

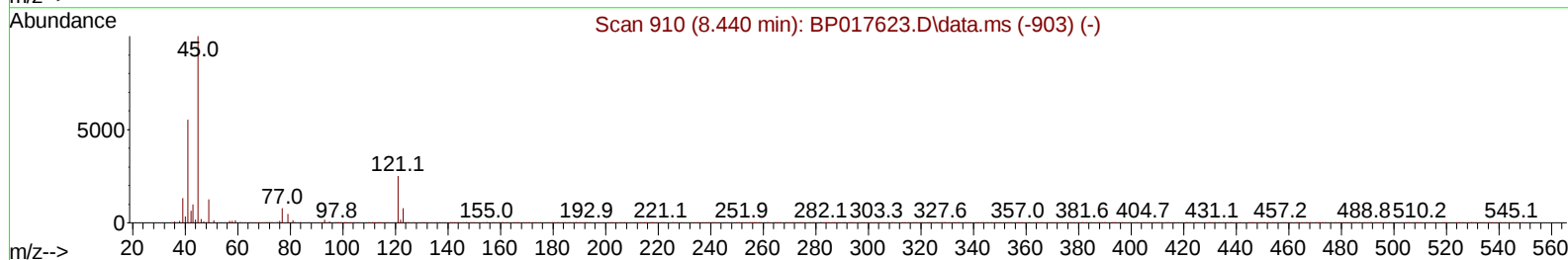
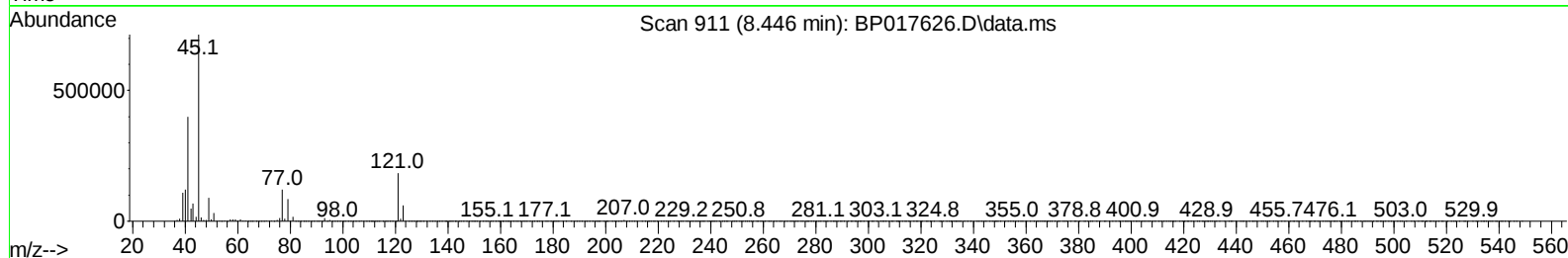
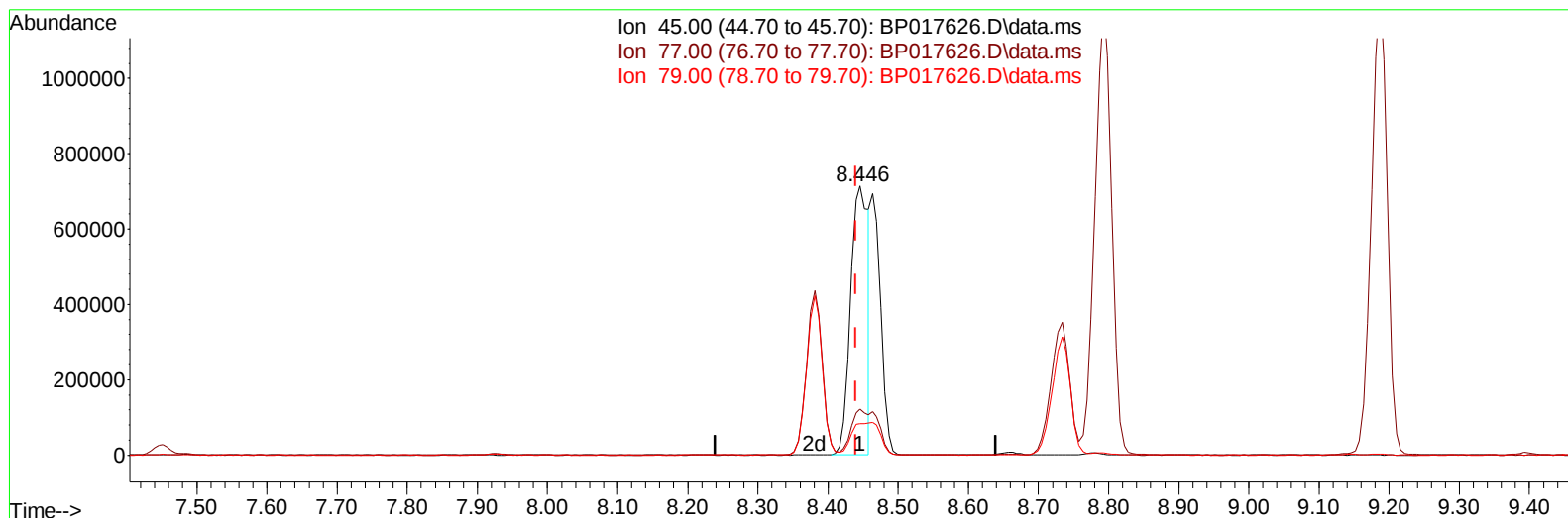
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
 Data File : BP017626.D  
 Acq On : 09 Oct 2023 20:46  
 Operator : MA/JU  
 Sample : SSTD16071  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160636

Manual IntegrationsAPPROVED

Quant Time: Oct 10 04:19:05 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 10 03:51:30 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/10/2023  
 Supervised By :mohammad ahmed 10/11/2023



TIC: BP017626.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.446min (+ 0.006) 94.49 ng/u1

response 1256006

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.80  | 17.12  |
| 79.00 | 13.00  | 11.78  |
| 0.00  | 0.00   | 0.00   |

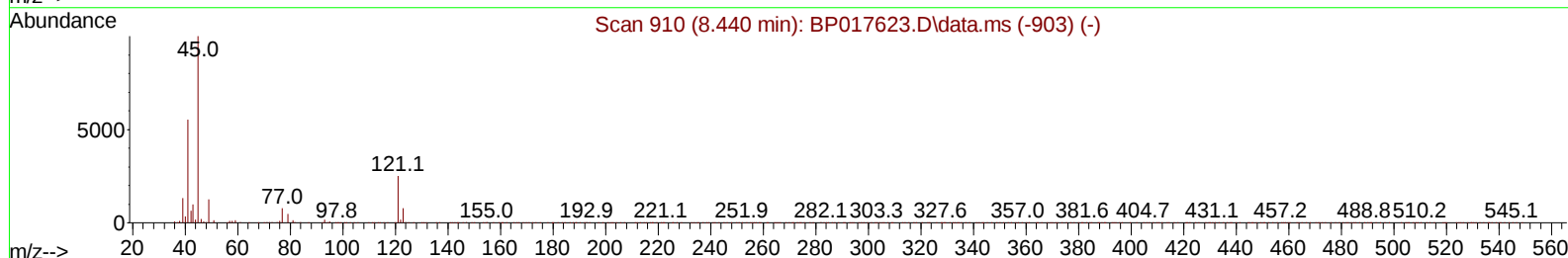
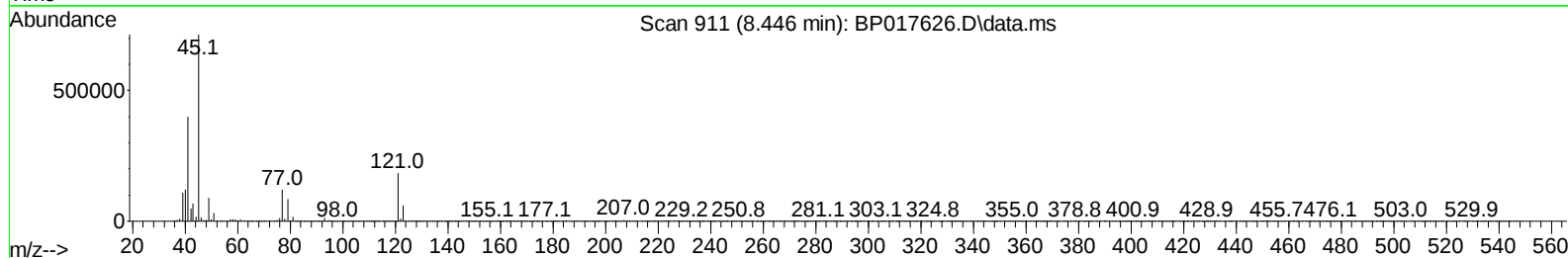
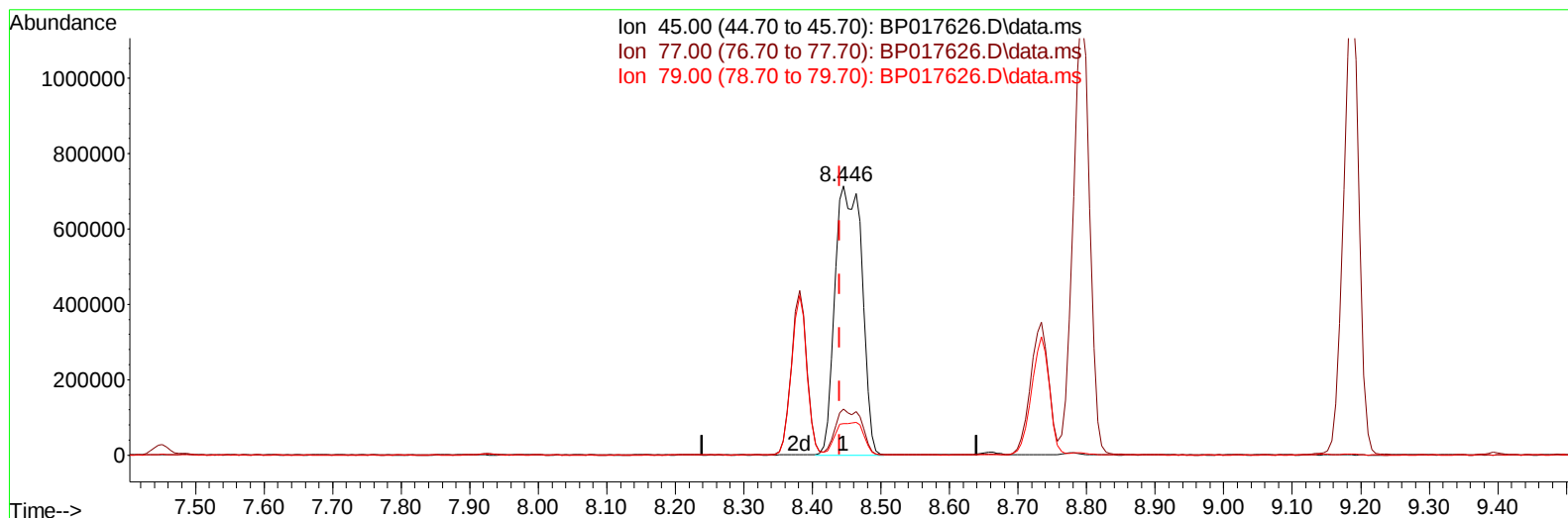
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 Data File : BP017626.D  
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Instrument :  
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Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 10/10/2023  
 Supervised By :mohammad ahmed 10/11/2023



TIC: BP017626.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.446min (+ 0.006) 146.27 ng/ul m

response 1944369

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.80  | 17.12  |
| 79.00 | 13.00  | 11.78  |
| 0.00  | 0.00   | 0.00   |

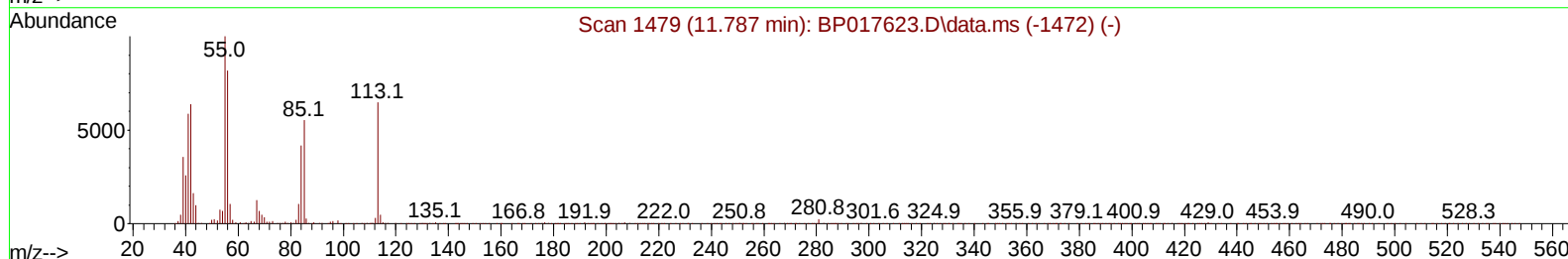
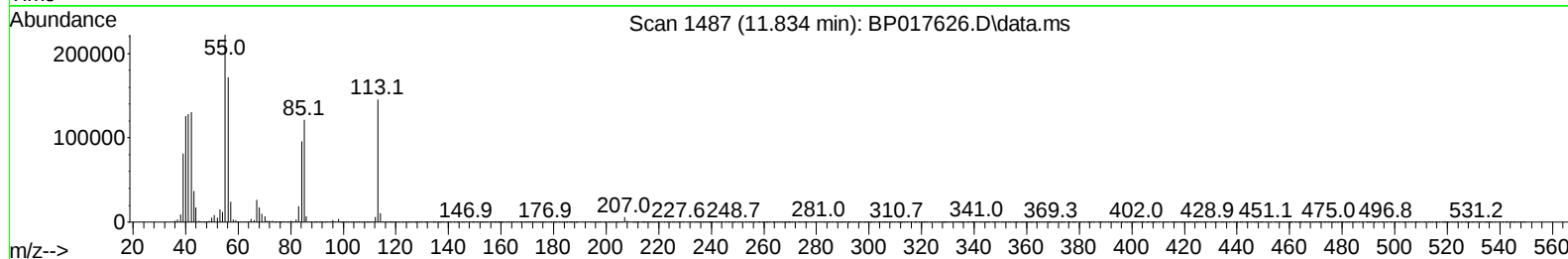
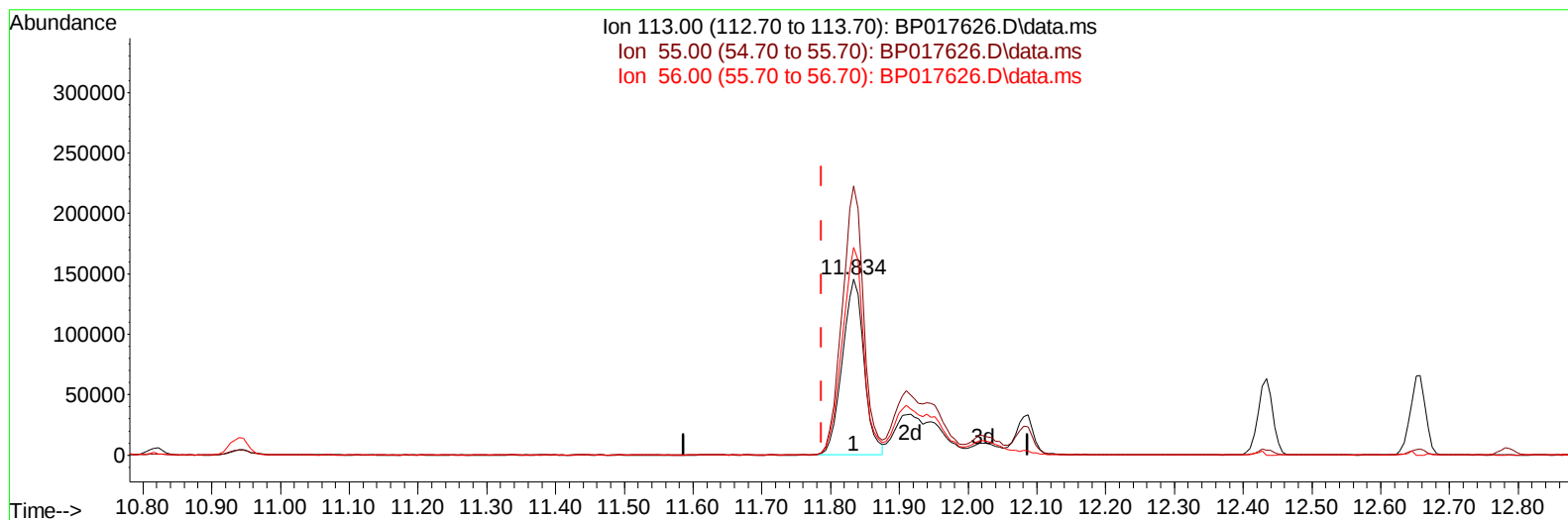
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
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Manual IntegrationsAPPROVED

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TIC: BP017626.D\data.ms

**(34) Caprolactam**

**11.834min (+ 0.047) 105.88 ng/ul**

**response 316727**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 155.20 | 152.57 |
| 56.00  | 126.50 | 117.73 |
| 0.00   | 0.00   | 0.00   |

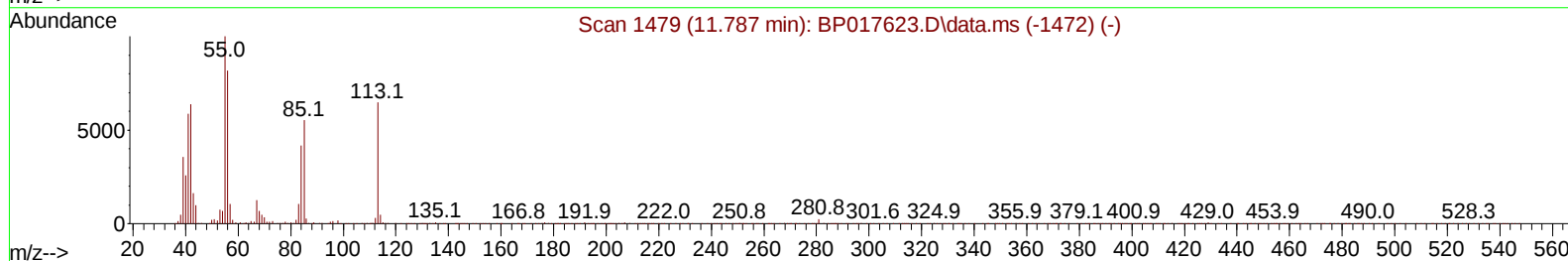
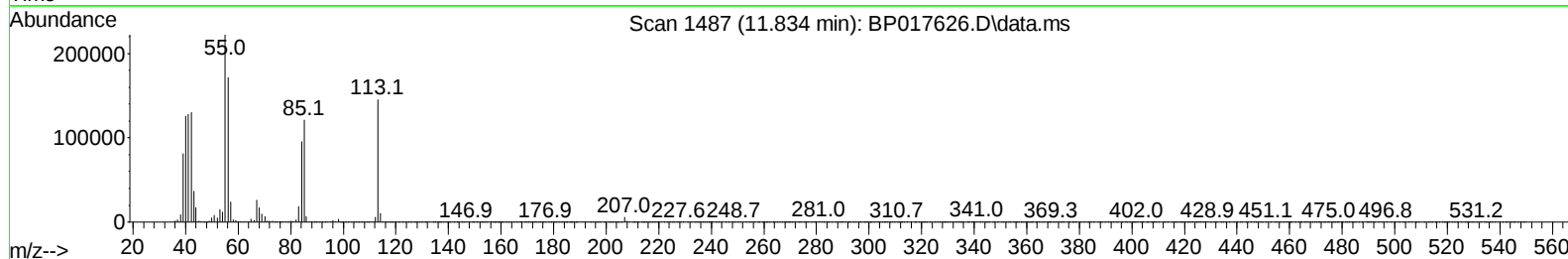
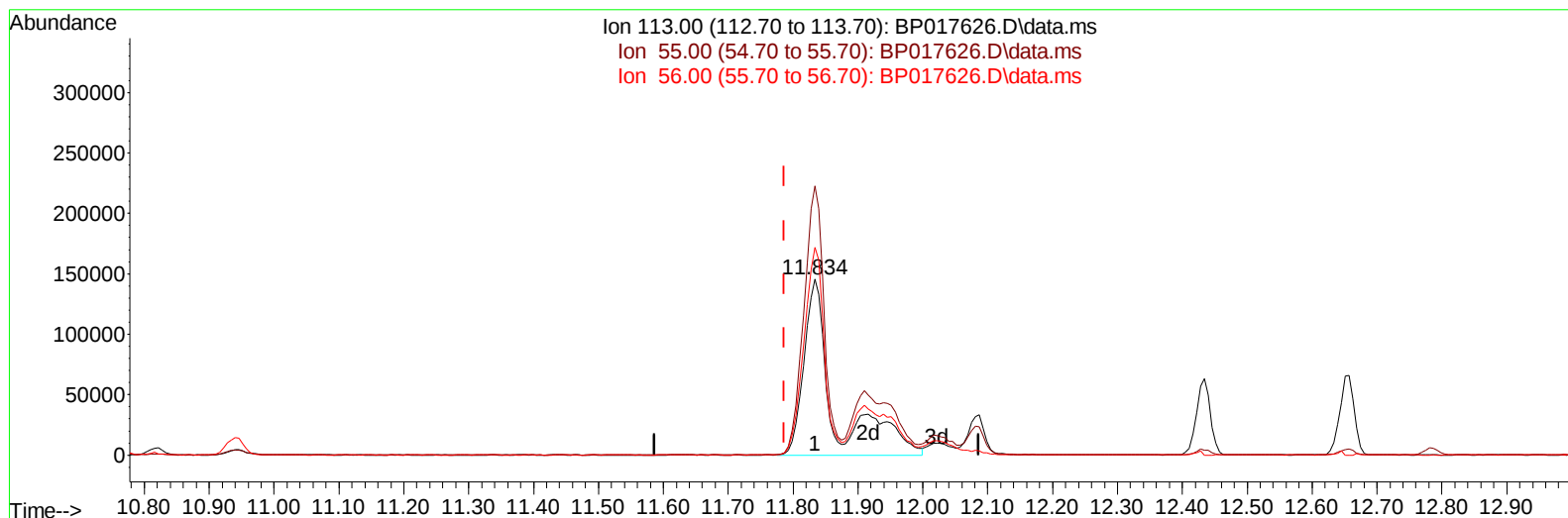
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
 Data File : BP017626.D  
 Acq On : 09 Oct 2023 20:46  
 Operator : MA/JU  
 Sample : SSTD16071  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160636

Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 10/10/2023  
 Supervised By :mohammad ahmed 10/11/2023



TIC: BP017626.D\data.ms

**(34) Caprolactam**

**11.834min (+ 0.047) 157.08 ng/ul m**

**response 469881**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 155.20 | 152.57 |
| 56.00  | 126.50 | 117.73 |
| 0.00   | 0.00   | 0.00   |

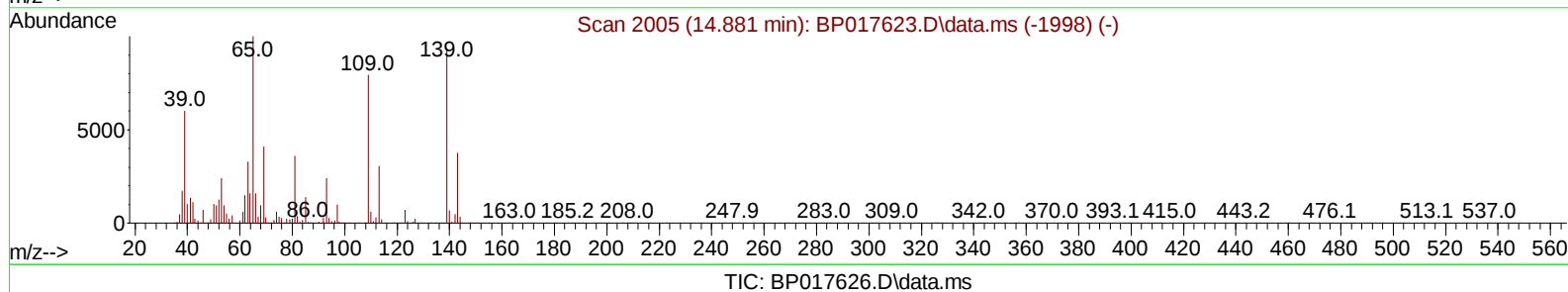
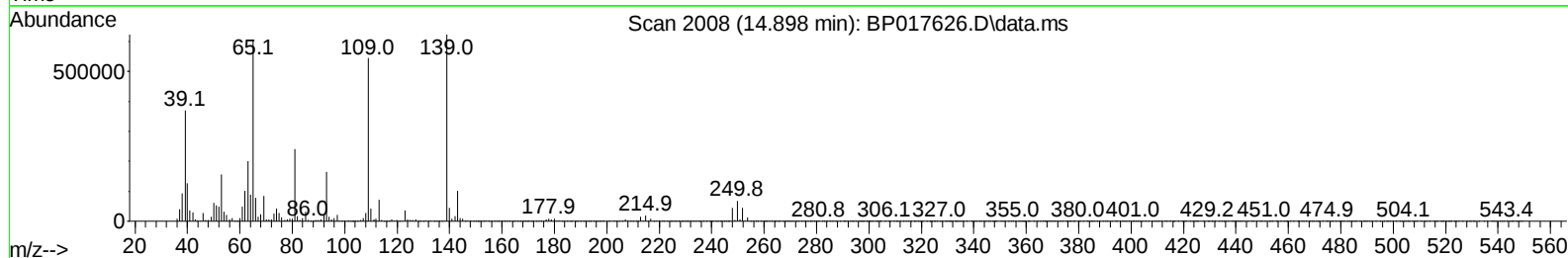
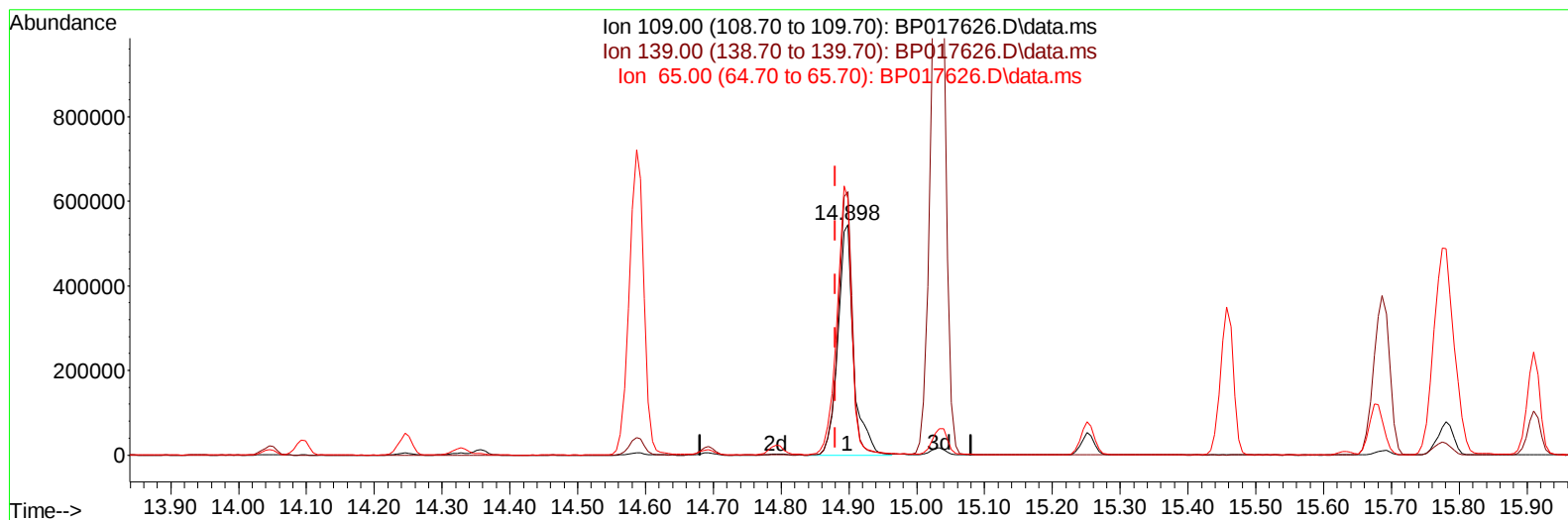
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
Data File : BP017626.D  
Acq On : 09 Oct 2023 20:46  
Operator : MA/JU  
Sample : SSTD16071  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD160636

Manual IntegrationsAPPROVED

Quant Time: Oct 10 04:29:42 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100923.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 10 03:51:30 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/10/2023  
Supervised By :mohammad ahmed 10/11/2023



(55) 4-Nitrophenol

14.898min (+ 0.018) 186.06 ng/u1

response 877883

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 109.00 | 100.00 | 100.00 |
| 139.00 | 118.30 | 114.52 |
| 65.00  | 127.20 | 111.43 |
| 0.00   | 0.00   | 0.00   |

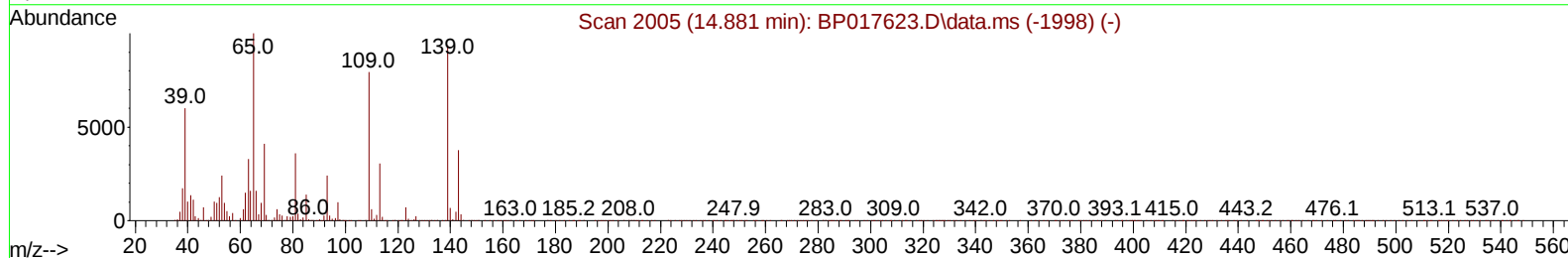
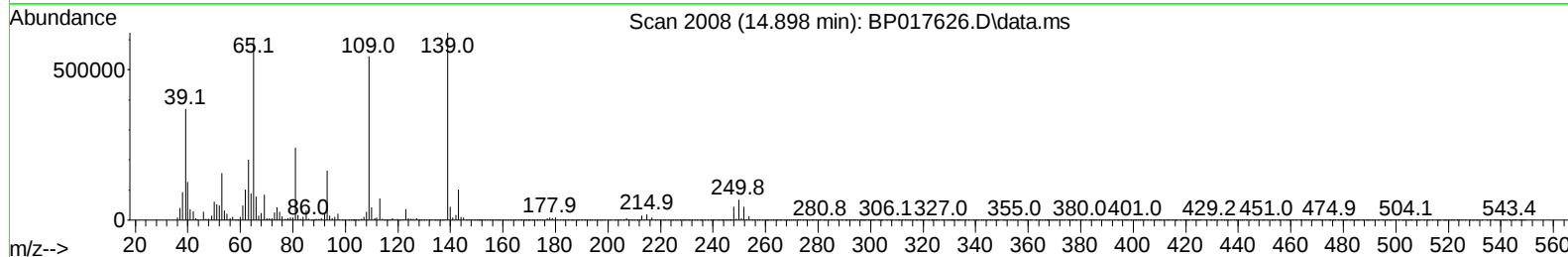
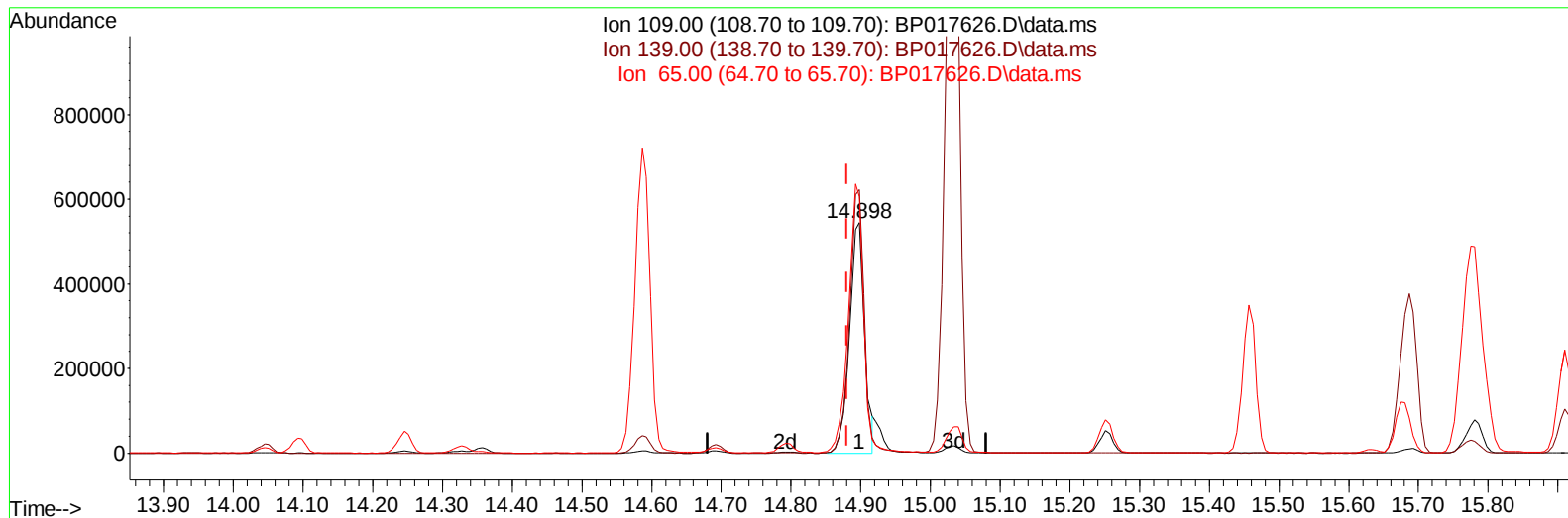
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
 Data File : BP017626.D  
 Acq On : 09 Oct 2023 20:46  
 Operator : MA/JU  
 Sample : SSTD16071  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160636

Manual IntegrationsAPPROVED

Quant Time: Oct 10 04:29:42 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 10 03:51:30 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/10/2023  
 Supervised By :mohammad ahmed 10/11/2023



TIC: BP017626.D\data.ms

**(55) 4-Nitrophenol**

**14.898min (+ 0.018) 172.33 ng/ul m**

**response 813131**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 109.00 | 100.00 | 100.00 |
| 139.00 | 118.30 | 114.52 |
| 65.00  | 127.20 | 111.43 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100923\  
 Data File : BP017626.D  
 Acq On : 09 Oct 2023 20:46  
 Operator : MA/JU  
 Sample : SSTD16071  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160636

Manual IntegrationsAPPROVED

Quant Time: Oct 10 04:30:17 2023  
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Reviewed By :Yogesh Patel 10/10/2023  
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| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.928  | 152  | 148627   | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 630412   | 20.000  | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.622 | 164  | 402949   | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.416 | 188  | 911970   | 20.000  | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.557 | 240  | 892953   | 20.000  | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.127 | 264  | 983049   | 20.000  | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 3.711  | 84   | 1584490  | 162.774 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.093  | 99   | 2004501  | 160.989 | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 1147066  | 150.585 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/ul  |          |
| 15) 4-Methylphenol-d8         | 8.663  | 113  | 1588037  | 159.662 | ng/ul  | 0.02     |
| 21) Nitrobenzene-d5           | 0.000  | 128  | 0d       | 0.000   | ng/ul  |          |
| 24) 2-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000   | ng/ul  |          |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165  | 0d       | 0.000   | ng/ul  |          |
| 31) 4-Chloroaniline-d4        | 10.940 | 131  | 2234382  | 154.873 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.000   | ng/ul  |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.000   | ng/ul  |          |
| 54) 4-Nitrophenol-d4          | 14.881 | 143  | 884749   | 169.116 | ng/ul  | 0.02     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/ul  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.781 | 200  | 968306   | 176.280 | ng/ul  | 0.01     |
| 73) Anthracene-d10            | 0.000  | 188  | 0d       | 0.000   | ng/ul  |          |
| 81) Pyrene-d10                | 0.000  | 212  | 0d       | 0.000   | ng/ul  |          |
| 92) Benzo(a)pyrene-d12        | 0.000  | 264  | 0d       | 0.000   | ng/ul  |          |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 5) Pyridine                   | 3.734  | 79   | 1608481  | 157.701 | ng/ul  | 97       |
| 6) Benzaldehyde               | 7.087  | 77   | 712112   | 113.574 | ng/ul  | 99       |
| 8) Phenol                     | 7.122  | 94   | 1977178  | 158.354 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.369  | 93   | 1525963  | 150.962 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.381  | 108  | 1501682  | 161.100 | ng/ul  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.446  | 45   | 1944369m | 146.272 | ng/ul  |          |
| 16) Acetophenone              | 8.793  | 105  | 2426175  | 154.897 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.734  | 108  | 1596865  | 157.854 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.963 | 127  | 2106412  | 152.905 | ng/ul  | 98       |
| 34) Caprolactam               | 11.834 | 113  | 469881m  | 157.083 | ng/ul  |          |
| 40) Hexachlorocyclopentadiene | 12.728 | 237  | 1337881  | 183.921 | ng/ul  | 100      |
| 51) 3-Nitroaniline            | 14.587 | 138  | 813342   | 147.270 | ng/ul  | 93       |
| 53) 2,4-Dinitrophenol         | 14.792 | 184  | 659273   | 201.502 | ng/ul  | 95       |
| 55) 4-Nitrophenol             | 14.898 | 109  | 813131m  | 172.333 | ng/ul  |          |
| 63) 4-Nitroaniline            | 15.775 | 138  | 731266   | 141.498 | ng/ul  | 88       |
| 66) 4,6-Dinitro-2-methylph... | 15.798 | 198  | 996897   | 169.094 | ng/ul# | 91       |
| 70) Atrazine                  | 16.881 | 200  | 1539360  | 156.168 | ng/ul  | 99       |
| 71) Pentachlorophenol         | 17.039 | 266  | 1203649  | 175.578 | ng/ul  | 96       |
| 77) Carbazole                 | 17.845 | 167  | 6642412  | 147.890 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.480 | 252  | 3054008  | 155.637 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.398 | 149  | 9698208  | 160.212 | ng/ul  | 100      |
| -----                         |        |      |          |         |        |          |

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 Sample : SSTD16071  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160636

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/10/2023  
 Supervised By :mohammad ahmed 10/11/2023

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 Response via : Initial Calibration

| Compound                                                                   | R.T. | Q | Ion | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|---|-----|----------|------|-------|----------|
| -----                                                                      |      |   |     |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |   |     |          |      |       |          |

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