

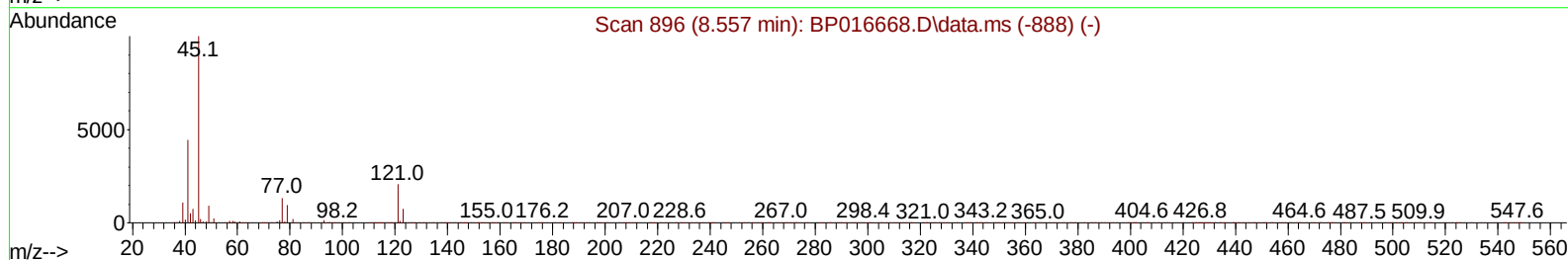
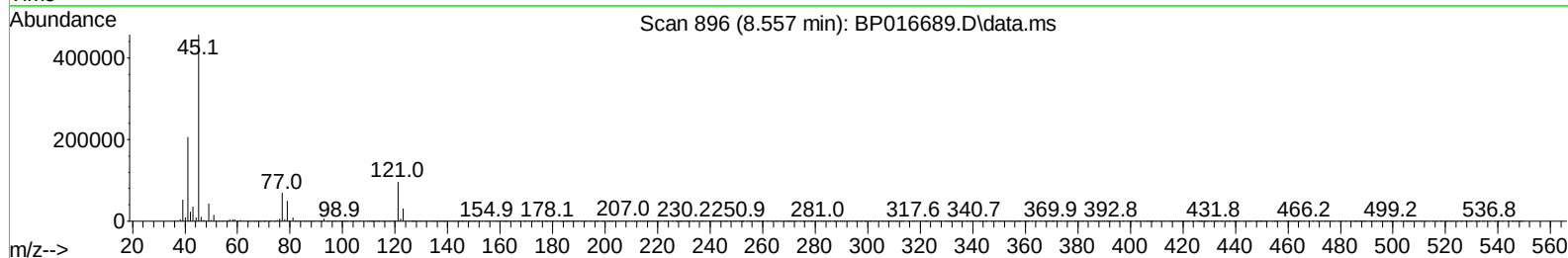
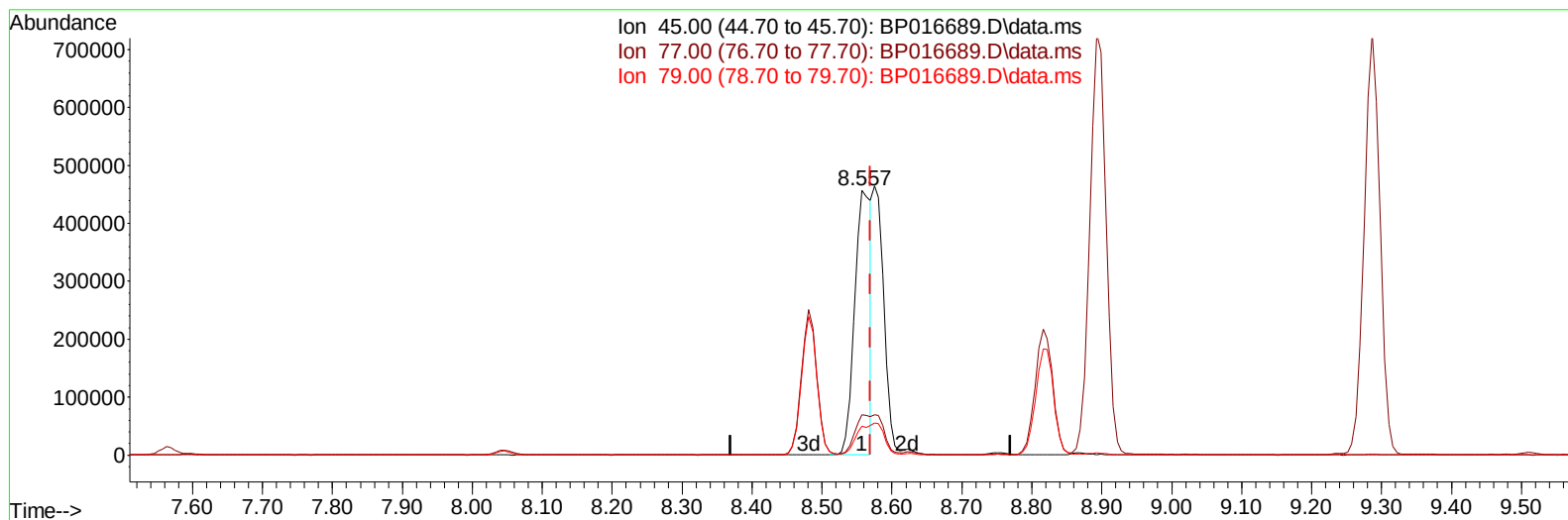
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082123\  
 Data File : BP016689.D  
 Acq On : 21 Aug 2023 23:38  
 Operator : MA/JU  
 Sample : PB154871BS  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS871

Manual IntegrationsAPPROVED

Quant Time: Aug 22 00:35:37 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 17 01:38:34 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/22/2023  
 Supervised By :mohammad ahmed 08/22/2023



TIC: BP016689.D\data.ms

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

8.557min (-0.012) 20.34 ng/uL

response 735563

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 15.10  | 15.22  |
| 79.00 | 10.30  | 10.87  |
| 0.00  | 0.00   | 0.00   |

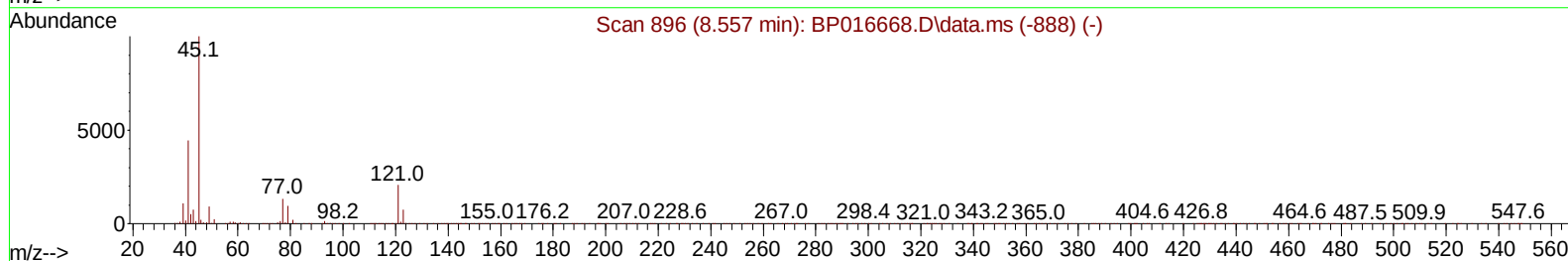
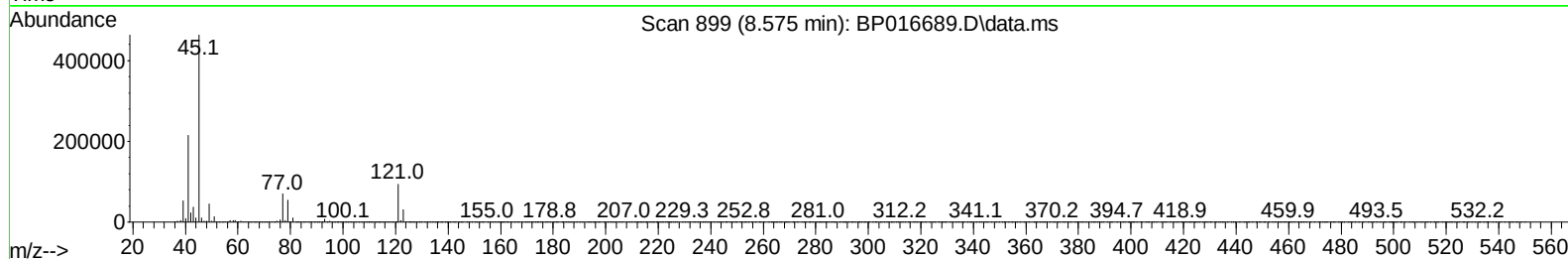
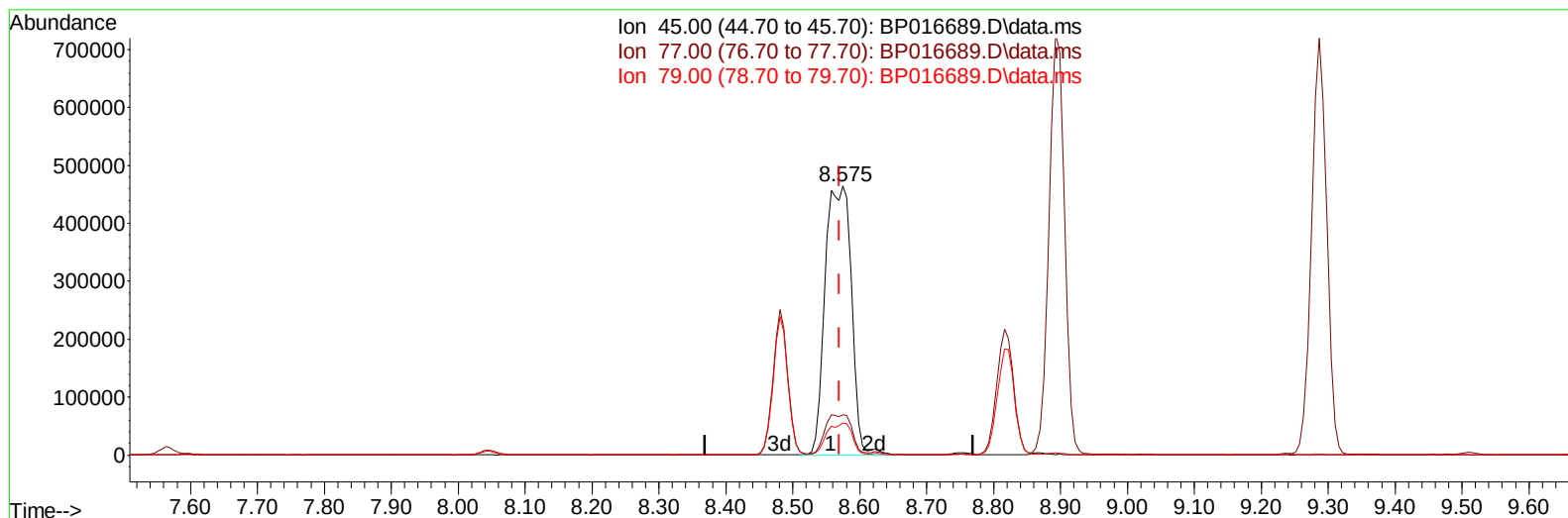
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 Data File : BP016689.D  
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 Operator : MA/JU  
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Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 08/22/2023  
 Supervised By :mohammad ahmed 08/22/2023



TIC: BP016689.D\data.ms

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

8.575min (+ 0.006) 34.91 ng/ul m

response 1262870

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 15.10  | 15.10  |
| 79.00 | 10.30  | 11.93  |
| 0.00  | 0.00   | 0.00   |

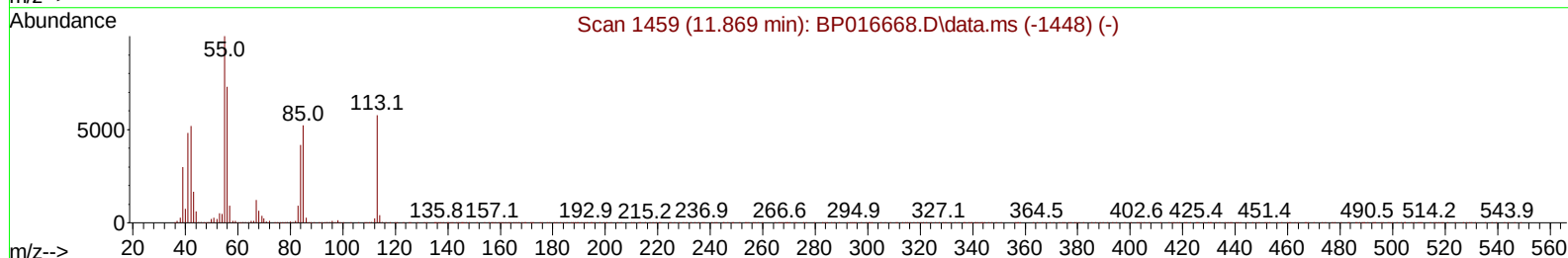
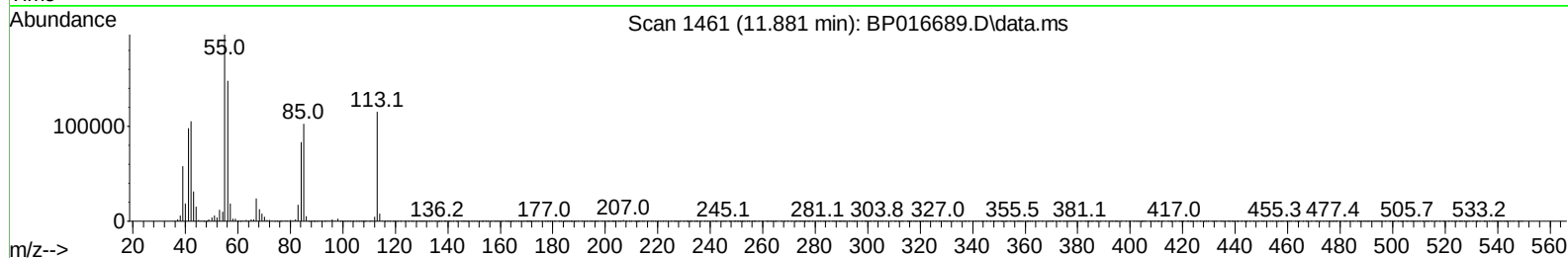
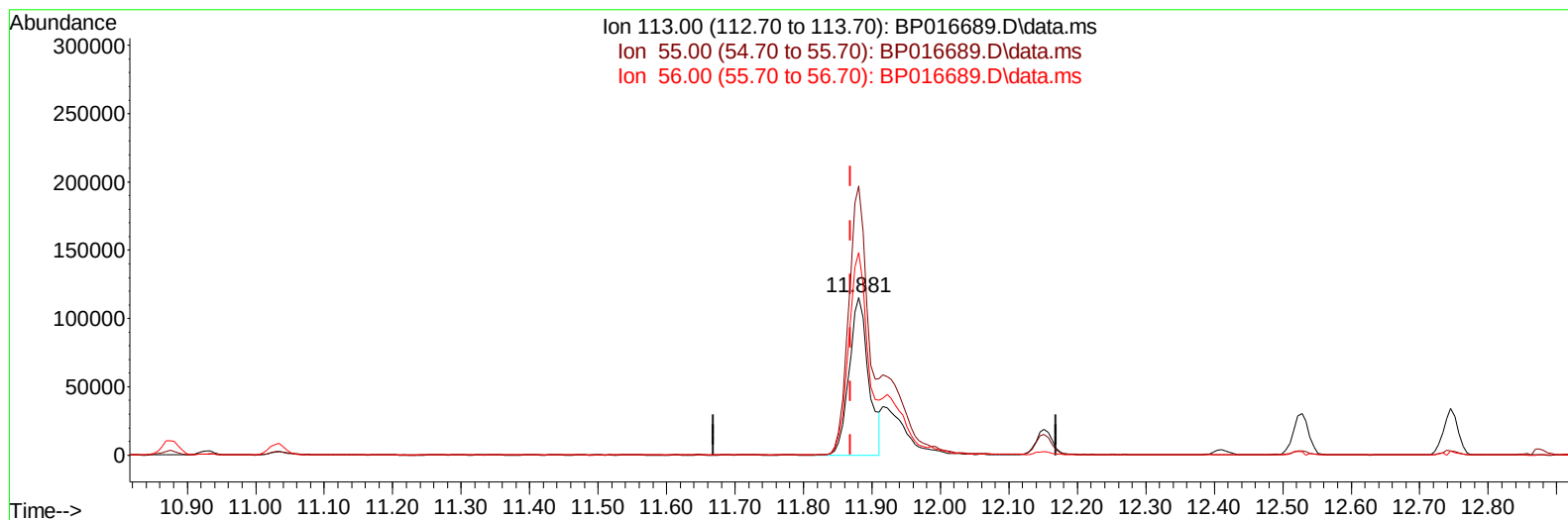
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082123\  
 Data File : BP016689.D  
 Acq On : 21 Aug 2023 23:38  
 Operator : MA/JU  
 Sample : PB154871BS  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
 SLCS871

Manual IntegrationsAPPROVED

Quant Time: Aug 22 00:35:37 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
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Reviewed By :Yogesh Patel 08/22/2023  
 Supervised By :mohammad ahmed 08/22/2023



TIC: BP016689.D\data.ms

**(34) Caprolactam**

**11.881min (+ 0.012) 29.18 ng/ul**

**response 228646**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 168.50 | 170.44 |
| 56.00  | 126.90 | 128.51 |
| 0.00   | 0.00   | 0.00   |

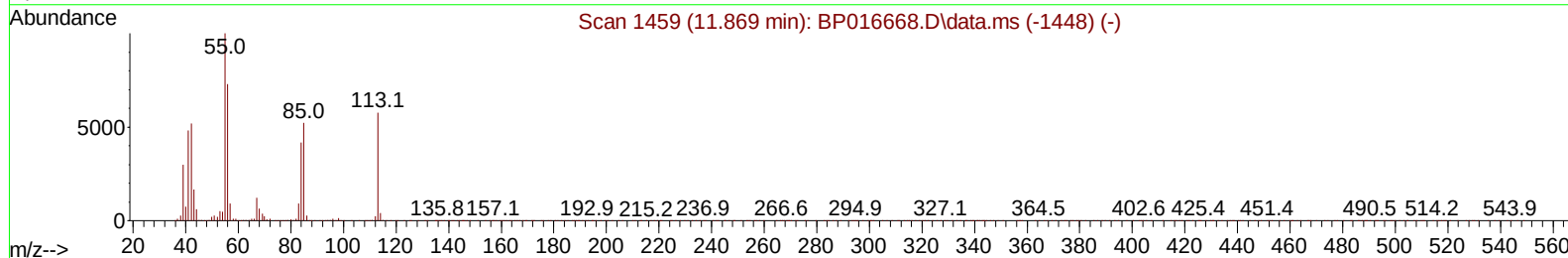
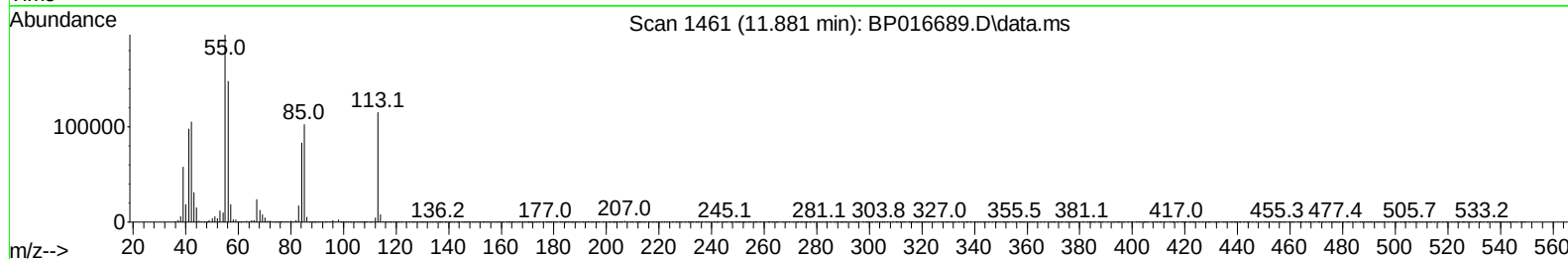
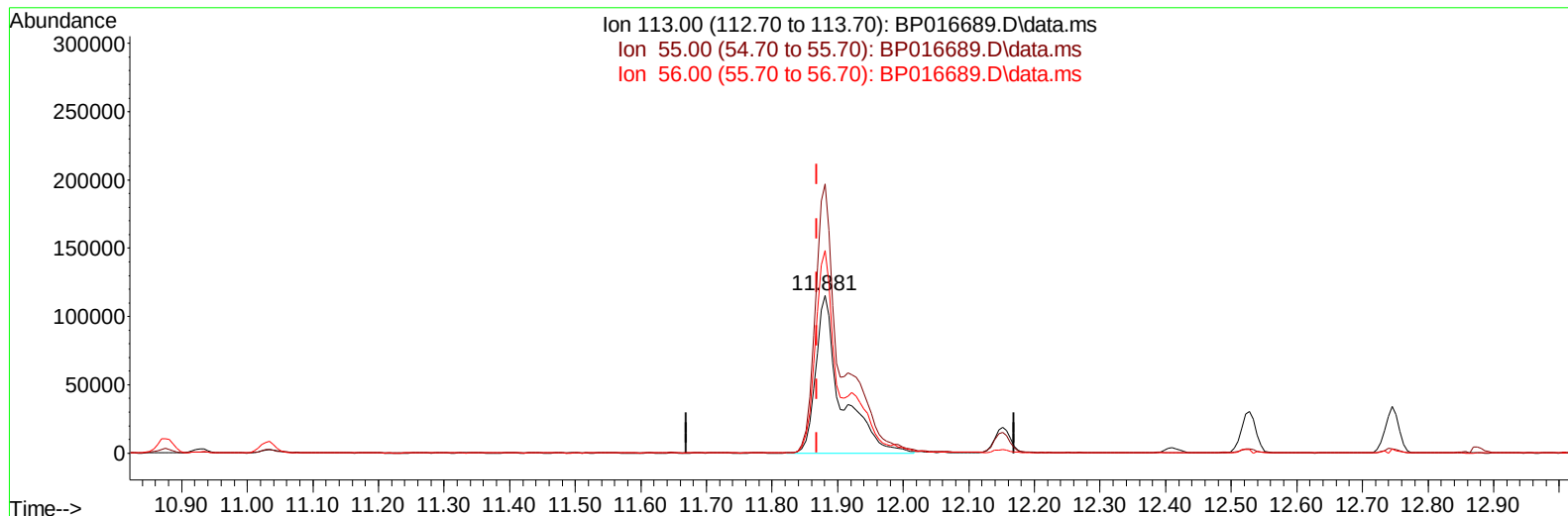
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082123\  
Data File : BP016689.D  
Acq On : 21 Aug 2023 23:38  
Operator : MA/JU  
Sample : PB154871BS  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS871

Manual IntegrationsAPPROVED

Quant Time: Aug 22 00:57:56 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 17 01:38:34 2023  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/22/2023  
Supervised By :mohammad ahmed 08/22/2023



TIC: BP016689.D\data.ms

(34) Caprolactam

11.881min (+ 0.012) 40.18 ng/ul m

response 314790

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 168.50 | 170.44 |
| 56.00  | 126.90 | 128.51 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082123\  
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 Acq On : 21 Aug 2023 23:38  
 Operator : MA/JU  
 Sample : PB154871BS  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS871

Manual IntegrationsAPPROVED

Quant Time: Aug 22 00:58:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 17 01:38:34 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/22/2023  
 Supervised By :mohammad ahmed 08/22/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.046  | 152  | 326056   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.875 | 136  | 1488305  | 20.000 | ng/ul  | -0.01    |
| 38) Acenaphthene-d10          | 14.698 | 164  | 911992   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.463 | 188  | 2006147  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.557 | 240  | 1849228  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.133 | 264  | 1867060  | 20.000 | ng/ul  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 70693    | 8.239  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 872694   | 33.428 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.187  | 99   | 1102501  | 36.677 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 675634   | 34.125 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.563  | 132  | 763141   | 35.346 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.751  | 113  | 852888   | 35.683 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.246  | 128  | 392979   | 37.526 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.957  | 143  | 420436   | 42.271 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 773219   | 39.503 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.034 | 131  | 1129831  | 33.642 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.110 | 166  | 2352337  | 35.661 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 2708021  | 36.372 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.904 | 143  | 498015   | 37.061 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.692 | 176  | 1999212  | 35.551 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.816 | 200  | 378905   | 39.701 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.563 | 188  | 3130350  | 36.074 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 3711185  | 38.301 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.968 | 264  | 3411523  | 38.128 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.428  | 88   | 134091   | 14.212 | ng/uL  | 99       |
| 5) Pyridine                   | 3.846  | 79   | 876587   | 32.363 | ng/ul  | 99       |
| 6) Benzaldehyde               | 7.199  | 77   | 698163   | 42.133 | ng/ul  | 98       |
| 8) Phenol                     | 7.216  | 94   | 1197291  | 37.266 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.475  | 93   | 929441   | 35.813 | ng/ul  | 97       |
| 12) 2-Chlorophenol            | 7.599  | 128  | 845374   | 37.212 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.481  | 108  | 879593   | 37.691 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.575  | 45   | 1262870m | 34.914 | ng/ul  |          |
| 16) Acetophenone              | 8.893  | 105  | 1432214  | 37.106 | ng/ul  | 98       |
| 17) N-Nitroso-di-n-propyla... | 8.869  | 70   | 802555   | 38.770 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.816  | 108  | 977502   | 38.111 | ng/ul  | 96       |
| 19) Hexachloroethane          | 9.116  | 117  | 336869   | 35.493 | ng/ul  | 96       |
| 22) Nitrobenzene              | 9.287  | 77   | 1140193  | 38.893 | ng/ul  | 99       |
| 23) Isophorone                | 9.804  | 82   | 2227402  | 40.227 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.993  | 139  | 491005   | 43.182 | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 10.028 | 107  | 957538   | 35.721 | ng/ul  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.287 | 93   | 1336160  | 37.645 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.510 | 162  | 812103   | 40.324 | ng/ul  | 99       |
| 30) Naphthalene               | 10.928 | 128  | 2822320  | 36.295 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 11.057 | 127  | 1170066  | 35.175 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 11.163 | 225  | 447456   | 36.858 | ng/ul  | 99       |
| 34) Caprolactam               | 11.881 | 113  | 314790m  | 40.178 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.151 | 107  | 1029442  | 41.819 | ng/ul  | 99       |

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 Sample : PB154871BS  
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Instrument :  
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 ClientSampleId :  
 SLCS871

Manual IntegrationsAPPROVED

Quant Time: Aug 22 00:58:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/22/2023  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.528 | 142  | 1897478  | 36.429 | ng/ul  | 100      |
| 37) 1-Methylnaphthalene       | 12.745 | 142  | 1908239  | 36.384 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.875 | 216  | 913213   | 37.437 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.828 | 237  | 471400   | 31.761 | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol     | 13.128 | 196  | 645784   | 42.538 | ng/ul  | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.192 | 196  | 721787   | 43.862 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.534 | 154  | 2542158  | 37.501 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.581 | 162  | 1958472  | 37.132 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.810 | 65   | 720703   | 45.192 | ng/ul  | 95       |
| 47) Dimethylphthalate         | 14.157 | 163  | 2593796  | 38.459 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene        | 14.298 | 165  | 535277   | 45.821 | ng/ul  | 100      |
| 50) Acenaphthylene            | 14.428 | 152  | 3160777  | 36.048 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.634 | 138  | 598705   | 45.708 | ng/ul  | 98       |
| 52) Acenaphthene              | 14.763 | 153  | 2165859  | 36.900 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.834 | 184  | 270467   | 41.589 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.916 | 109  | 440830   | 40.353 | ng/ul  | 98       |
| 56) Dibenzofuran              | 15.098 | 168  | 2979299  | 37.203 | ng/ul  | 100      |
| 57) 2,4-Dinitrotoluene        | 15.075 | 165  | 784834   | 44.864 | ng/ul  | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.310 | 232  | 593845   | 43.071 | ng/ul  | 98       |
| 59) Diethylphthalate          | 15.510 | 149  | 2648474  | 38.407 | ng/ul  | 99       |
| 61) Fluorene                  | 15.751 | 166  | 2465071  | 37.540 | ng/ul  | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.739 | 204  | 1150646  | 37.329 | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 15.804 | 138  | 614792   | 49.046 | ng/ul  | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.828 | 198  | 423147   | 39.987 | ng/ul# | 98       |
| 67) N-Nitrosodiphenylamine    | 15.963 | 169  | 2119755  | 38.981 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.639 | 248  | 706145   | 39.423 | ng/ul  | 99       |
| 69) Hexachlorobenzene         | 16.722 | 284  | 805025   | 38.464 | ng/ul  | 97       |
| 70) Atrazine                  | 16.916 | 200  | 705231   | 36.047 | ng/ul  | 99       |
| 71) Pentachlorophenol         | 17.086 | 266  | 511661   | 39.707 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.510 | 178  | 4003176  | 37.876 | ng/ul  | 100      |
| 74) Anthracene                | 17.598 | 178  | 3993252  | 37.615 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.487 | 216  | 902796   | 35.411 | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 14.986 | 250  | 887172   | 38.783 | ng/uL  | 99       |
| 77) Carbazole                 | 17.880 | 167  | 3891728  | 39.779 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.386 | 149  | 4737586  | 41.832 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.463 | 202  | 4796696  | 41.477 | ng/ul  | 98       |
| 82) Pyrene                    | 19.822 | 202  | 4923243  | 40.315 | ng/ul  | 100      |
| 83) Butylbenzylphthalate      | 20.680 | 149  | 2320175  | 47.163 | ng/ul  | 94       |
| 84) 3,3'-Dichlorobenzidine    | 21.480 | 252  | 1642664  | 40.793 | ng/ul  | 100      |
| 85) Benzo(a)anthracene        | 21.539 | 228  | 4769330  | 38.376 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.416 | 149  | 3377863  | 47.311 | ng/ul  | 98       |
| 87) Chrysene                  | 21.598 | 228  | 4497730  | 38.675 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.398 | 149  | 5814659  | 49.911 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.333 | 252  | 4587864  | 41.269 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.386 | 252  | 4538218  | 41.149 | ng/ul  | 100      |
| 93) Benzo(a)pyrene            | 24.021 | 252  | 3911732  | 37.301 | ng/ul  | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 26.880 | 276  | 4523062  | 36.544 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.909 | 278  | 3748975  | 36.397 | ng/ul  | 100      |
| 96) Benzo(g,h,i)perylene      | 27.739 | 276  | 3577335  | 36.316 | ng/ul  | 97       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS871

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 08/22/2023

Supervised By :mohammad ahmed 08/22/2023

