

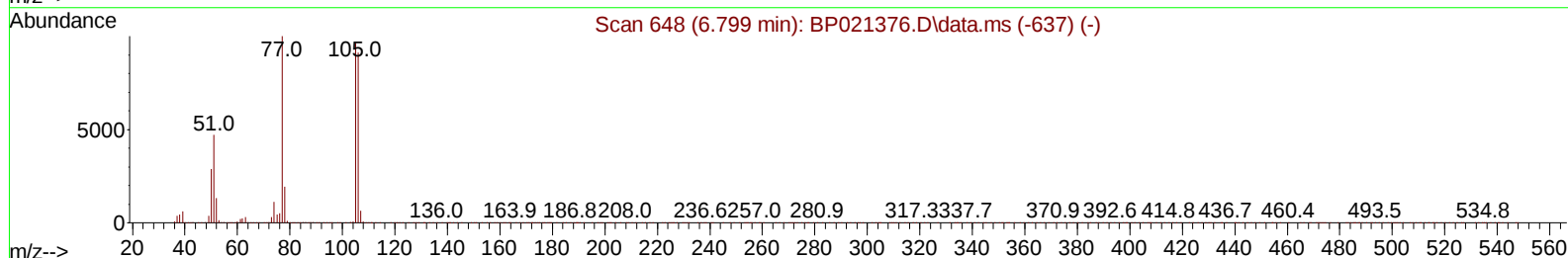
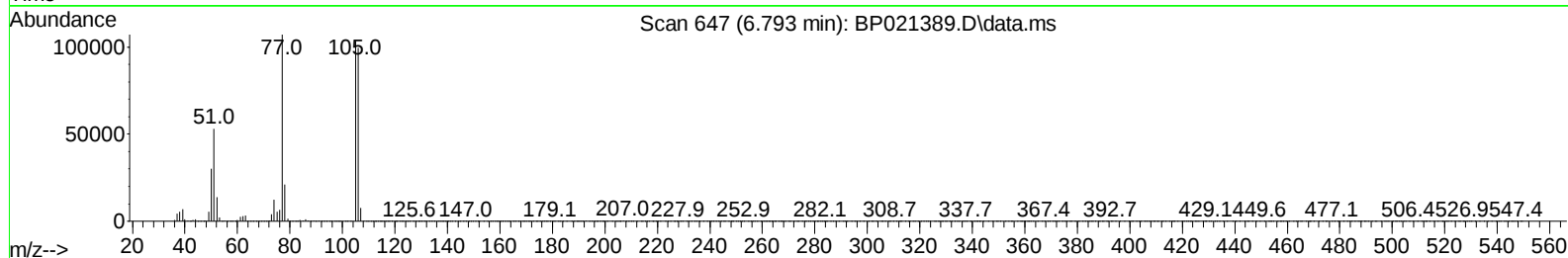
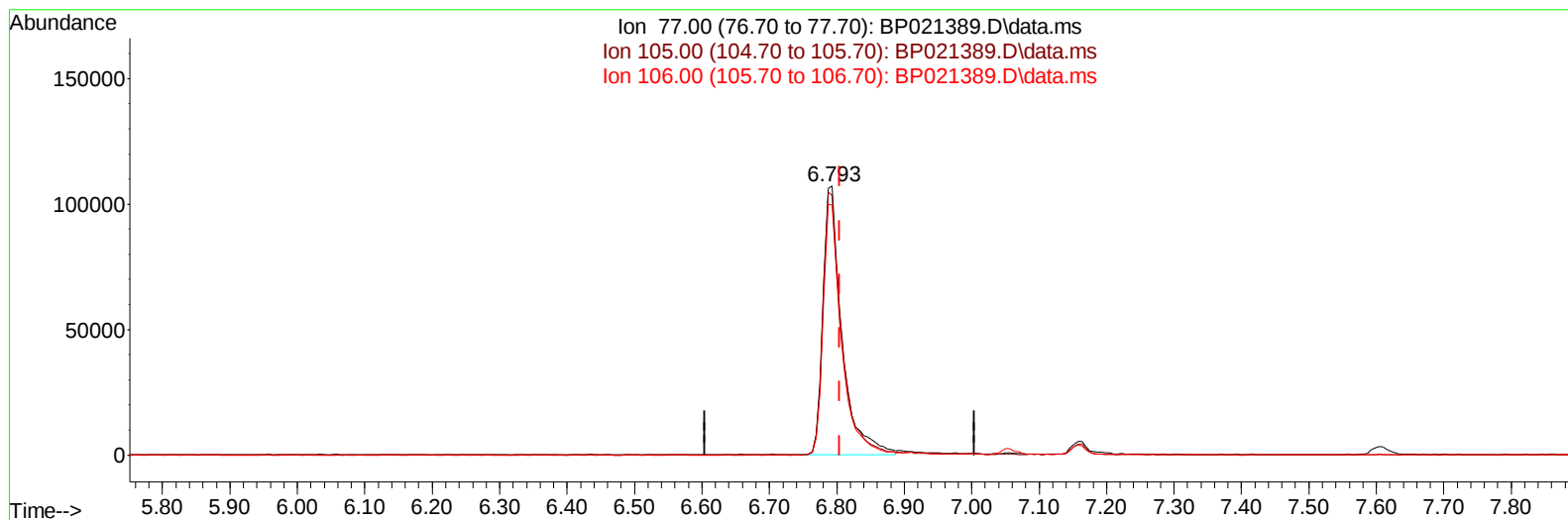
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/06/2024  
Supervised By :mohammad ahmed 08/07/2024

Quant Time: Aug 06 04:21:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration



TIC: BP021389.D\data.ms

## (6) Benzaldehyde

6.793min (-0.011) 29.25 ng/u1

response 209742

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 96.45  |
| 106.00 | 93.30  | 92.98  |
| 0.00   | 0.00   | 0.00   |

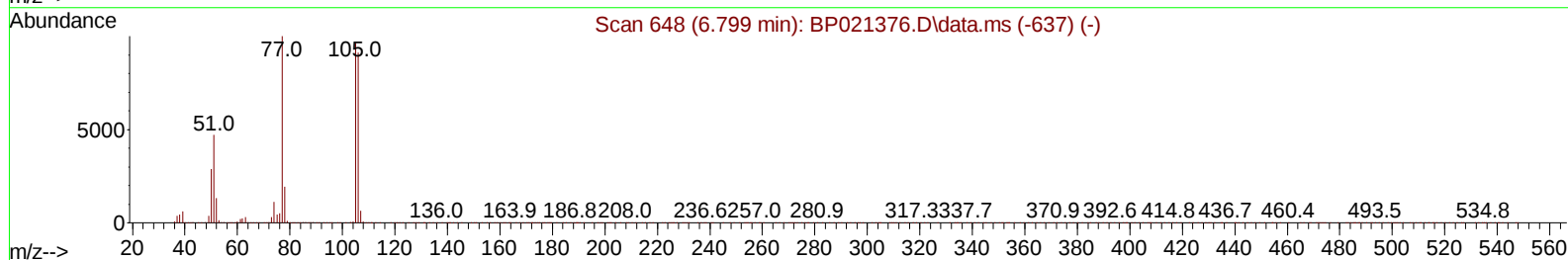
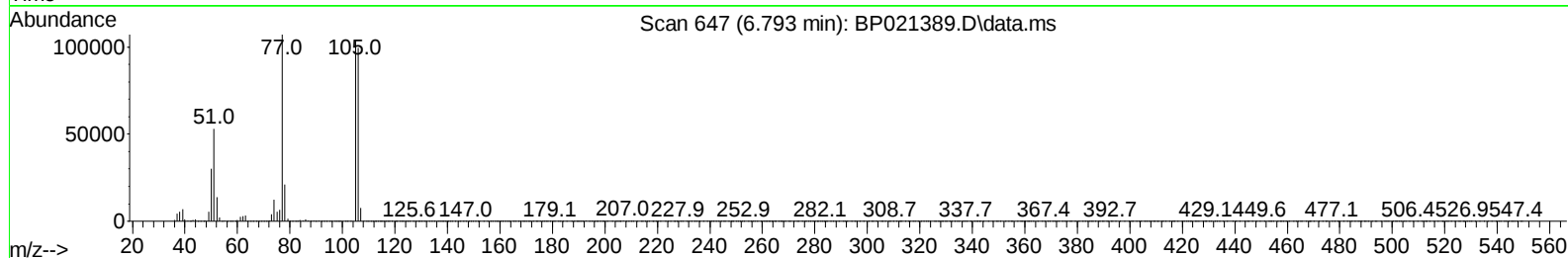
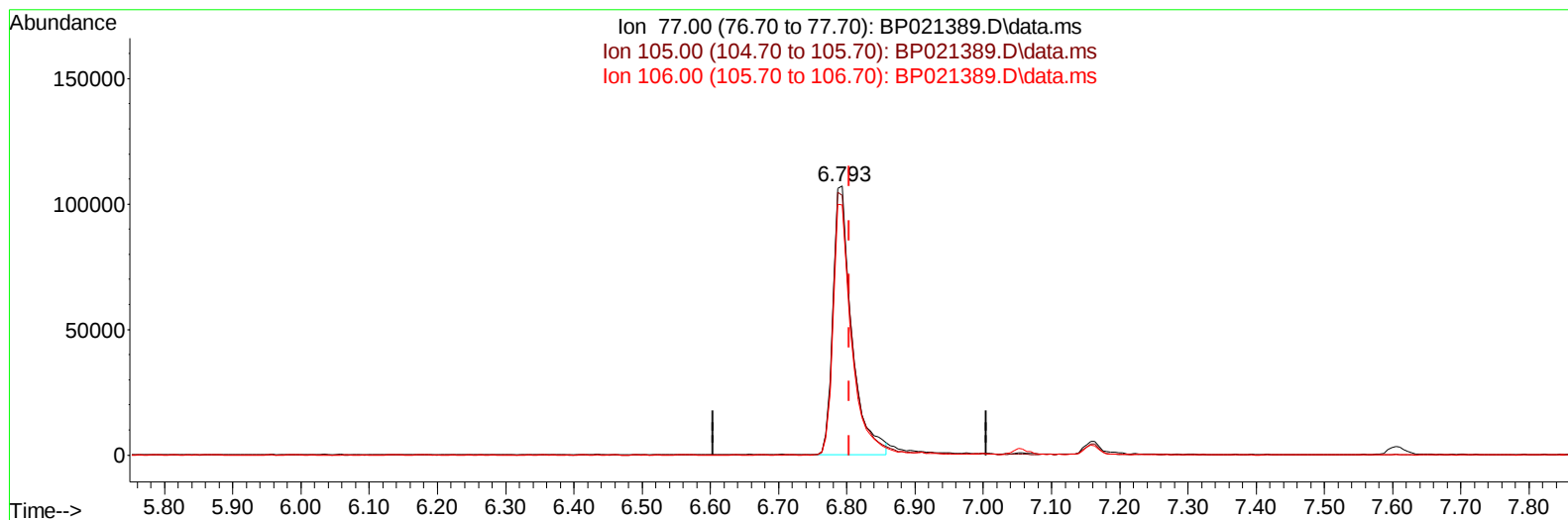
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Manual IntegrationsAPPROVED

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TIC: BP021389.D\data.ms

**(6) Benzaldehyde**

6.793min (-0.011) 28.72 ng/ul m

response 205921

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 96.45  |
| 106.00 | 93.30  | 92.98  |
| 0.00   | 0.00   | 0.00   |

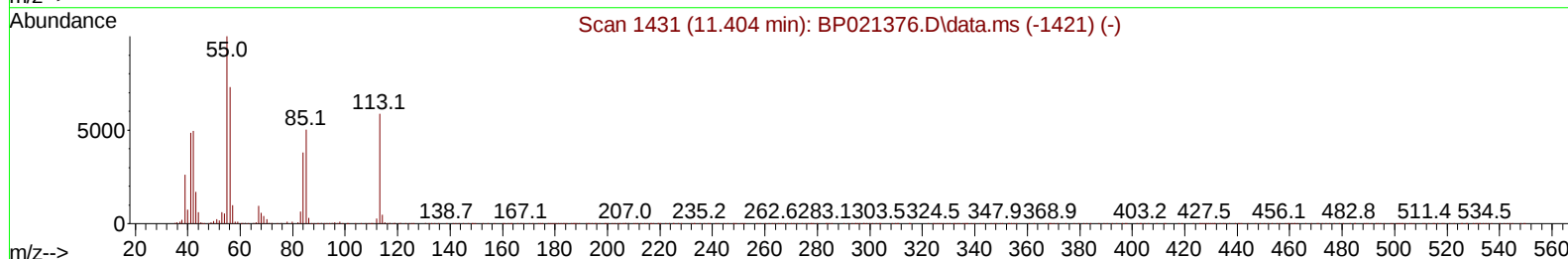
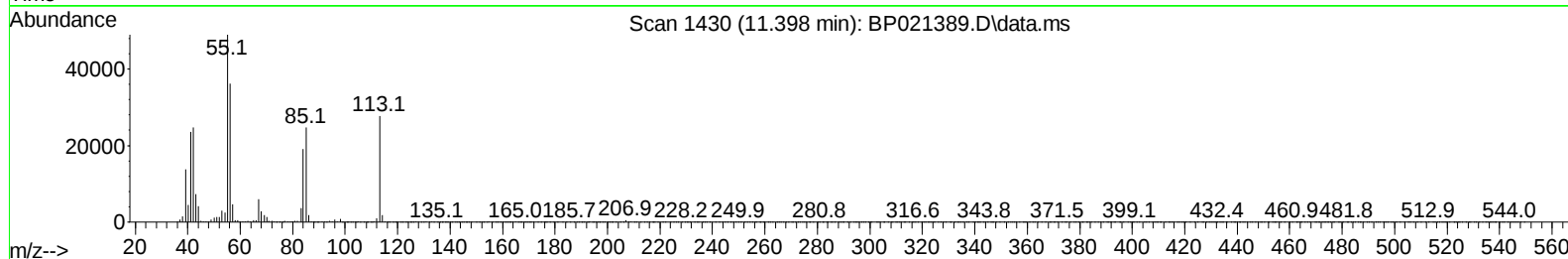
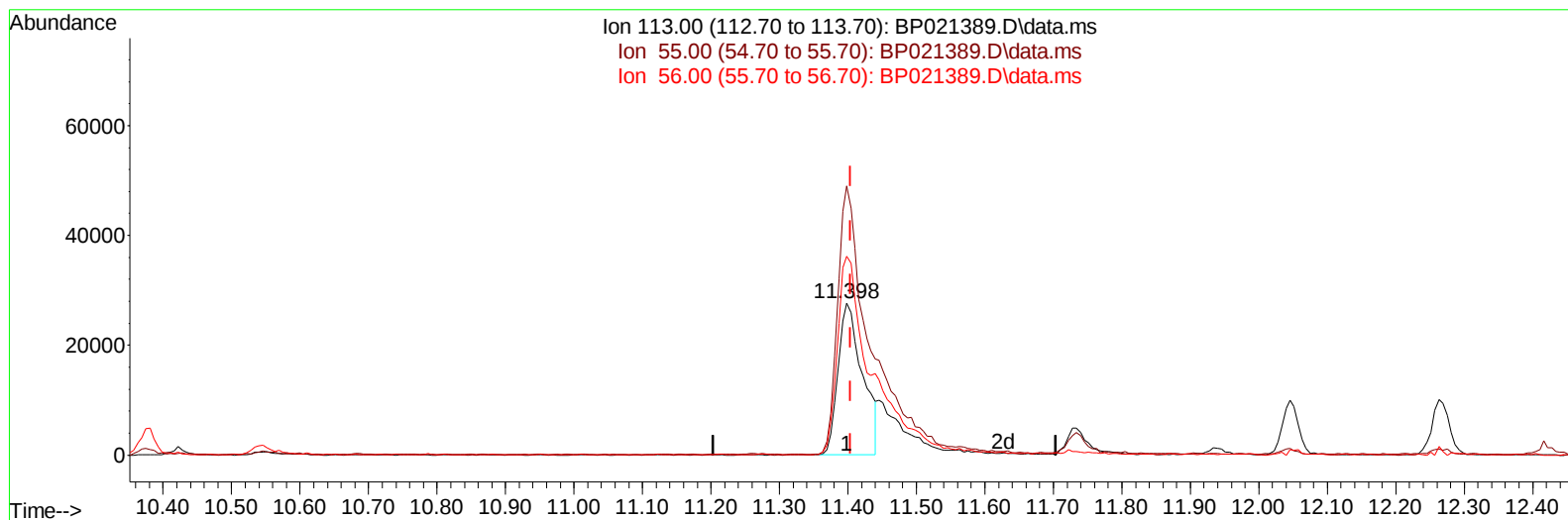
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

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Quant Title : SVOA CALIBRATION  
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TIC: BP021389.D\data.ms

(34) Caprolactam

11.398min (-0.006) 22.92 ng/ul

response 68683

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 176.73 |
| 56.00  | 131.90 | 130.72 |
| 0.00   | 0.00   | 0.00   |

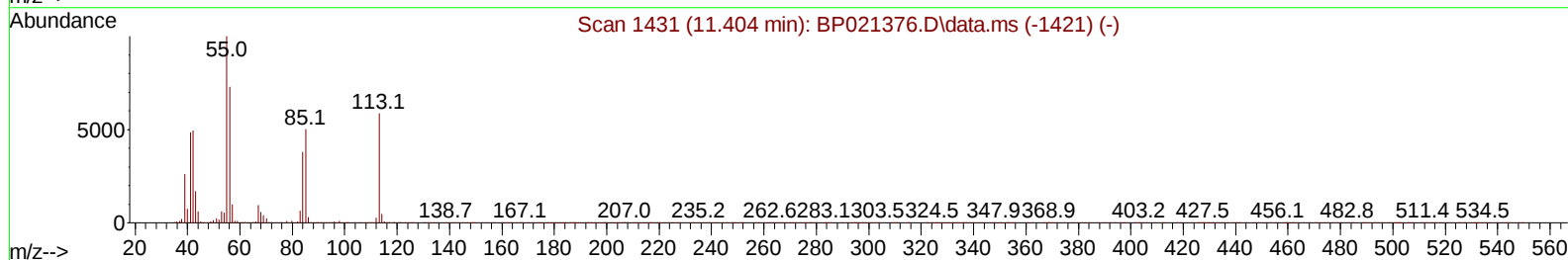
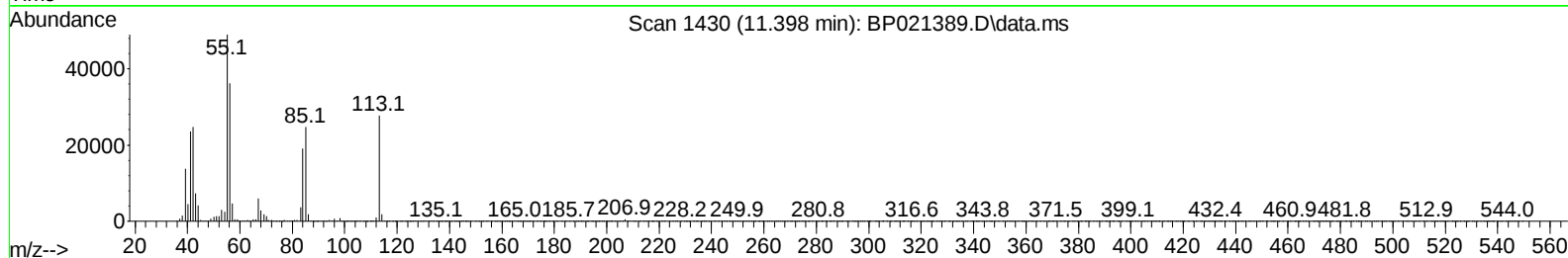
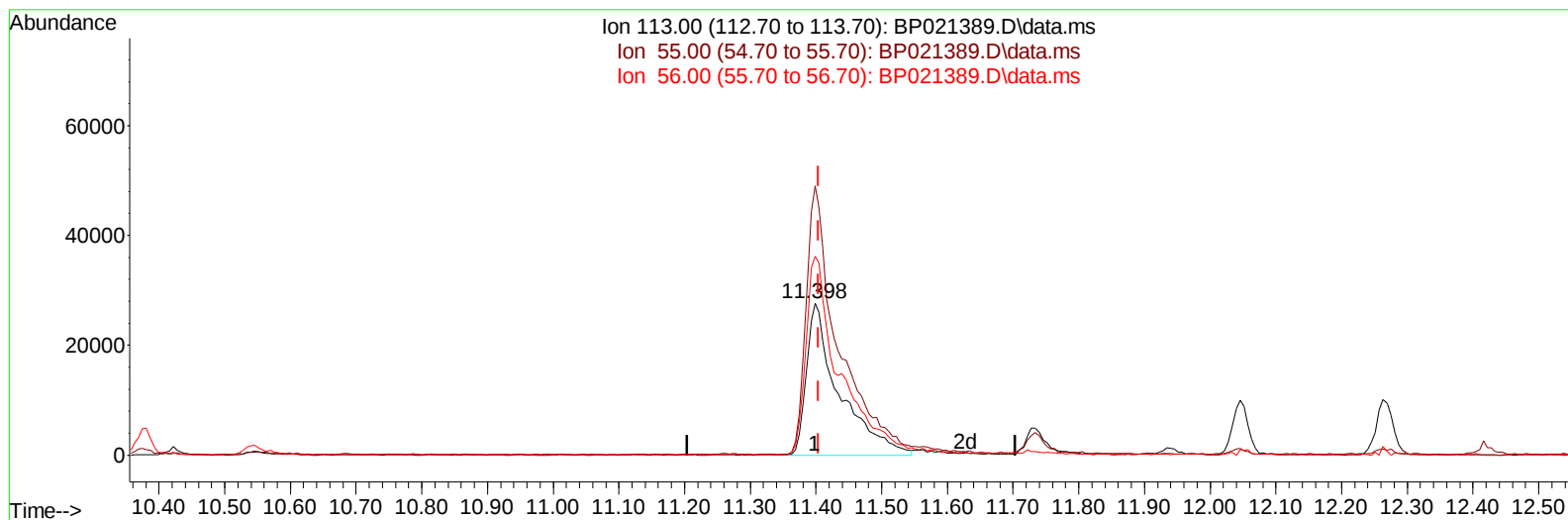
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Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:21:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
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Reviewed By :Jagrut Upadhyay 08/06/2024  
Supervised By :mohammad ahmed 08/07/2024



TIC: BP021389.D\data.ms

(34) Caprolactam

11.398min (-0.006) 31.79 ng/ul m

response 95247

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 176.73 |
| 56.00  | 131.90 | 130.72 |
| 0.00   | 0.00   | 0.00   |

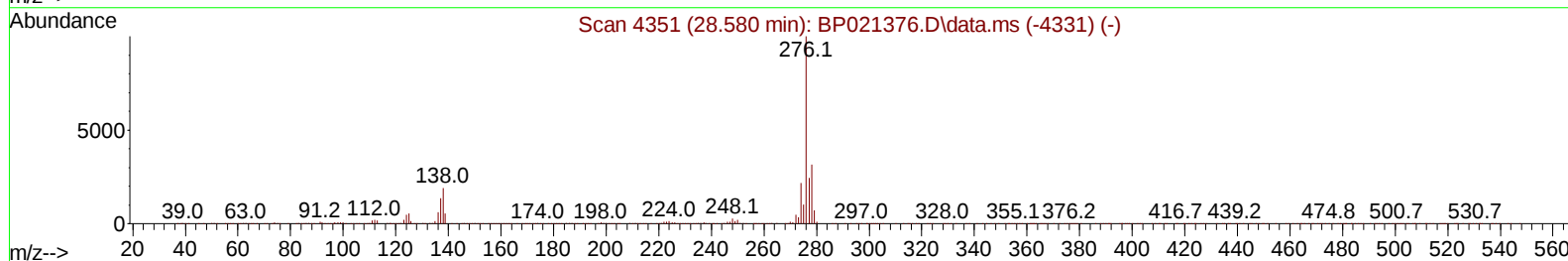
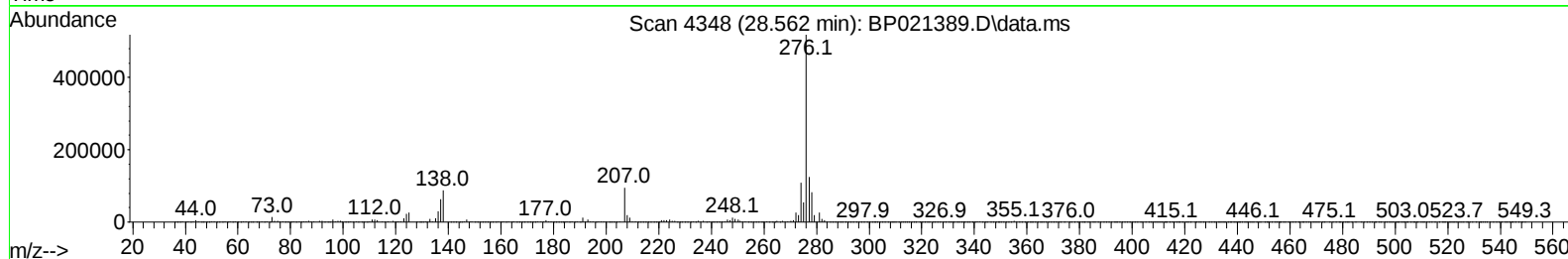
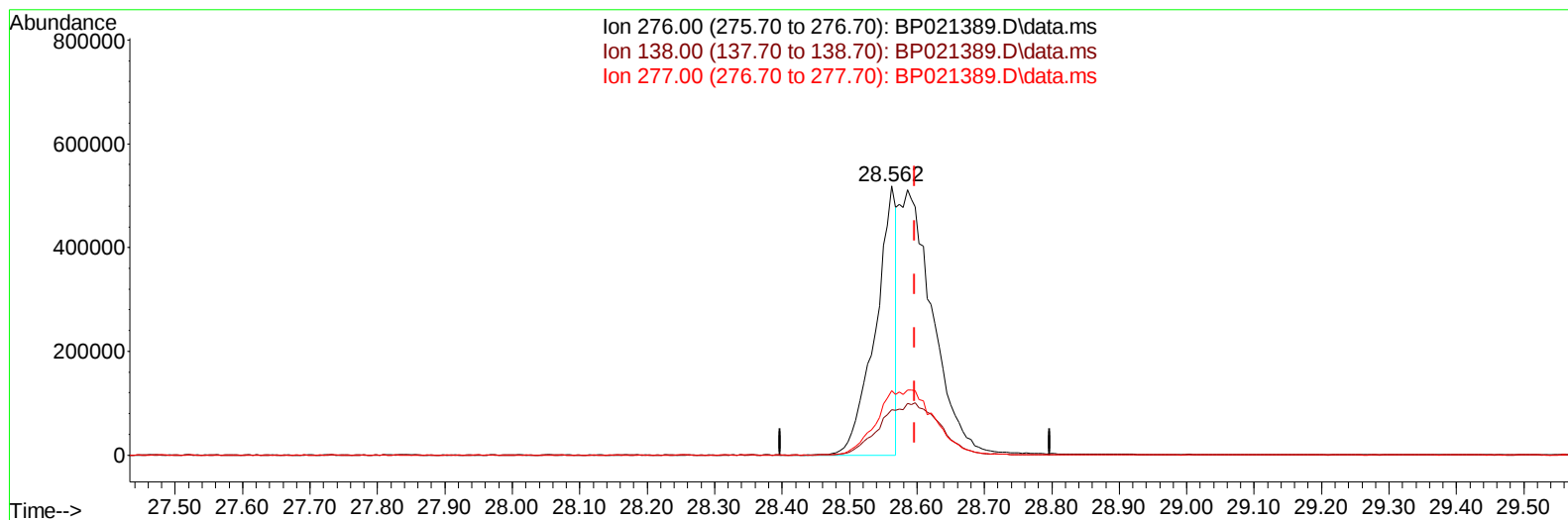
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:22:12 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
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TIC: BP021389.D\data.ms

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

**28.562min (-0.035) 12.67 ng/u1**

**response 1112330**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 16.90  |
| 277.00 | 24.00  | 24.02  |
| 0.00   | 0.00   | 0.00   |

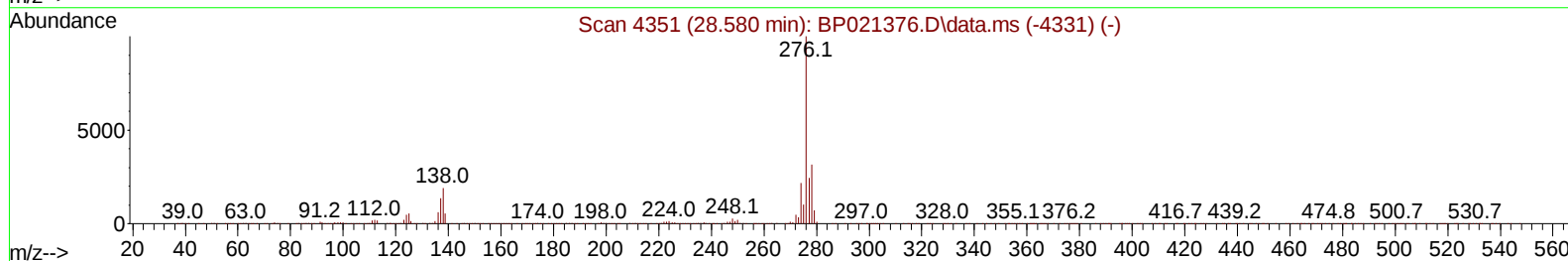
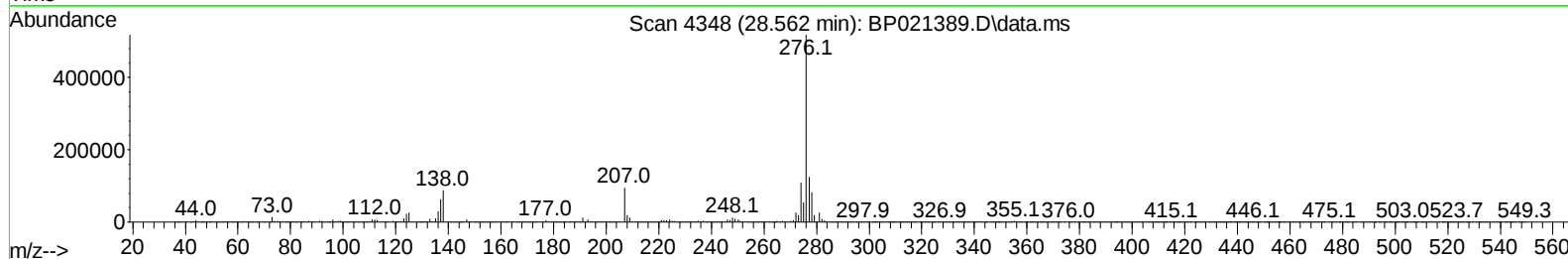
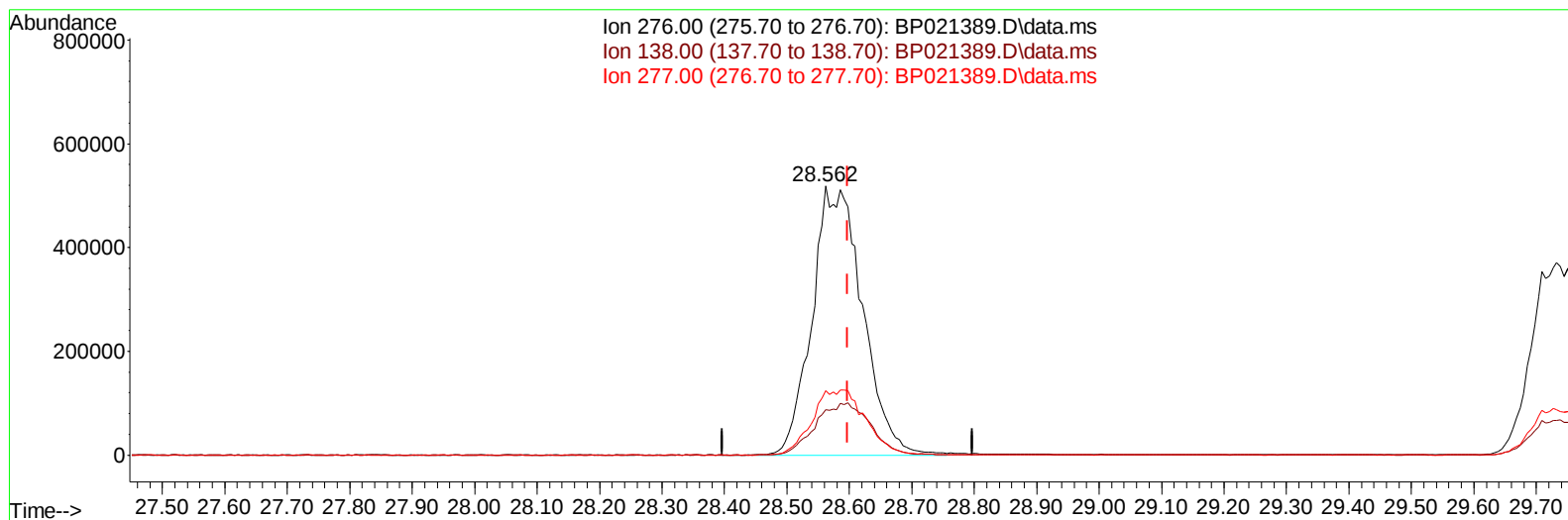
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:22:12 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
Supervised By :mohammad ahmed 08/07/2024



TIC: BP021389.D\data.ms

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

28.562min (-0.035) 32.97 ng/ul m

response 2893762

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 16.90  |
| 277.00 | 24.00  | 24.02  |
| 0.00   | 0.00   | 0.00   |

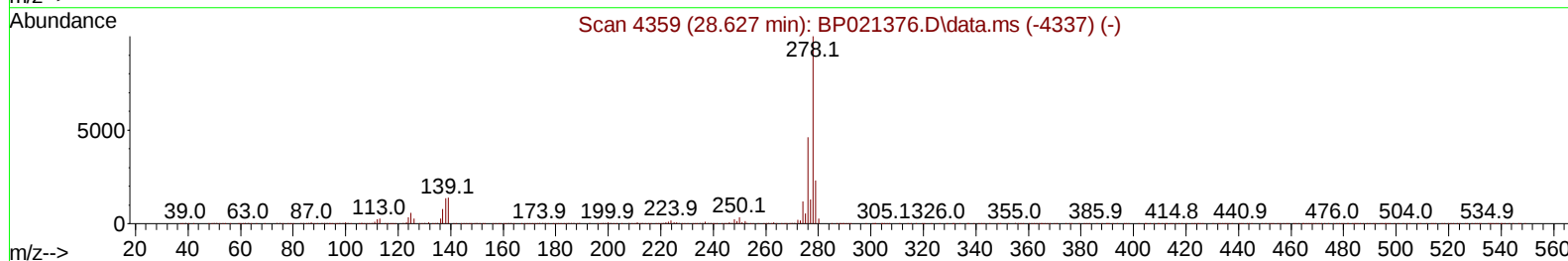
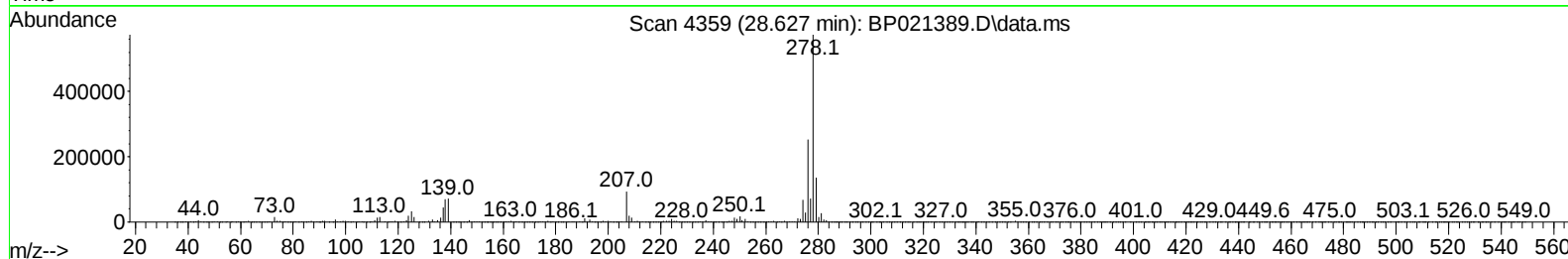
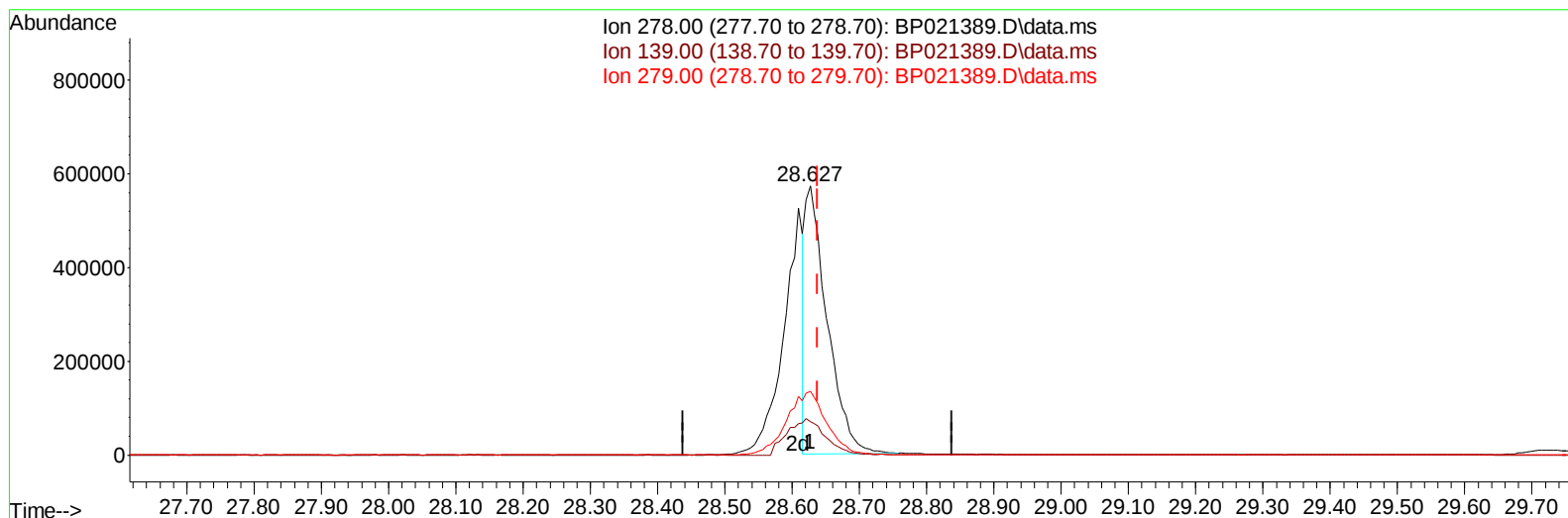
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
 Data File : BP021389.D  
 Acq On : 06 Aug 2024 03:56  
 Operator : CG/JU  
 Sample : PB162219BS  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:22:12 2024  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 29 12:46:18 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
 Supervised By :mohammad ahmed 08/07/2024



TIC: BP021389.D\data.ms

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

**28.627min (-0.012) 17.73 ng/u1**

**response 1288513**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 13.60  | 12.47  |
| 279.00 | 24.20  | 23.75  |
| 0.00   | 0.00   | 0.00   |

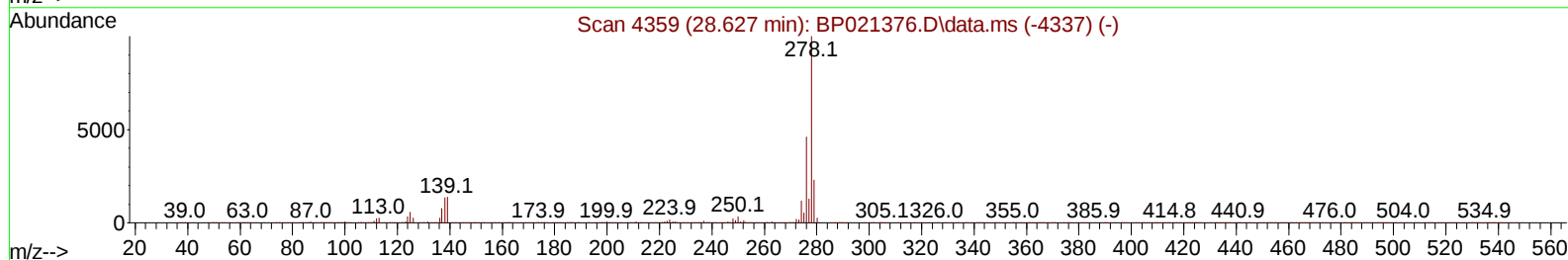
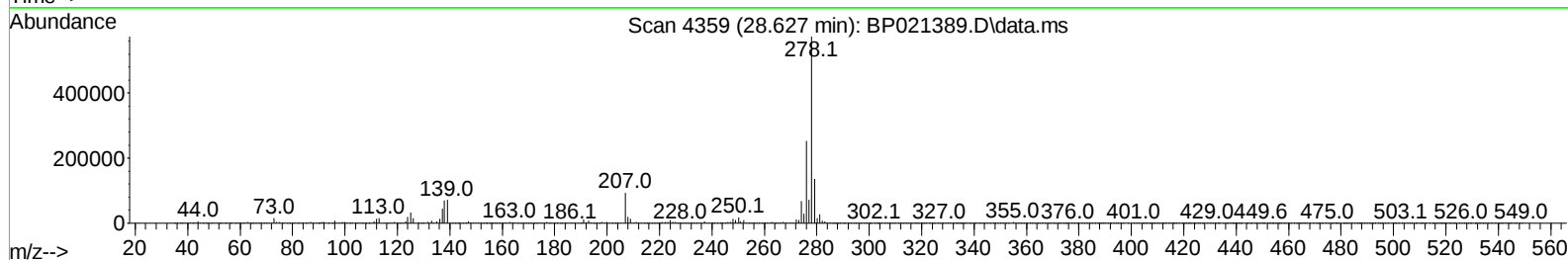
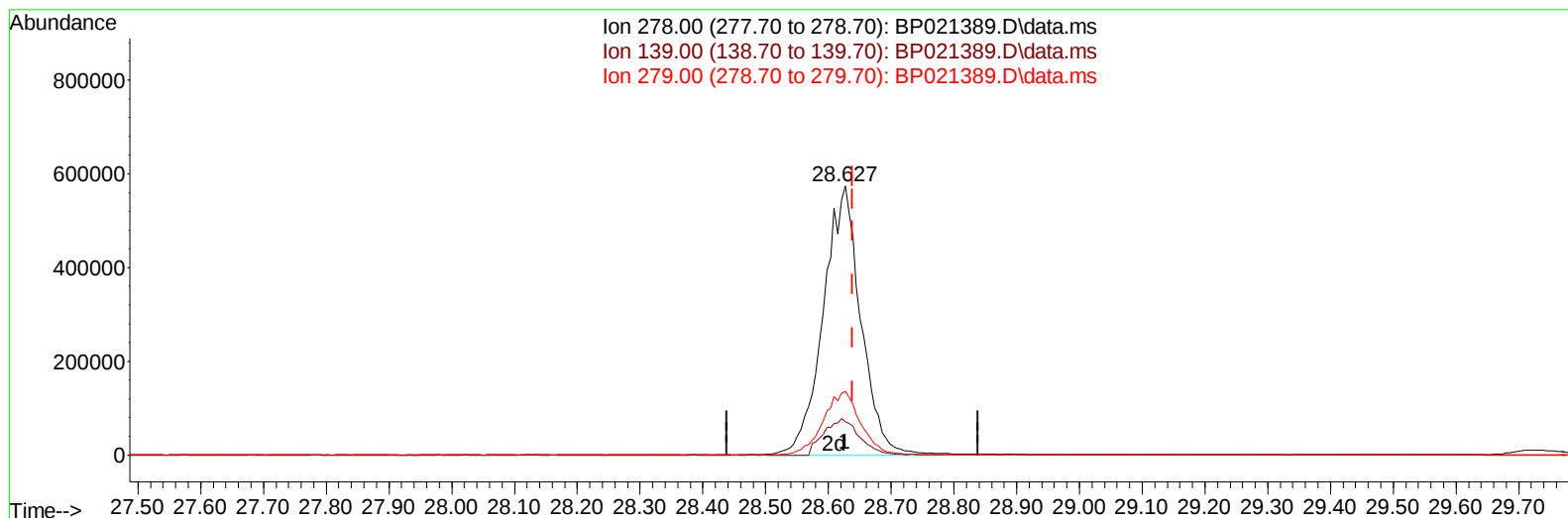
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:22:12 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
Supervised By :mohammad ahmed 08/07/2024



TIC: BP021389.D\data.ms

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

**28.627min (-0.012) 32.84 ng/ul m**

**response 2386441**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 13.60  | 12.47  |
| 279.00 | 24.20  | 23.75  |
| 0.00   | 0.00   | 0.00   |

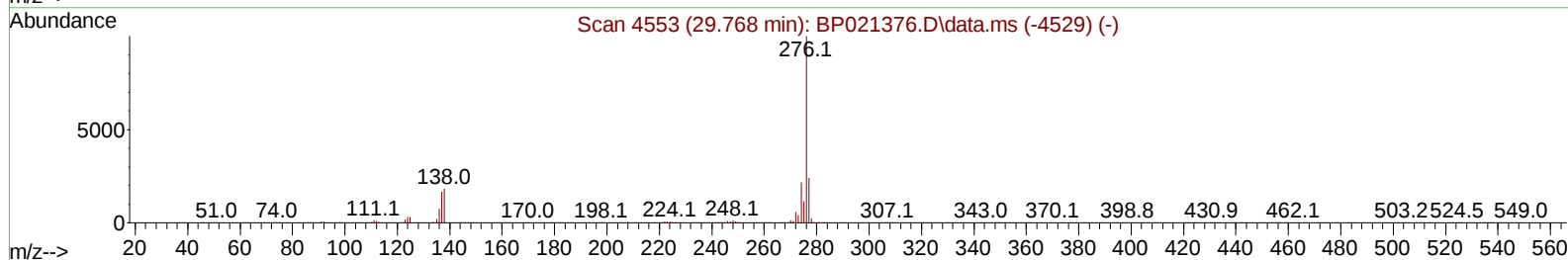
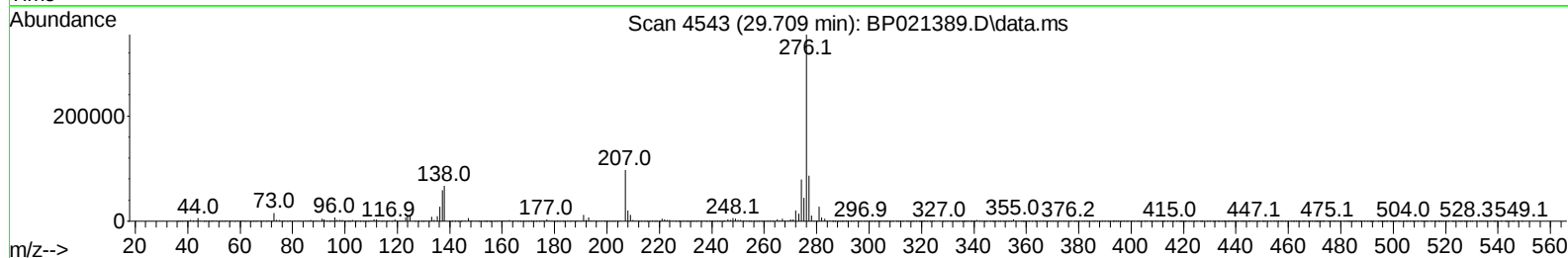
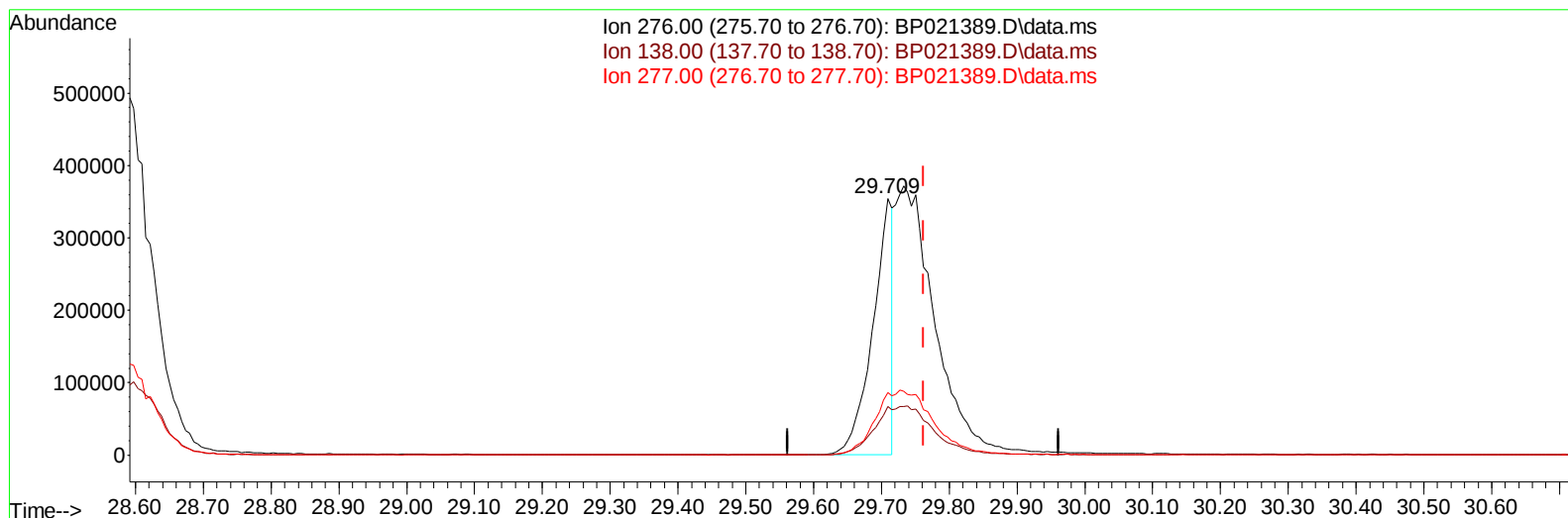
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Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

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TIC: BP021389.D\data.ms

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

29.709min (-0.053) 10.05 ng/u1

response 718598

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 18.95  |
| 277.00 | 22.70  | 24.28  |
| 0.00   | 0.00   | 0.00   |

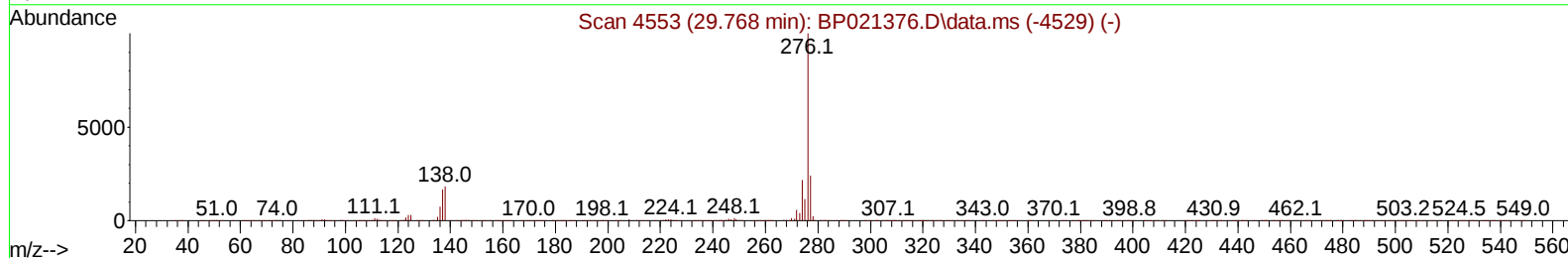
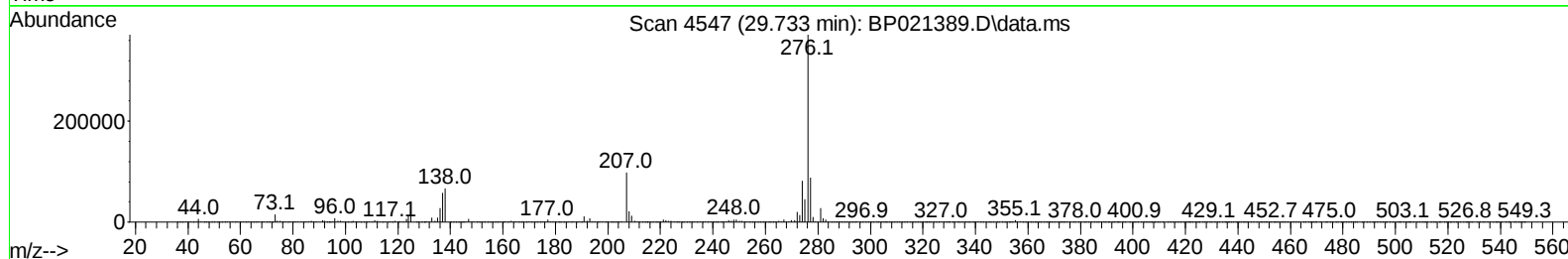
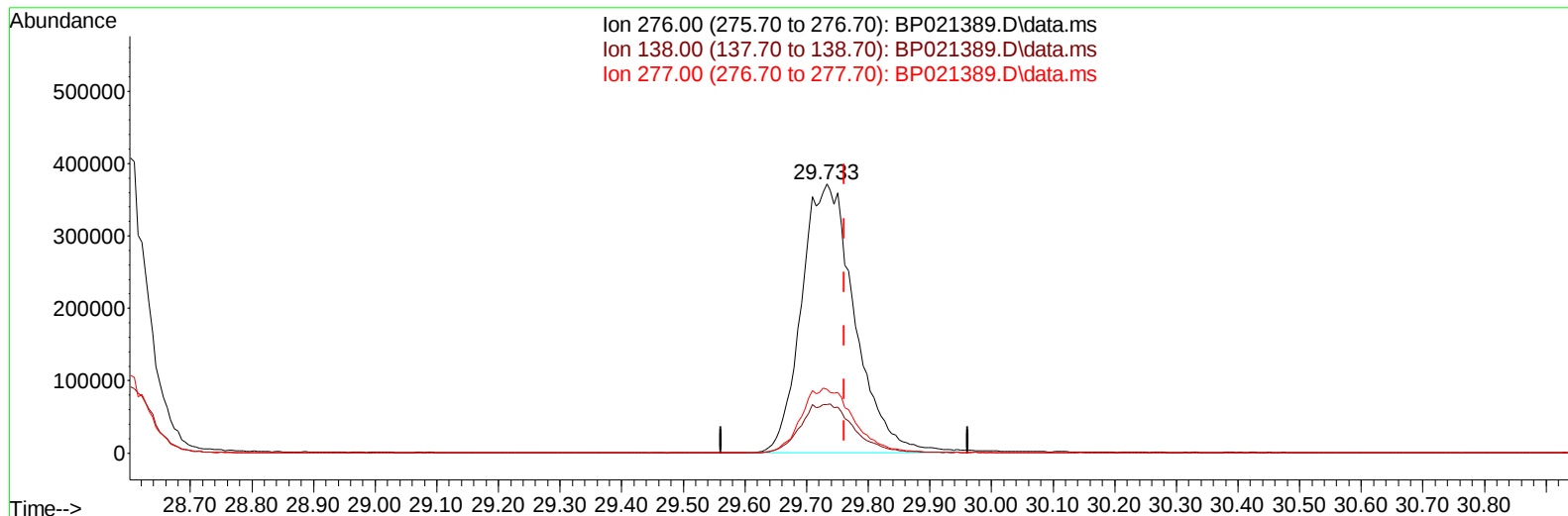
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Data File : BP021389.D  
Acq On : 06 Aug 2024 03:56  
Operator : CG/JU  
Sample : PB162219BS  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:22:12 2024  
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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
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TIC: BP021389.D\data.ms

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

29.733min (-0.029) 31.16 ng/ul m

response 2227009

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 18.07  |
| 277.00 | 22.70  | 23.66  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
 Data File : BP021389.D  
 Acq On : 06 Aug 2024 03:56  
 Operator : CG/JU  
 Sample : PB162219BS  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:23:45 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 29 12:46:18 2024  
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Reviewed By :Jagrut Upadhyay 08/06/2024  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.604  | 152  | 174465   | 20.000 | ng/ul  | -0.03    |
| 20) Naphthalene-d8            | 10.381 | 136  | 613898   | 20.000 | ng/ul  | -0.01    |
| 38) Acenaphthene-d10          | 14.257 | 164  | 379686   | 20.000 | ng/ul  | -0.01    |
| 64) Phenanthrene-d10          | 17.080 | 188  | 840876   | 20.000 | ng/ul  | -0.01    |
| 79) Chrysene-d12              | 21.527 | 240  | 1001749  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.815 | 264  | 1187669  | 20.000 | ng/ul  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 28450    | 7.421  | ng/uL  | -0.01    |
| 4) Pyridine-d5                | 3.581  | 84   | 318714   | 28.714 | ng/ul  | -0.01    |
| 7) Phenol-d5                  | 6.828  | 99   | 394513   | 27.736 | ng/ul  | -0.02    |
| 9) Bis-(2-Chloroethyl)eth...  | 6.957  | 67   | 240867   | 27.943 | ng/ul  | -0.02    |
| 11) 2-Chlorophenol-d4         | 7.157  | 132  | 329515   | 28.118 | ng/ul  | -0.02    |
| 15) 4-Methylphenol-d8         | 8.340  | 113  | 296941   | 25.136 | ng/ul  | -0.02    |
| 21) Nitrobenzene-d5           | 8.775  | 128  | 153108   | 32.110 | ng/ul  | -0.02    |
| 24) 2-Nitrophenol-d4          | 9.487  | 143  | 170927   | 31.320 | ng/ul  | -0.02    |
| 28) 2,4-Dichlorophenol-d3     | 10.034 | 165  | 300921   | 30.493 | ng/ul  | -0.02    |
| 31) 4-Chloroaniline-d4        | 10.546 | 131  | 367035   | 28.075 | ng/ul  | -0.01    |
| 46) Dimethylphthalate-d6      | 13.669 | 166  | 871601   | 31.578 | ng/ul  | -0.01    |
| 49) Acenaphthylene-d8         | 13.951 | 160  | 1006577  | 32.478 | ng/ul  | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 178960   | 45.571 | ng/ul  | -0.01    |
| 60) Fluorene-d10              | 15.275 | 176  | 749219   | 33.087 | ng/ul  | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 158369   | 31.128 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.175 | 188  | 1221991  | 32.719 | ng/ul  | -0.02    |
| 81) Pyrene-d10                | 19.569 | 212  | 1672337  | 27.003 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.586 | 264  | 1960362  | 33.467 | ng/ul  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.193  | 88   | 51120    | 12.124 | ng/uL  | 96       |
| 5) Pyridine                   | 3.599  | 79   | 330355   | 28.856 | ng/ul  | 99       |
| 6) Benzaldehyde               | 6.793  | 77   | 205921m  | 28.717 | ng/ul  |          |
| 8) Phenol                     | 6.857  | 94   | 413743   | 28.746 | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.052  | 93   | 330301   | 27.780 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.187  | 128  | 347707   | 29.603 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.081  | 108  | 306653   | 27.549 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.122  | 45   | 477706   | 27.810 | ng/ul  | 96       |
| 16) Acetophenone              | 8.440  | 105  | 491247   | 26.943 | ng/ul  | 93       |
| 17) N-Nitroso-di-n-propyla... | 8.410  | 70   | 240416   | 25.166 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.410  | 108  | 321279   | 26.942 | ng/ul  | 96       |
| 19) Hexachloroethane          | 8.651  | 117  | 152563   | 28.905 | ng/ul  | 98       |
| 22) Nitrobenzene              | 8.816  | 77   | 377687   | 33.066 | ng/ul  | 99       |
| 23) Isophorone                | 9.328  | 82   | 658590   | 29.003 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.516  | 139  | 185926   | 32.665 | ng/ul  | 100      |
| 26) 2,4-Dimethylphenol        | 9.593  | 107  | 273351   | 24.354 | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 9.804  | 93   | 408702   | 29.610 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.069 | 162  | 300280   | 30.875 | ng/ul  | 99       |
| 30) Naphthalene               | 10.422 | 128  | 1033528  | 32.006 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.575 | 127  | 344229   | 27.574 | ng/ul  | 97       |
| 33) Hexachlorobutadiene       | 10.687 | 225  | 233443   | 32.012 | ng/ul  | 99       |
| 34) Caprolactam               | 11.398 | 113  | 95247m   | 31.791 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.728 | 107  | 332033   | 31.408 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080524\  
 Data File : BP021389.D  
 Acq On : 06 Aug 2024 03:56  
 Operator : CG/JU  
 Sample : PB162219BS  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS219

Manual IntegrationsAPPROVED

Quant Time: Aug 06 04:23:45 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 29 12:46:18 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024  
 Supervised By :mohammad ahmed 08/07/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.045 | 142  | 677110   | 30.547 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.263 | 142  | 677993   | 29.740 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.422 | 216  | 402069   | 32.607 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.369 | 237  | 53079    | 8.229  | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.687 | 196  | 218409   | 28.032 | ng/ul | 98       |
| 42) 2,4,5-Trichlorophenol     | 12.798 | 196  | 243077   | 29.486 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.075 | 154  | 882870   | 31.817 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.122 | 162  | 708329   | 31.976 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.363 | 65   | 214288   | 34.981 | ng/ul | 98       |
| 47) Dimethylphthalate         | 13.722 | 163  | 874719   | 31.233 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 13.863 | 165  | 184121   | 32.884 | ng/ul | 98       |
| 50) Acenaphthylene            | 13.981 | 152  | 1128560  | 32.162 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.216 | 138  | 184711   | 37.910 | ng/ul | 99       |
| 52) Acenaphthene              | 14.328 | 153  | 741322   | 31.830 | ng/ul | 98       |
| 53) 2,4-Dinitrophenol         | 14.463 | 184  | 92015    | 31.293 | ng/ul | 98       |
| 55) 4-Nitrophenol             | 14.592 | 109  | 133840   | 41.898 | ng/ul | 90       |
| 56) Dibenzofuran              | 14.669 | 168  | 1042741  | 32.694 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.698 | 165  | 282439   | 36.746 | ng/ul | 92       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.922 | 232  | 189479   | 26.841 | ng/ul | 100      |
| 59) Diethylphthalate          | 15.110 | 149  | 894617   | 31.963 | ng/ul | 99       |
| 61) Fluorene                  | 15.333 | 166  | 868430   | 33.368 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.322 | 204  | 449315   | 31.816 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.422 | 138  | 204080   | 50.578 | ng/ul | 95       |
| 66) 4,6-Dinitro-2-methylph... | 15.475 | 198  | 171005   | 31.695 | ng/ul | 98       |
| 67) N-Nitrosodiphenylamine    | 15.557 | 169  | 744516   | 30.405 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.245 | 248  | 293209   | 29.749 | ng/ul | 97       |
| 69) Hexachlorobenzene         | 16.357 | 284  | 356386   | 31.536 | ng/ul | 99       |
| 70) Atrazine                  | 16.545 | 200  | 28767    | 3.224  | ng/ul | 97       |
| 71) Pentachlorophenol         | 16.745 | 266  | 137022   | 22.324 | ng/ul | 95       |
| 72) Phenanthrene              | 17.116 | 178  | 1454840  | 33.352 | ng/ul | 99       |
| 74) Anthracene                | 17.216 | 178  | 1451402  | 32.638 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.034 | 216  | 399007   | 29.647 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.581 | 250  | 385753   | 28.845 | ng/uL | 100      |
| 77) Carbazole                 | 17.510 | 167  | 1380245  | 37.968 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.075 | 149  | 1619521  | 32.290 | ng/ul | 100      |
| 80) Fluoranthene              | 19.222 | 202  | 1941478  | 25.958 | ng/ul | 99       |
| 82) Pyrene                    | 19.604 | 202  | 2096689  | 27.549 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.533 | 149  | 874127   | 27.566 | ng/ul | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.427 | 252  | 797198   | 34.827 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.510 | 228  | 2245852  | 32.004 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.404 | 149  | 1216347  | 26.655 | ng/ul | 100      |
| 87) Chrysene                  | 21.574 | 228  | 2153108  | 33.020 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.651 | 149  | 2249373  | 28.670 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.768 | 252  | 2391439  | 33.181 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.839 | 252  | 2370082  | 32.951 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.662 | 252  | 2145496  | 31.501 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 28.562 | 276  | 2893762m | 32.972 | ng/ul |          |
| 95) Dibenzo(a,h)anthracene    | 28.627 | 278  | 2386441m | 32.837 | ng/ul |          |
| 96) Benzo(g,h,i)perylene      | 29.733 | 276  | 2227009m | 31.157 | ng/ul |          |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS219

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 08/06/2024

Supervised By :mohammad ahmed 08/07/2024

