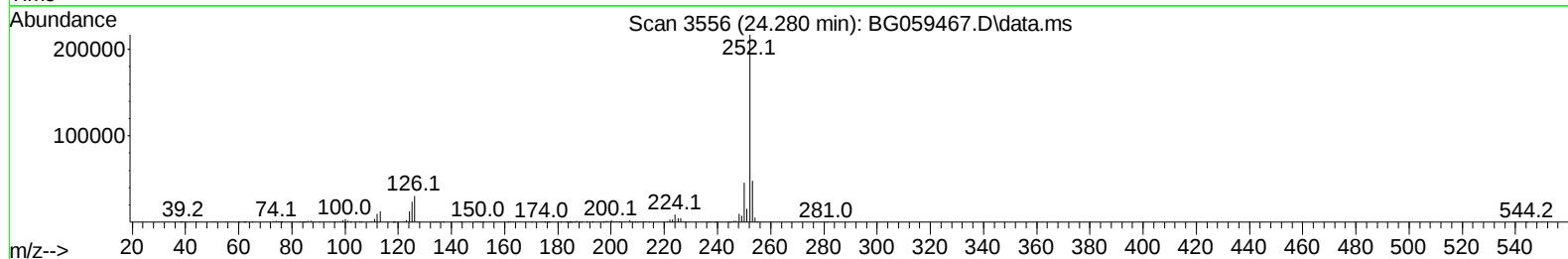
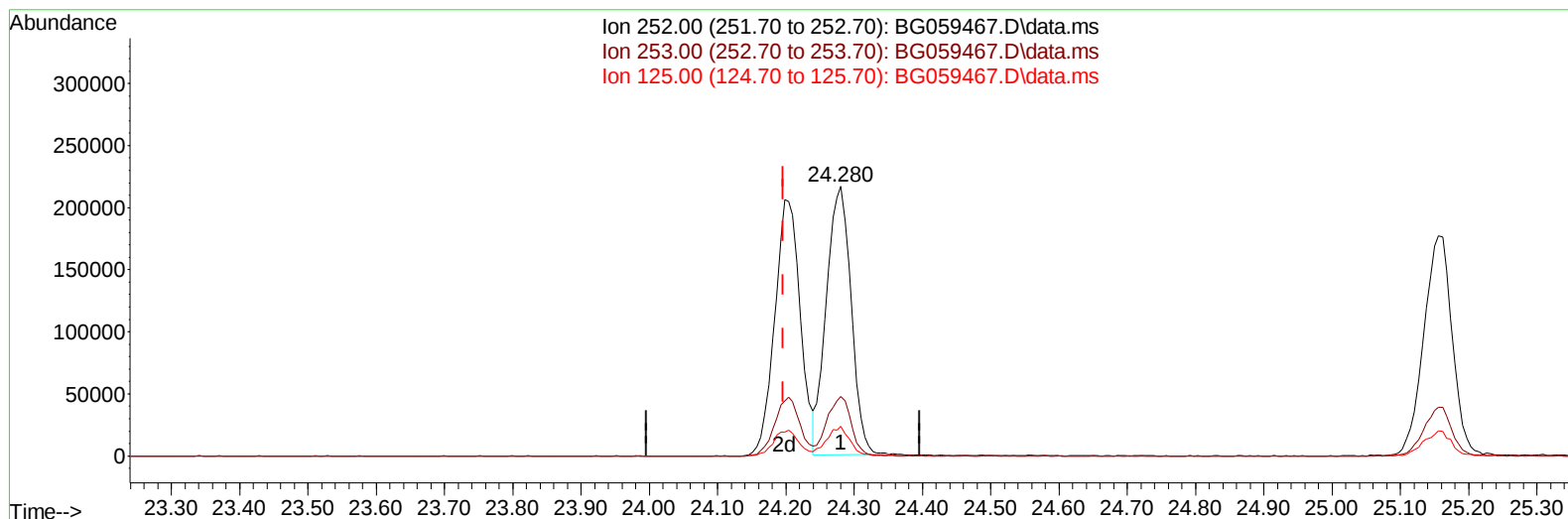
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
Data File : BG059467.D  
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ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD040416

Manual IntegrationsAPPROVED

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QLast Update : Wed Oct 18 03:51:17 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

**(90) Benzo(b)fluoranthene**

**24.280min (+ 0.085) 43.86 ng/u1**

**response 544558**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.50  | 22.11  |
| 125.00 | 12.20  | 11.00  |
| 0.00   | 0.00   | 0.00   |

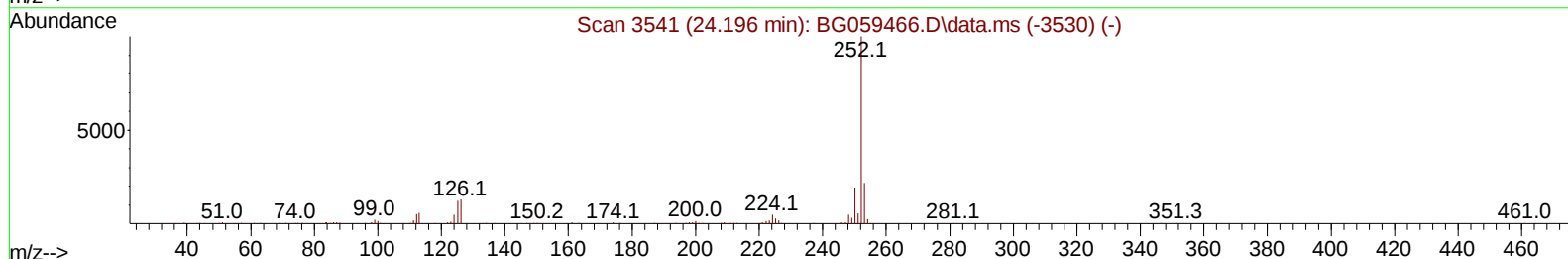
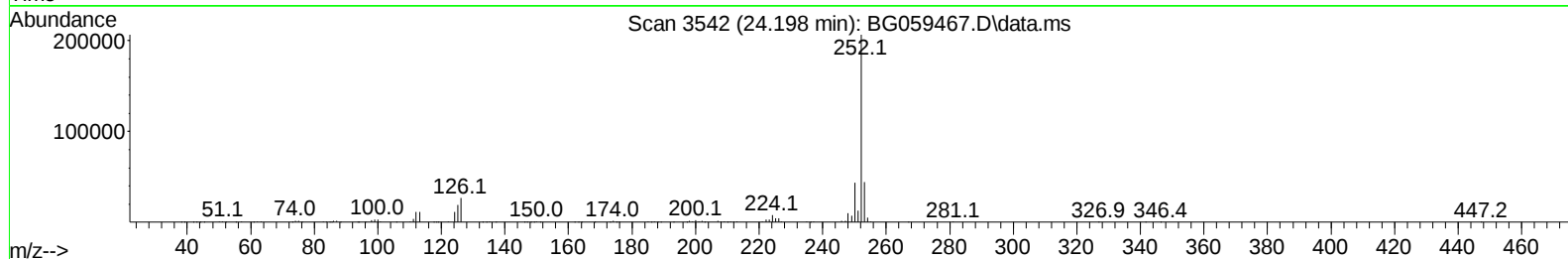
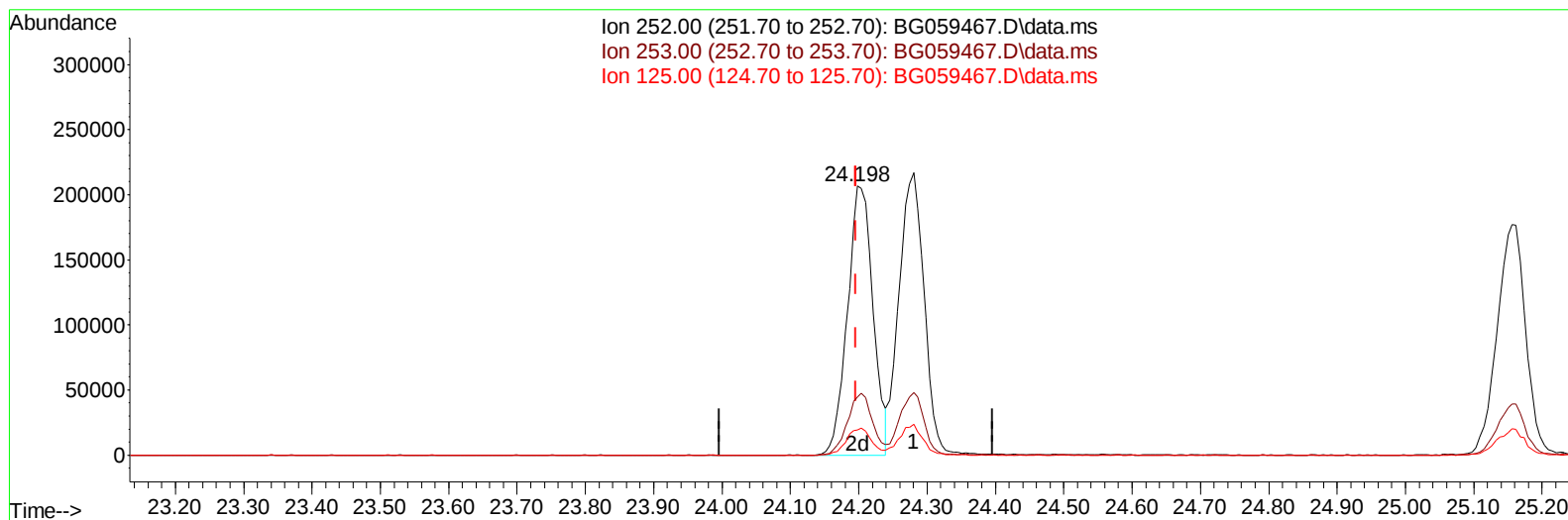
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
Data File : BG059467.D  
Acq On : 17 Oct 2023 21:02  
Operator : MA/JU  
Sample : SSTD04052  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 18 03:51:17 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

(90) Benzo(b)fluoranthene

24.198min (+ 0.002) 43.52 ng/ul m

response 540325

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.50  | 21.49  |
| 125.00 | 12.20  | 9.30#  |
| 0.00   | 0.00   | 0.00   |

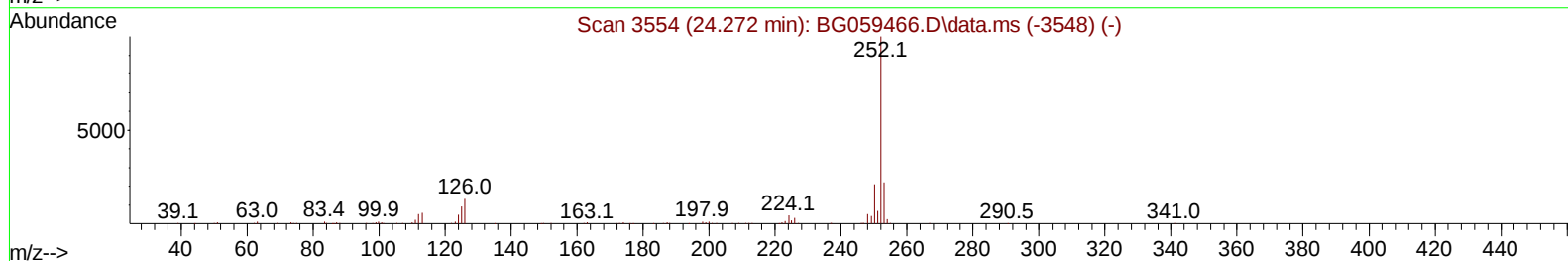
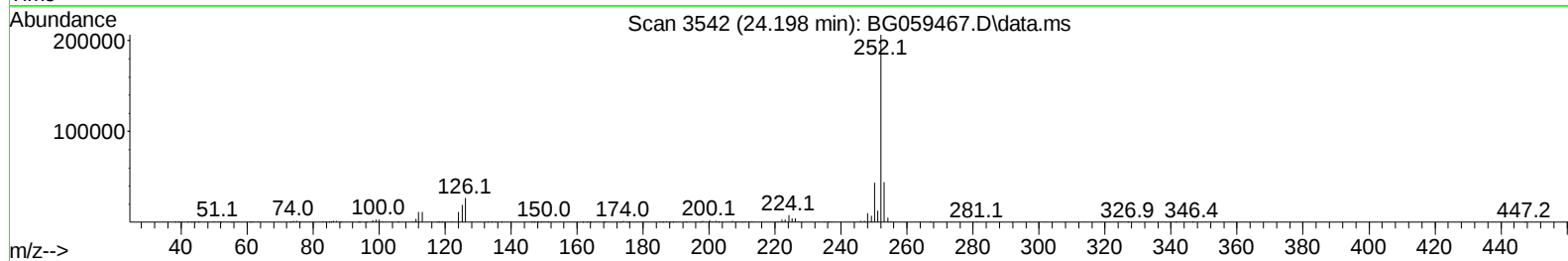
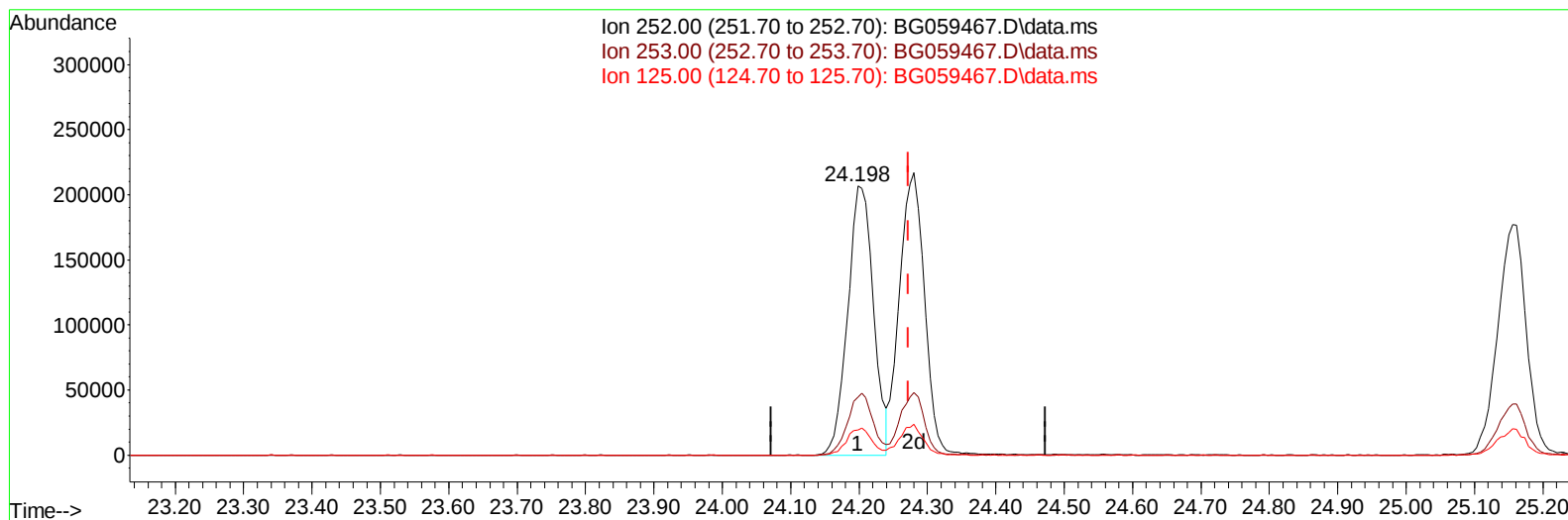
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
Data File : BG059467.D  
Acq On : 17 Oct 2023 21:02  
Operator : MA/JU  
Sample : SST04052  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 18 03:51:17 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

(91) Benzo(k)fluoranthene

24.198min (-0.074) 42.80 ng/u1

response 540325

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 22.00  | 21.49  |
| 125.00 | 9.00   | 9.30   |
| 0.00   | 0.00   | 0.00   |

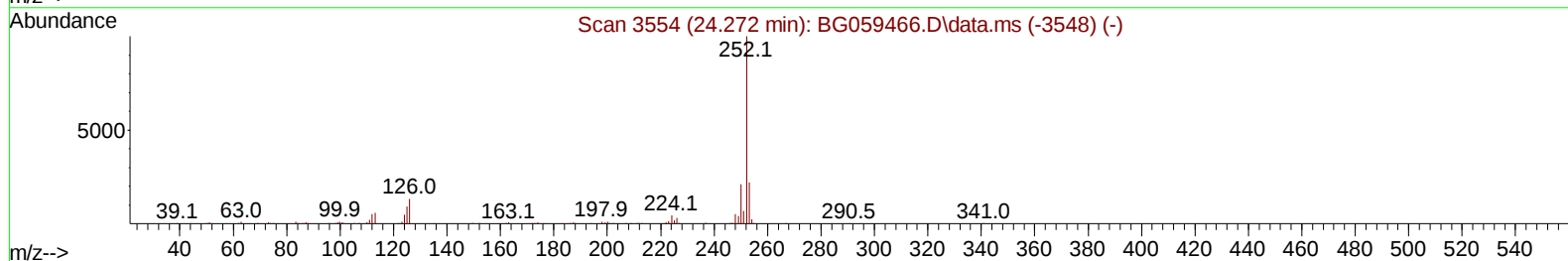
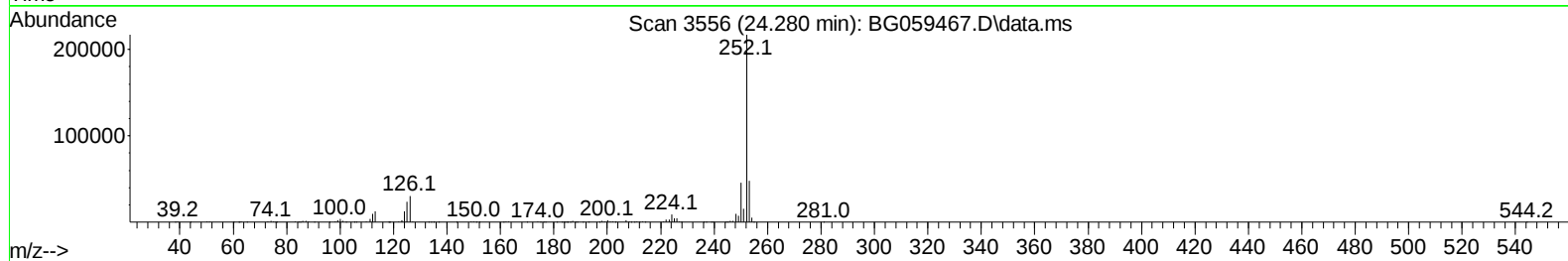
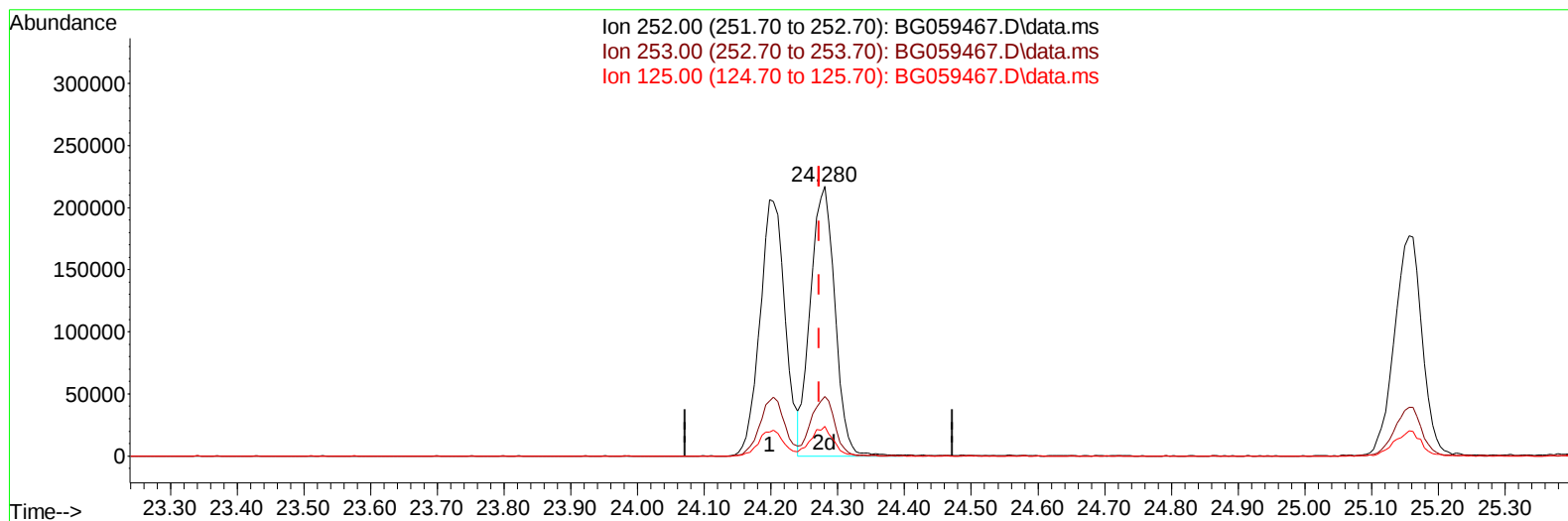
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
 Data File : BG059467.D  
 Acq On : 17 Oct 2023 21:02  
 Operator : MA/JU  
 Sample : SST04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 18 03:51:17 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
 Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

(91) Benzo(k)fluoranthene

24.280min (+ 0.008) 43.93 ng/ul m

response 554548

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 22.00  | 22.11  |
| 125.00 | 9.00   | 11.00# |
| 0.00   | 0.00   | 0.00   |

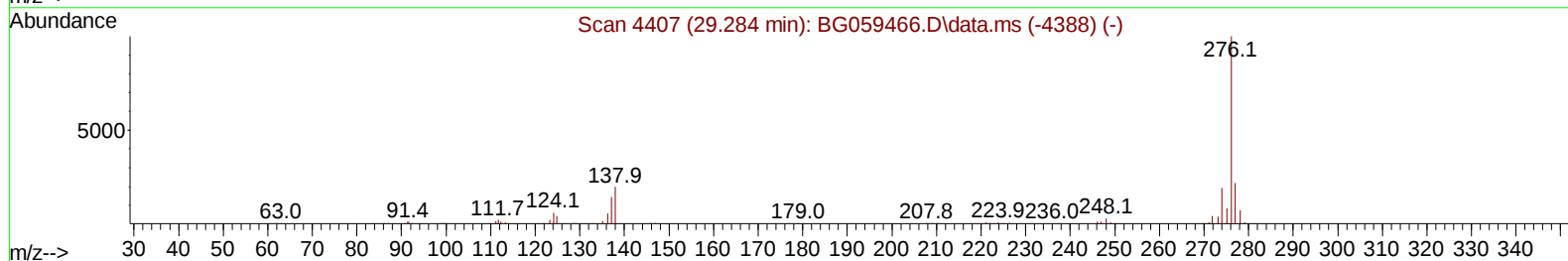
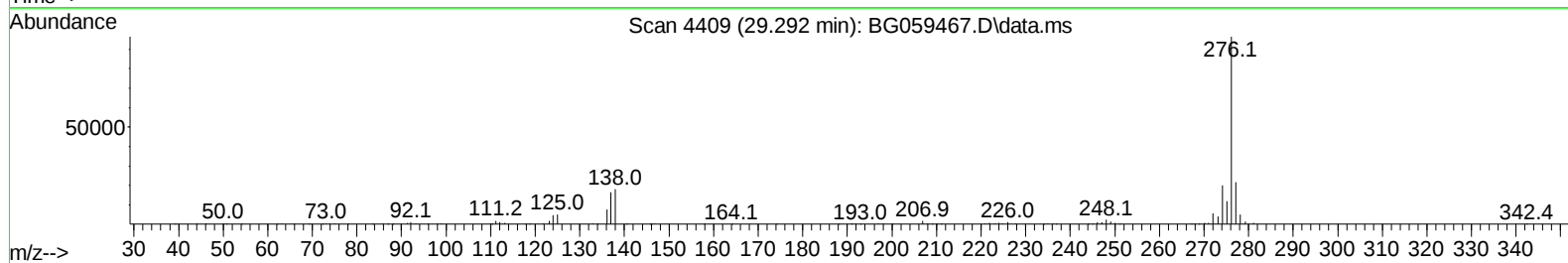
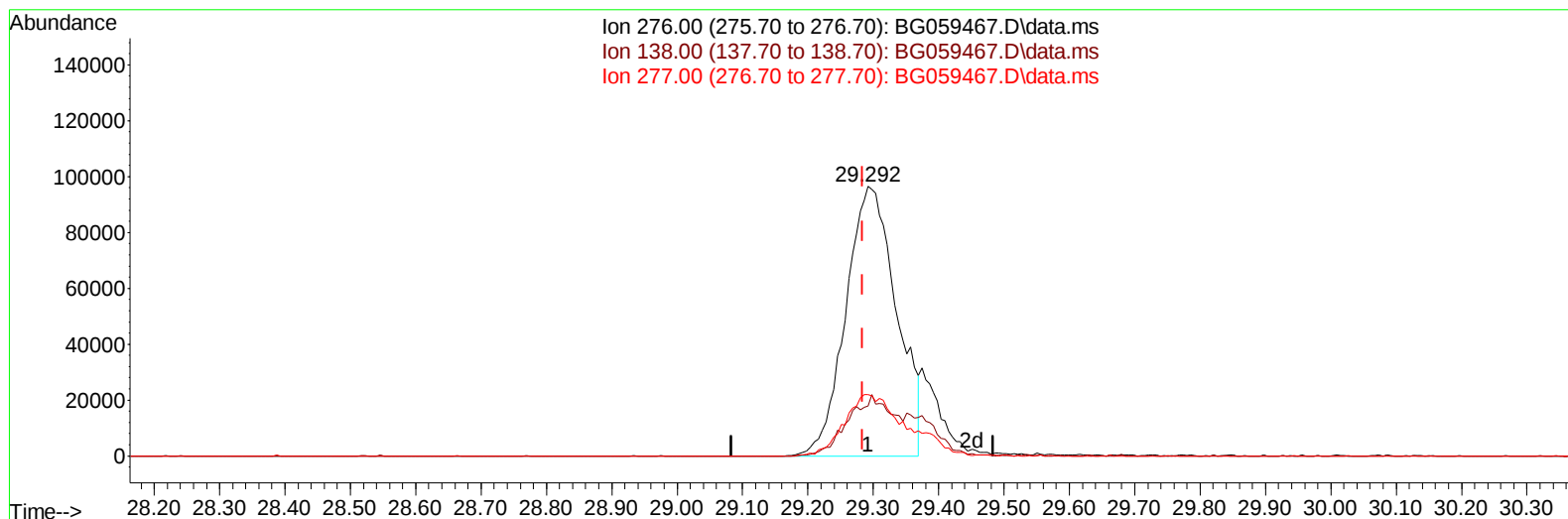
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
 Data File : BG059467.D  
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 Operator : MA/JU  
 Sample : SSTD04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
 Quant Title : SVOA CALIBRATION  
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TIC: BG059467.D\data.ms

**(94) Indeno(1,2,3-cd)pyrene**

29.292min (+ 0.008) 38.47 ng/u1

response 520276

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 20.10  | 18.63  |
| 277.00 | 21.80  | 22.67  |
| 0.00   | 0.00   | 0.00   |

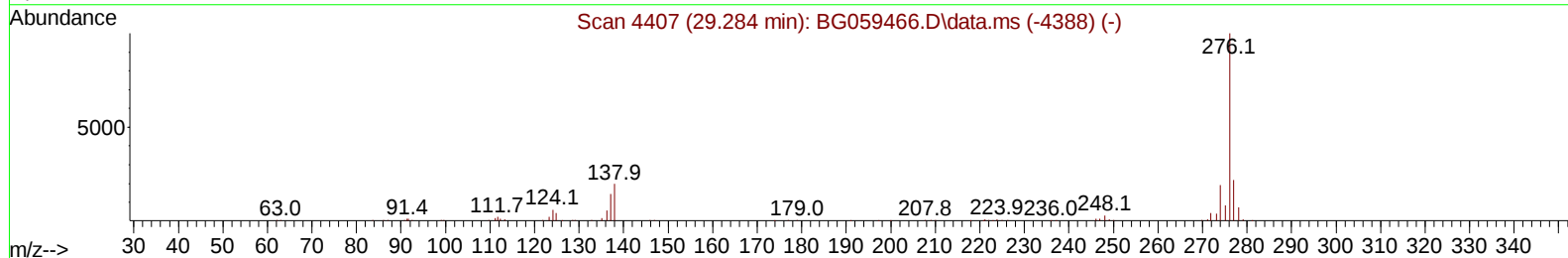
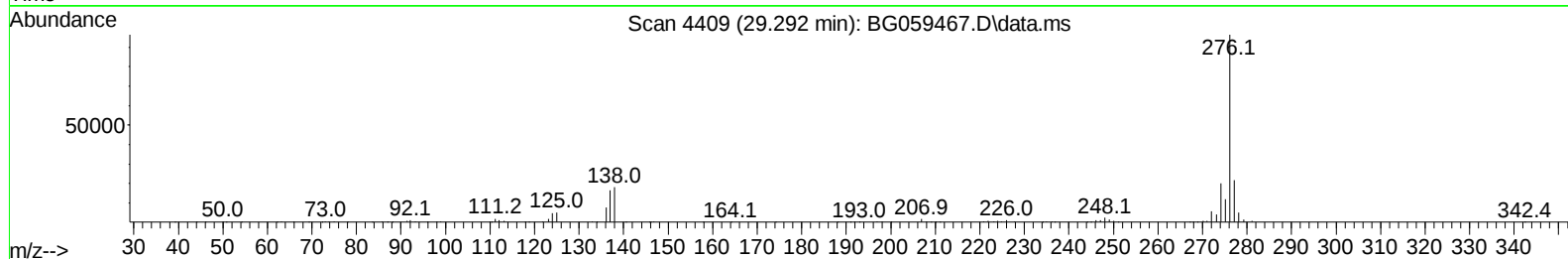
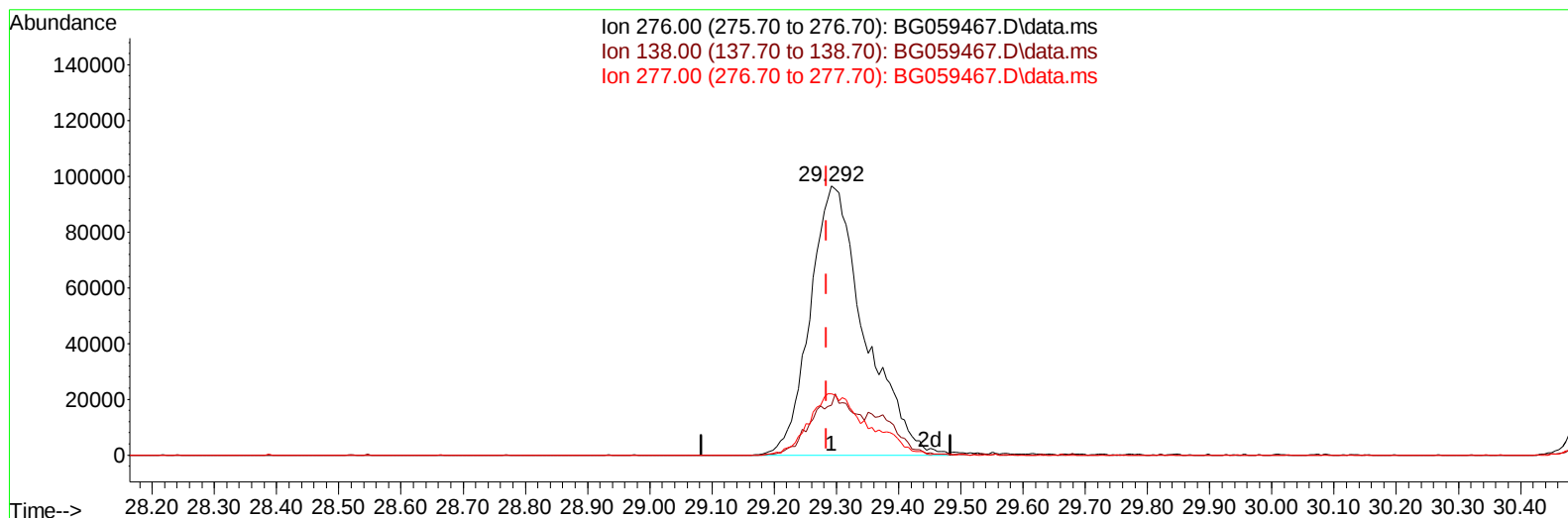
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
 Data File : BG059467.D  
 Acq On : 17 Oct 2023 21:02  
 Operator : MA/JU  
 Sample : SSTD04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 03:57:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG101723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 18 03:51:17 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023  
 Supervised By :mohammad ahmed 10/19/2023



TIC: BG059467.D\data.ms

**(94) Indeno(1,2,3-cd)pyrene**

29.292min (+ 0.008) 43.47 ng/ul m

response 587981

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 20.10  | 18.63  |
| 277.00 | 21.80  | 22.67  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG101723\  
 Data File : BG059467.D  
 Acq On : 17 Oct 2023 21:02  
 Operator : MA/JU  
 Sample : SST04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST040416

Manual IntegrationsAPPROVED

Quant Time: Oct 18 04:19:18 2023  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.193  | 152  | 27812    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 11.031 | 136  | 181621   | 20.000 | ng/ul  | # 0.00   |
| 38) Acenaphthene-d10          | 14.827 | 164  | 78187    | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.577 | 188  | 166653   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.871 | 240  | 153236   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 25.320 | 264  | 183602   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.481  | 96   | 13341    | 15.663 | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.916  | 84   | 135531   | 52.622 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.318  | 99   | 126421   | 41.804 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.512  | 67   | 71012    | 41.545 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.712  | 132  | 95849    | 44.305 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.887  | 113  | 95978    | 42.059 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.392  | 128  | 44944    | 30.078 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.109 | 143  | 65663    | 44.874 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.638 | 165  | 130025   | 42.355 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.178 | 131  | 173586   | 36.179 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.227 | 166  | 269656   | 40.759 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.533 | 160  | 335802   | 41.741 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.027 | 143  | 48481    | 41.883 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.820 | 176  | 250263   | 39.778 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.937 | 200  | 42378    | 41.876 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.676 | 188  | 378445   | 44.025 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.956 | 212  | 420850   | 44.577 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.079 | 264  | 432967   | 42.502 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        |        | Qvalue   |
| 2) 1,4-Dioxane                | 3.517  | 88   | 14630    | 13.659 | ng/uL  | 98       |
| 5) Pyridine                   | 3.940  | 79   | 141673   | 56.338 | ng/ul  | 89       |
| 6) Benzaldehyde               | 7.336  | 77   | 59718    | 45.132 | ng/ul  | 91       |
| 8) Phenol                     | 7.347  | 94   | 123086   | 41.138 | ng/ul  | 93       |
| 10) Bis(2-Chloroethyl)ether   | 7.612  | 93   | 96198    | 40.391 | ng/ul  | 93       |
| 12) 2-Chlorophenol            | 7.741  | 128  | 97878    | 43.542 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.622  | 108  | 94926    | 43.208 | ng/ul  | 93       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.699  | 45   | 203117   | 41.513 | ng/ul  | 98       |
| 16) Acetophenone              | 9.034  | 105  | 143798   | 41.729 | ng/ul  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.998  | 70   | 65640    | 41.369 | ng/ul  | 92       |
| 18) 4-Methylphenol            | 8.957  | 108  | 99322    | 41.587 | ng/ul  | 89       |
| 19) Hexachloroethane          | 9.274  | 117  | 37440    | 43.825 | ng/ul# | 81       |
| 22) Nitrobenzene              | 9.433  | 77   | 100532   | 30.046 | ng/ul  | 99       |
| 23) Isophorone                | 9.938  | 82   | 324833   | 46.226 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 10.144 | 139  | 66357    | 41.149 | ng/ul  | 92       |
| 26) 2,4-Dimethylphenol        | 10.168 | 107  | 126868   | 38.401 | ng/ul  | 94       |
| 27) Bis(2-Chloroethoxy)met... | 10.426 | 93   | 210004   | 46.466 | ng/ul  | 95       |
| 29) 2,4-Dichlorophenol        | 10.667 | 162  | 127630   | 43.018 | ng/ul# | 83       |
| 30) Naphthalene               | 11.084 | 128  | 455803   | 41.347 | ng/ul  | 97       |
| 32) 4-Chloroaniline           | 11.208 | 127  | 161382   | 35.336 | ng/ul  | 92       |
| 33) Hexachlorobutadiene       | 11.313 | 225  | 59106    | 30.216 | ng/ul  | 96       |
| 34) Caprolactam               | 11.983 | 113  | 26124m   | 25.531 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.289 | 107  | 87065    | 29.332 | ng/ul  | 97       |

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 Operator : MA/JU  
 Sample : SST04052  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

## Instrument :

BNA\_G

ClientSampleId :

SSTD040416

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/18/2023

Supervised By :mohammad ahmed 10/19/2023

Quant Time: Oct 18 04:19:18 2023  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 18 03:51:17 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.670 | 142  | 208509   | 29.637 | ng/ul  | 94       |
| 37) 1-Methylnaphthalene       | 12.888 | 142  | 209927   | 29.031 | ng/ul# | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 13.017 | 216  | 110228   | 42.980 | ng/ul  | 97       |
| 40) Hexachlorocyclopentadiene | 12.970 | 237  | 65945    | 44.389 | ng/ul  | 92       |
| 41) 2,4,6-Trichlorophenol     | 13.264 | 196  | 68706    | 43.924 | ng/ul  | 93       |
| 42) 2,4,5-Trichlorophenol     | 13.334 | 196  | 71978    | 42.311 | ng/ul  | 97       |
| 43) 1,1'-Biphenyl             | 13.663 | 154  | 292193   | 41.326 | ng/ul  | 100      |
| 44) 2-Chloronaphthalene       | 13.722 | 162  | 223475   | 41.136 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.940 | 65   | 58761    | 40.999 | ng/ul  | 94       |
| 47) Dimethylphthalate         | 14.275 | 163  | 267040   | 40.351 | ng/ul  | 97       |
| 48) 2,6-Dinitrotoluene        | 14.421 | 165  | 51940    | 42.847 | ng/ul  | 97       |
| 50) Acenaphthylene            | 14.562 | 152  | 352559   | 41.634 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.762 | 138  | 54809    | 41.749 | ng/ul# | 80       |
| 52) Acenaphthene              | 14.897 | 153  | 236267   | 40.520 | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 14.956 | 184  | 27803    | 42.237 | ng/ul  | 95       |
| 55) 4-Nitrophenol             | 15.038 | 109  | 34367    | 41.363 | ng/ul  | 89       |
| 56) Dibenzofuran              | 15.226 | 168  | 318798   | 40.219 | ng/ul  | 100      |
| 57) 2,4-Dinitrotoluene        | 15.197 | 165  | 75088    | 42.831 | ng/ul# | 96       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.438 | 232  | 66442    | 43.100 | ng/ul# | 92       |
| 59) Diethylphthalate          | 15.620 | 149  | 252538   | 39.348 | ng/ul  | 98       |
| 61) Fluorene                  | 15.873 | 166  | 262084   | 39.987 | ng/ul  | 94       |
| 62) 4-Chlorophenyl-phenyle... | 15.855 | 204  | 136025   | 40.756 | ng/ul  | 91       |
| 63) 4-Nitroaniline            | 15.926 | 138  | 54339    | 42.117 | ng/ul  | 92       |
| 66) 4,6-Dinitro-2-methylph... | 15.949 | 198  | 45834    | 42.771 | ng/ul# | 80       |
| 67) N-Nitrosodiphenylamine    | 16.078 | 169  | 216711   | 43.961 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.754 | 248  | 80508    | 44.305 | ng/ul  | 93       |
| 69) Hexachlorobenzene         | 16.848 | 284  | 88352    | 42.548 | ng/ul  | 97       |
| 70) Atrazine                  | 17.012 | 200  | 78875    | 41.098 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.206 | 266  | 51895    | 43.926 | ng/ul  | 85       |
| 72) Phenanthrene              | 17.624 | 178  | 441876   | 42.533 | ng/ul  | 98       |
| 74) Anthracene                | 17.718 | 178  | 454108   | 42.630 | ng/ul  | 97       |
| 75) 1,2,3,4-Tetrachloroben... | 13.628 | 216  | 111460   | 46.322 | ng/uL  | 94       |
| 76) Pentachlorobenzene        | 15.121 | 250  | 106480   | 45.408 | ng/uL  | 95       |
| 77) Carbazole                 | 17.994 | 167  | 388383   | 42.162 | ng/ul  | 98       |
| 78) Di-n-butylphthalate       | 18.493 | 149  | 446628   | 43.009 | ng/ul  | 97       |
| 80) Fluoranthene              | 19.615 | 202  | 517863   | 46.024 | ng/ul  | 98       |
| 82) Pyrene                    | 19.985 | 202  | 537046   | 44.444 | ng/ul  | 97       |
| 83) Butylbenzylphthalate      | 20.832 | 149  | 195146   | 47.931 | ng/ul  | 95       |
| 84) 3,3'-Dichlorobenzidine    | 21.766 | 252  | 187627   | 46.783 | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.848 | 228  | 529038   | 43.663 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.683 | 149  | 286127   | 48.213 | ng/ul  | 96       |
| 87) Chrysene                  | 21.924 | 228  | 500326   | 43.772 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.958 | 149  | 484517   | 45.050 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 24.198 | 252  | 540325m  | 43.516 | ng/ul  |          |
| 91) Benzo(k)fluoranthene      | 24.280 | 252  | 554548m  | 43.929 | ng/ul  |          |
| 93) Benzo(a)pyrene            | 25.156 | 252  | 510746   | 42.414 | ng/ul  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.292 | 276  | 587981m  | 43.471 | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 29.374 | 278  | 472706   | 42.004 | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene      | 30.561 | 276  | 475933   | 43.056 | ng/ul# | 94       |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTD040416

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/18/2023

Supervised By :mohammad ahmed 10/19/2023

