

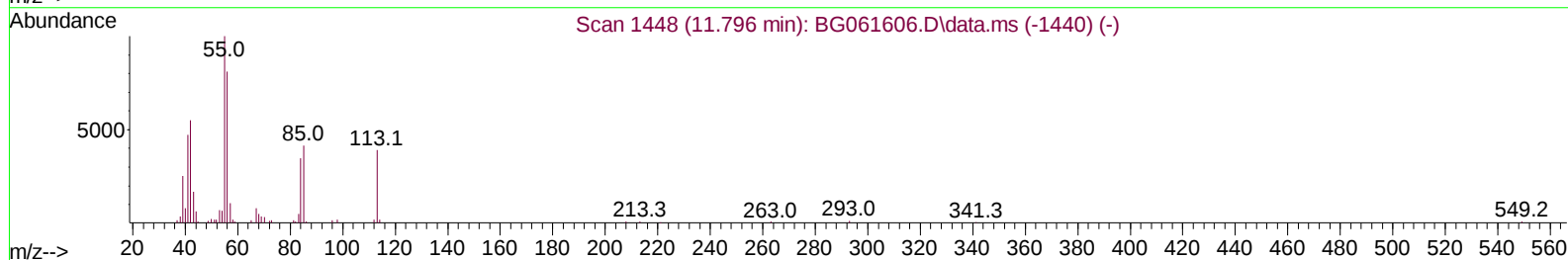
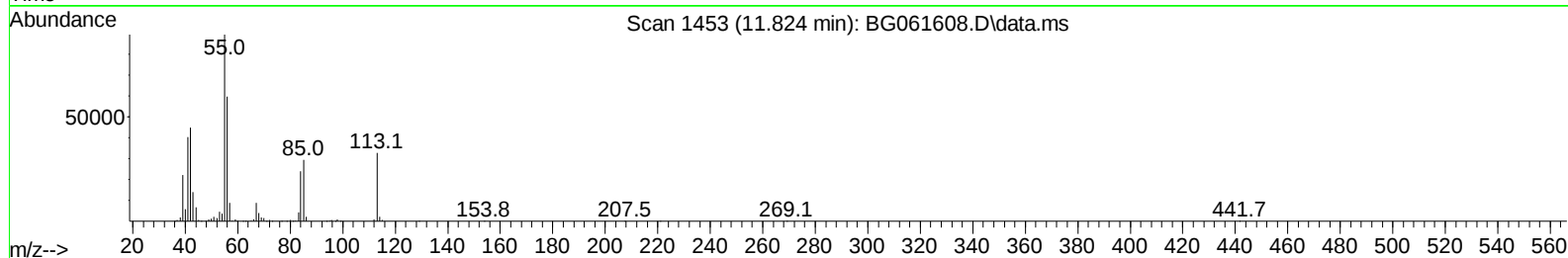
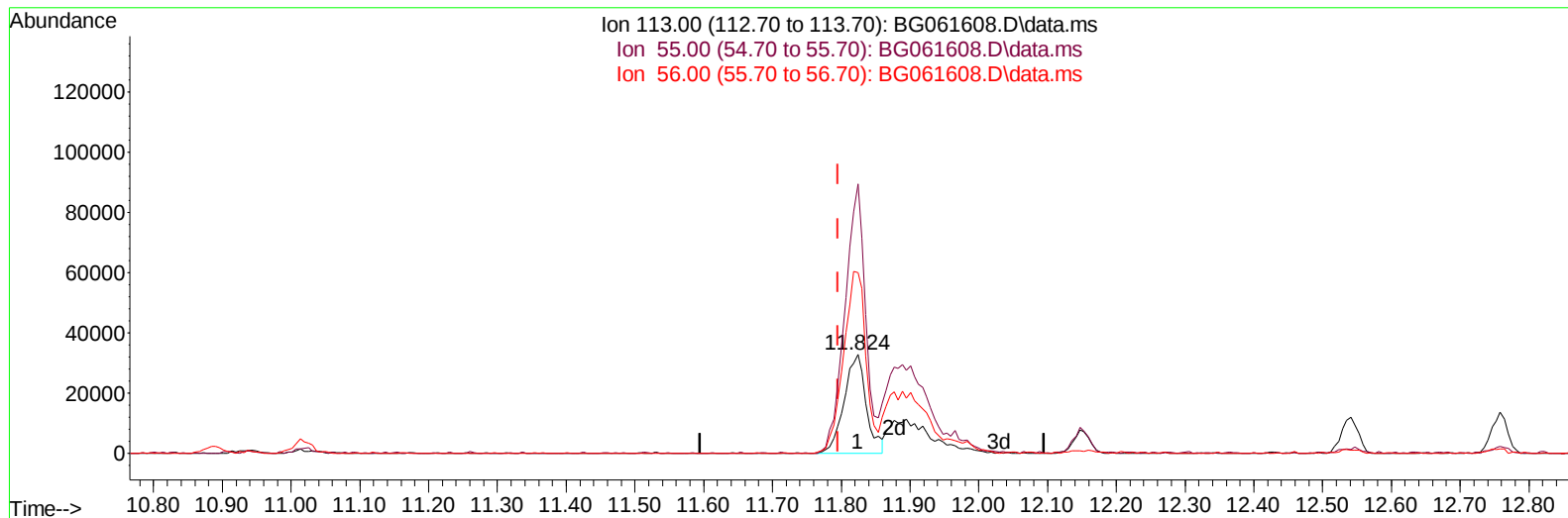
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062024\  
Data File : BG061608.D  
Acq On : 20 Jun 2024 16:27  
Operator : MA/JU  
Sample : SSTD08014  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD080443

Manual IntegrationsAPPROVED

Quant Time: Jun 20 17:04:02 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 20 16:29:55 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2024  
Supervised By :mohammad ahmed 06/21/2024



TIC: BG061608.D\data.ms

(34) Caprolactam

11.824min (+ 0.029) 55.27 ng/ul

response 73665

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 259.30 | 272.33 |
| 56.00  | 209.00 | 182.46 |
| 0.00   | 0.00   | 0.00   |

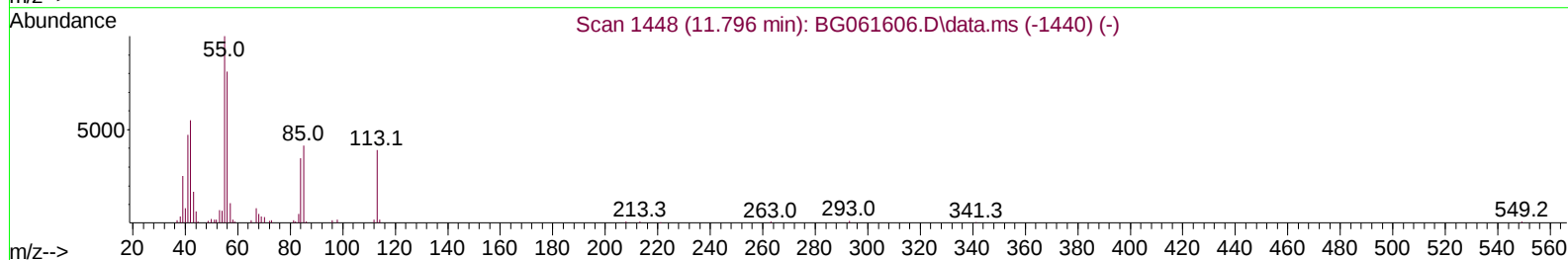
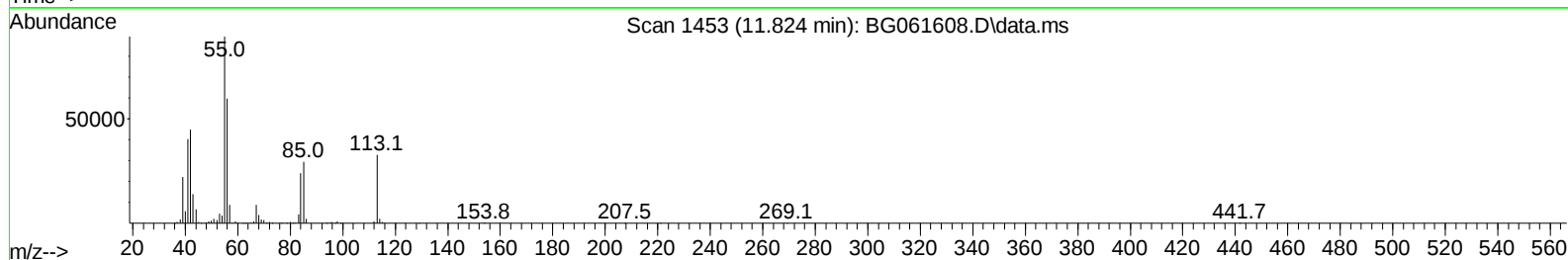
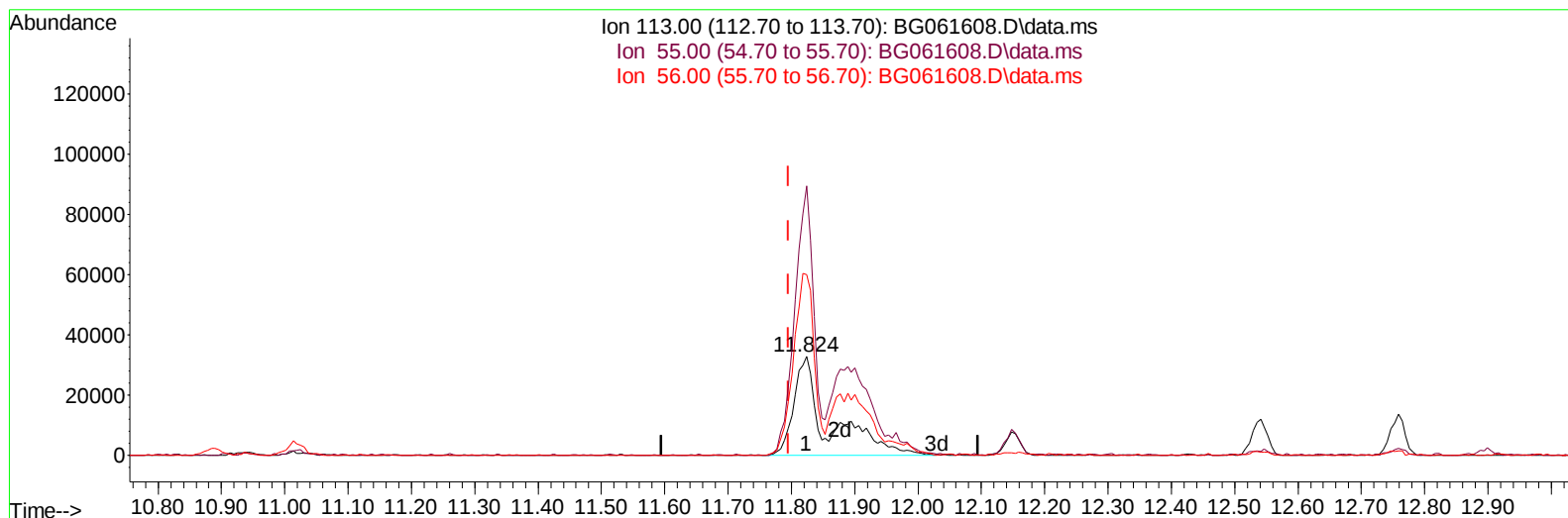
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062024\  
Data File : BG061608.D  
Acq On : 20 Jun 2024 16:27  
Operator : MA/JU  
Sample : SSTD08014  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD080443

Manual IntegrationsAPPROVED

Quant Time: Jun 20 17:04:02 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 20 16:29:55 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2024  
Supervised By :mohammad ahmed 06/21/2024



TIC: BG061608.D\data.ms

(34) Caprolactam

11.824min (+ 0.029) 90.97 ng/ul m

response 121247

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 259.30 | 272.33 |
| 56.00  | 209.00 | 182.46 |
| 0.00   | 0.00   | 0.00   |

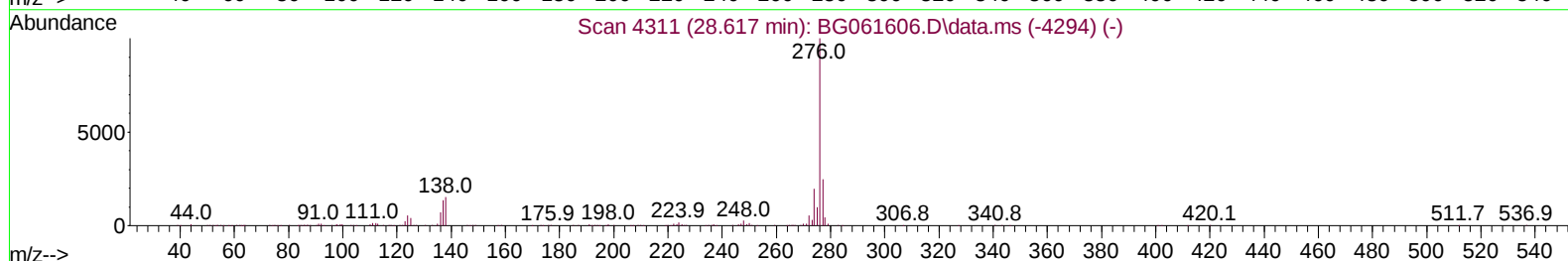
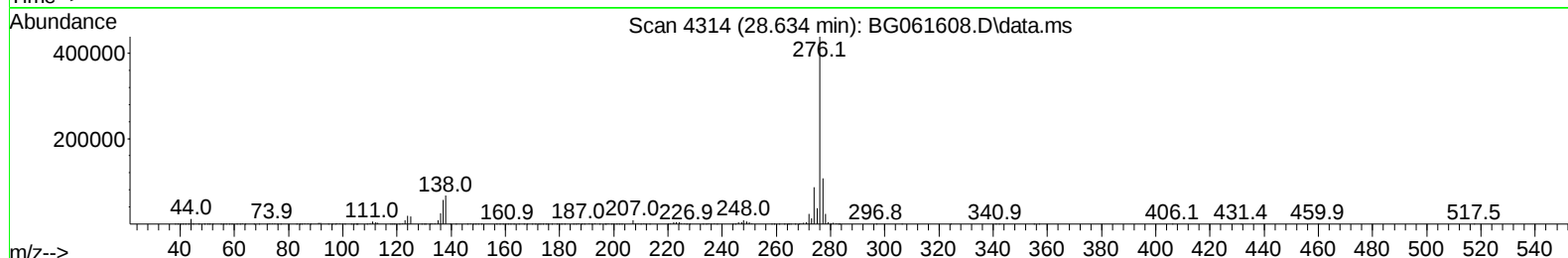
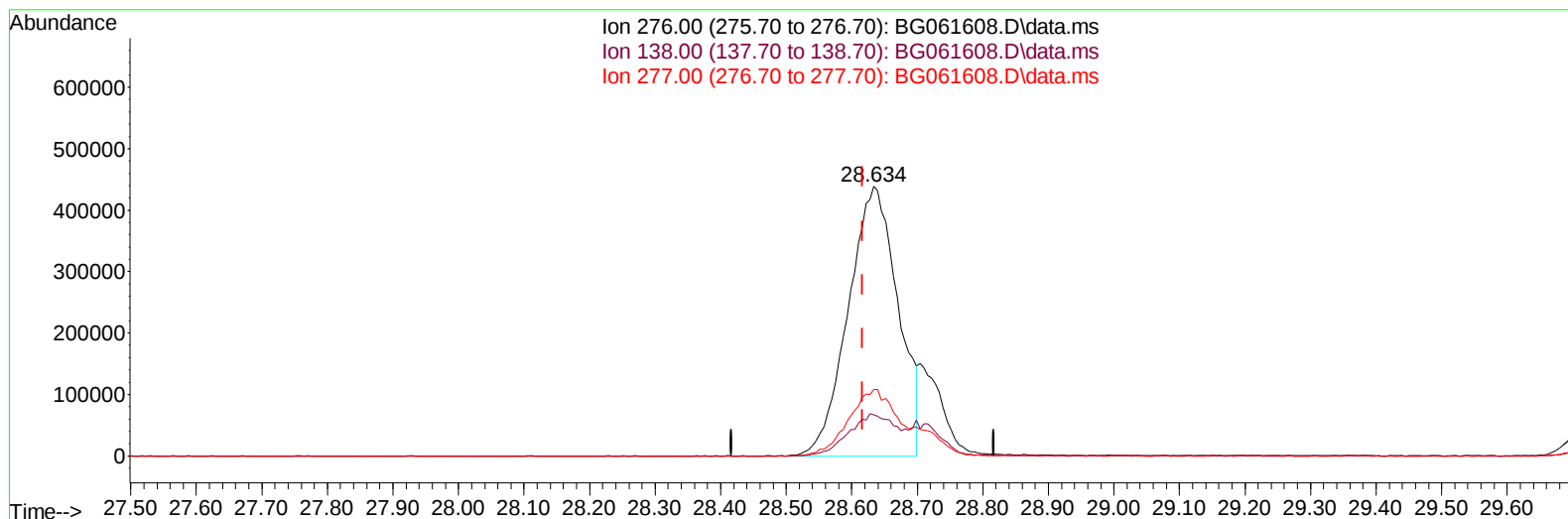
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062024\  
Data File : BG061608.D  
Acq On : 20 Jun 2024 16:27  
Operator : MA/JU  
Sample : SST08014  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD080443

Manual IntegrationsAPPROVED

Quant Time: Jun 20 17:04:02 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 20 16:29:55 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2024  
Supervised By :mohammad ahmed 06/21/2024



TIC: BG061608.D\data.ms

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

**28.634min (+ 0.017) 72.80 ng/u1**

**response 2305911**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 15.30  | 15.18  |
| 277.00 | 24.60  | 24.59  |
| 0.00   | 0.00   | 0.00   |

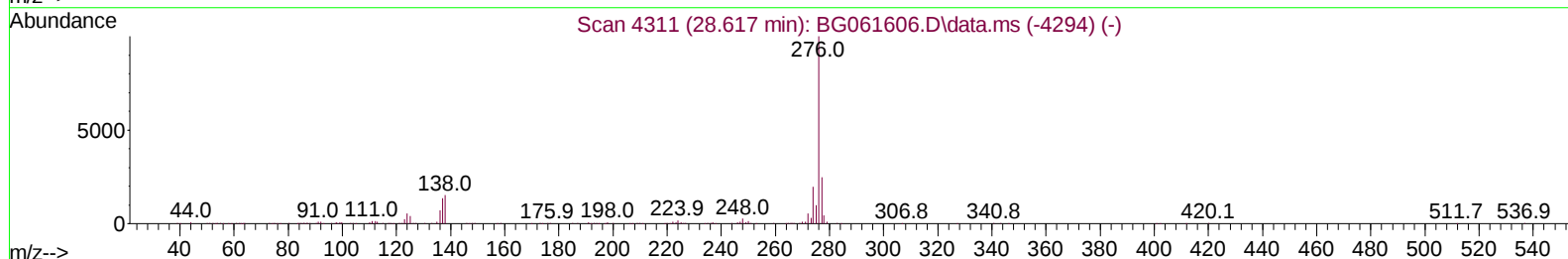
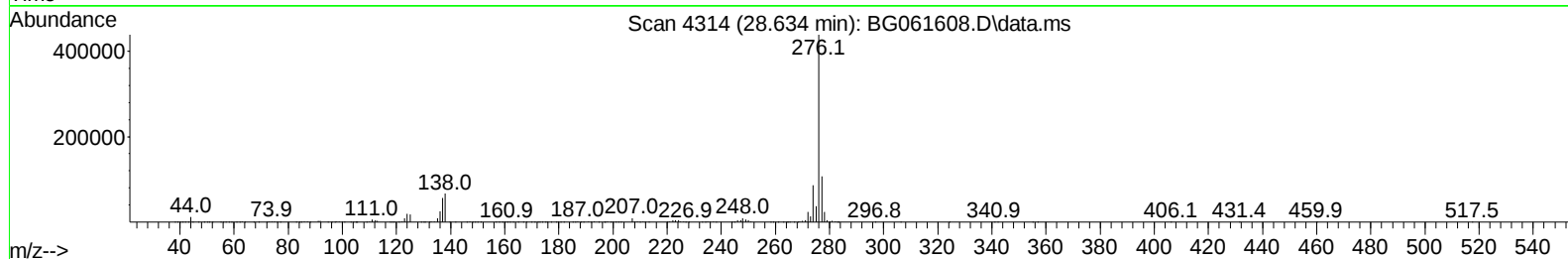
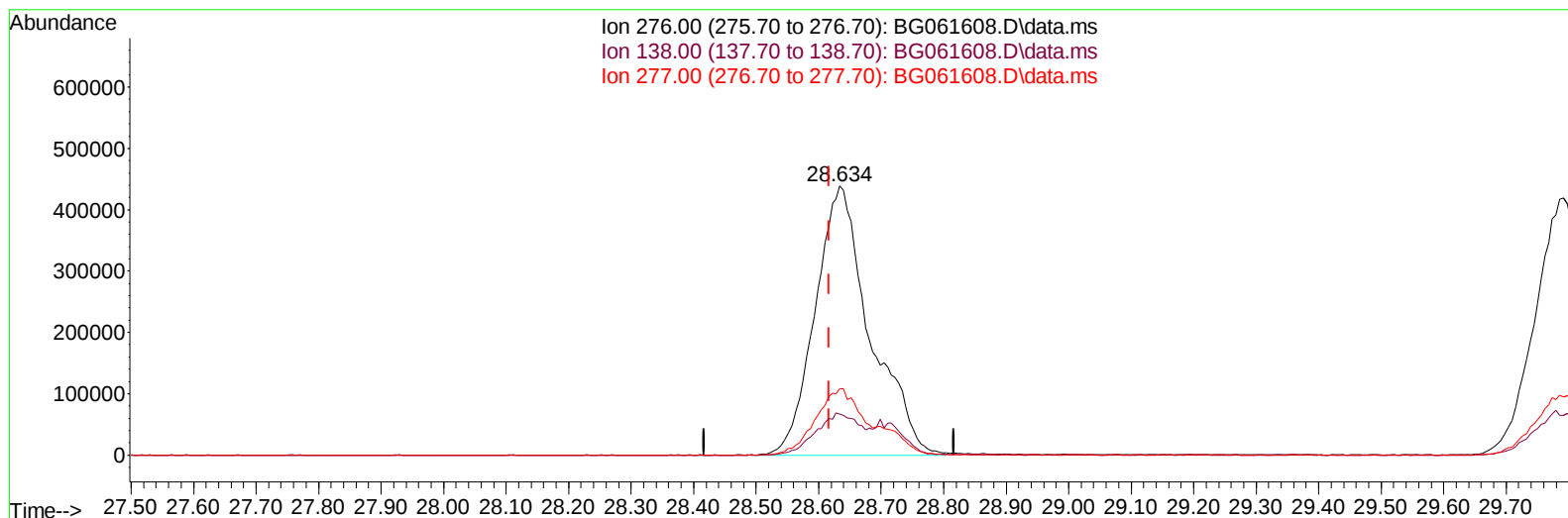
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062024\  
Data File : BG061608.D  
Acq On : 20 Jun 2024 16:27  
Operator : MA/JU  
Sample : SSTD08014  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD080443

Manual IntegrationsAPPROVED

Quant Time: Jun 20 17:04:02 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 20 16:29:55 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2024  
Supervised By :mohammad ahmed 06/21/2024



TIC: BG061608.D\data.ms

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

28.634min (+ 0.017) 84.43 ng/ul m

response 2674426

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 15.30  | 15.18  |
| 277.00 | 24.60  | 24.59  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062024\  
 Data File : BG061608.D  
 Acq On : 20 Jun 2024 16:27  
 Operator : MA/JU  
 Sample : SST08014  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST080443

Manual IntegrationsAPPROVED

Quant Time: Jun 20 17:05:33 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 20 16:29:55 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2024  
 Supervised By :mohammad ahmed 06/21/2024

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.076  | 152  | 46401    | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.890 | 136  | 233267   | 20.000  | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.703 | 164  | 180476   | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.447 | 188  | 421818   | 20.000  | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.713 | 240  | 383085   | 20.000  | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.938 | 264  | 458406   | 20.000  | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.469  | 96   | 36480    | 26.418  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.886  | 84   | 305621   | 69.417  | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.218  | 99   | 410512   | 87.265  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.400  | 67   | 243450   | 69.928  | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.600  | 132  | 280178   | 116.758 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.775  | 113  | 318406   | 90.156  | ng/ul  | 0.01     |
| 21) Nitrobenzene-d5           | 9.239  | 128  | 139975   | 123.854 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.962  | 143  | 157219   | 152.188 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.496 | 165  | 319639   | 103.817 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.019 | 131  | 476039   | 88.672  | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.116 | 166  | 1071823  | 93.882  | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 1196609  | 82.148  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.879 | 143  | 195017   | 114.295 | ng/ul  | 0.01     |
| 60) Fluorene-d10              | 15.696 | 176  | 920254   | 78.220  | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.802 | 200  | 164729   | 128.844 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.547 | 188  | 1463561  | 78.672  | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.827 | 212  | 1616919  | 77.237  | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.727 | 264  | 1773430  | 81.732  | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 2) 1,4-Dioxane                | 3.505  | 88   | 41922    | 25.218  | ng/uL# | 91       |
| 5) Pyridine                   | 3.904  | 79   | 318751   | 70.715  | ng/ul  | 98       |
| 6) Benzaldehyde               | 7.212  | 77   | 178391   | 60.749  | ng/ul  | 94       |
| 8) Phenol                     | 7.241  | 94   | 432268   | 87.156  | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.494  | 93   | 308933   | 79.144  | ng/ul# | 97       |
| 12) 2-Chlorophenol            | 7.635  | 128  | 280714   | 107.336 | ng/ul  | 94       |
| 13) 2-Methylphenol            | 8.505  | 108  | 299941   | 86.465  | ng/ul  | 94       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.599  | 45   | 634658   | 75.128  | ng/ul  | 97       |
| 16) Acetophenone              | 8.904  | 105  | 502741   | 75.545  | ng/ul  | 96       |
| 17) N-Nitroso-di-n-propyla... | 8.892  | 70   | 271754   | 68.011  | ng/ul# | 97       |
| 18) 4-Methylphenol            | 8.839  | 108  | 334092   | 87.959  | ng/ul  | 97       |
| 19) Hexachloroethane          | 9.163  | 117  | 109050   | 93.671  | ng/ul  | 91       |
| 22) Nitrobenzene              | 9.280  | 77   | 395451   | 86.619  | ng/ul  | 99       |
| 23) Isophorone                | 9.815  | 82   | 849490   | 68.294  | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.997  | 139  | 164838   | 141.804 | ng/ul# | 80       |
| 26) 2,4-Dimethylphenol        | 10.044 | 107  | 378560   | 84.017  | ng/ul  | 92       |
| 27) Bis(2-Chloroethoxy)met... | 10.297 | 93   | 461475   | 72.708  | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.526 | 162  | 288706   | 99.346  | ng/ul  | 96       |
| 30) Naphthalene               | 10.937 | 128  | 1000816  | 80.221  | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 11.043 | 127  | 463761   | 87.979  | ng/ul  | 94       |
| 33) Hexachlorobutadiene       | 11.219 | 225  | 189985   | 65.060  | ng/ul  | 92       |
| 34) Caprolactam               | 11.824 | 113  | 121247m  | 90.966  | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.153 | 107  | 401204   | 90.030  | ng/ul  | 98       |

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 Data File : BG061608.D  
 Acq On : 20 Jun 2024 16:27  
 Operator : MA/JU  
 Sample : SST08014  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :

BNA\_G

ClientSampleId :

SSTD080443

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/21/2024  
 Supervised By :mohammad ahmed 06/21/2024

Quant Time: Jun 20 17:05:33 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 20 16:29:55 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.541 | 142  | 743480   | 86.643  | ng/ul  | 98       |
| 37) 1-Methylnaphthalene       | 12.758 | 142  | 765660   | 86.368  | ng/ul  | 94       |
| 39) 1,2,4,5-Tetrachloroben... | 12.899 | 216  | 405434   | 76.460  | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.882 | 237  | 244535   | 102.099 | ng/ul# | 97       |
| 41) 2,4,6-Trichlorophenol     | 13.134 | 196  | 271894   | 101.781 | ng/ul  | 94       |
| 42) 2,4,5-Trichlorophenol     | 13.205 | 196  | 299039   | 101.407 | ng/ul  | 91       |
| 43) 1,1'-Biphenyl             | 13.546 | 154  | 1006365  | 82.146  | ng/ul  | 98       |
| 44) 2-Chloronaphthalene       | 13.587 | 162  | 790404   | 83.147  | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.781 | 65   | 308789   | 110.954 | ng/ul  | 99       |
| 47) Dimethylphthalate         | 14.163 | 163  | 1041465  | 86.727  | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.280 | 165  | 211791   | 137.268 | ng/ul  | 95       |
| 50) Acenaphthylene            | 14.427 | 152  | 1291483  | 78.953  | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.603 | 138  | 192565   | 100.562 | ng/ul# | 97       |
| 52) Acenaphthene              | 14.768 | 153  | 870119   | 79.650  | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 14.803 | 184  | 101702   | 117.567 | ng/ul  | 87       |
| 55) 4-Nitrophenol             | 14.897 | 109  | 197245   | 90.455  | ng/ul  | 97       |
| 56) Dibenzofuran              | 15.103 | 168  | 1232584  | 82.246  | ng/ul  | 91       |
| 57) 2,4-Dinitrotoluene        | 15.062 | 165  | 305304   | 132.840 | ng/ul  | 93       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.320 | 232  | 278340   | 107.002 | ng/ul  | 99       |
| 59) Diethylphthalate          | 15.526 | 149  | 1069143  | 92.908  | ng/ul  | 97       |
| 61) Fluorene                  | 15.749 | 166  | 994259   | 78.278  | ng/ul  | 94       |
| 62) 4-Chlorophenyl-phenyle... | 15.743 | 204  | 492538   | 72.365  | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 15.767 | 138  | 195304   | 92.450  | ng/ul  | 93       |
| 66) 4,6-Dinitro-2-methylph... | 15.820 | 198  | 181539   | 123.530 | ng/ul# | 94       |
| 67) N-Nitrosodiphenylamine    | 15.955 | 169  | 860248   | 79.694  | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.636 | 248  | 342199   | 83.293  | ng/ul  | 97       |
| 69) Hexachlorobenzene         | 16.742 | 284  | 405474   | 86.006  | ng/ul  | 98       |
| 70) Atrazine                  | 16.907 | 200  | 335399   | 94.421  | ng/ul  | 97       |
| 71) Pentachlorophenol         | 17.083 | 266  | 267276   | 110.293 | ng/ul  | 95       |
| 72) Phenanthrene              | 17.488 | 178  | 1625940  | 77.094  | ng/ul  | 100      |
| 74) Anthracene                | 17.582 | 178  | 1664078  | 76.835  | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.505 | 216  | 400020   | 77.423  | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 15.015 | 250  | 462615   | 90.019  | ng/uL  | 94       |
| 77) Carbazole                 | 17.847 | 167  | 1489837  | 82.676  | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.416 | 149  | 1722182  | 97.192  | ng/ul  | 98       |
| 80) Fluoranthene              | 19.492 | 202  | 1871165  | 74.050  | ng/ul  | 98       |
| 82) Pyrene                    | 19.856 | 202  | 1948822  | 73.143  | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.749 | 149  | 787504   | 118.743 | ng/ul  | 97       |
| 84) 3,3'-Dichlorobenzidine    | 21.607 | 252  | 625163   | 85.381  | ng/ul  | 97       |
| 85) Benzo(a)anthracene        | 21.695 | 228  | 2034590  | 77.241  | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.619 | 149  | 1140741  | 109.317 | ng/ul  | 97       |
| 87) Chrysene                  | 21.760 | 228  | 1914696  | 76.751  | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.852 | 149  | 2041584  | 95.400  | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.922 | 252  | 2217314  | 81.275  | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene      | 23.992 | 252  | 2157384  | 76.703  | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.797 | 252  | 2101122  | 79.693  | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.634 | 276  | 2674426m | 84.433  | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 28.716 | 278  | 2245157  | 86.415  | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene      | 29.791 | 276  | 2150932  | 83.603  | ng/ul  | 99       |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTD080443

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 06/21/2024

Supervised By :mohammad ahmed 06/21/2024

