

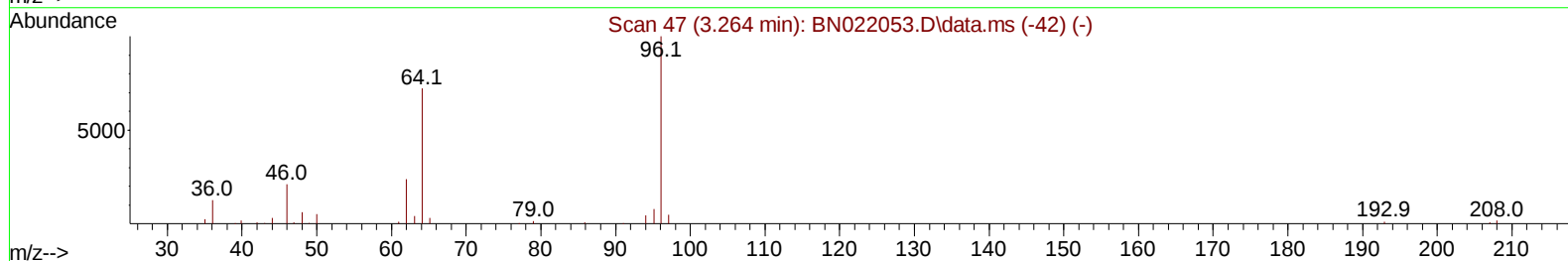
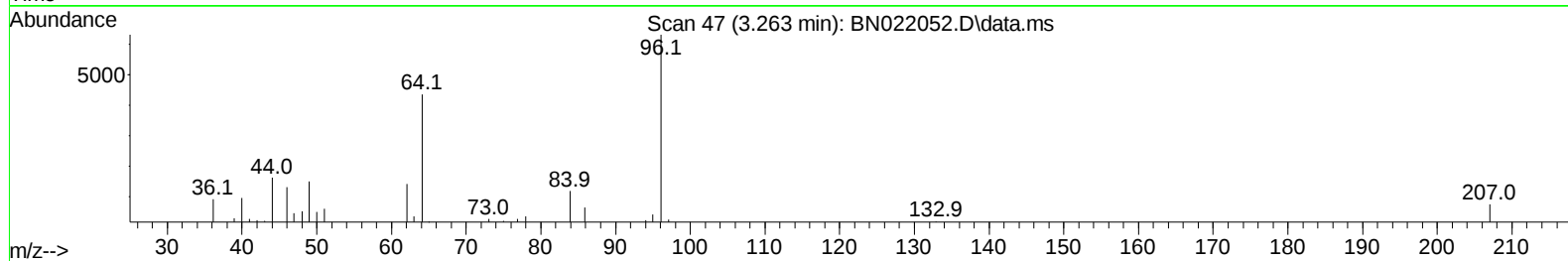
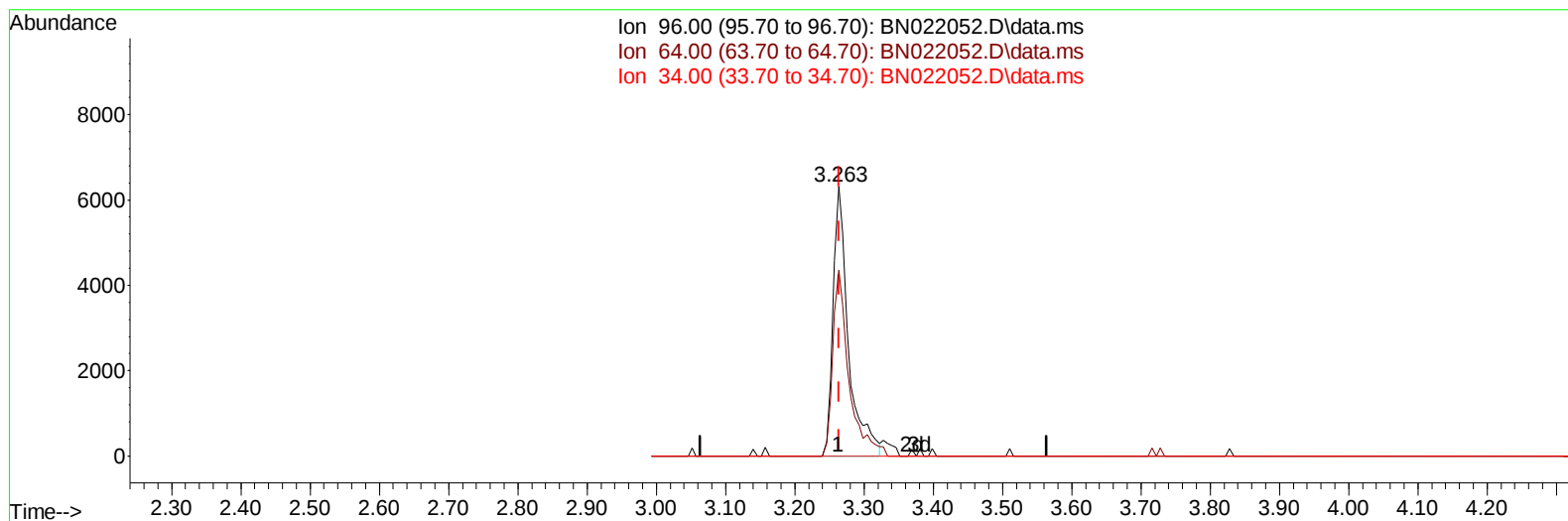
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SSTD01030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 03:00:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
 Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

**(3) 1,4-Dioxane-d8 (S)**

**3.263min (-0.000) 4.04 ng/uL**

**response 9734**

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 72.10  | 69.04  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

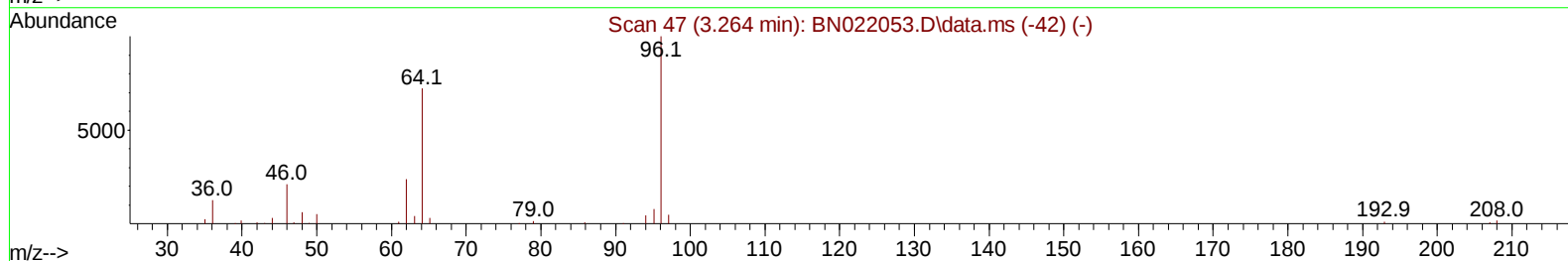
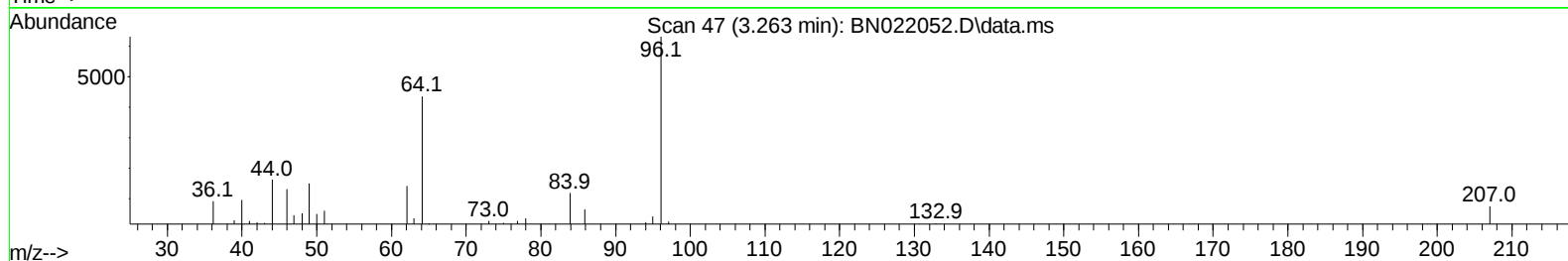
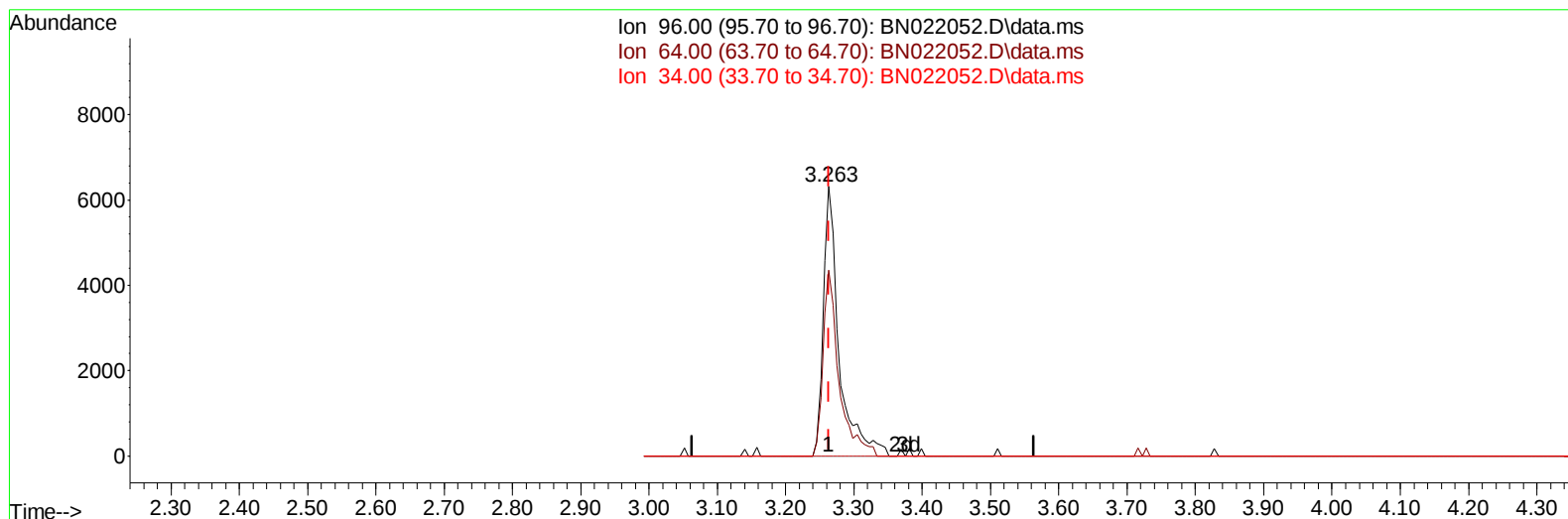
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 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SST01030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 03:00:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
 Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

**(3) 1,4-Dioxane-d8 (S)**

3.263min (-0.000) 4.20 ng/uL m

response 10123

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 72.10  | 69.04  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

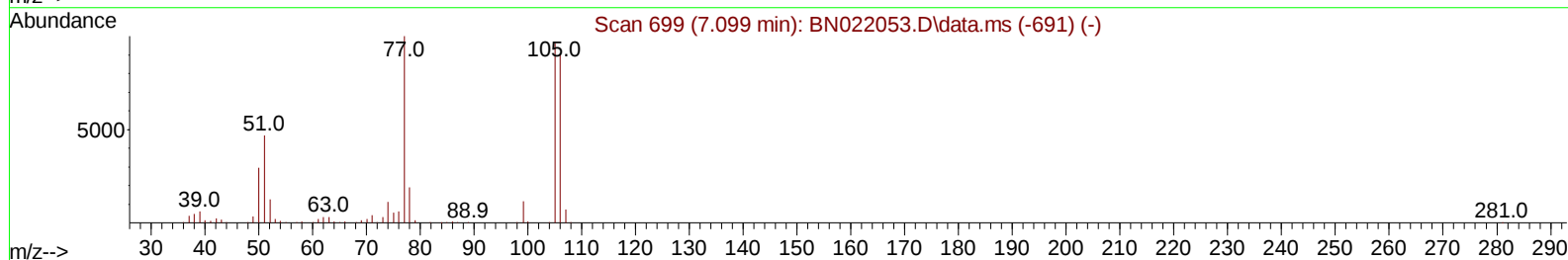
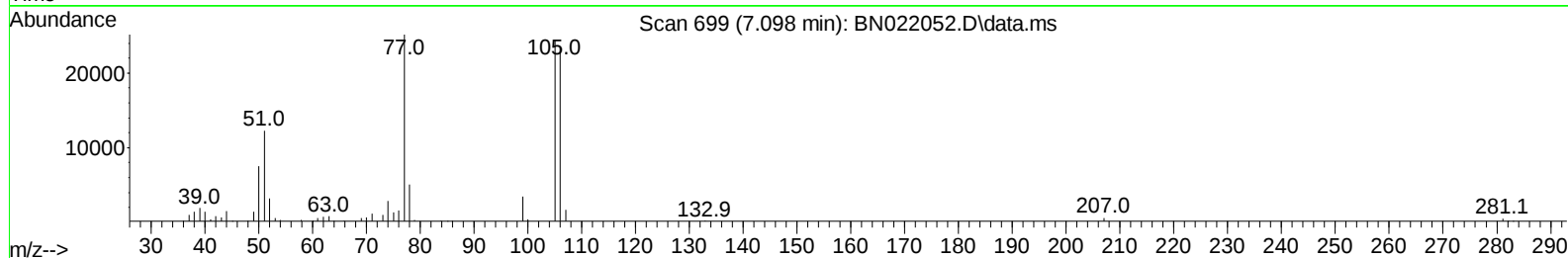
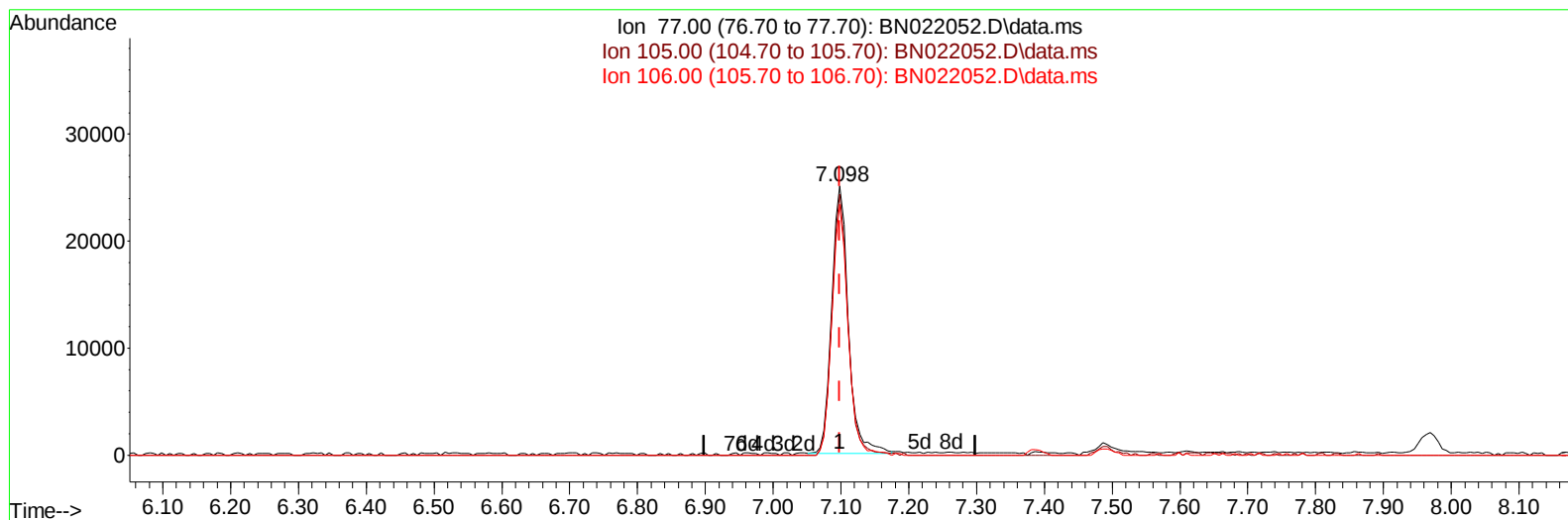
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SSTD01030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 03:00:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
 Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

**(6) Benzaldehyde**

7.098min (-0.000) 11.95 ng/u1

response 42609

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 96.30  | 97.07  |
| 106.00 | 90.40  | 93.29  |
| 0.00   | 0.00   | 0.00   |

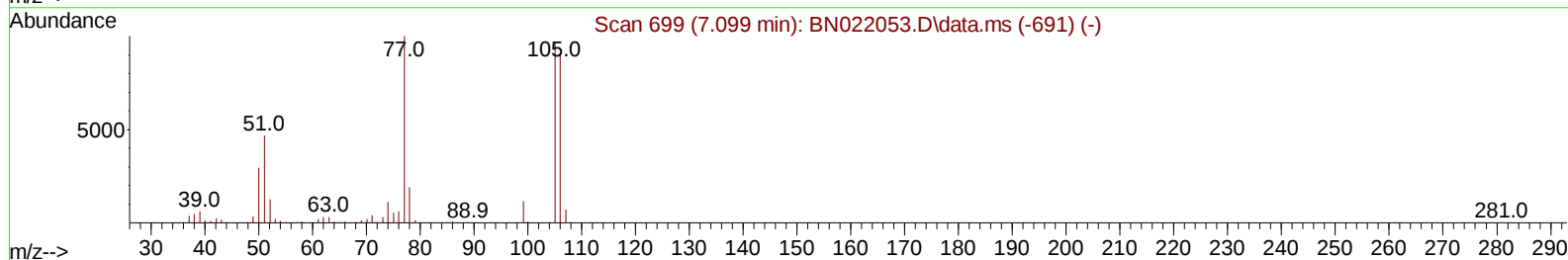
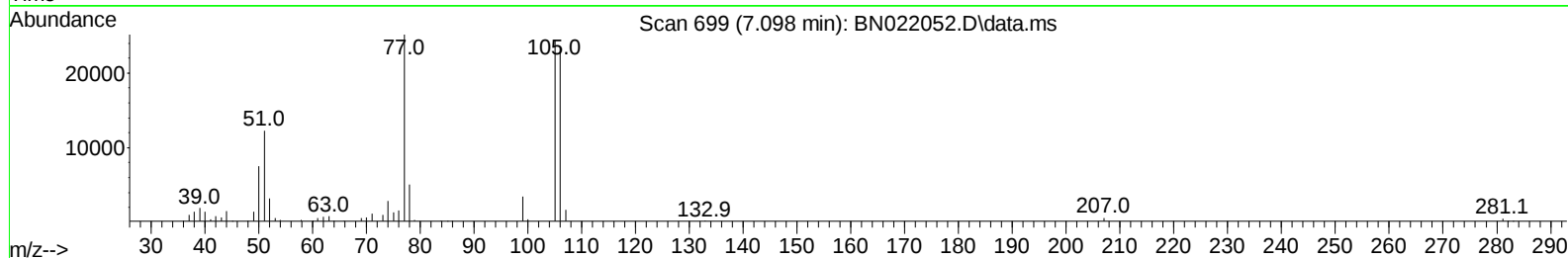
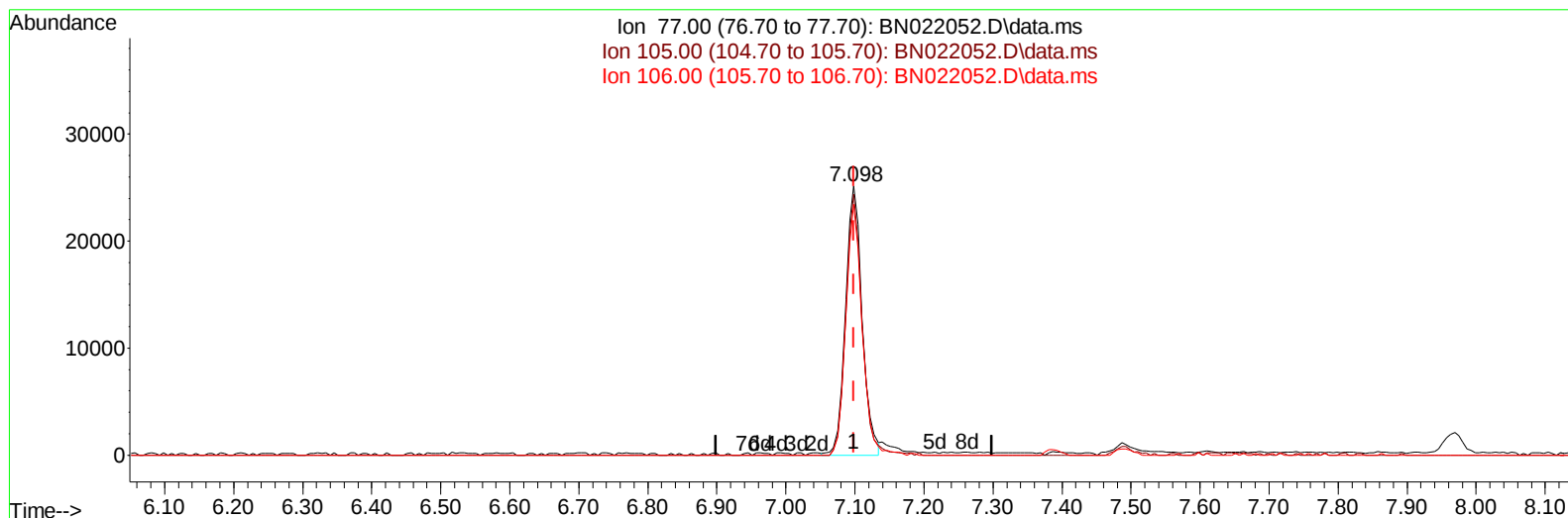
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
Data File : BN022052.D  
Acq On : 11 Oct 2022 15:01  
Operator : CG/JU  
Sample : SSTD01030  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 12 03:00:49 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

(6) Benzaldehyde

7.098min (-0.000) 11.78 ng/u1 m

response 42000

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 96.30  | 97.07  |
| 106.00 | 90.40  | 93.29  |
| 0.00   | 0.00   | 0.00   |

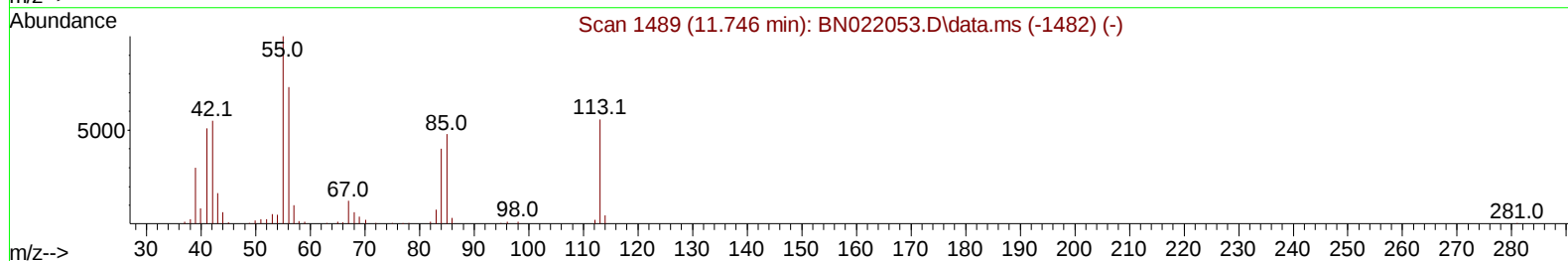
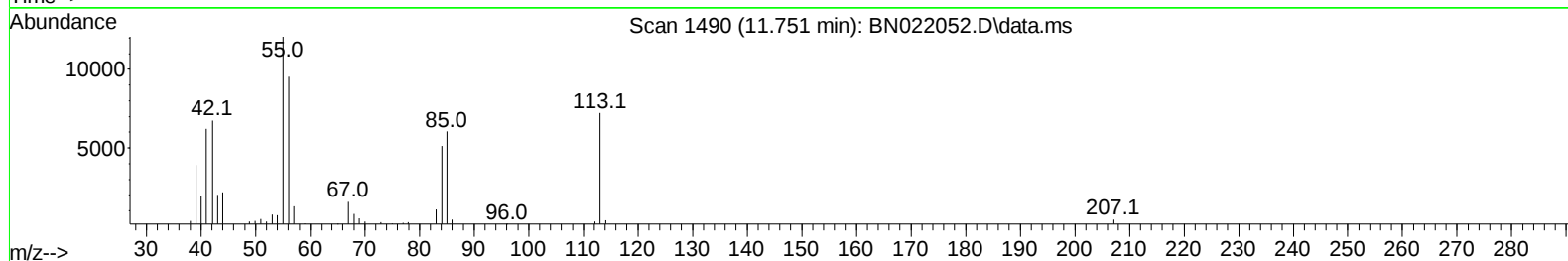
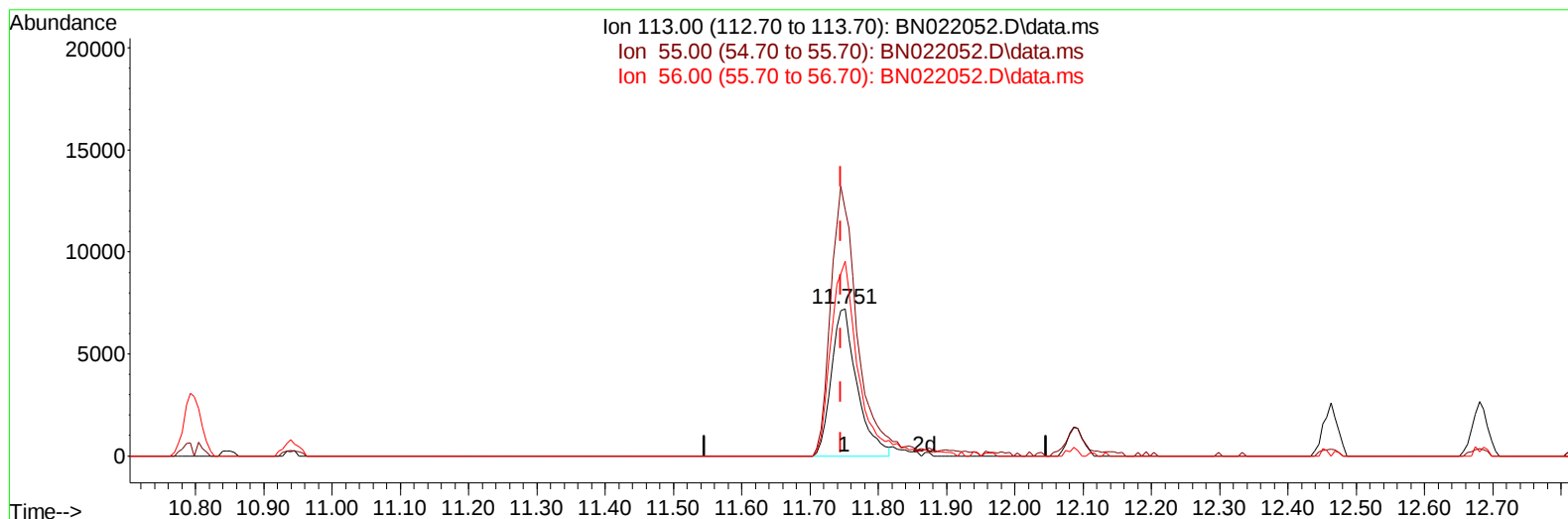
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
Data File : BN022052.D  
Acq On : 11 Oct 2022 15:01  
Operator : CG/JU  
Sample : SST01030  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SST010239

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022

Quant Time: Oct 12 03:06:53 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 12 03:00:49 2022  
Response via : Initial Calibration



TIC: BN022052.D\data.ms

## (34) Caprolactam

11.751min (+ 0.006) 8.35 ng/ul

response 18909

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 179.20 | 167.22 |
| 56.00  | 130.80 | 131.98 |
| 0.00   | 0.00   | 0.00   |

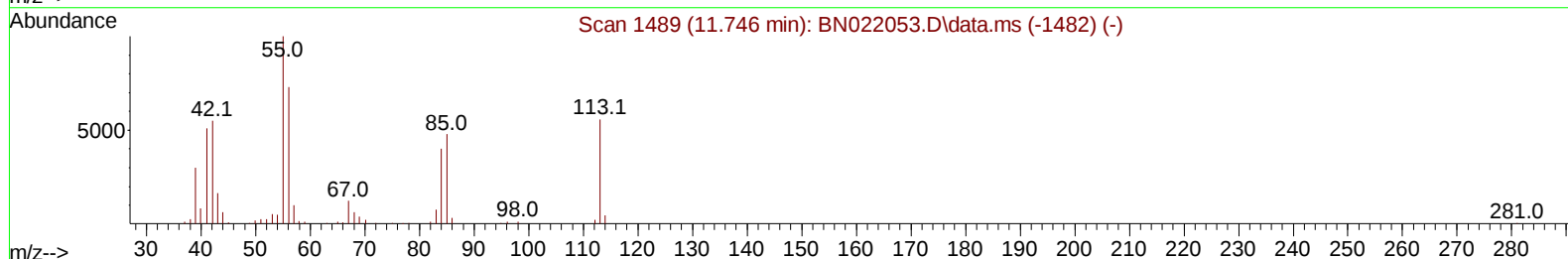
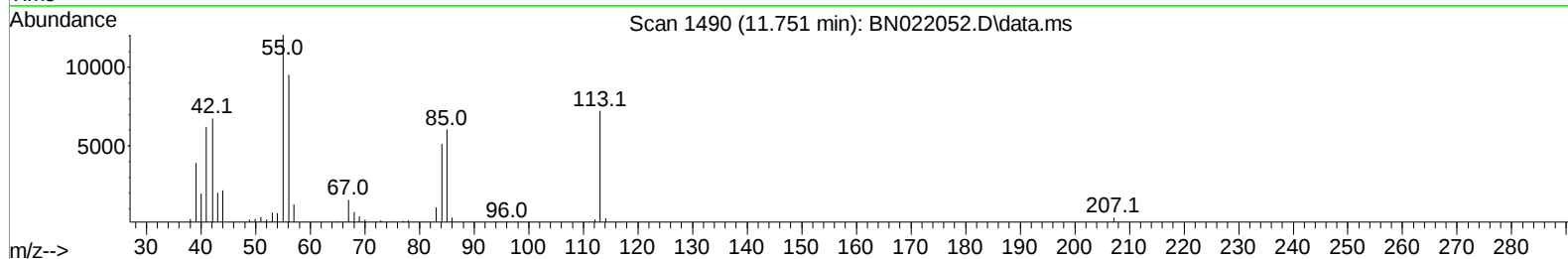
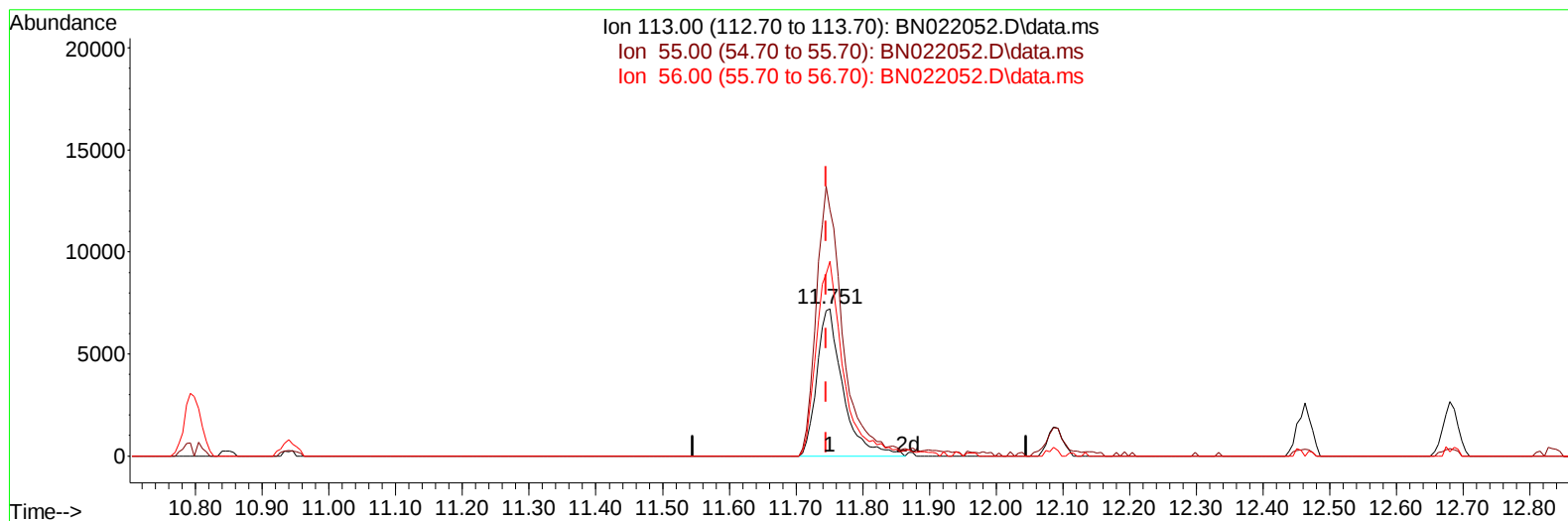
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
Data File : BN022052.D  
Acq On : 11 Oct 2022 15:01  
Operator : CG/JU  
Sample : SST01030  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SST010239

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022

Quant Time: Oct 12 03:06:53 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 12 03:00:49 2022  
Response via : Initial Calibration



TIC: BN022052.D\data.ms

## (34) Caprolactam

11.751min (+ 0.006) 8.65 ng/ul m

response 19597

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 179.20 | 167.22 |
| 56.00  | 130.80 | 131.98 |
| 0.00   | 0.00   | 0.00   |

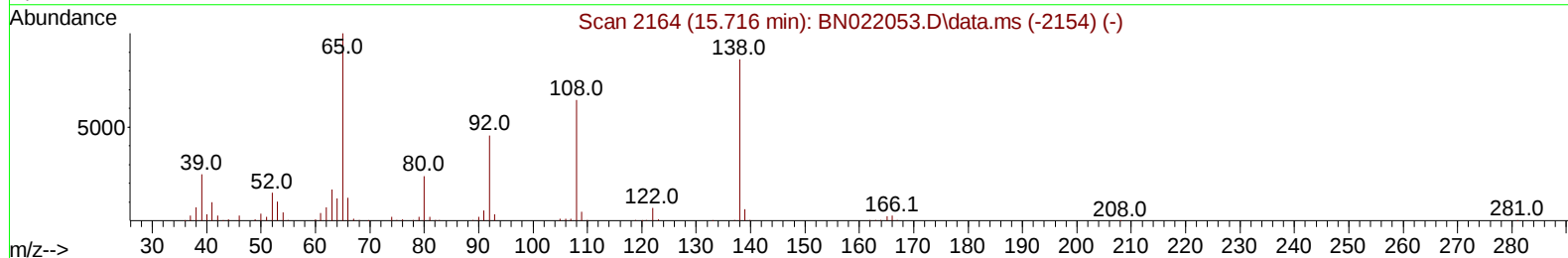
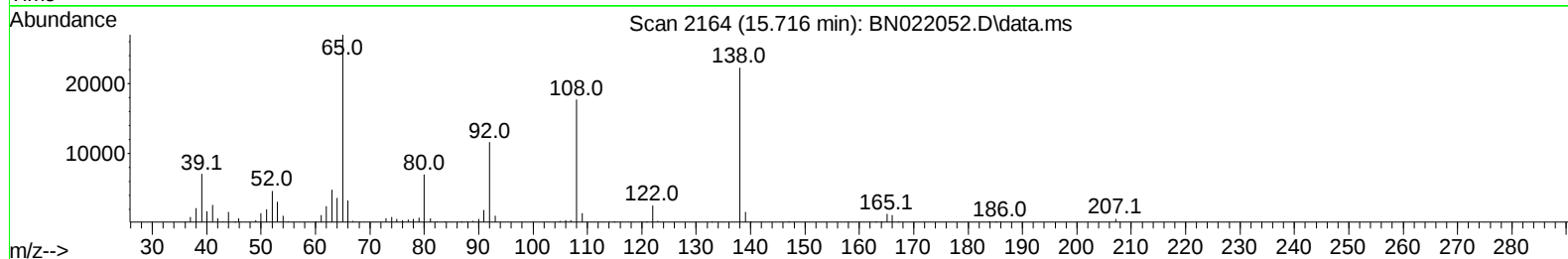
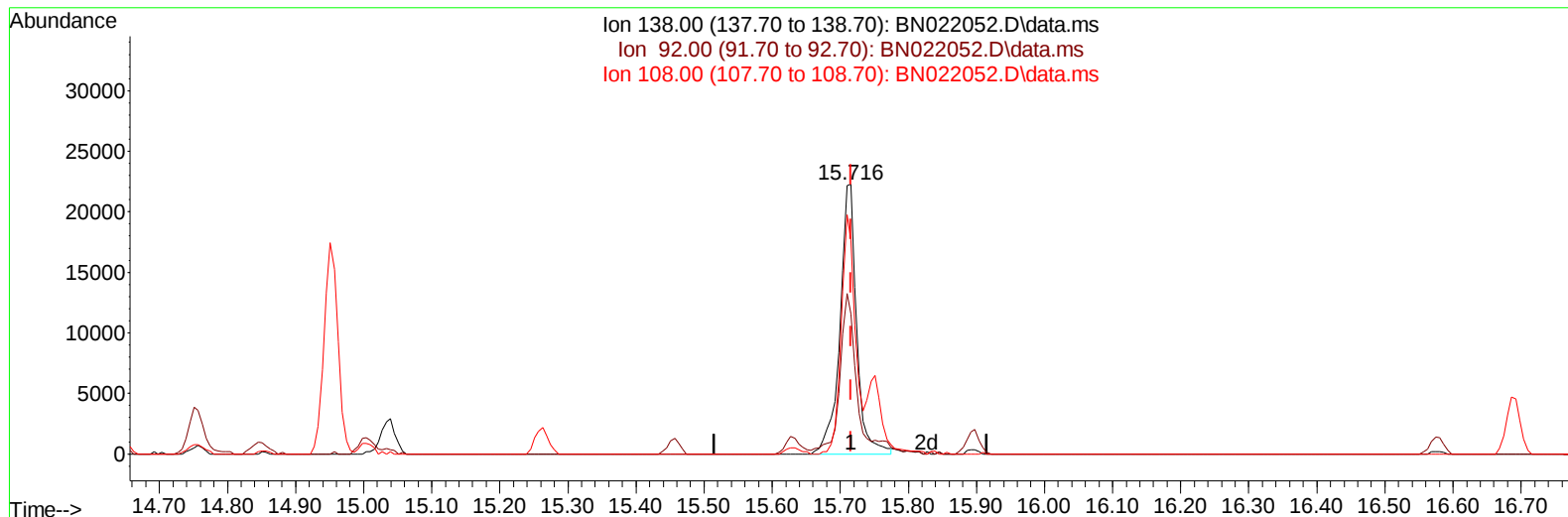
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
Data File : BN022052.D  
Acq On : 11 Oct 2022 15:01  
Operator : CG/JU  
Sample : SSTD01030  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 12 03:00:49 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

**(63) 4-Nitroaniline**

**15.716min (-0.000) 10.06 ng/ul**

**response 38021**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 53.50  | 52.02  |
| 108.00 | 75.10  | 79.40  |
| 0.00   | 0.00   | 0.00   |



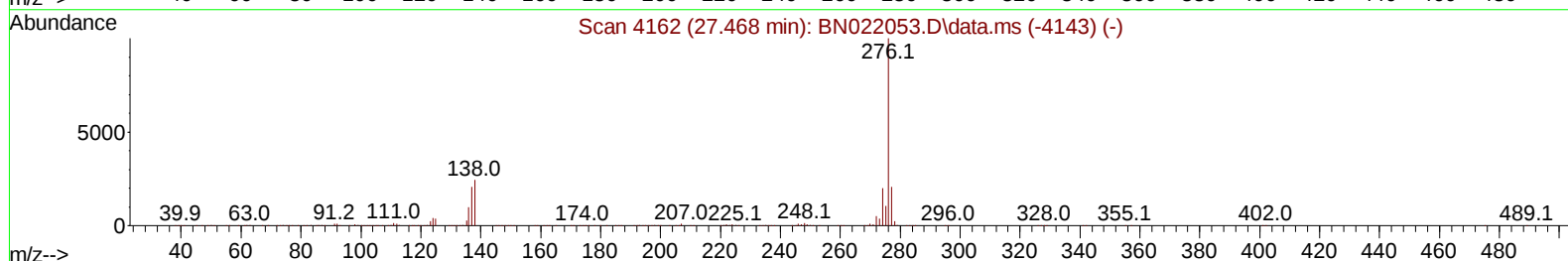
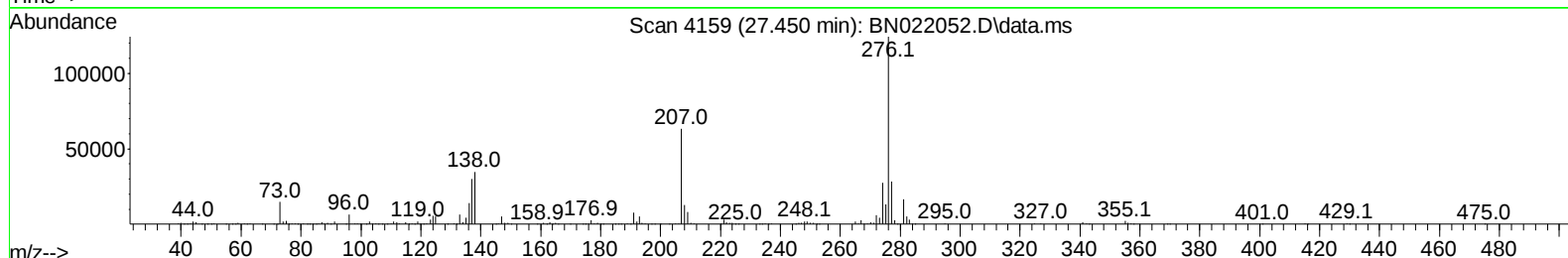
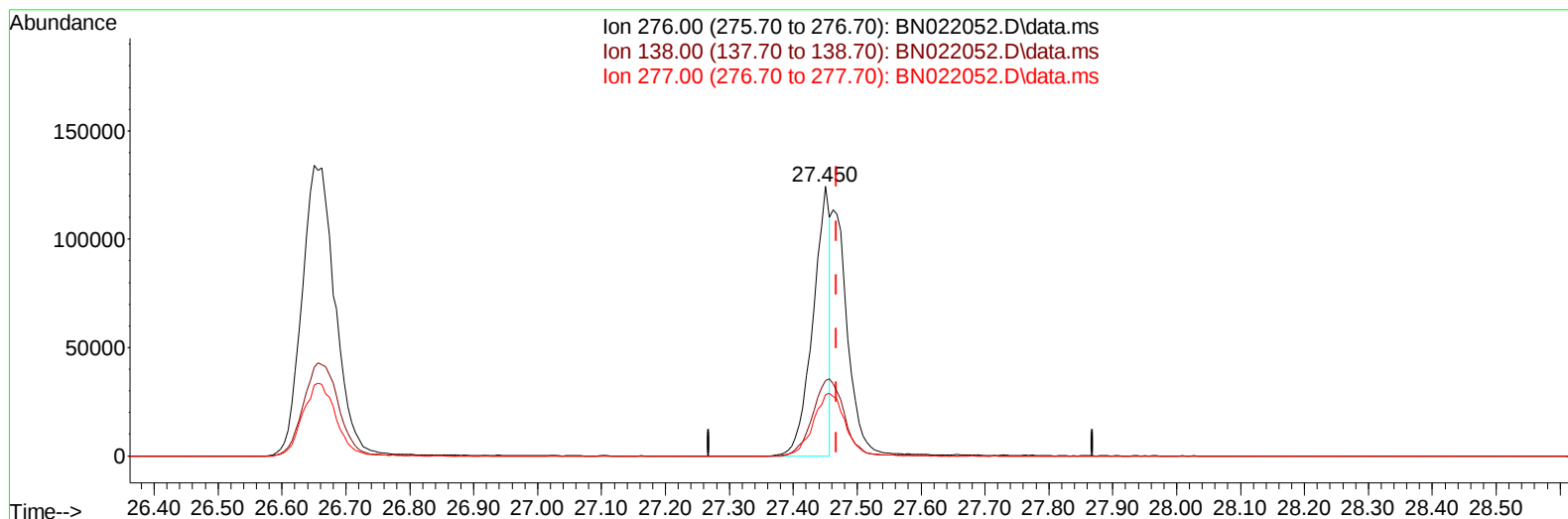
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SSTD01030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 03:00:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
 Supervised By :mohammad ahmed 10/14/2022



TIC: BN022052.D\data.ms

**(96) Benzo(g,h,i)perylene**

27.450min (-0.018) 5.38 ng/u1

response 226523

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 24.30  | 28.00  |
| 277.00 | 20.60  | 22.81  |
| 0.00   | 0.00   | 0.00   |

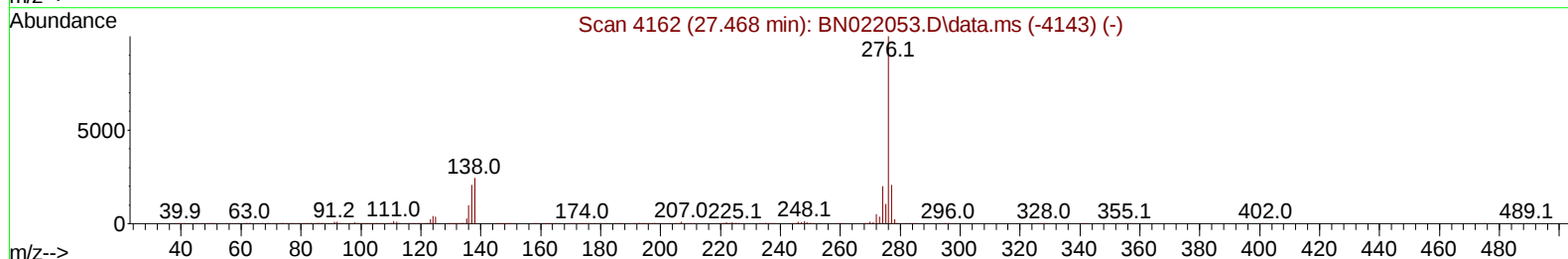
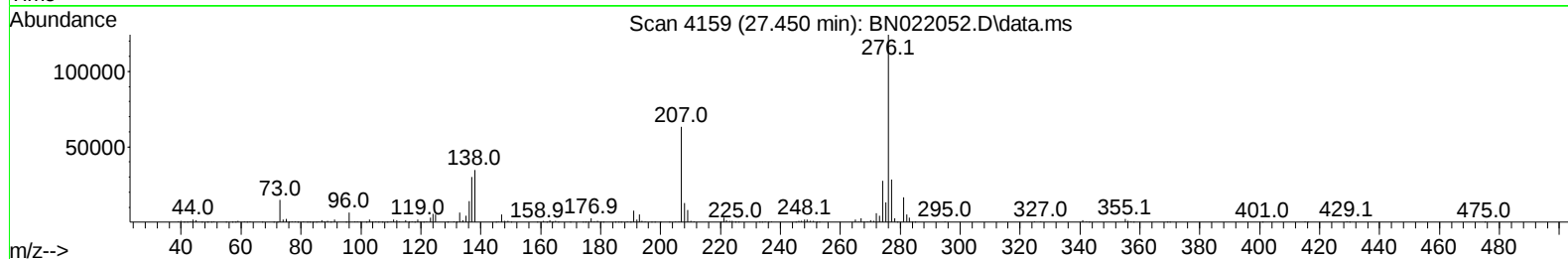
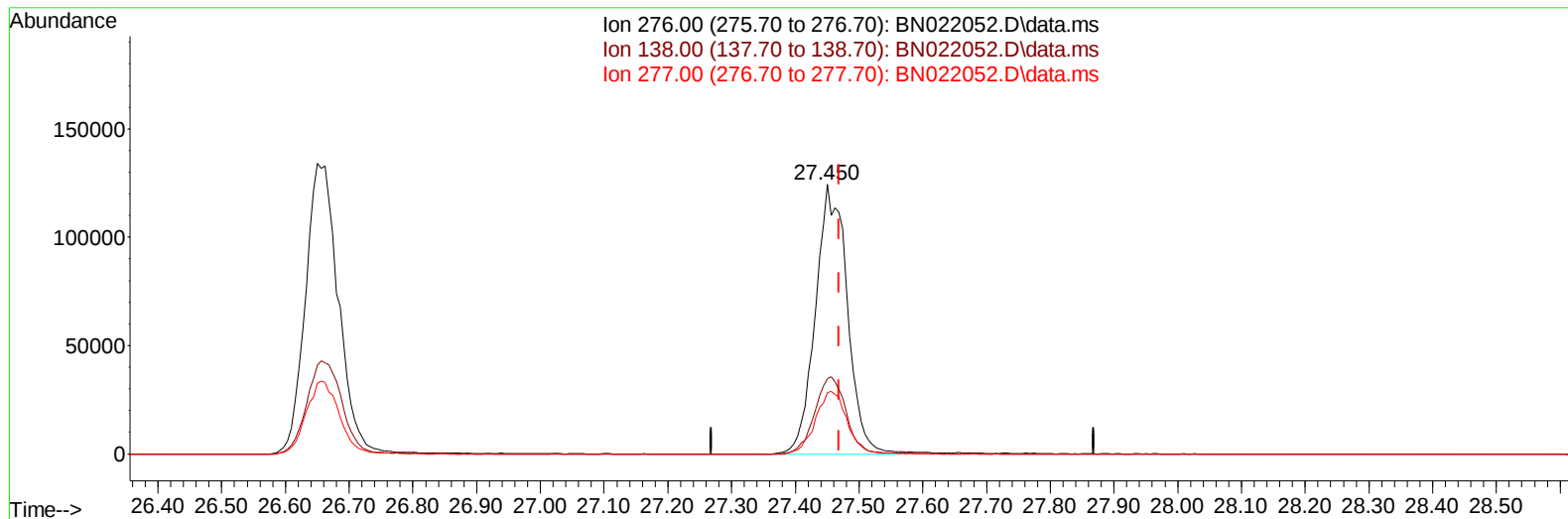
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
Data File : BN022052.D  
Acq On : 11 Oct 2022 15:01  
Operator : CG/JU  
Sample : SSTD01030  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022

Quant Time: Oct 12 03:06:53 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Oct 12 03:00:49 2022  
Response via : Initial Calibration



TIC: BN022052.D\data.ms

**(96) Benzo(g,h,i)perylene**

27.450min (-0.018) 10.14 ng/ul m

response 427245

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 24.30  | 28.00  |
| 277.00 | 20.60  | 22.81  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SST001030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.969  | 152  | 90445    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.798 | 136  | 419464   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.639 | 164  | 281917   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.392 | 188  | 614760   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.592 | 240  | 626810   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.062 | 264  | 675719   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.263  | 96   | 10123m   | 4.202  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.710  | 84   | 64878    | 9.882  | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.122  | 99   | 75928    | 9.256  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.293  | 67   | 54222    | 10.920 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.492  | 132  | 55369    | 9.277  | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.687  | 113  | 59146    | 9.189  | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.145  | 128  | 27273    | 9.543  | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.875  | 143  | 26948    | 10.557 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.416 | 165  | 54600    | 9.658  | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.939 | 131  | 95115    | 9.907  | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 200973   | 10.430 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.333 | 160  | 210435   | 9.368  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.833 | 143  | 36103    | 9.416  | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.627 | 176  | 170120   | 10.146 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 25715    | 10.458 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.492 | 188  | 274537   | 10.134 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 309620   | 10.387 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.903 | 264  | 320569   | 10.009 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        |        | Qvalue   |
| 2) 1,4-Dioxane                | 3.299  | 88   | 11026    | 4.444  | ng/uL  | 93       |
| 5) Pyridine                   | 3.728  | 79   | 67978    | 10.290 | ng/ul  | 99       |
| 6) Benzaldehyde               | 7.098  | 77   | 42000m   | 11.784 | ng/ul  |          |
| 8) Phenol                     | 7.151  | 94   | 83905    | 9.898  | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.387  | 93   | 71615    | 10.596 | ng/ul  | 97       |
| 12) 2-Chlorophenol            | 7.522  | 128  | 63150    | 9.863  | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.416  | 108  | 60516    | 9.396  | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.504  | 45   | 110242   | 12.931 | ng/ul  | 99       |
| 16) Acetophenone              | 8.804  | 105  | 110704   | 10.875 | ng/ul  | 98       |
| 17) N-Nitroso-di-n-propyla... | 8.787  | 70   | 55215    | 10.893 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.745  | 108  | 68526    | 9.686  | ng/ul  | 98       |
| 19) Hexachloroethane          | 9.057  | 117  | 26194    | 10.264 | ng/ul  | 92       |
| 22) Nitrobenzene              | 9.187  | 77   | 80830    | 11.477 | ng/ul  | 98       |
| 23) Isophorone                | 9.716  | 82   | 151893   | 10.474 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.904  | 139  | 32592    | 10.900 | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.963  | 107  | 76152    | 10.285 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.204 | 93   | 102851   | 11.325 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.439 | 162  | 59843    | 10.103 | ng/ul  | 97       |
| 30) Naphthalene               | 10.845 | 128  | 232842   | 10.661 | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.963 | 127  | 103301   | 10.382 | ng/ul  | 96       |
| 33) Hexachlorobutadiene       | 11.128 | 225  | 34191    | 10.020 | ng/ul# | 92       |
| 34) Caprolactam               | 11.751 | 113  | 19597m   | 8.650  | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.086 | 107  | 71162    | 10.660 | ng/ul  | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101122\  
 Data File : BN022052.D  
 Acq On : 11 Oct 2022 15:01  
 Operator : CG/JU  
 Sample : SST001030  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010239

Manual IntegrationsAPPROVED

Quant Time: Oct 12 03:06:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 03:00:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022  
 Supervised By :mohammad ahmed 10/14/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.463 | 142  | 153528   | 10.391 | ng/ul  | 96       |
| 37) 1-Methylnaphthalene       | 12.680 | 142  | 158453   | 10.671 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachloroben... | 12.828 | 216  | 72196    | 9.987  | ng/ul  | 98       |
| 40) Hexachlorocyclopentadiene | 12.804 | 237  | 33433    | 7.768  | ng/ul  | 89       |
| 41) 2,4,6-Trichlorophenol     | 13.075 | 196  | 45550    | 10.155 | ng/ul  | 98       |
| 42) 2,4,5-Trichlorophenol     | 13.139 | 196  | 50333    | 10.227 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.475 | 154  | 205132   | 9.957  | ng/ul  | 98       |
| 44) 2-Chloronaphthalene       | 13.516 | 162  | 163275   | 10.352 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.727 | 65   | 43599    | 11.021 | ng/ul  | 97       |
| 47) Dimethylphthalate         | 14.092 | 163  | 217574   | 11.039 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.222 | 165  | 32628    | 9.649  | ng/ul  | 95       |
| 50) Acenaphthylene            | 14.363 | 152  | 251832   | 9.711  | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.551 | 138  | 37612    | 9.985  | ng/ul  | 97       |
| 52) Acenaphthene              | 14.704 | 153  | 183051   | 10.924 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol         | 14.757 | 184  | 18245    | 11.022 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.851 | 109  | 31955    | 10.396 | ng/ul  | 99       |
| 56) Dibenzofuran              | 15.033 | 168  | 254008   | 10.551 | ng/ul  | 100      |
| 57) 2,4-Dinitrotoluene        | 15.004 | 165  | 52652    | 10.721 | ng/ul  | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.263 | 232  | 43424    | 11.495 | ng/ul  | 99       |
| 59) Diethylphthalate          | 15.457 | 149  | 220883   | 11.113 | ng/ul  | 97       |
| 61) Fluorene                  | 15.686 | 166  | 211962   | 11.163 | ng/ul  | 91       |
| 62) 4-Chlorophenyl-phenyle... | 15.680 | 204  | 96212    | 10.992 | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 15.716 | 138  | 38348m   | 10.144 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.763 | 198  | 30282    | 11.625 | ng/ul# | 95       |
| 67) N-Nitrosodiphenylamine    | 15.898 | 169  | 181759   | 10.549 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.580 | 248  | 54891    | 10.040 | ng/ul  | 99       |
| 69) Hexachlorobenzene         | 16.692 | 284  | 68000    | 10.448 | ng/ul  | 99       |
| 70) Atrazine                  | 16.851 | 200  | 58535    | 9.601  | ng/ul  | 96       |
| 71) Pentachlorophenol         | 17.039 | 266  | 34619    | 9.554  | ng/ul# | 87       |
| 72) Phenanthrene              | 17.433 | 178  | 344219   | 10.600 | ng/ul  | 98       |
| 74) Anthracene                | 17.527 | 178  | 346494   | 10.716 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.439 | 216  | 71997    | 9.202  | ng/uL  | 95       |
| 76) Pentachlorobenzene        | 14.957 | 250  | 81314    | 10.909 | ng/uL  | 98       |
| 77) Carbazole                 | 17.798 | 167  | 318444   | 10.422 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.363 | 149  | 361434   | 10.492 | ng/ul  | 100      |
| 80) Fluoranthene              | 19.457 | 202  | 390669   | 10.865 | ng/ul  | 100      |
| 82) Pyrene                    | 19.821 | 202  | 428580   | 11.437 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.721 | 149  | 156480   | 11.032 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.509 | 252  | 128996   | 10.793 | ng/ul  | 92       |
| 85) Benzo(a)anthracene        | 21.574 | 228  | 409315   | 10.845 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.498 | 149  | 233664   | 10.760 | ng/ul  | 98       |
| 87) Chrysene                  | 21.627 | 228  | 418991   | 11.467 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.445 | 149  | 403864   | 10.203 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.303 | 252  | 433021   | 10.533 | ng/ul  | 94       |
| 91) Benzo(k)fluoranthene      | 23.356 | 252  | 427761   | 11.632 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 23.950 | 252  | 360086   | 9.155  | ng/ul# | 91       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.650 | 276  | 481685   | 9.975  | ng/ul  | 98       |
| 95) Dibenzo(a,h)anthracene    | 26.680 | 278  | 431462   | 10.588 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 27.450 | 276  | 427245m  | 10.139 | ng/ul  |          |

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

**Instrument :**

BNA\_N

**ClientSampleId :**

SSTD010239

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 10/12/2022  
Supervised By :mohammad ahmed 10/14/2022

