

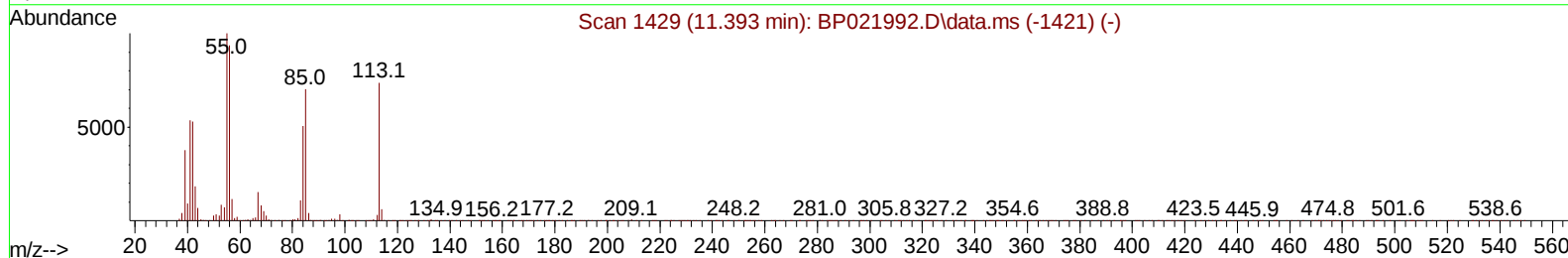
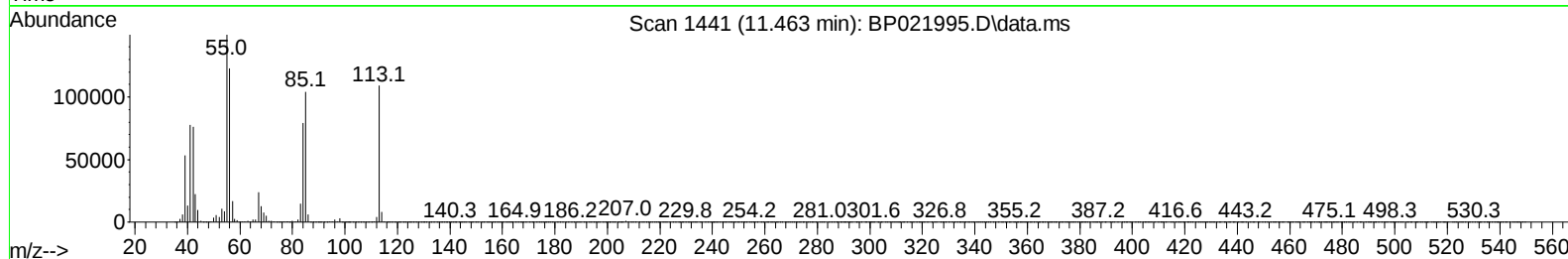
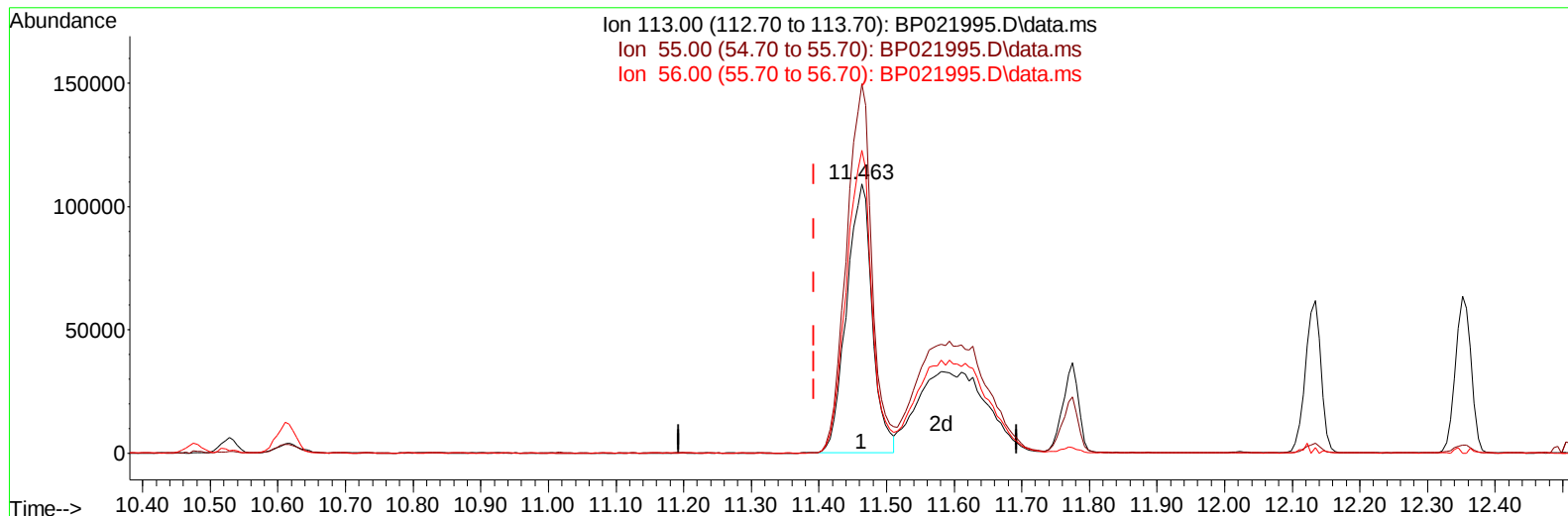
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092424\  
Data File : BP021995.D  
Acq On : 24 Sep 2024 14:18  
Operator : RC/JU  
Sample : SSTD16042  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD160620

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/25/2024  
Supervised By :mohammad ahmed 09/28/2024

Quant Time: Sep 24 15:18:44 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092424.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Sep 24 15:01:38 2024  
Response via : Initial Calibration



TIC: BP021995.D\data.ms

## (34) Caprolactam

11.463min (+ 0.071) 94.23 ng/ul

response 287415

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 136.10 | 137.36 |
| 56.00  | 127.30 | 112.61 |
| 0.00   | 0.00   | 0.00   |

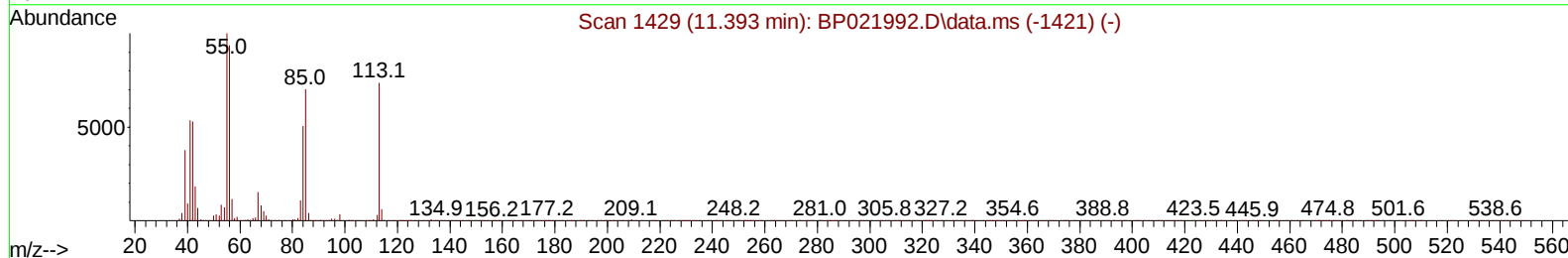
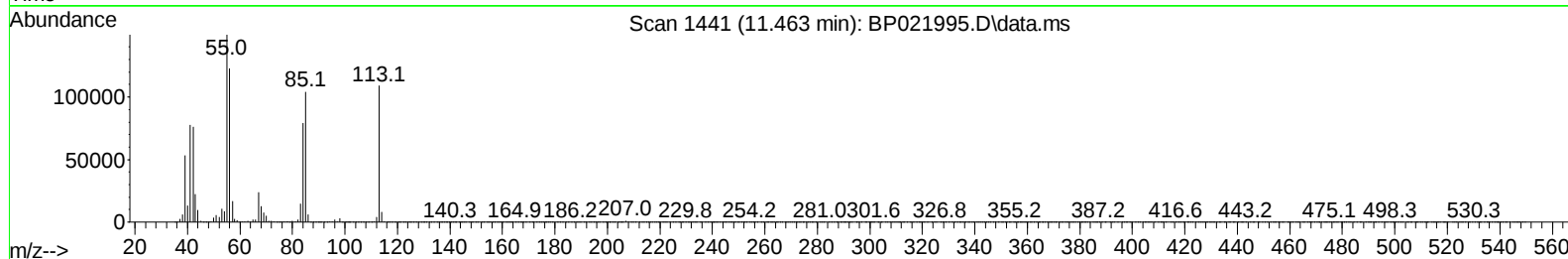
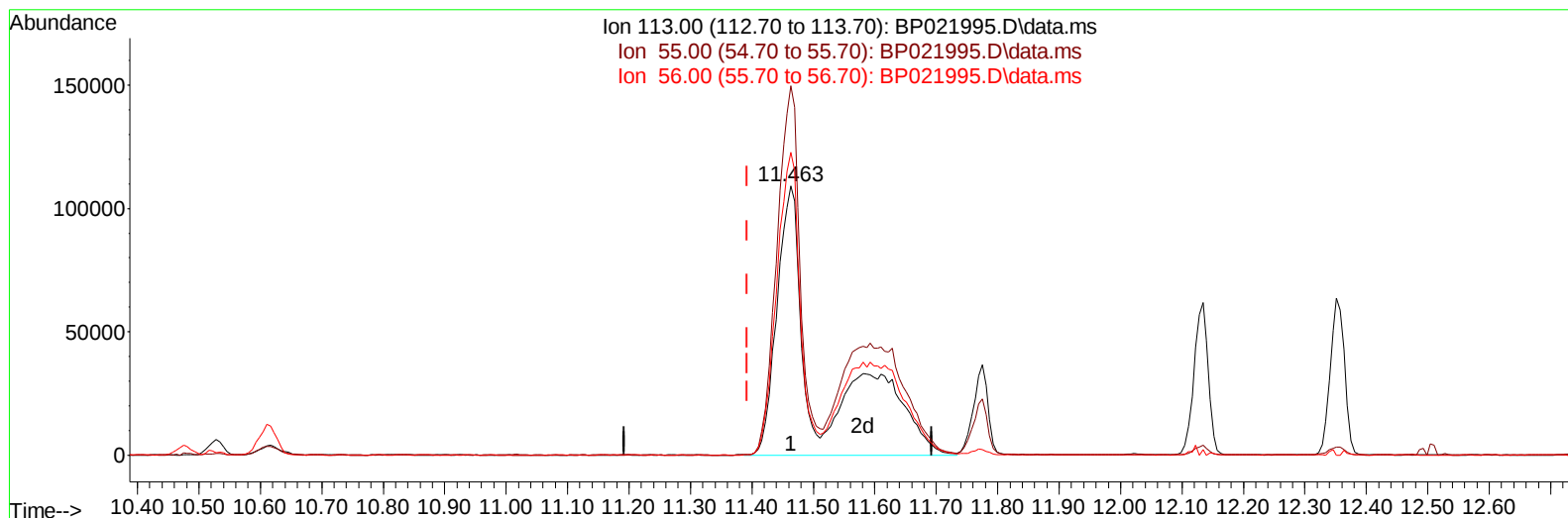
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TIC: BP021995.D\data.ms

(34) Caprolactam

11.463min (+ 0.071) 172.79 ng/u1 m

response 527044

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 136.10 | 137.36 |
| 56.00  | 127.30 | 112.61 |
| 0.00   | 0.00   | 0.00   |

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Manual IntegrationsAPPROVED

Quant Time: Sep 24 15:21:53 2024  
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 Response via : Initial Calibration

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| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.711  | 152  | 141883   | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.475 | 136  | 549904   | 20.000  | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.322 | 164  | 415215   | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.122 | 188  | 1001677  | 20.000  | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.569 | 240  | 1081994  | 20.000  | ng/ul  | 0.03     |
| 88) Perylene-d12              | 24.845 | 264  | 1220716  | 20.000  | ng/ul  | 0.02     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 3.634  | 84   | 1583709  | 152.956 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.905  | 99   | 1980452  | 170.638 | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.058  | 67   | 1119026  | 156.581 | ng/ul  | 0.01     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/ul  |          |
| 15) 4-Methylphenol-d8         | 8.428  | 113  | 1540769  | 167.534 | 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        | 10.616 | 131  | 1906901  | 157.697 | ng/ul  | 0.01     |
| 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          | 14.557 | 143  | 919091   | 167.408 | ng/ul  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/ul  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.463 | 200  | 1145803  | 184.246 | ng/ul  | 0.02     |
| 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                   | 3.658  | 79   | 1595248  | 151.650 | ng/ul  | 96       |
| 6) Benzaldehyde               | 6.869  | 77   | 741705   | 112.970 | ng/ul  | 99       |
| 8) Phenol                     | 6.934  | 94   | 2037355  | 168.209 | ng/ul  | 96       |
| 10) Bis(2-Chloroethyl)ether   | 7.152  | 93   | 1477561  | 159.386 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.152  | 108  | 1486445  | 169.216 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.228  | 45   | 1325516  | 154.134 | ng/ul  | 98       |
| 16) Acetophenone              | 8.534  | 105  | 2433610  | 159.082 | ng/ul  | 97       |
| 18) 4-Methylphenol            | 8.499  | 108  | 1591180  | 166.106 | ng/ul  | 96       |
| 32) 4-Chloroaniline           | 10.640 | 127  | 1873199  | 156.411 | ng/ul  | 99       |
| 34) Caprolactam               | 11.463 | 113  | 527044m  | 172.790 | ng/ul  |          |
| 40) Hexachlorocyclopentadiene | 12.487 | 237  | 1598462  | 165.074 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.234 | 138  | 943208   | 167.632 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.440 | 184  | 834257   | 198.392 | ng/ul  | 97       |
| 55) 4-Nitrophenol             | 14.575 | 109  | 985055   | 158.593 | ng/ul  | 96       |
| 63) 4-Nitroaniline            | 15.428 | 138  | 821114   | 143.384 | ng/ul  | 96       |
| 66) 4,6-Dinitro-2-methylph... | 15.481 | 198  | 1216219  | 179.634 | ng/ul# | 95       |
| 70) Atrazine                  | 16.598 | 200  | 1701508  | 144.575 | ng/ul  | 99       |
| 71) Pentachlorophenol         | 16.769 | 266  | 1725198  | 177.034 | ng/ul  | 99       |
| 77) Carbazole                 | 17.545 | 167  | 7368884  | 156.436 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.469 | 252  | 3773456  | 152.317 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.739 | 149  | 10064312 | 166.894 | ng/ul  | 100      |
| -----                         |        |      |          |         |        |          |

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

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