

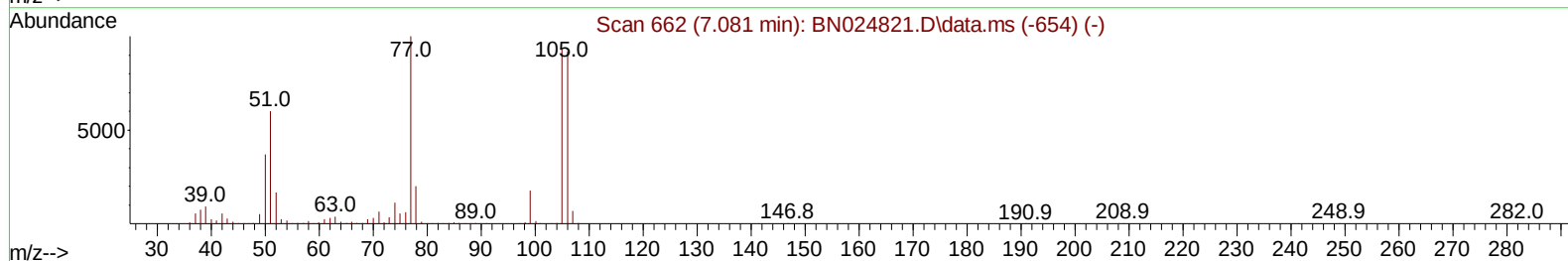
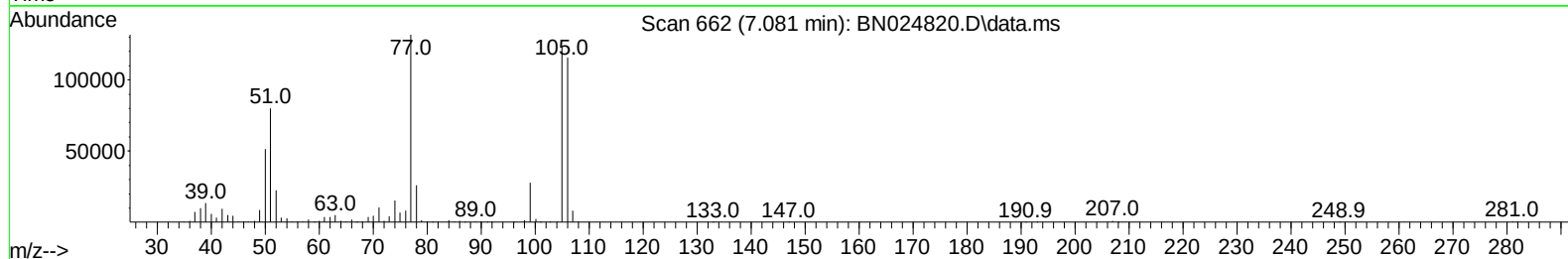
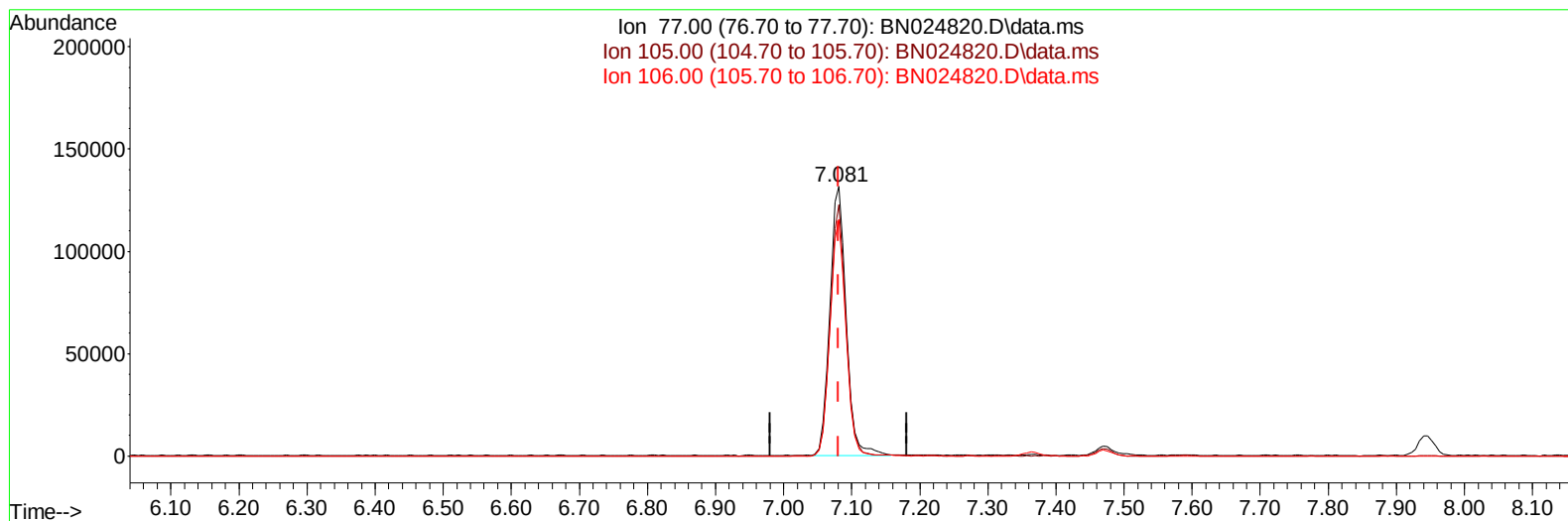
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

## Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/11/2023  
Supervised By :Jagrut Upadhyay 04/11/2023

Quant Time: Apr 11 02:39:34 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration



TIC: BN024820.D\data.ms

**(6) Benzaldehyde****7.081min (+ 0.000) 11.80 ng/ul****response 219682**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 92.90  | 93.36  |
| 106.00 | 91.60  | 87.76  |
| 0.00   | 0.00   | 0.00   |

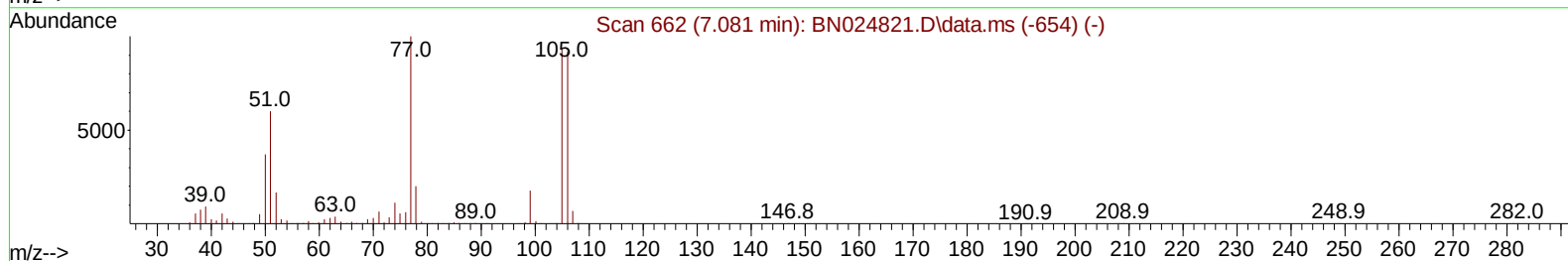
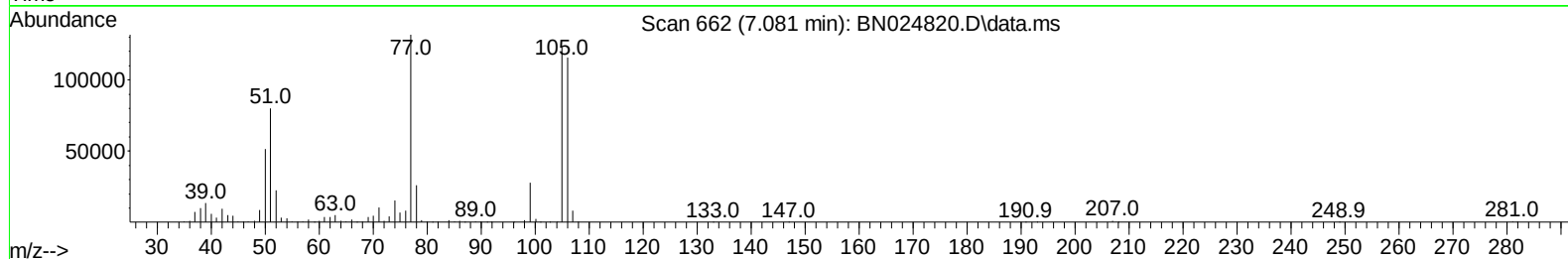
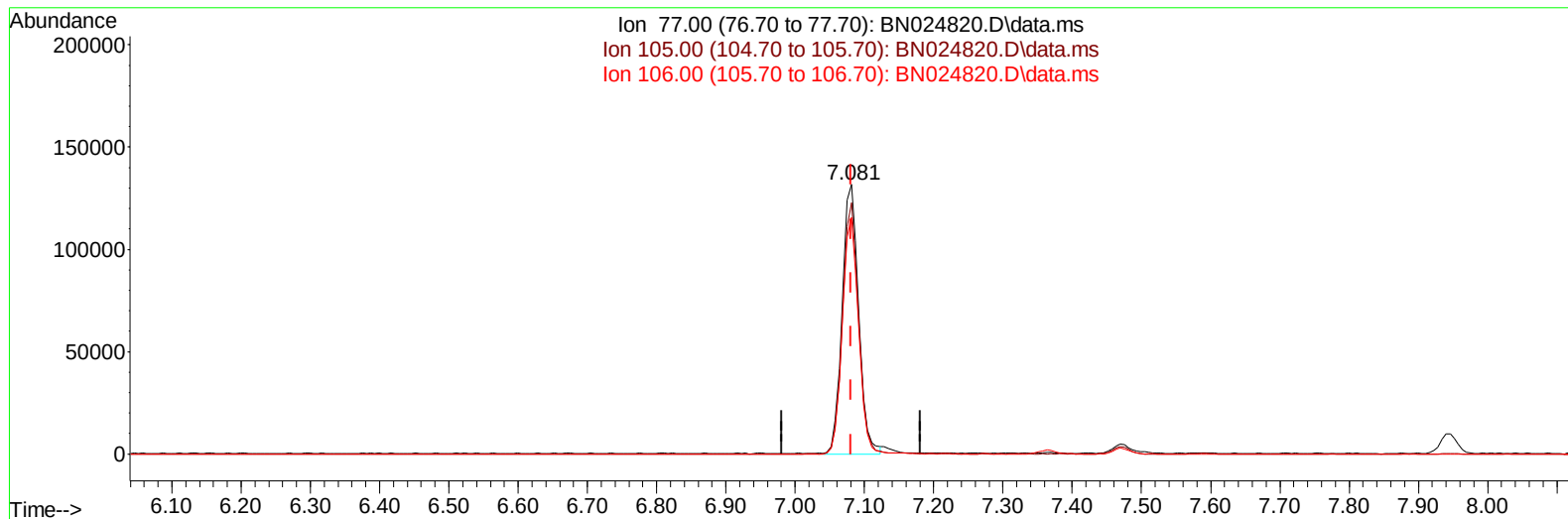
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
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ALS Vial : 3 Sample Multiplier: 1

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Response via : Initial Calibration



TIC: BN024820.D\data.ms

## (6) Benzaldehyde

7.081min (+ 0.000) 11.71 ng/ul m

response 217998

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 92.90  | 93.36  |
| 106.00 | 91.60  | 87.76  |
| 0.00   | 0.00   | 0.00   |

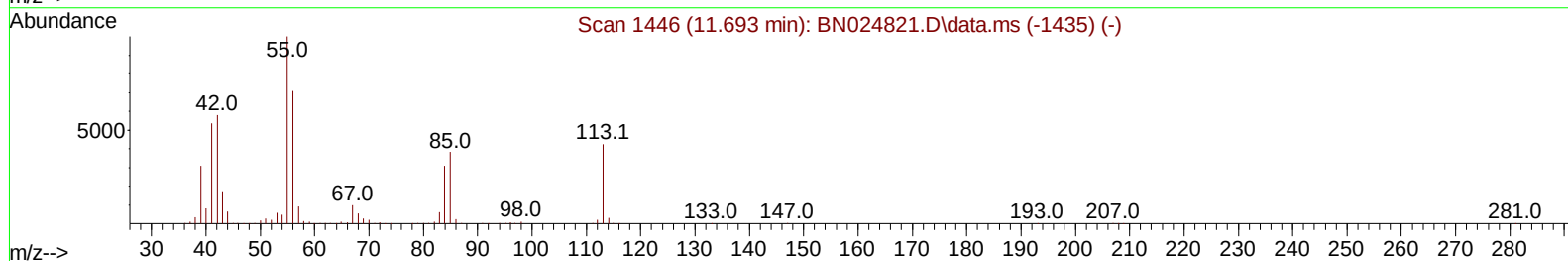
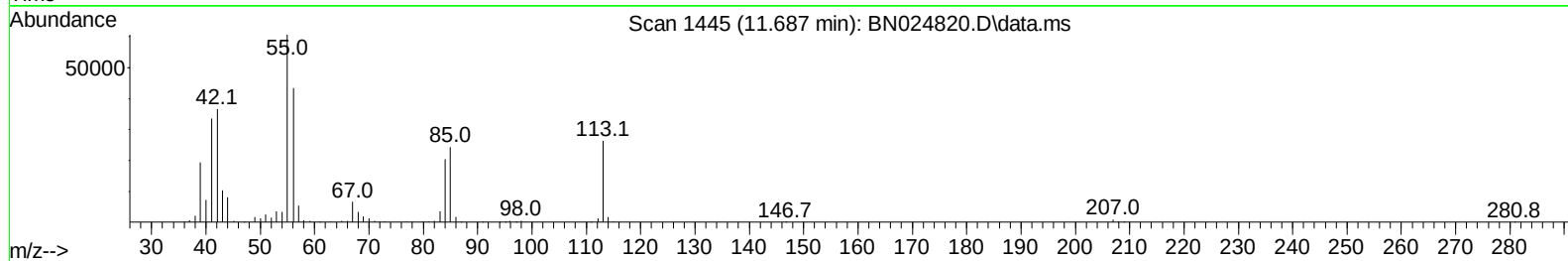
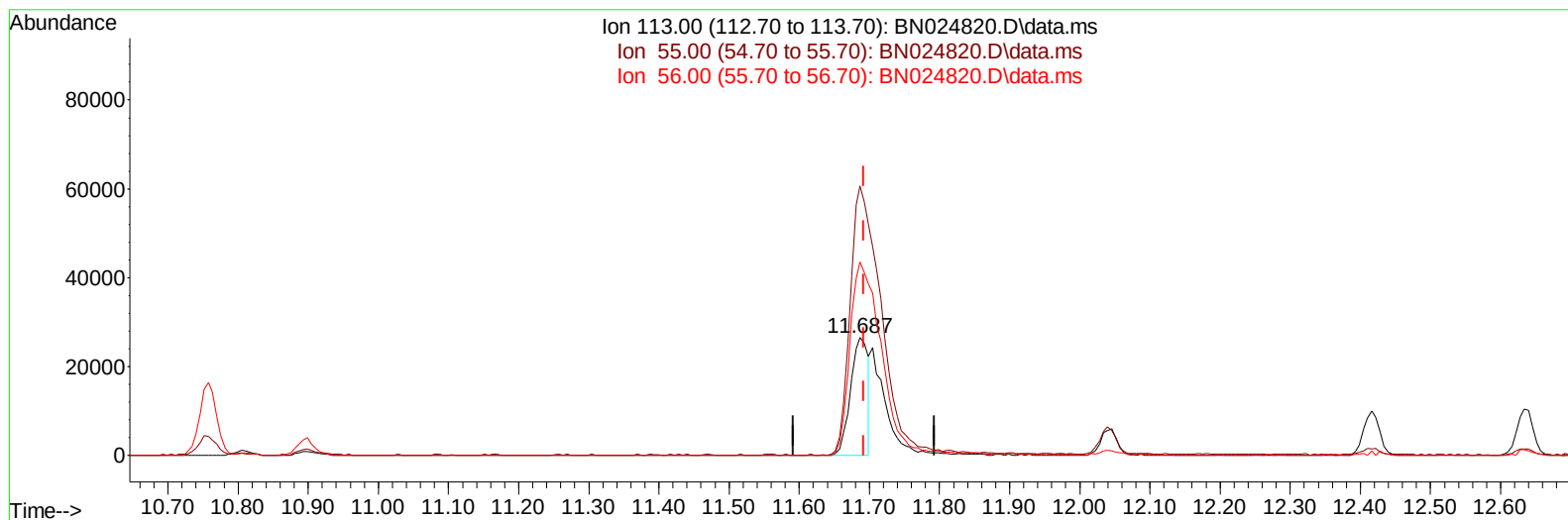
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
 Data File : BN024820.D  
 Acq On : 10 Apr 2023 11:55  
 Operator : CG/JU  
 Sample : SST01090  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST010212

Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 10 15:00:02 2023  
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
 Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BN024820.D\data.ms

**(34) Caprolactam**

**11.687min (-0.006) 4.82 ng/ul**

**response 46314**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 235.40 | 229.10 |
| 56.00  | 166.90 | 164.34 |
| 0.00   | 0.00   | 0.00   |

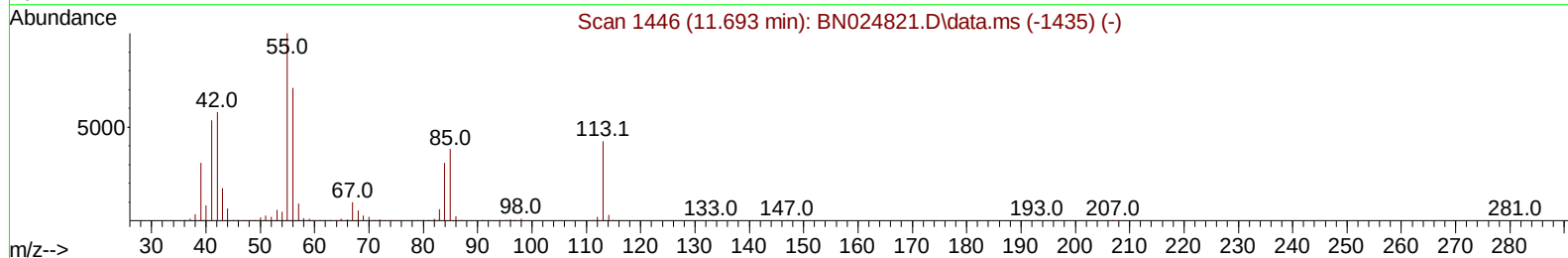
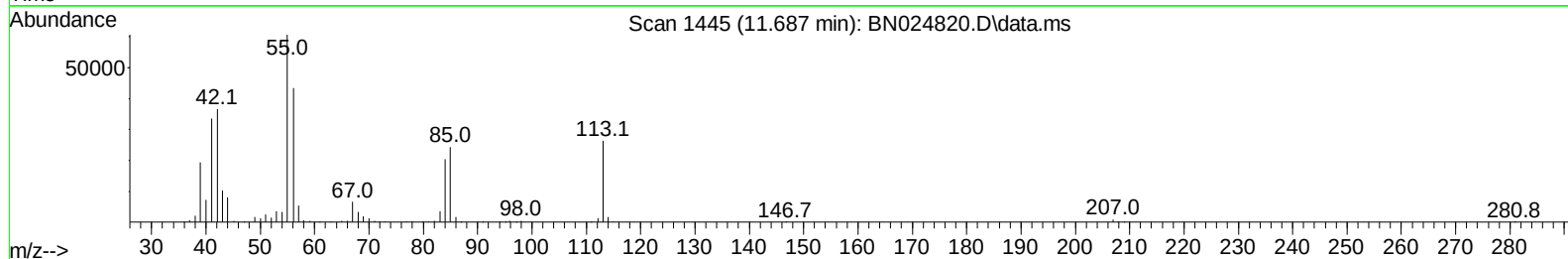
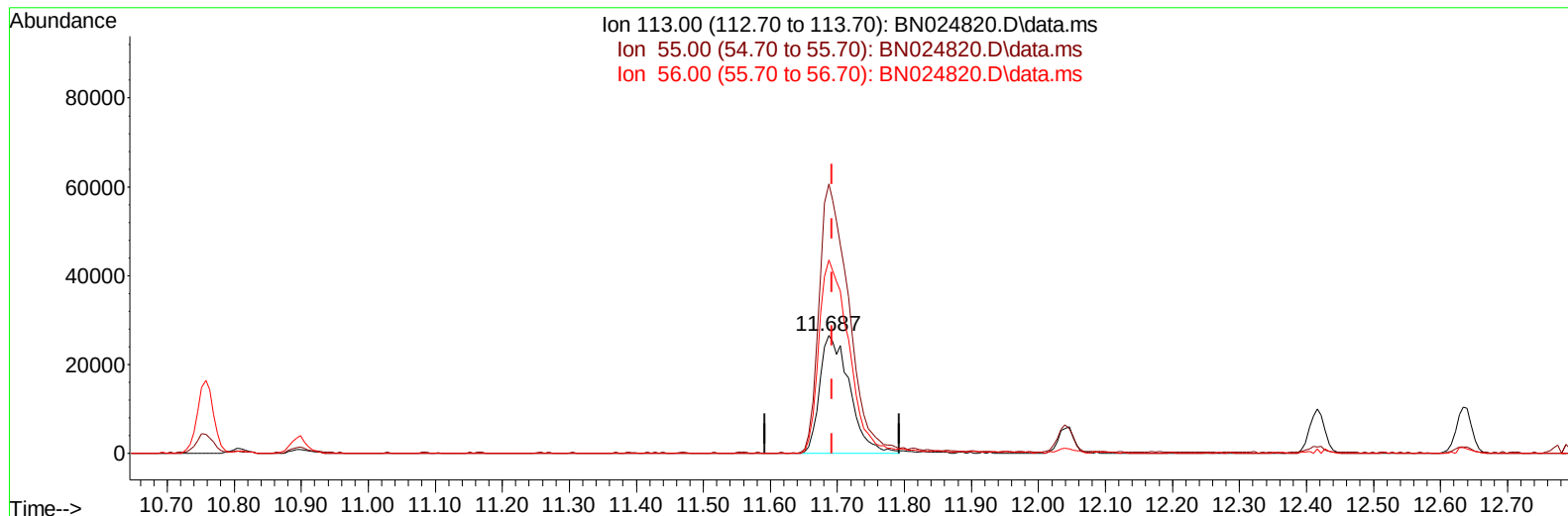
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
 Data File : BN024820.D  
 Acq On : 10 Apr 2023 11:55  
 Operator : CG/JU  
 Sample : SST01090  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
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 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
 Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BN024820.D\data.ms

**(34) Caprolactam**

11.687min (-0.006) 8.50 ng/ul m

response 81746

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 235.40 | 229.10 |
| 56.00  | 166.90 | 164.34 |
| 0.00   | 0.00   | 0.00   |

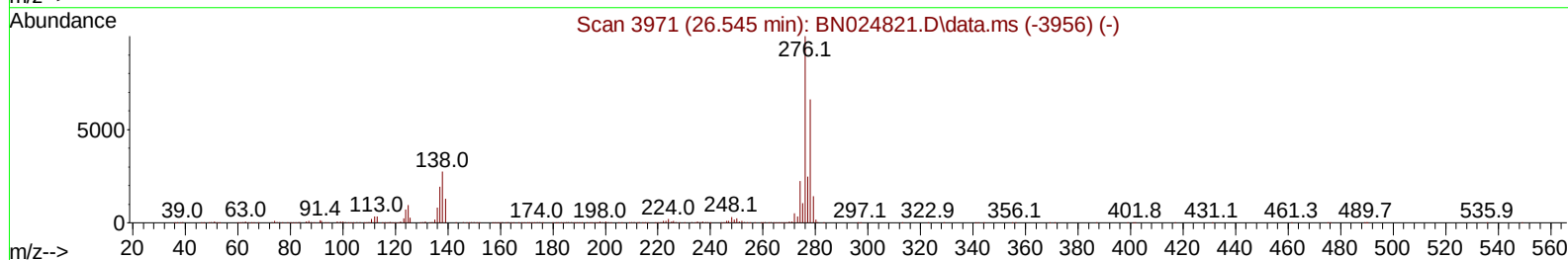
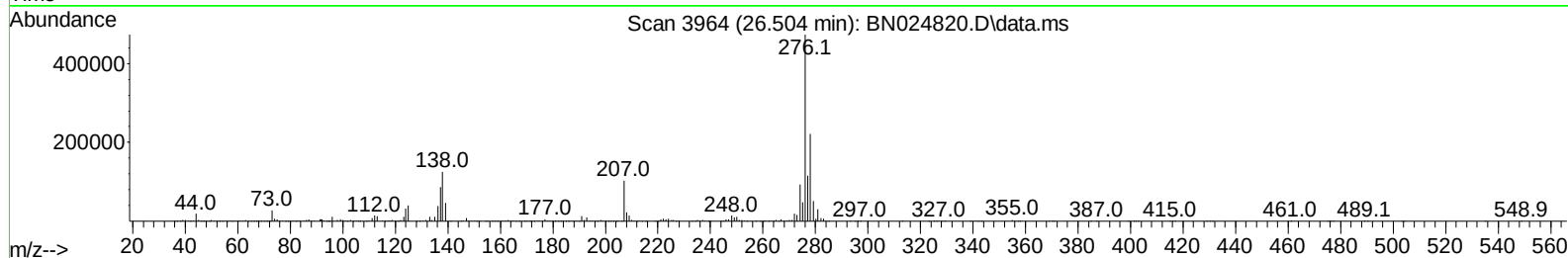
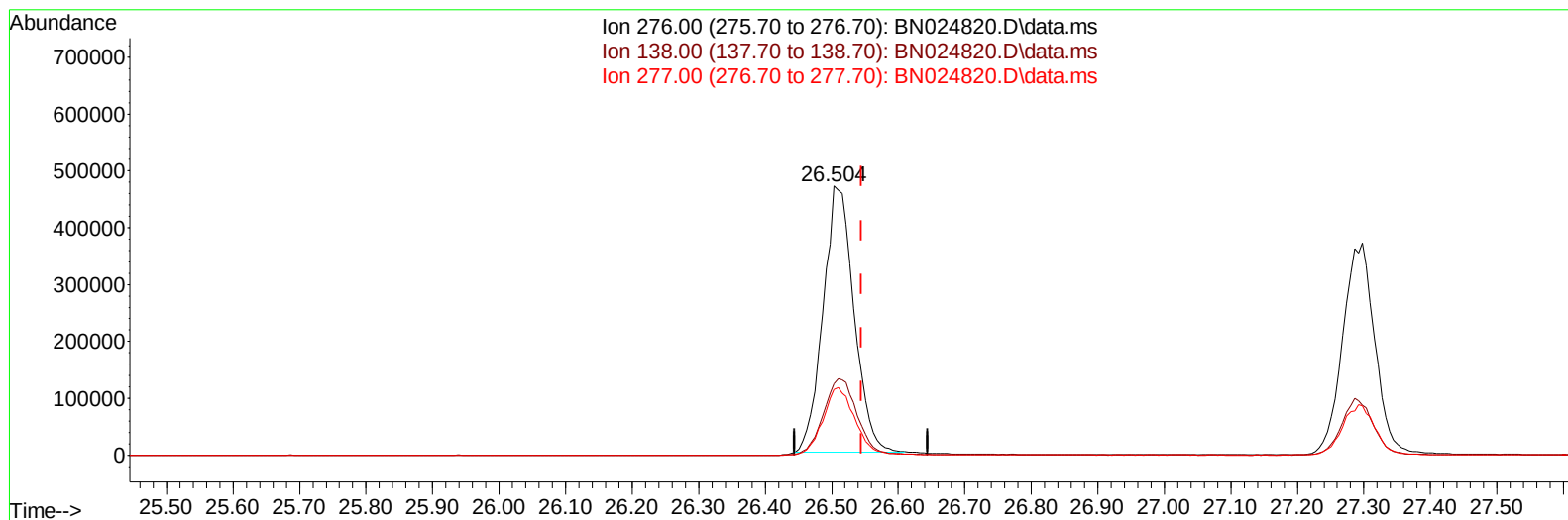
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

## Manual IntegrationsAPPROVED

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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration



TIC: BN024820.D\data.ms

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

26.504min (-0.041) 9.57 ng/uL

response 1509721

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 27.40  | 26.43  |
| 277.00 | 24.70  | 24.43  |
| 0.00   | 0.00   | 0.00   |

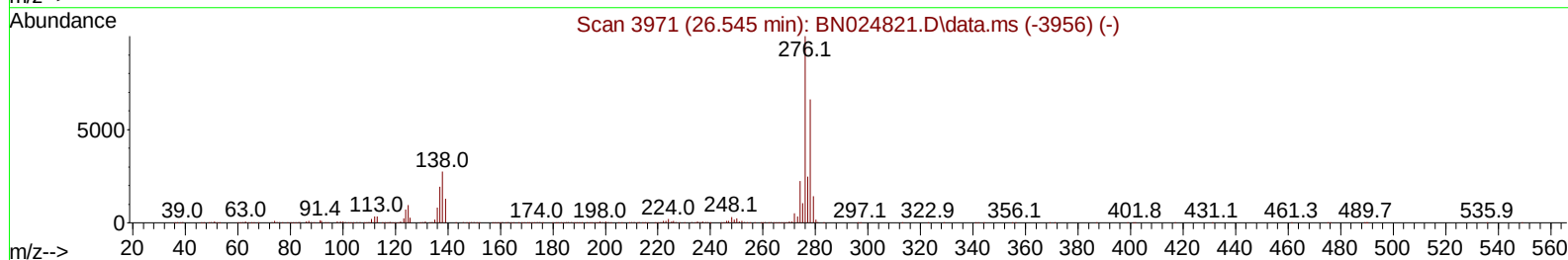
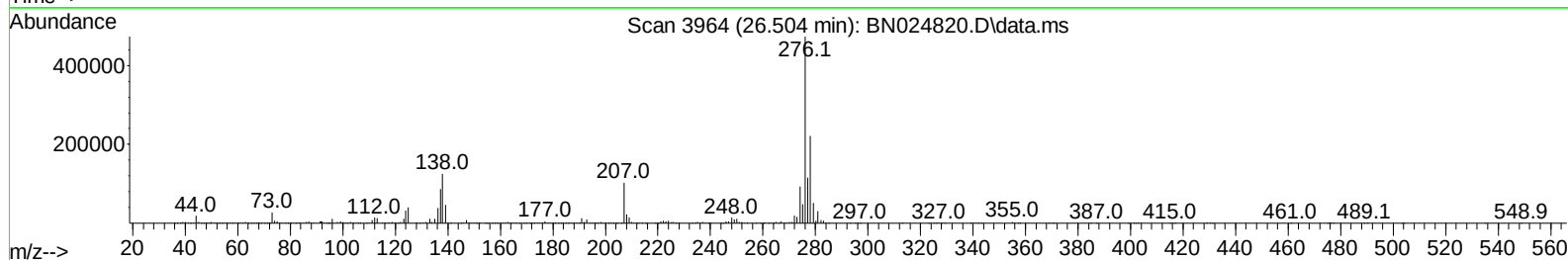
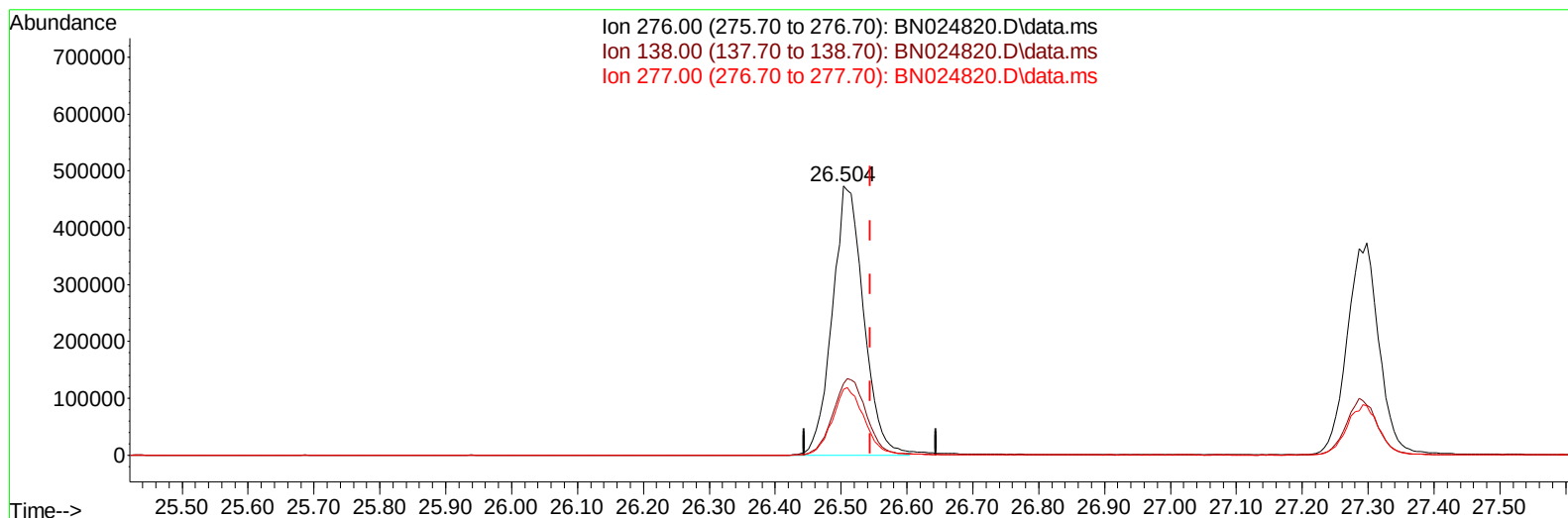
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BN024820.D\data.ms

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

26.504min (-0.041) 9.90 ng/ul m

response 1562054

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 27.40  | 26.43  |
| 277.00 | 24.70  | 24.43  |
| 0.00   | 0.00   | 0.00   |

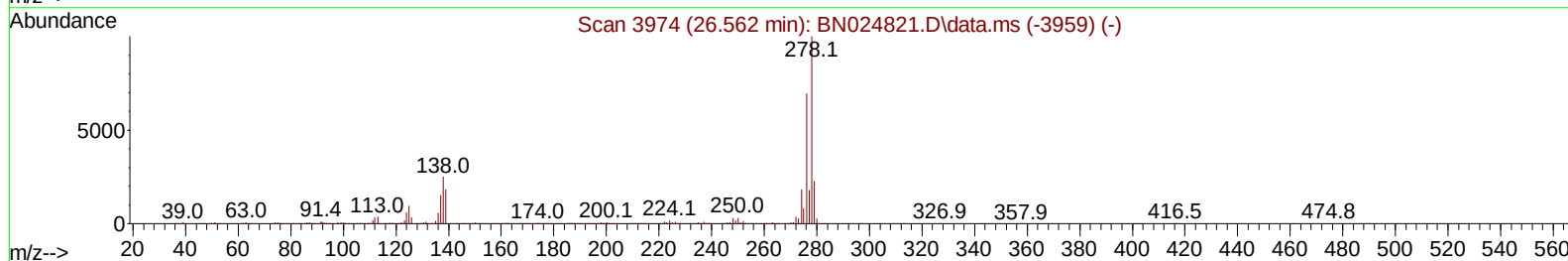
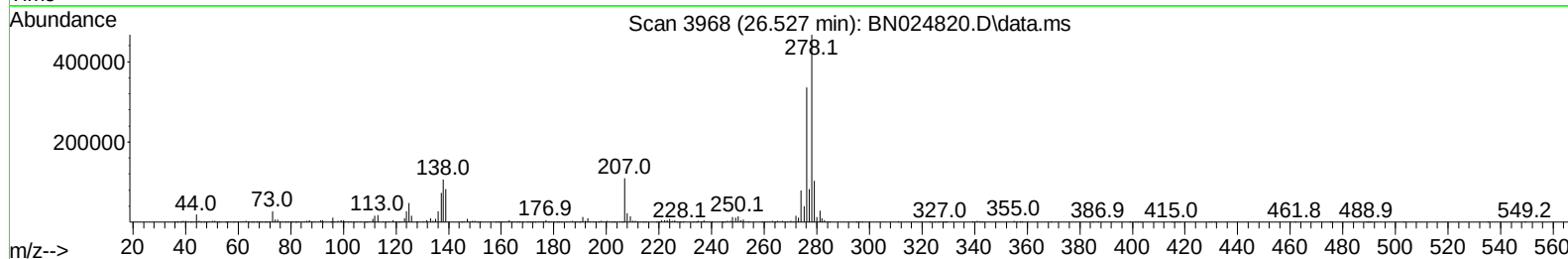
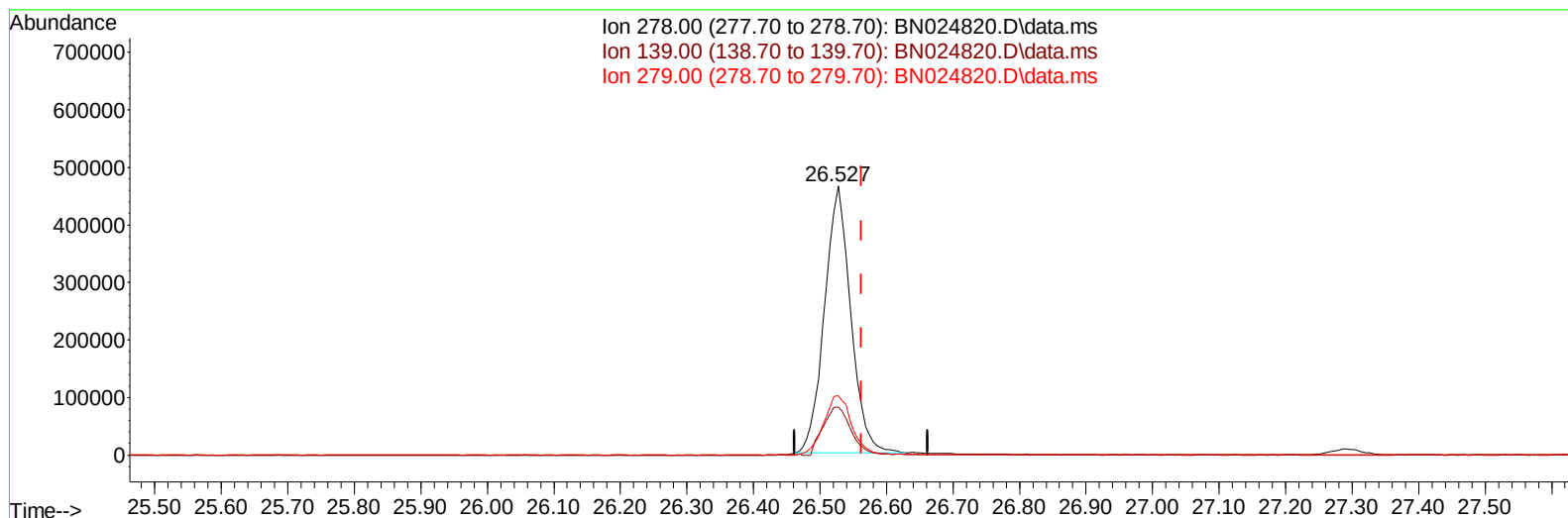
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

## Manual IntegrationsAPPROVED

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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
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TIC: BN024820.D\data.ms

**(95) Dibenzo(a,h)anthracene**

26.527min (-0.035) 9.73 ng/uI

response 1270054

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 18.20  | 17.83  |
| 279.00 | 22.80  | 22.16  |
| 0.00   | 0.00   | 0.00   |

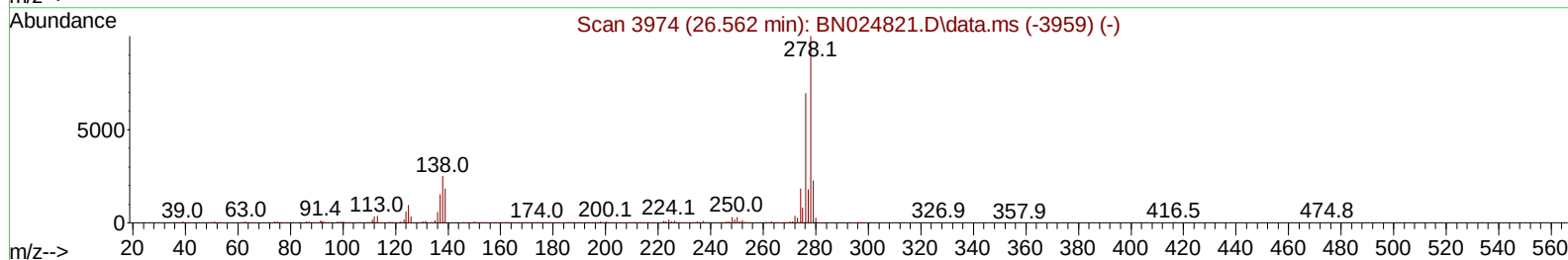
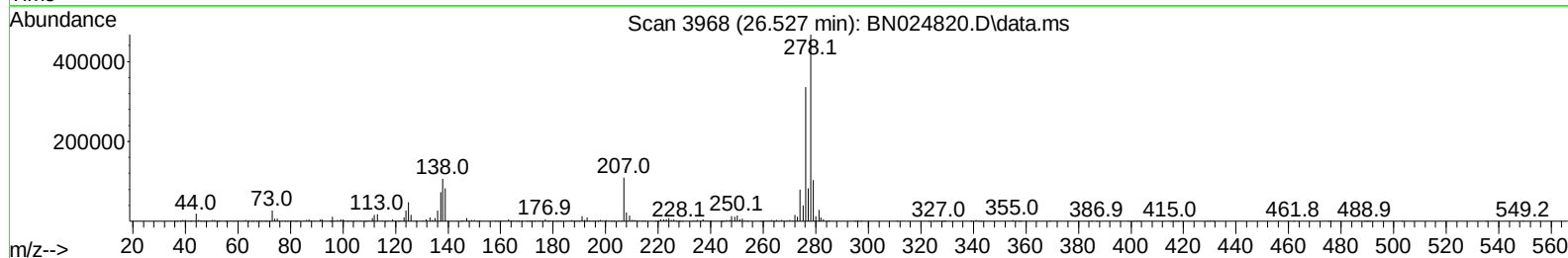
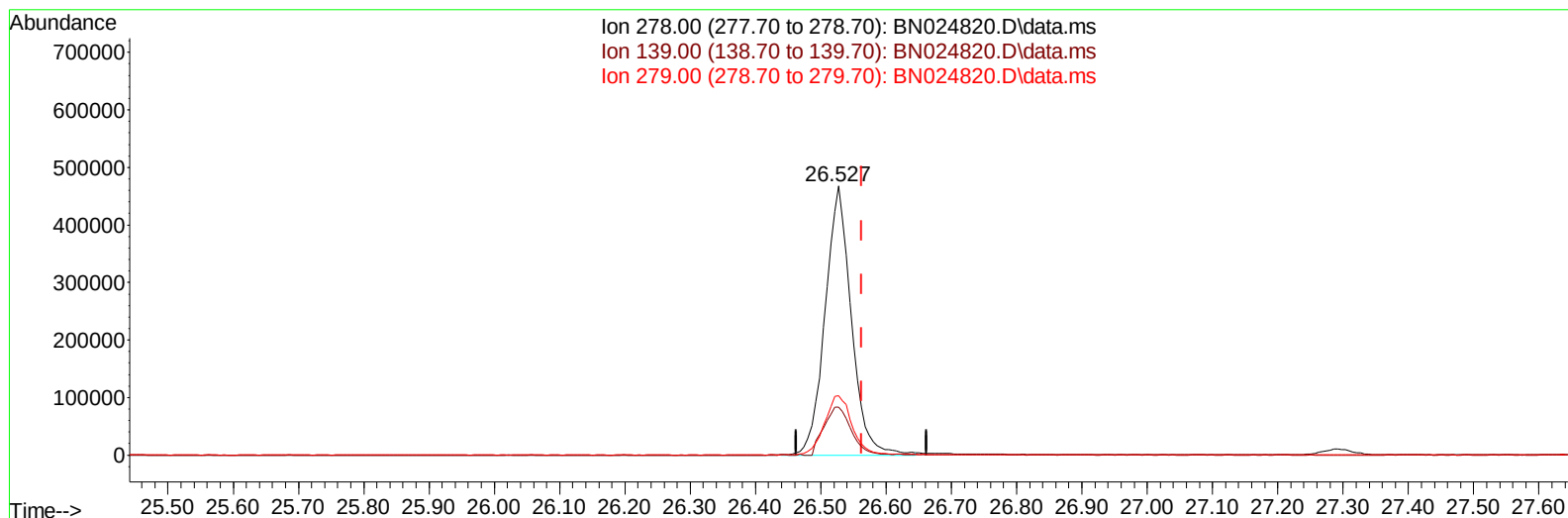
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BN024820.D\data.ms

(95) Dibenzo(a,h)anthracene

26.527min (-0.035) 10.06 ng/ul m

response 1314144

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 18.20  | 17.83  |
| 279.00 | 22.80  | 22.16  |
| 0.00   | 0.00   | 0.00   |



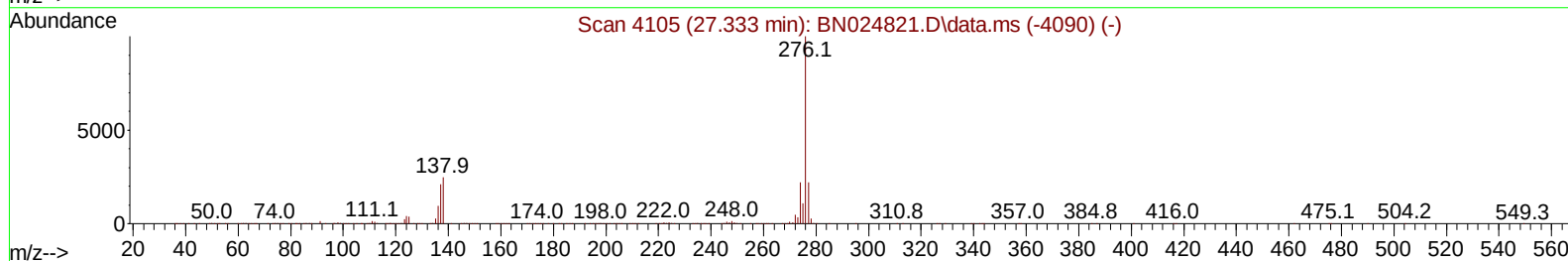
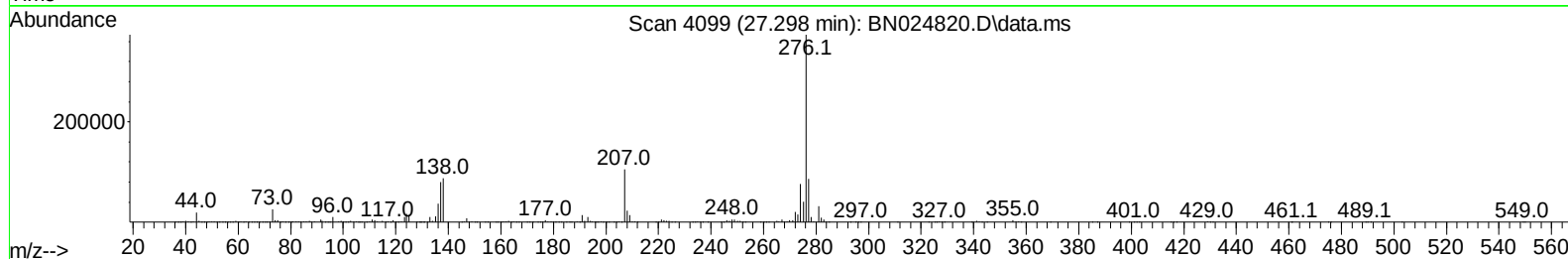
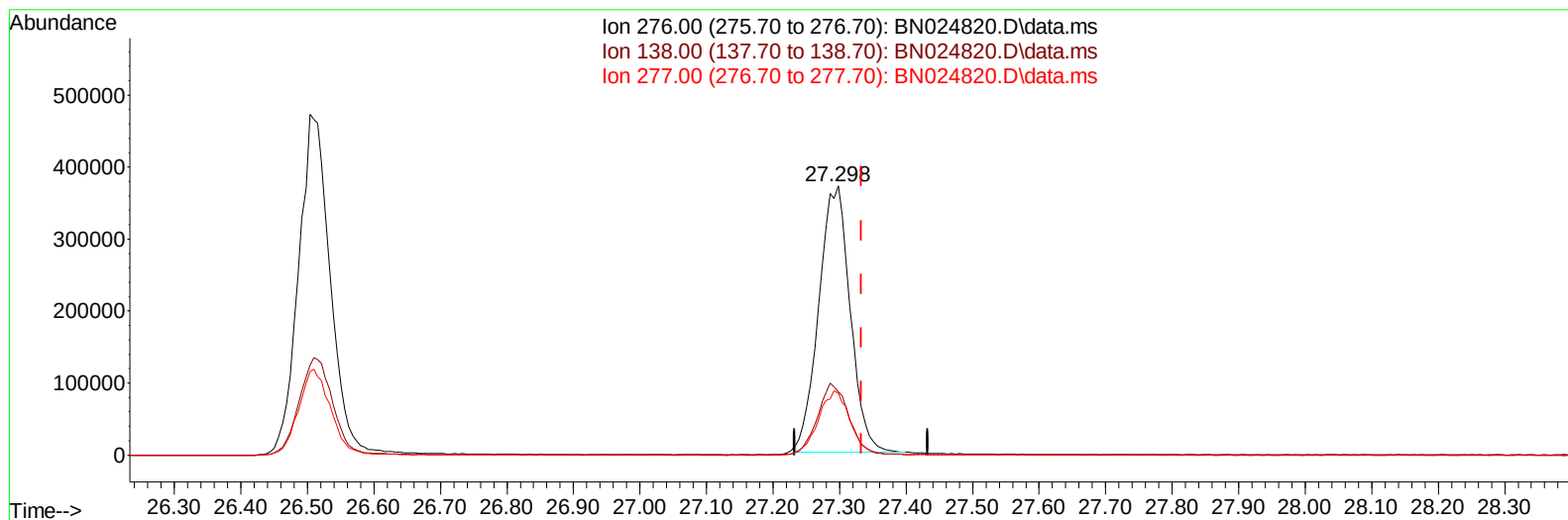
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration



TIC: BN024820.D\data.ms

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

27.298min (-0.035) 9.50 ng/u1

response 1210825

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 24.80  | 23.55  |
| 277.00 | 22.00  | 23.08  |
| 0.00   | 0.00   | 0.00   |

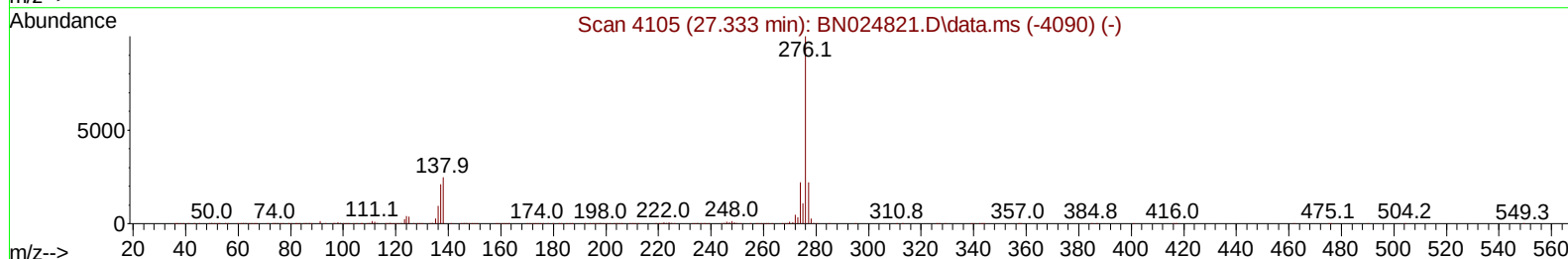
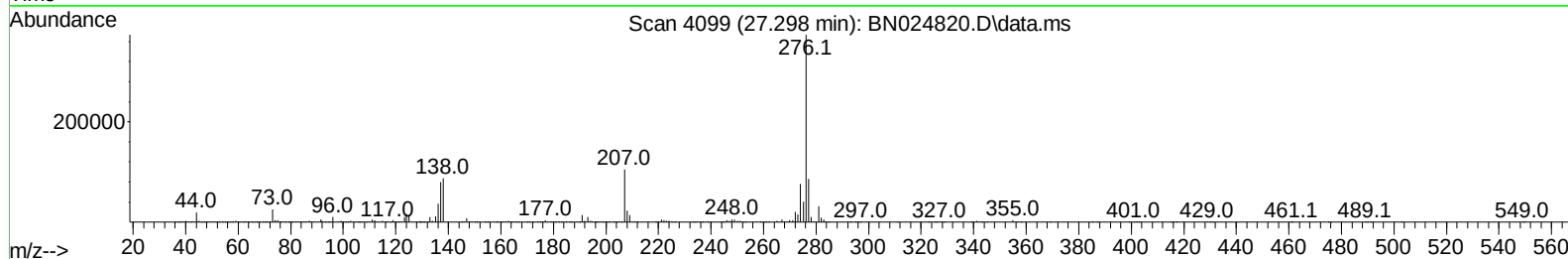
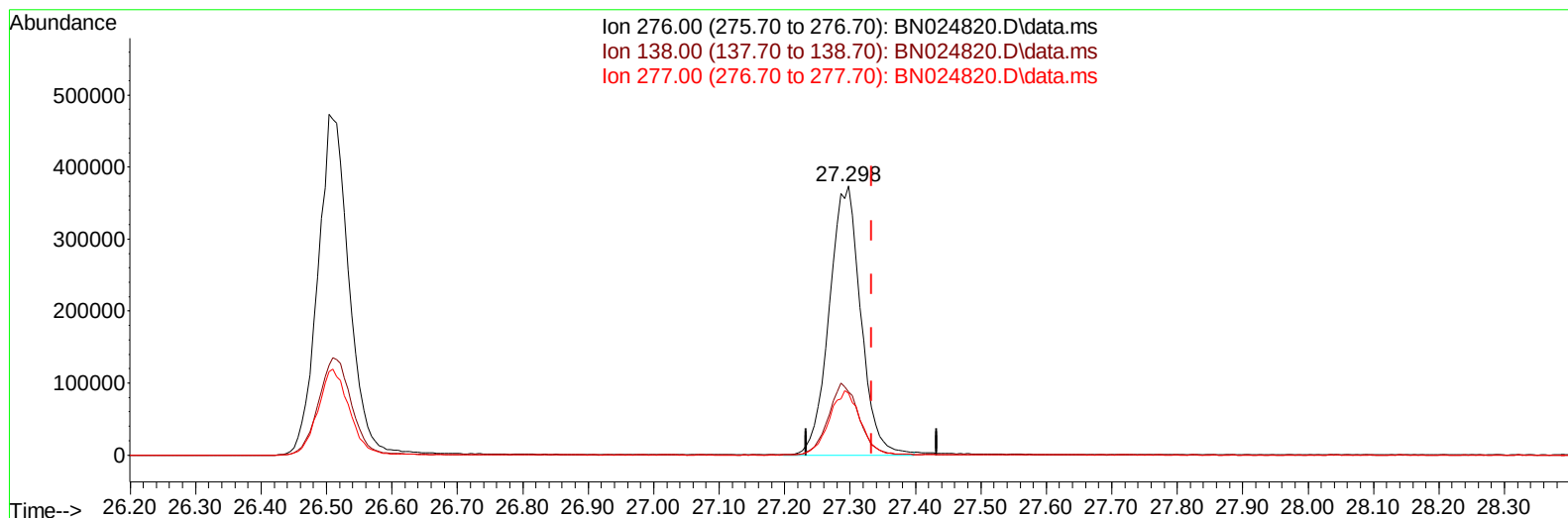
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
Data File : BN024820.D  
Acq On : 10 Apr 2023 11:55  
Operator : CG/JU  
Sample : SSTD01090  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010212

## Manual IntegrationsAPPROVED

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Quant Time: Apr 11 02:39:34 2023  
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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 10 15:00:02 2023  
Response via : Initial Calibration



TIC: BN024820.D\data.ms

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

27.298min (-0.035) 9.89 ng/ul m

response 1260189

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 24.80  | 23.55  |
| 277.00 | 22.00  | 23.08  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
 Data File : BN024820.D  
 Acq On : 10 Apr 2023 11:55  
 Operator : CG/JU  
 Sample : SST001090  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 390253   | 20.000 | ng/uL  | 0.00     |
| 20) Naphthalene-d8            | 10.757 | 136  | 1745639  | 20.000 | ng/uL  | 0.00     |
| 38) Acenaphthene-d10          | 14.593 | 164  | 1088625  | 20.000 | ng/uL  | 0.00     |
| 64) Phenanthrene-d10          | 17.334 | 188  | 2277425  | 20.000 | ng/uL  | 0.00     |
| 79) Chrysene-d12              | 21.527 | 240  | 2124389  | 20.000 | ng/uL  | 0.00     |
| 88) Perylene-d12              | 23.974 | 264  | 2310070  | 20.000 | ng/uL  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.311  | 96   | 39867    | 3.912  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.740  | 84   | 290114   | 9.449  | ng/uL  | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 352214   | 9.257  | ng/uL  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 249920   | 9.821  | ng/uL  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.469  | 132  | 269183   | 9.627  | ng/uL  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 279875   | 9.266  | ng/uL  | 0.00     |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 131603   | 9.311  | ng/uL  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 142264   | 9.293  | ng/uL  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 265269   | 9.694  | ng/uL  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.899 | 131  | 395490   | 9.524  | ng/uL  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 827676   | 9.807  | ng/uL  | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 943486   | 9.756  | ng/uL  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.781 | 143  | 151599   | 8.763  | ng/uL  | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 712537   | 10.000 | ng/uL  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.698 | 200  | 121515   | 8.199  | ng/uL  | 0.00     |
| 73) Anthracene-d10            | 17.434 | 188  | 1067274  | 9.978  | ng/uL  | 0.00     |
| 81) Pyrene-d10                | 19.728 | 212  | 1207324  | 9.787  | ng/uL  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.810 | 264  | 1203862  | 9.967  | ng/uL  | -0.02    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.346  | 88   | 43679    | 3.947  | ng/uL  | 98       |
| 5) Pyridine                   | 3.764  | 79   | 295836   | 9.326  | ng/uL  | 94       |
| 6) Benzaldehyde               | 7.081  | 77   | 217998m  | 11.706 | ng/uL  |          |
| 8) Phenol                     | 7.128  | 94   | 373855   | 9.435  | ng/uL  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.364  | 93   | 303378   | 9.752  | ng/uL  | 98       |
| 12) 2-Chlorophenol            | 7.505  | 128  | 282136   | 9.661  | ng/uL  | 98       |
| 13) 2-Methylphenol            | 8.387  | 108  | 269807   | 9.274  | ng/uL  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.469  | 45   | 609363   | 9.976  | ng/uL  | 98       |
| 16) Acetophenone              | 8.775  | 105  | 472001   | 9.883  | ng/uL  | 98       |
| 17) N-Nitroso-di-n-propyla... | 8.758  | 70   | 249350   | 9.679  | ng/uL  | 99       |
| 18) 4-Methylphenol            | 8.716  | 108  | 304490   | 9.412  | ng/uL  | 92       |
| 19) Hexachloroethane          | 9.028  | 117  | 112246   | 9.524  | ng/uL  | 100      |
| 22) Nitrobenzene              | 9.158  | 77   | 370772   | 9.762  | ng/uL  | 98       |
| 23) Isophorone                | 9.681  | 82   | 660661   | 9.476  | ng/uL  | 100      |
| 25) 2-Nitrophenol             | 9.869  | 139  | 156447   | 9.473  | ng/uL  | 97       |
| 26) 2,4-Dimethylphenol        | 9.928  | 107  | 337462   | 9.705  | ng/uL  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.163 | 93   | 428129   | 9.868  | ng/uL  | 99       |
| 29) 2,4-Dichlorophenol        | 10.405 | 162  | 266919   | 9.719  | ng/uL  | 98       |
| 30) Naphthalene               | 10.810 | 128  | 964314   | 9.891  | ng/uL  | 99       |
| 32) 4-Chloroaniline           | 10.922 | 127  | 407622   | 9.696  | ng/uL  | 97       |
| 33) Hexachlorobutadiene       | 11.093 | 225  | 158399   | 9.895  | ng/uL  | 95       |
| 34) Caprolactam               | 11.687 | 113  | 81746m   | 8.499  | ng/uL  |          |
| 35) 4-Chloro-3-methylphenol   | 12.040 | 107  | 309482   | 9.611  | ng/uL  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
 Data File : BN024820.D  
 Acq On : 10 Apr 2023 11:55  
 Operator : CG/JU  
 Sample : SST001090  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 10 15:00:02 2023  
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
 Supervised By :Jagrut Upadhyay 04/11/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.416 | 142  | 633044   | 9.825  | ng/ul  | 99       |
| 37) 1-Methylnaphthalene       | 12.634 | 142  | 638201   | 9.922  | ng/ul  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.787 | 216  | 314069   | 9.891  | ng/ul  | 96       |
| 40) Hexachlorocyclopentadiene | 12.763 | 237  | 173643   | 8.968  | ng/ul  | 96       |
| 41) 2,4,6-Trichlorophenol     | 13.022 | 196  | 209251   | 9.645  | ng/ul  | 96       |
| 42) 2,4,5-Trichlorophenol     | 13.093 | 196  | 230242   | 9.697  | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.422 | 154  | 859910   | 9.914  | ng/ul  | 97       |
| 44) 2-Chloronaphthalene       | 13.469 | 162  | 662035   | 9.830  | ng/ul  | 98       |
| 45) 2-Nitroaniline            | 13.675 | 65   | 215232   | 9.152  | ng/ul  | 100      |
| 47) Dimethylphthalate         | 14.046 | 163  | 837324   | 9.861  | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 14.169 | 165  | 150001   | 9.050  | ng/ul  | 96       |
| 50) Acenaphthylene            | 14.310 | 152  | 1059350  | 9.973  | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.498 | 138  | 161159   | 9.253  | ng/ul  | 98       |
| 52) Acenaphthene              | 14.651 | 153  | 727656   | 9.909  | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.698 | 184  | 72678    | 6.820  | ng/ul  | 90       |
| 55) 4-Nitrophenol             | 14.798 | 109  | 127886   | 9.083  | ng/ul  | 97       |
| 56) Dibenzofuran              | 14.987 | 168  | 1002765  | 9.907  | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 14.951 | 165  | 224534   | 9.359  | ng/ul# | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.210 | 232  | 193965   | 9.921  | ng/ul  | 99       |
| 59) Diethylphthalate          | 15.404 | 149  | 835465   | 9.792  | ng/ul  | 98       |
| 61) Fluorene                  | 15.634 | 166  | 818365   | 10.052 | ng/ul  | 96       |
| 62) 4-Chlorophenyl-phenyle... | 15.628 | 204  | 385338   | 10.169 | ng/ul  | 96       |
| 63) 4-Nitroaniline            | 15.657 | 138  | 157072   | 9.295  | ng/ul  | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.716 | 198  | 124163   | 8.552  | ng/ul# | 95       |
| 67) N-Nitrosodiphenylamine    | 15.845 | 169  | 695464   | 10.066 | ng/ul  | 96       |
| 68) 4-Bromophenyl-phenylether | 16.522 | 248  | 229464   | 10.024 | ng/ul  | 94       |
| 69) Hexachlorobenzene         | 16.639 | 284  | 258070   | 9.957  | ng/ul  | 97       |
| 70) Atrazine                  | 16.792 | 200  | 222589   | 9.468  | ng/ul  | 98       |
| 71) Pentachlorophenol         | 16.987 | 266  | 147228   | 9.139  | ng/ul  | 91       |
| 72) Phenanthrene              | 17.381 | 178  | 1275576  | 9.925  | ng/ul  | 99       |
| 74) Anthracene                | 17.469 | 178  | 1308060  | 10.101 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.393 | 216  | 322728   | 9.963  | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 14.904 | 250  | 309314   | 9.981  | ng/uL  | 97       |
| 77) Carbazole                 | 17.739 | 167  | 1173547  | 10.036 | ng/ul  | 100      |
| 78) Di-n-butylphthalate       | 18.298 | 149  | 1326064  | 9.554  | ng/ul  | 99       |
| 80) Fluoranthene              | 19.392 | 202  | 1431076  | 9.679  | ng/ul  | 99       |
| 82) Pyrene                    | 19.757 | 202  | 1559321  | 10.020 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.657 | 149  | 589196   | 9.119  | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.445 | 252  | 435617   | 8.937  | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.516 | 228  | 1510368  | 9.899  | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.433 | 149  | 905310   | 9.386  | ng/ul  | 99       |
| 87) Chrysene                  | 21.569 | 228  | 1434883  | 9.873  | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.374 | 149  | 1493729  | 9.409  | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.227 | 252  | 1495026  | 9.868  | ng/ul  | 95       |
| 91) Benzo(k)fluoranthene      | 23.274 | 252  | 1502941  | 10.082 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 23.863 | 252  | 1290801  | 9.732  | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.504 | 276  | 1562054m | 9.897  | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 26.527 | 278  | 1314144m | 10.064 | ng/ul  |          |
| 96) Benzo(g,h,i)perylene      | 27.298 | 276  | 1260189m | 9.892  | ng/ul  |          |

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

**Instrument :**

BNA\_N

**ClientSampleId :**

SSTD010212

**Manual IntegrationsAPPROVED**

Reviewed By :Christian Giraldo 04/11/2023  
Supervised By :Jagrut Upadhyay 04/11/2023

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041023\  
 Data File : BN024820.D  
 Acq On : 10 Apr 2023 11:55  
 Operator : CG/JU  
 Sample : SSTD01090  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010212

Manual IntegrationsAPPROVED

Quant Time: Apr 11 02:39:34 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 10 15:00:02 2023  
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023  
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