

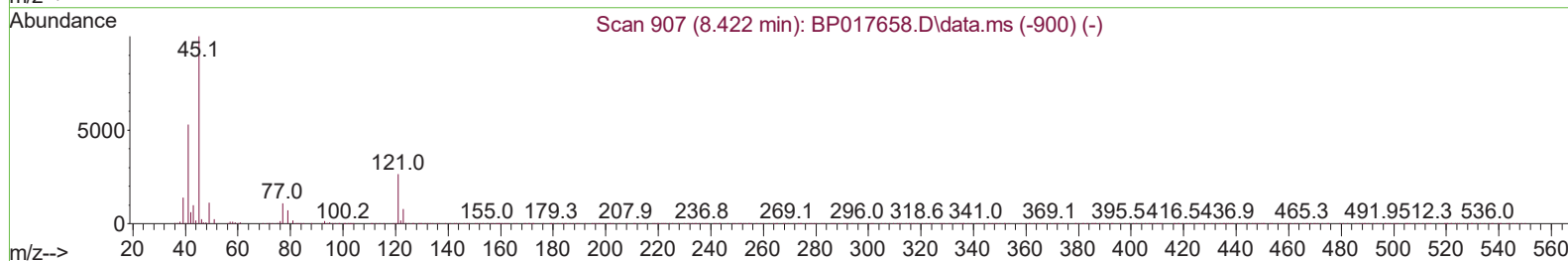
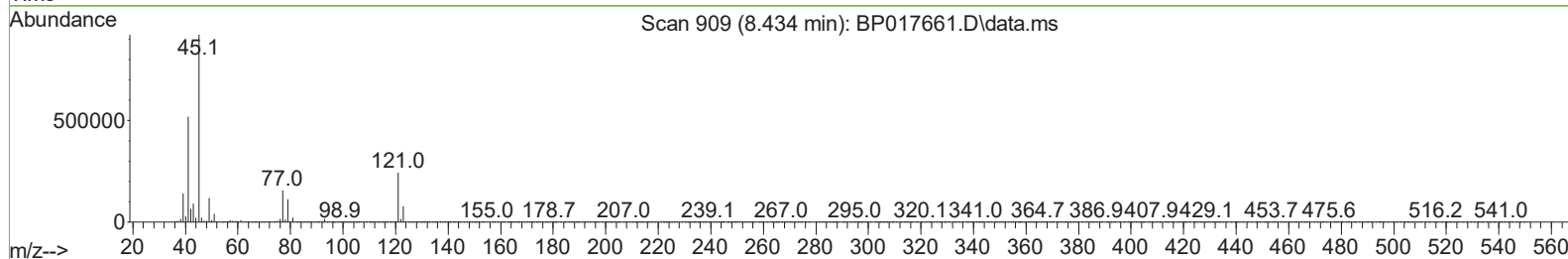
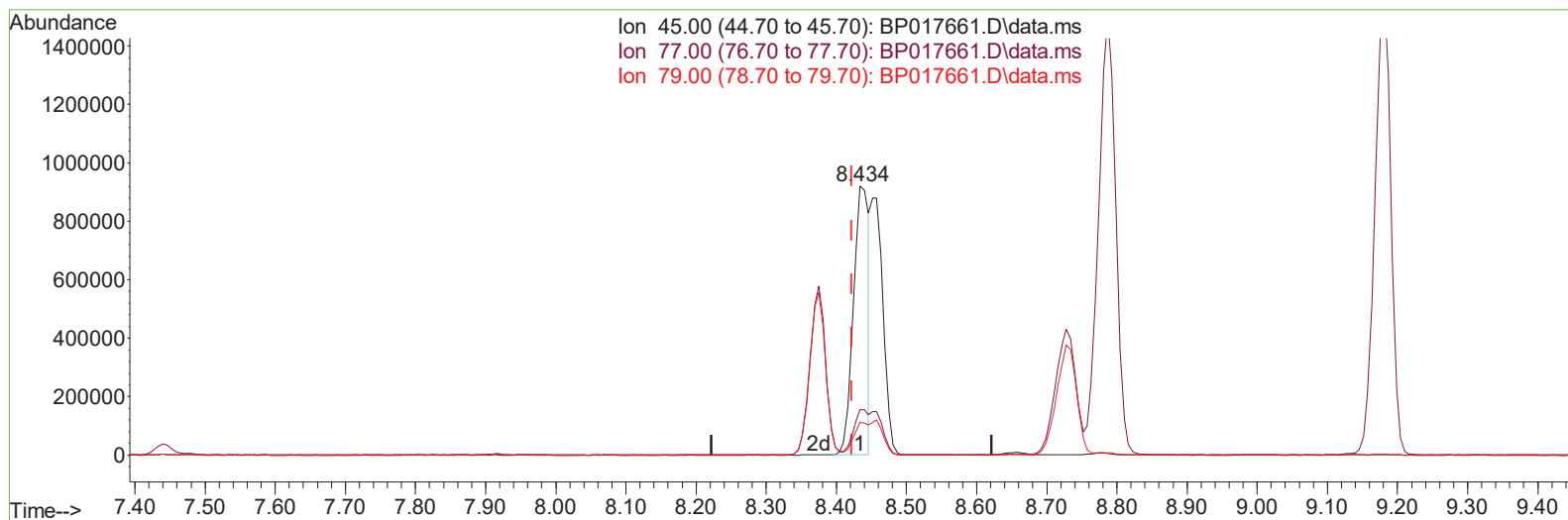
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
 Data File : BP017661.D  
 Acq On : 12 Oct 2023 15:37  
 Operator : MA/JU  
 Sample : SSTD16033  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160643

Manual IntegrationsAPPROVED

Quant Time: Oct 12 16:08:53 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 12 15:43:48 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/13/2023  
 Supervised By :mohammad ahmed 10/13/2023



TIC: BP017661.D\data.ms

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

8.434min (+ 0.012) 94.68 ng/u1

response 1408034

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.30  | 16.88  |
| 79.00 | 11.90  | 12.15  |
| 0.00  | 0.00   | 0.00   |

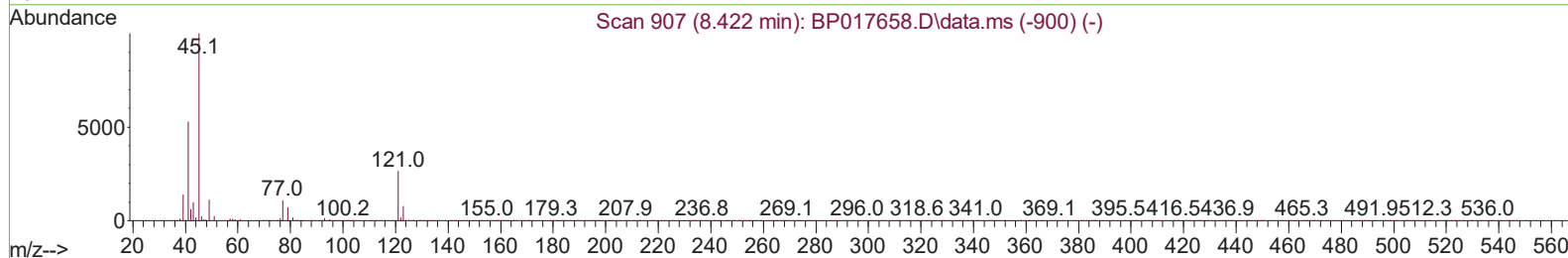
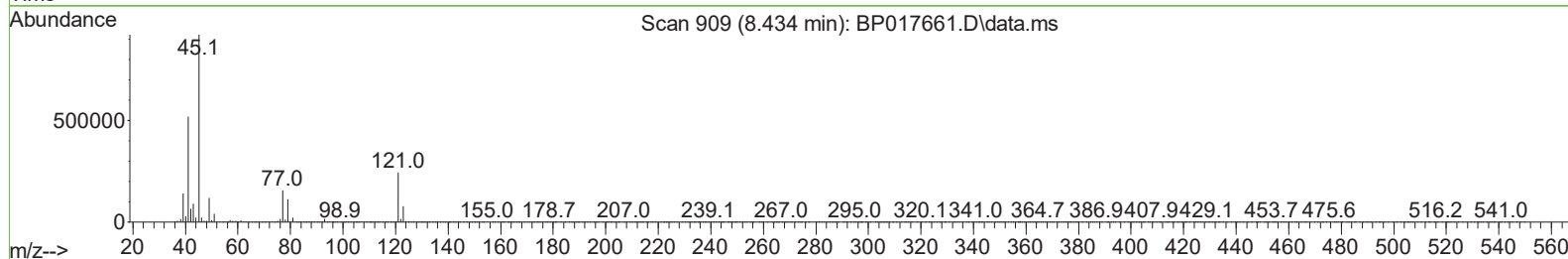
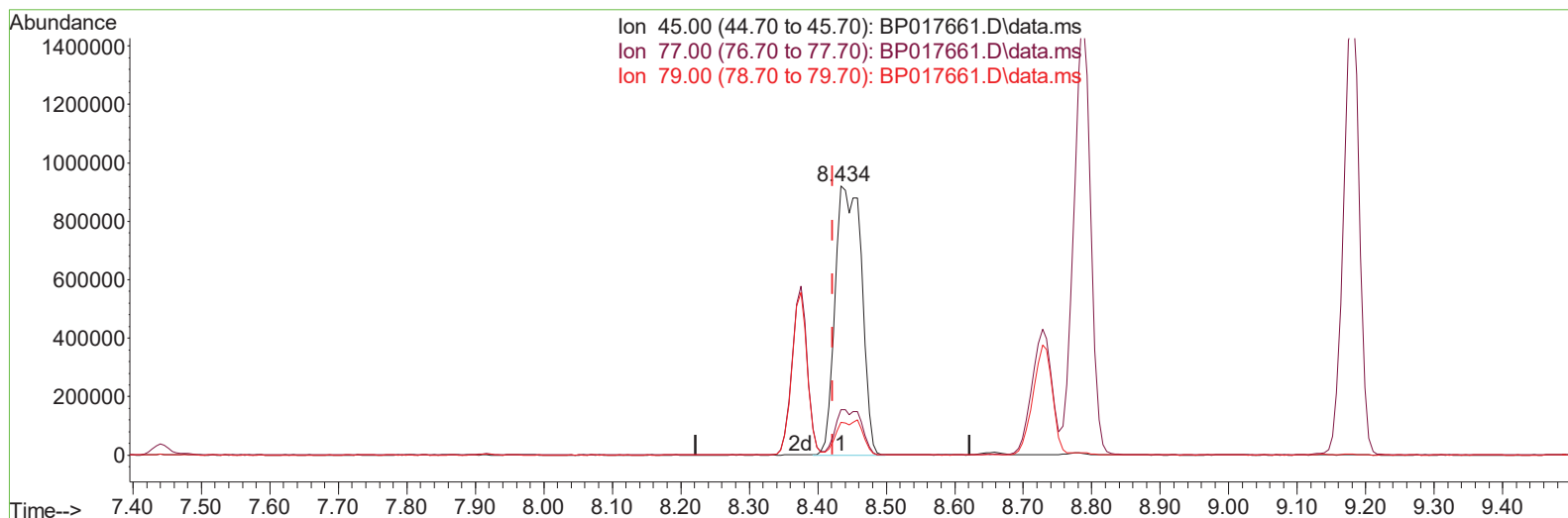
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
 Data File : BP017661.D  
 Acq On : 12 Oct 2023 15:37  
 Operator : MA/JU  
 Sample : SSTD16033  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160643

Manual IntegrationsAPPROVED

Quant Time: Oct 12 16:08:53 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 12 15:43:48 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/13/2023  
 Supervised By :mohammad ahmed 10/13/2023



TIC: BP017661.D\data.ms

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

8.434min (+ 0.012) 166.36 ng/ul m

response 2474184

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.30  | 16.88  |
| 79.00 | 11.90  | 12.15  |
| 0.00  | 0.00   | 0.00   |

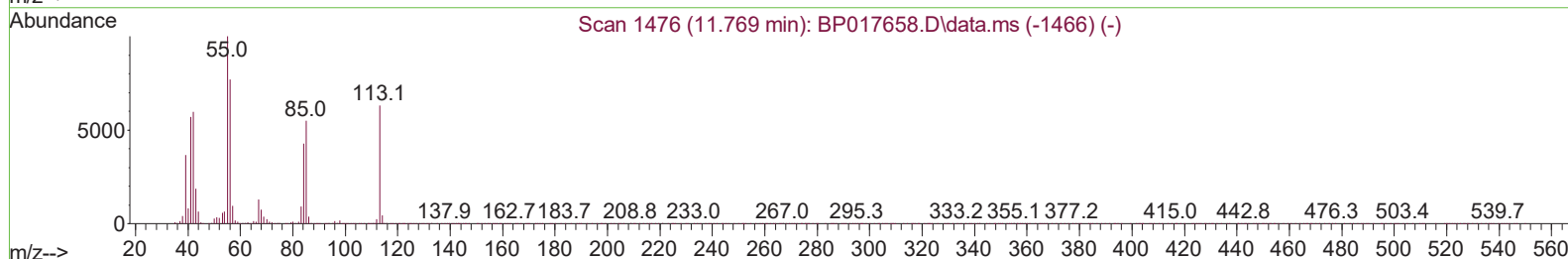
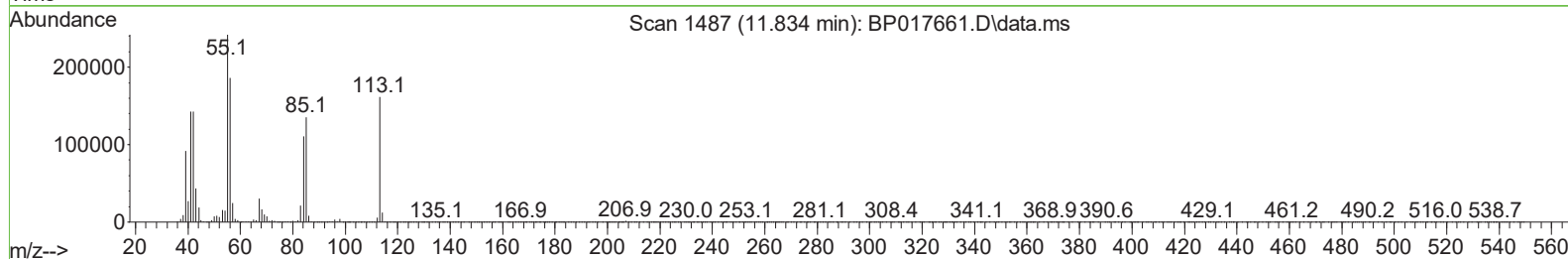
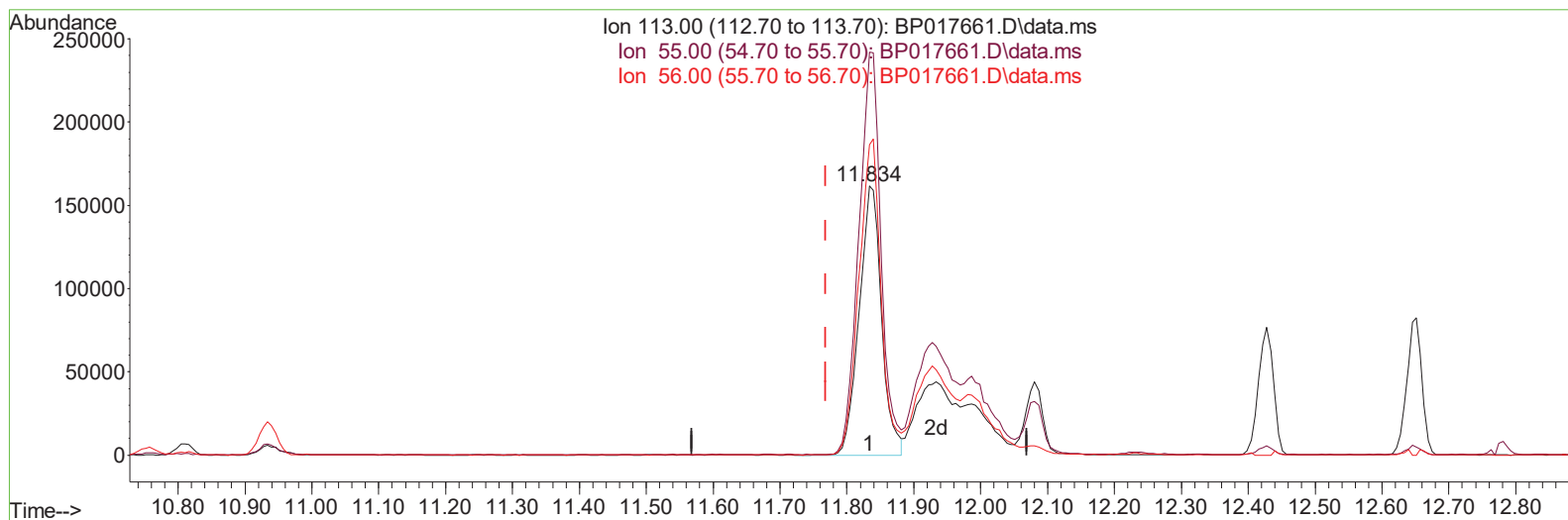
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
Data File : BP017661.D  
Acq On : 12 Oct 2023 15:37  
Operator : MA/JU  
Sample : SSTD16033  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD160643

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/13/2023  
Supervised By :mohammad ahmed 10/13/2023

Quant Time: Oct 12 16:08:53 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Oct 12 15:43:48 2023  
Response via : Initial Calibration



TIC: BP017661.D\data.ms

## (34) Caprolactam

11.834min (+ 0.065) 113.23 ng/ul

response 374690

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 158.60 | 149.76 |
| 56.00  | 122.40 | 115.29 |
| 0.00   | 0.00   | 0.00   |

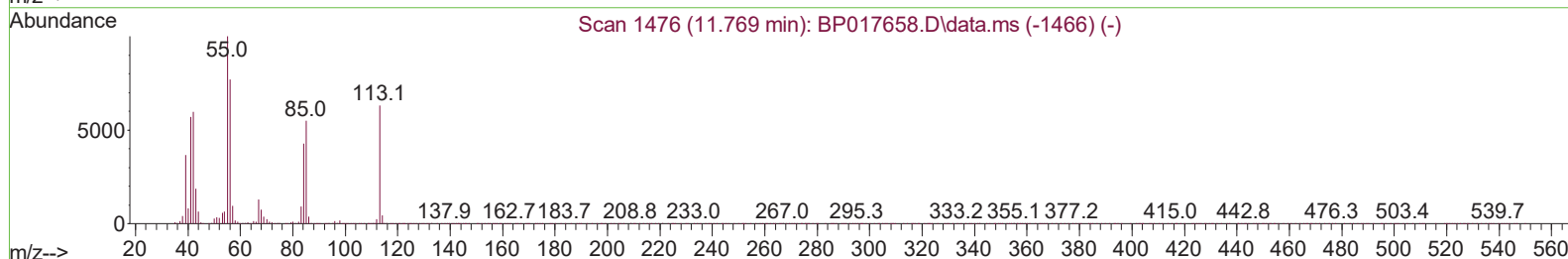
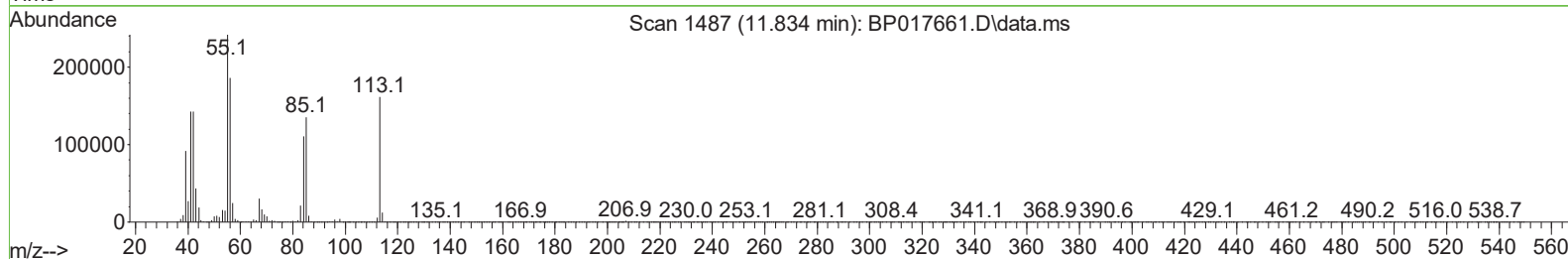
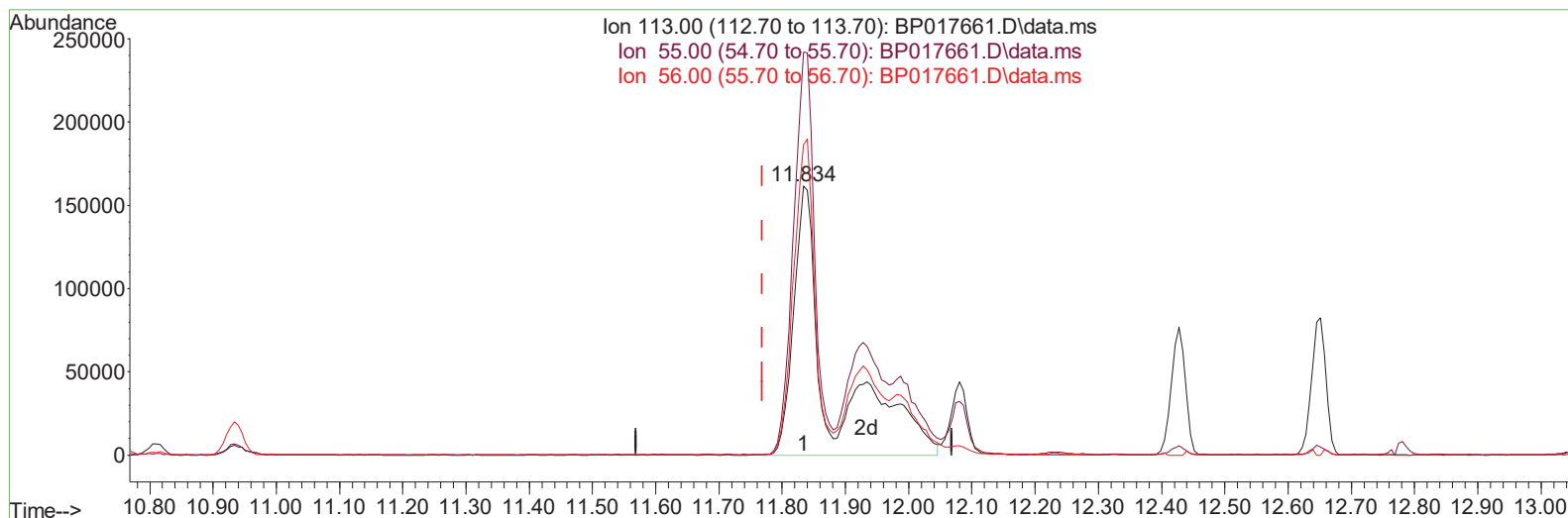
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
 Data File : BP017661.D  
 Acq On : 12 Oct 2023 15:37  
 Operator : MA/JU  
 Sample : SSTD16033  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160643

Manual IntegrationsAPPROVED

Quant Time: Oct 12 16:08:53 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 12 15:43:48 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/13/2023  
 Supervised By :mohammad ahmed 10/13/2023



TIC: BP017661.D\data.ms

**(34) Caprolactam**

11.834min (+ 0.065) 192.53 ng/ul m

response 637093

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 158.60 | 149.76 |
| 56.00  | 122.40 | 115.29 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
 Data File : BP017661.D  
 Acq On : 12 Oct 2023 15:37  
 Operator : MA/JU  
 Sample : SSTD16033  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160643

Manual IntegrationsAPPROVED

Quant Time: Oct 12 16:11:48 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 12 15:43:48 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/13/2023  
 Supervised By :mohammad ahmed 10/13/2023

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 166197   | 20.000  | ng/ul  | 0.01     |
| 20) Naphthalene-d8            | 10.757 | 136  | 686827   | 20.000  | ng/ul  | 0.02     |
| 38) Acenaphthene-d10          | 14.622 | 164  | 454773   | 20.000  | ng/ul  | 0.02     |
| 64) Phenanthrene-d10          | 17.416 | 188  | 1049345  | 20.000  | ng/ul  | 0.02     |
| 79) Chrysene-d12              | 21.563 | 240  | 1094735  | 20.000  | ng/ul  | 0.02     |
| 88) Perylene-d12              | 24.145 | 264  | 1227020  | 20.000  | ng/ul  | 0.03     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 3.699  | 84   | 2078833  | 181.467 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.081  | 99   | 2688079  | 192.239 | ng/ul  | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 1510195  | 174.024 | ng/ul  | 0.02     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/ul  |          |
| 15) 4-Methylphenol-d8         | 8.657  | 113  | 2117845  | 187.448 | ng/ul  | 0.03     |
| 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        | 10.934 | 131  | 2945705  | 184.137 | ng/ul  | 0.02     |
| 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          | 14.887 | 143  | 1299286  | 229.956 | ng/ul  | 0.04     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/ul  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.786 | 200  | 1341623  | 208.162 | ng/ul  | 0.03     |
| 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                   | 3.722  | 79   | 2117637  | 179.458 | ng/ul  | 96       |
| 6) Benzaldehyde               | 7.075  | 77   | 802521   | 114.532 | ng/ul  | 99       |
| 8) Phenol                     | 7.116  | 94   | 2630264  | 185.463 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.363  | 93   | 1987131  | 174.913 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.375  | 108  | 1964626  | 186.141 | ng/ul  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.434  | 45   | 2474184m | 166.365 | ng/ul  |          |
| 16) Acetophenone              | 8.787  | 105  | 3137326  | 176.360 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.728  | 108  | 2103941  | 182.671 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.963 | 127  | 2796087  | 180.832 | ng/ul  | 98       |
| 34) Caprolactam               | 11.834 | 113  | 637093m  | 192.526 | ng/ul  |          |
| 40) Hexachlorocyclopentadiene | 12.722 | 237  | 1473303  | 192.232 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.587 | 138  | 1085545  | 172.872 | ng/ul  | 95       |
| 53) 2,4-Dinitrophenol         | 14.798 | 184  | 970364   | 221.173 | ng/ul  | 92       |
| 55) 4-Nitrophenol             | 14.904 | 109  | 1215652  | 237.797 | ng/ul  | 95       |
| 63) 4-Nitroaniline            | 15.786 | 138  | 1017878  | 170.106 | ng/ul# | 85       |
| 66) 4,6-Dinitro-2-methylph... | 15.804 | 198  | 1375260  | 202.219 | ng/ul# | 92       |
| 70) Atrazine                  | 16.886 | 200  | 1995014  | 175.058 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.045 | 266  | 1652819  | 199.068 | ng/ul  | 99       |
| 77) Carbazole                 | 17.851 | 167  | 8940920  | 172.628 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.492 | 252  | 4344878  | 182.719 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.410 | 149  | 12798900 | 174.420 | ng/ul  | 100      |
| -----                         |        |      |          |         |        |          |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
 Data File : BP017661.D  
 Acq On : 12 Oct 2023 15:37  
 Operator : MA/JU  
 Sample : SSTD16033  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160643

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/13/2023  
 Supervised By :mohammad ahmed 10/13/2023

Quant Time: Oct 12 16:11:48 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 12 15:43:48 2023  
 Response via : Initial Calibration

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101223\  
Data File : BP017661.D  
Acq On : 12 Oct 2023 15:37  
Operator : MA/JU  
Sample : SSTD16033  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD160643

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/13/2023  
Supervised By :mohammad ahmed 10/13/2023

Quant Time: Oct 12 16:11:48 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP101223.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Oct 12 15:43:48 2023  
Response via : Initial Calibration

