

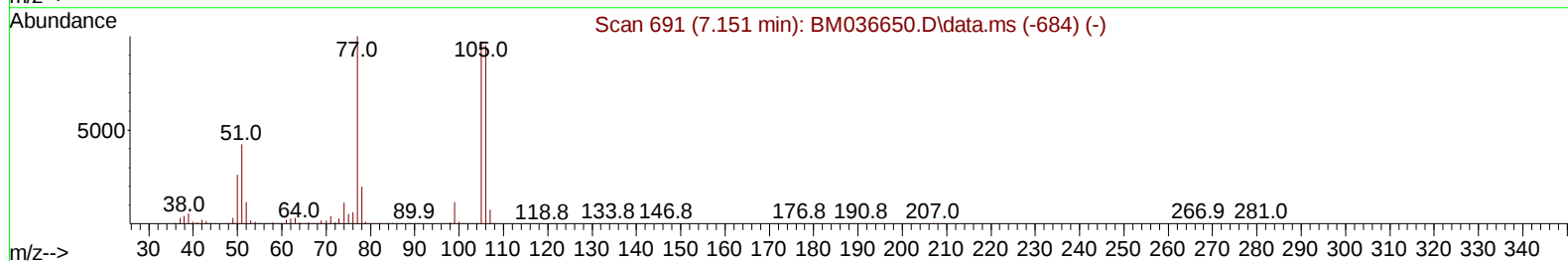
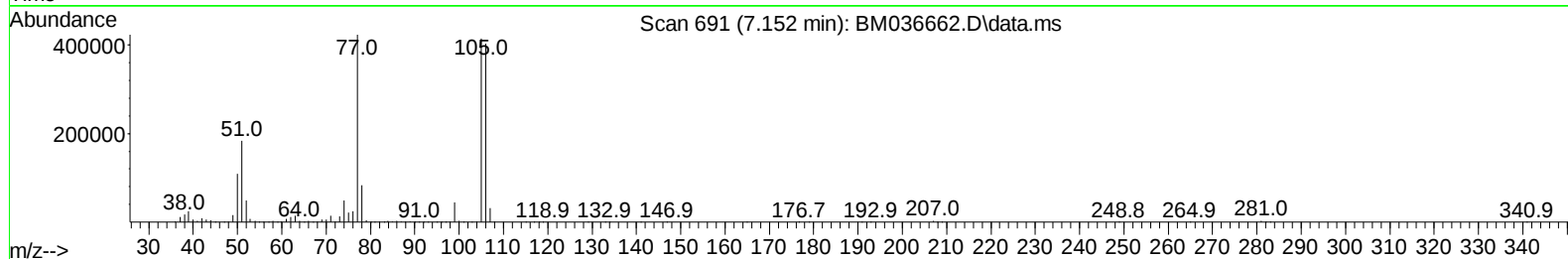
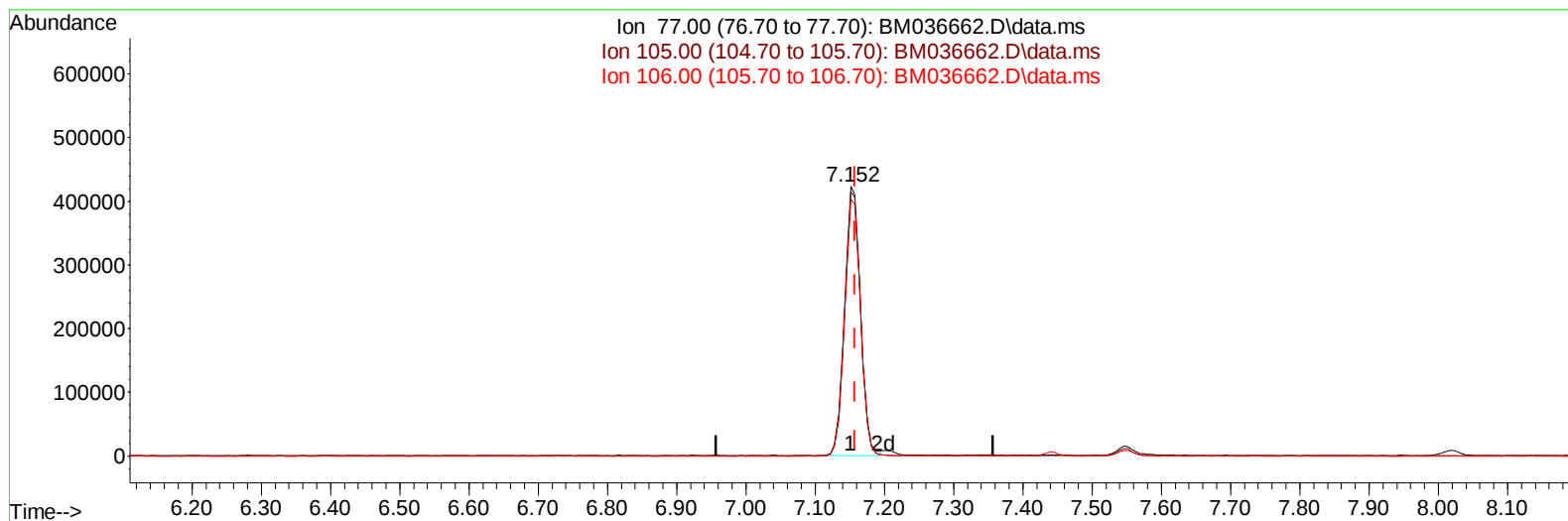
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
Data File : BM036662.D  
Acq On : 22 Sep 2022 14:44  
Operator : CG/JU  
Sample : PB147686BS  
Misc :  
ALS Vial : 48 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Sep 21 22:42:41 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
Supervised By :mohammad ahmed 09/23/2022



TIC: BM036662.D\data.ms

(6) Benzaldehyde

7.152min (-0.006) 43.54 ng/u1

response 673418

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 93.90  | 98.02  |
| 106.00 | 91.30  | 95.31  |
| 0.00   | 0.00   | 0.00   |

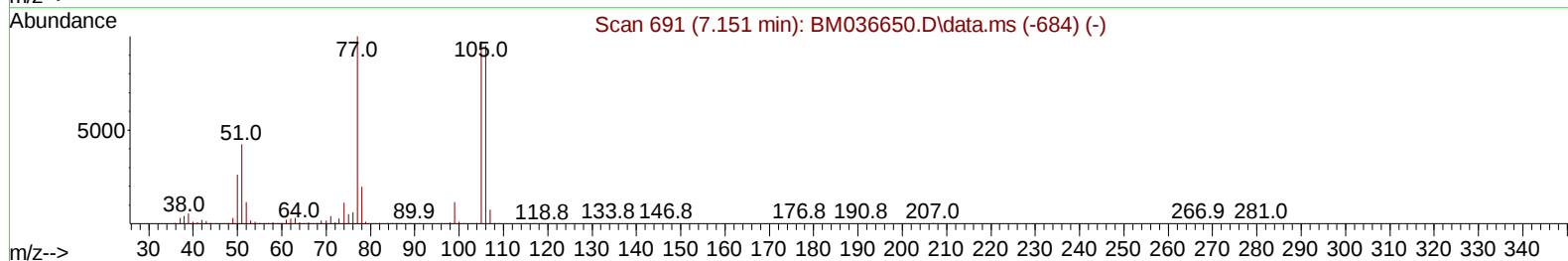
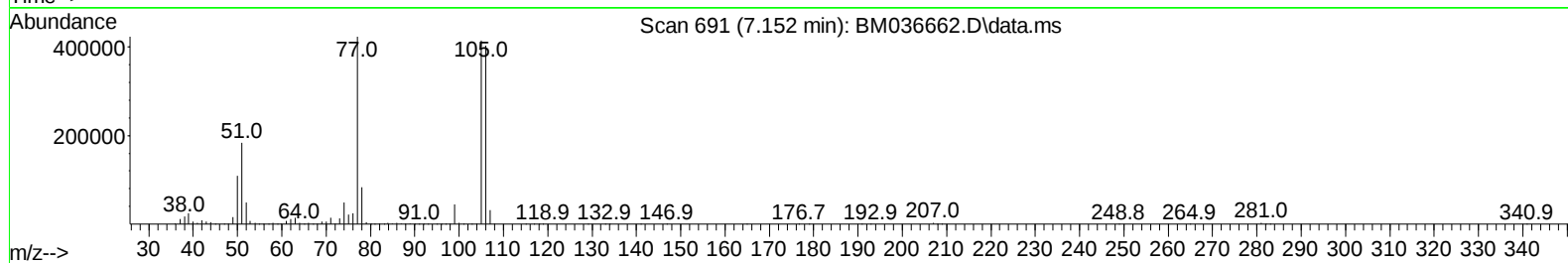
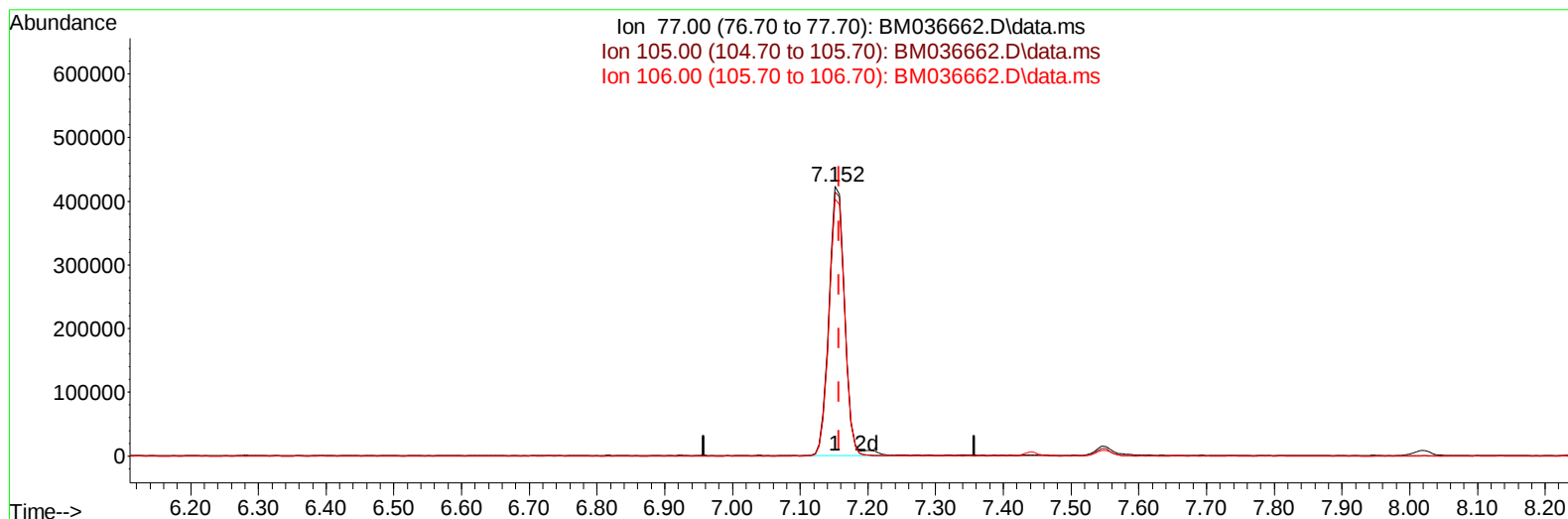
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
 Data File : BM036662.D  
 Acq On : 22 Sep 2022 14:44  
 Operator : CG/JU  
 Sample : PB147686BS  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 21 22:42:41 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
 Supervised By :mohammad ahmed 09/23/2022



TIC: BM036662.D\data.ms

**(6) Benzaldehyde**

7.152min (-0.006) 44.36 ng/ul m

response 686088

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 93.90  | 98.02  |
| 106.00 | 91.30  | 95.31  |
| 0.00   | 0.00   | 0.00   |

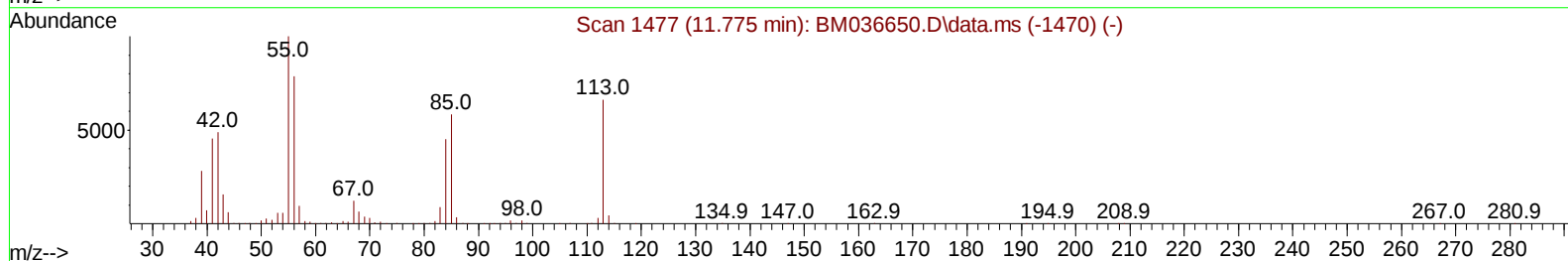
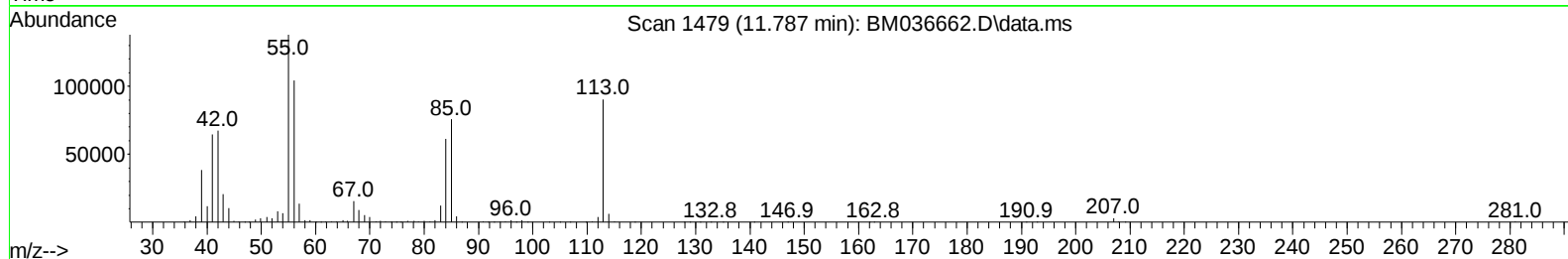
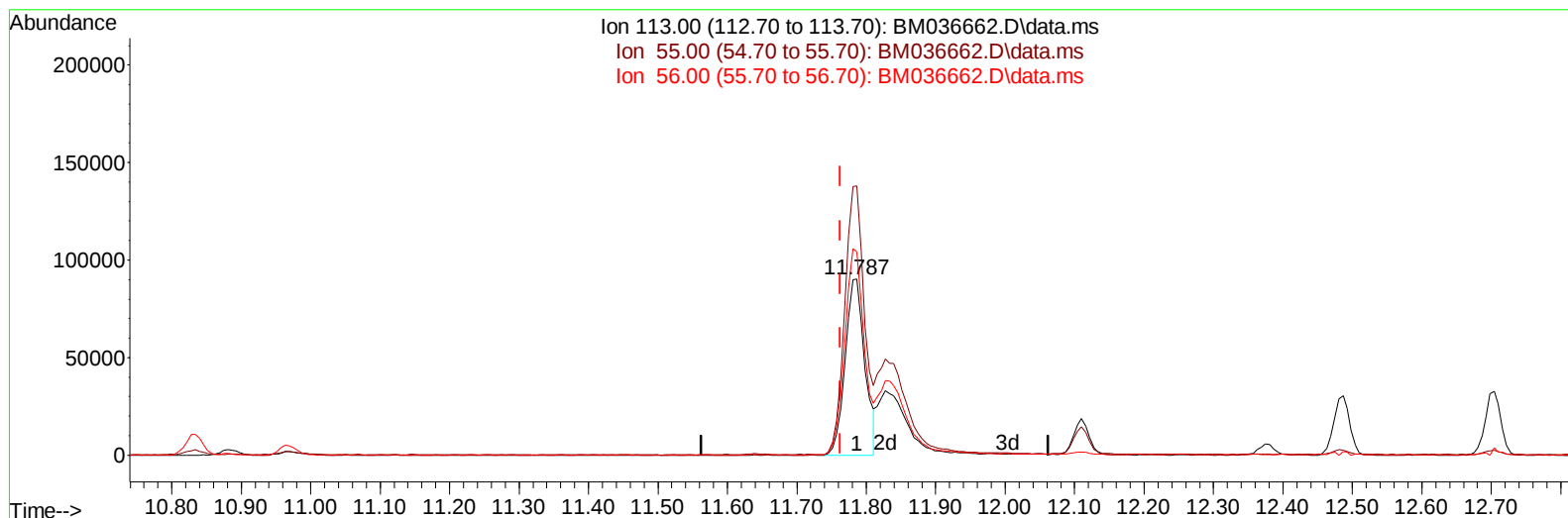
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
 Data File : BM036662.D  
 Acq On : 22 Sep 2022 14:44  
 Operator : CG/JU  
 Sample : PB147686BS  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 21 22:42:41 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
 Supervised By :mohammad ahmed 09/23/2022



TIC: BM036662.D\data.ms

**(34) Caprolactam**

11.787min (+ 0.024) 22.68 ng/ul

response 176548

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 159.40 | 152.65 |
| 56.00  | 118.30 | 115.31 |
| 0.00   | 0.00   | 0.00   |

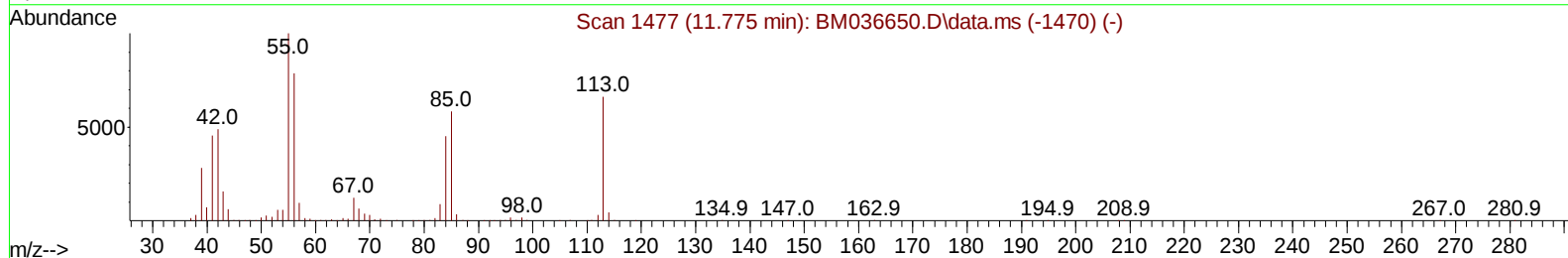
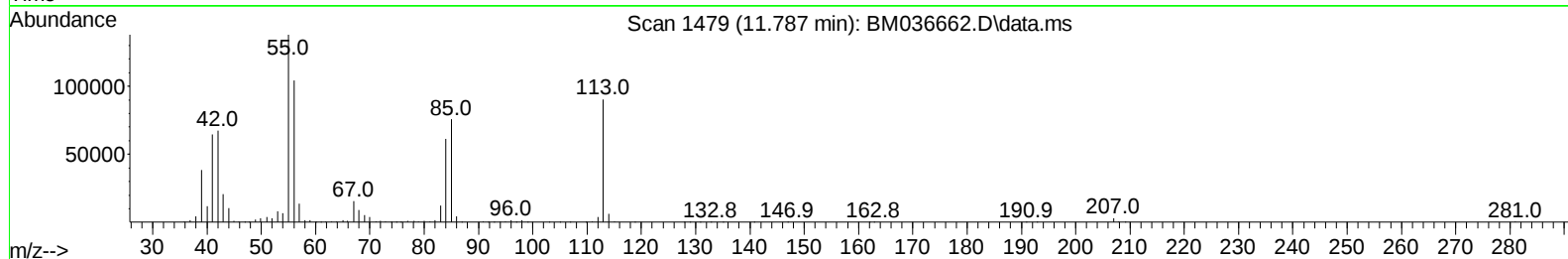
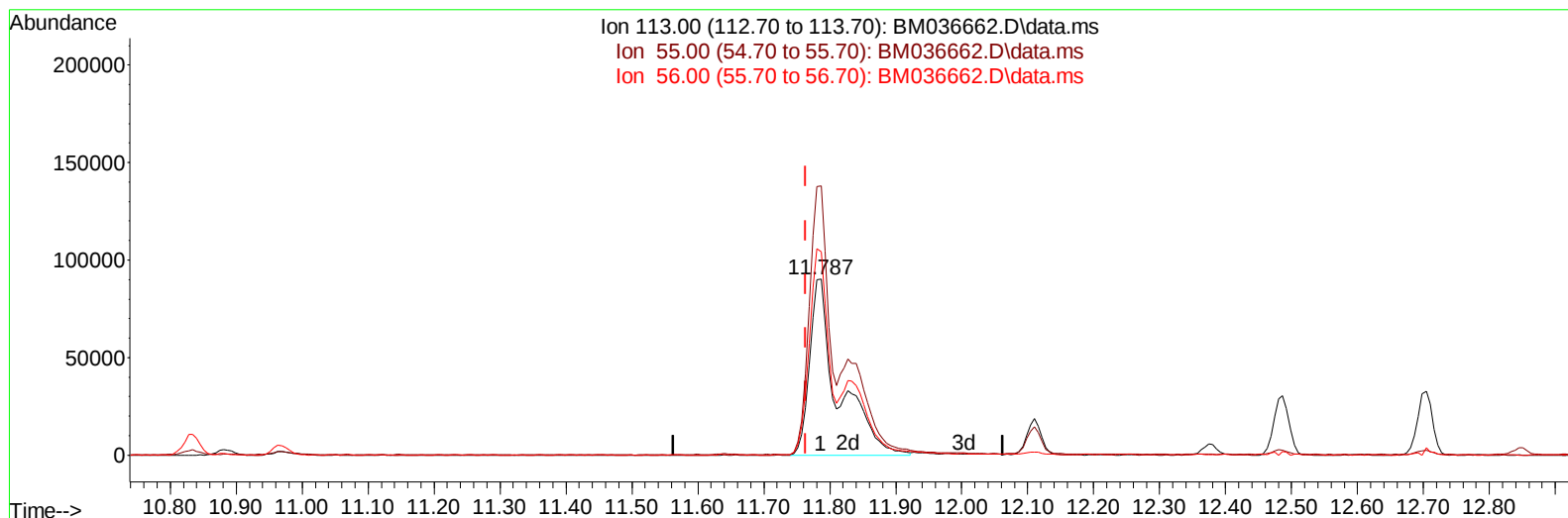
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
Data File : BM036662.D  
Acq On : 22 Sep 2022 14:44  
Operator : CG/JU  
Sample : PB147686BS  
Misc :  
ALS Vial : 48 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Sep 21 22:42:41 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
Supervised By :mohammad ahmed 09/23/2022



TIC: BM036662.D\data.ms

(34) Caprolactam

11.787min (+ 0.024) 34.78 ng/ul m

response 270711

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 159.40 | 152.65 |
| 56.00  | 118.30 | 115.31 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
 Data File : BM036662.D  
 Acq On : 22 Sep 2022 14:44  
 Operator : CG/JU  
 Sample : PB147686BS  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 21 22:42:41 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
 Supervised By :mohammad ahmed 09/23/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.016  | 152  | 357374   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.834 | 136  | 1624380  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 1013886  | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.380 | 188  | 2003313  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.551 | 240  | 1549857  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 1246555  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 58352    | 6.121  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.817  | 84   | 678169   | 24.392 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.175  | 99   | 1182702  | 37.033 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.346  | 67   | 696582   | 35.916 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.546  | 132  | 913078   | 39.711 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.734  | 113  | 941110   | 39.881 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.187  | 128  | 459006   | 41.748 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.910  | 143  | 460259   | 46.736 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.451 | 165  | 895933   | 39.256 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 778685   | 20.521 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.057 | 166  | 2571709  | 37.609 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.339 | 160  | 3144514  | 38.768 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.839 | 143  | 495084   | 37.328 | ng/ul  | 0.01     |
| 60) Fluorene-d10              | 15.633 | 176  | 2281117  | 37.225 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 340526   | 42.963 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.480 | 188  | 3354106  | 36.618 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 3528787  | 46.281 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.839 | 264  | 2397461  | 39.366 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.428  | 88   | 115334   | 11.231 | ng/uL  | 95       |
| 5) Pyridine                   | 3.840  | 79   | 815244   | 29.186 | ng/ul  | 93       |
| 6) Benzaldehyde               | 7.152  | 77   | 686088m  | 44.364 | ng/ul  |          |
| 8) Phenol                     | 7.205  | 94   | 1172975  | 34.221 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.440  | 93   | 892606   | 33.470 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.581  | 128  | 899667   | 36.172 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.463  | 108  | 877875   | 34.844 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.557  | 45   | 1049237  | 33.195 | ng/ul  | 98       |
| 16) Acetophenone              | 8.851  | 105  | 1396150  | 34.109 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.840  | 70   | 705056   | 35.699 | ng/ul  | 95       |
| 18) 4-Methylphenol            | 8.793  | 108  | 966210   | 35.465 | ng/ul  | 99       |
| 19) Hexachloroethane          | 9.104  | 117  | 347678   | 34.511 | ng/ul  | 94       |
| 22) Nitrobenzene              | 9.228  | 77   | 1034009  | 35.422 | ng/ul  | 97       |
| 23) Isophorone                | 9.763  | 82   | 1973772  | 33.162 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.945  | 139  | 481444   | 39.580 | ng/ul  | 99       |
| 26) 2,4-Dimethylphenol        | 10.004 | 107  | 994703   | 33.156 | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.245 | 93   | 1193902  | 33.654 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.475 | 162  | 858965   | 35.800 | ng/ul  | 99       |
| 30) Naphthalene               | 10.881 | 128  | 2950739  | 33.345 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.993 | 127  | 1179783  | 30.115 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 11.169 | 225  | 481130   | 32.787 | ng/ul  | 99       |
| 34) Caprolactam               | 11.787 | 113  | 270711m  | 34.782 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.110 | 107  | 951659   | 35.931 | ng/ul  | 98       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092122\  
 Data File : BM036662.D  
 Acq On : 22 Sep 2022 14:44  
 Operator : CG/JU  
 Sample : PB147686BS  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS686

Manual IntegrationsAPPROVED

Quant Time: Sep 23 01:01:04 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 21 22:42:41 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/23/2022  
 Supervised By :mohammad ahmed 09/23/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.487 | 142  | 1994527  | 33.829 | ng/ul  | 100      |
| 37) 1-Methylnaphthalene       | 12.704 | 142  | 2022169  | 34.117 | ng/ul  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.845 | 216  | 967284   | 33.707 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.828 | 237  | 351490   | 23.503 | ng/ul  | 100      |
| 41) 2,4,6-Trichlorophenol     | 13.086 | 196  | 636970   | 36.251 | ng/ul  | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.157 | 196  | 687414   | 36.490 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.486 | 154  | 2553486  | 34.165 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.528 | 162  | 1986132  | 33.865 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.728 | 65   | 587203   | 40.678 | ng/ul  | 98       |
| 47) Dimethylphthalate         | 14.110 | 163  | 2440959  | 33.652 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.222 | 165  | 479970   | 40.699 | ng/ul  | 98       |
| 50) Acenaphthylene            | 14.369 | 152  | 3139990  | 33.860 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.551 | 138  | 549047   | 38.818 | ng/ul# | 95       |
| 52) Acenaphthene              | 14.710 | 153  | 2189964  | 33.751 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.751 | 184  | 208634   | 39.870 | ng/ul  | 96       |
| 55) 4-Nitrophenol             | 14.851 | 109  | 376074   | 33.170 | ng/ul  | 96       |
| 56) Dibenzofuran              | 15.045 | 168  | 2956689  | 33.930 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 15.004 | 165  | 687890   | 40.123 | ng/ul  | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.263 | 232  | 562782   | 35.688 | ng/ul# | 95       |
| 59) Diethylphthalate          | 15.463 | 149  | 2465040  | 33.948 | ng/ul  | 99       |
| 61) Fluorene                  | 15.692 | 166  | 2530293  | 34.093 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.680 | 204  | 1184140  | 33.362 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 15.710 | 138  | 567070   | 41.265 | ng/ul  | 97       |
| 66) 4,6-Dinitro-2-methylph... | 15.763 | 198  | 332522   | 37.178 | ng/ul  | 99       |
| 67) N-Nitrosodiphenylamine    | 15.892 | 169  | 2058116  | 34.882 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.574 | 248  | 656971   | 33.272 | ng/ul  | 99       |
| 69) Hexachlorobenzene         | 16.686 | 284  | 762635   | 32.655 | ng/ul  | 98       |
| 70) Atrazine                  | 16.845 | 200  | 561211   | 26.954 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.027 | 266  | 447261   | 32.241 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.427 | 178  | 3789638  | 34.344 | ng/ul  | 99       |
| 74) Anthracene                | 17.516 | 178  | 3788673  | 33.780 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.451 | 216  | 1005706  | 36.705 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.963 | 250  | 901679   | 30.426 | ng/uL  | 100      |
| 77) Carbazole                 | 17.786 | 167  | 3458592  | 33.306 | ng/ul  | 100      |
| 78) Di-n-butylphthalate       | 18.351 | 149  | 4178867  | 35.505 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.433 | 202  | 4070446  | 43.565 | ng/ul  | 99       |
| 82) Pyrene                    | 19.792 | 202  | 4187552  | 42.341 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.686 | 149  | 1683309  | 44.917 | ng/ul  | 93       |
| 84) 3,3'-Dichlorobenzidine    | 21.468 | 252  | 1059154  | 34.975 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.533 | 228  | 3443987  | 35.466 | ng/ul  | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.468 | 149  | 2514344  | 45.929 | ng/ul  | 99       |
| 87) Chrysene                  | 21.592 | 228  | 3190998  | 34.620 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.409 | 149  | 3816208  | 50.780 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.251 | 252  | 2963029  | 37.484 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.298 | 252  | 2837651  | 35.926 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 23.886 | 252  | 2486018  | 35.774 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.544 | 276  | 2782838  | 34.292 | ng/ul  | 96       |
| 95) Dibenzo(a,h)anthracene    | 26.562 | 278  | 2530202  | 34.380 | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene      | 27.327 | 276  | 2411170  | 34.273 | ng/ul  | 99       |

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

**Instrument :**

BNA\_M

**ClientSampleId :**

SLCS686

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 09/23/2022  
Supervised By :mohammad ahmed 09/23/2022

