

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM072319\

Data File : BM021611.D

Acq On : 23 Jul 2019 12:26

Operator : HP/JU

Sample : SSTD16034

Misc :

ALS Vial : 8 Sample Multiplier: 1

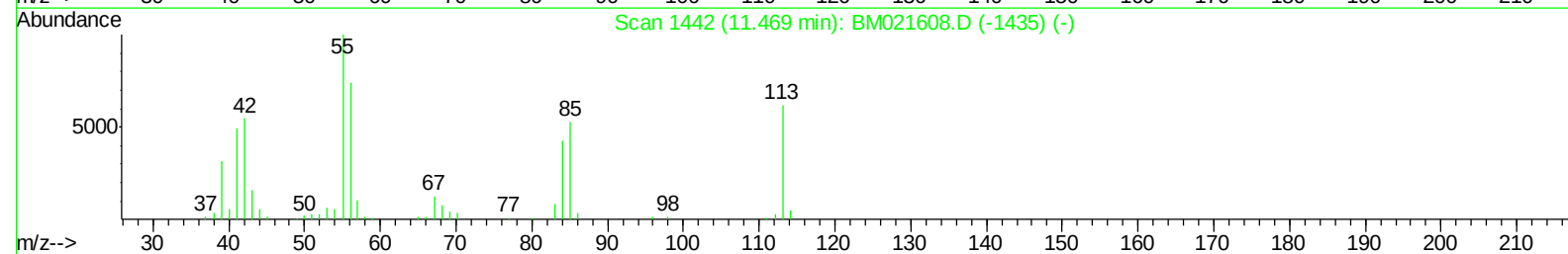
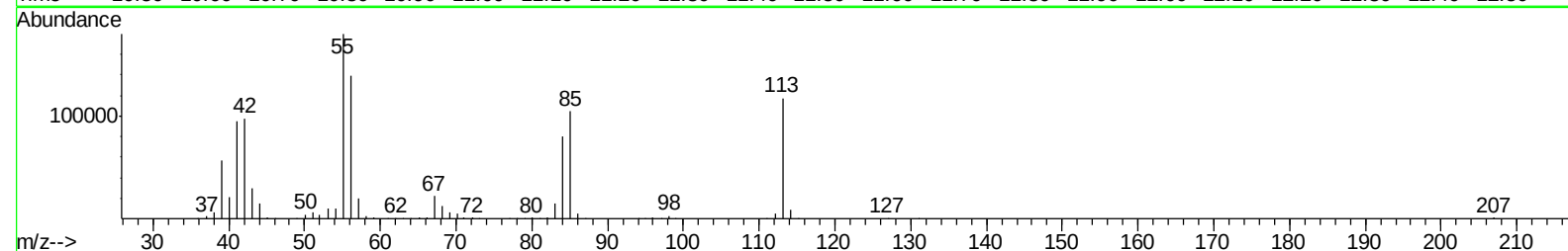
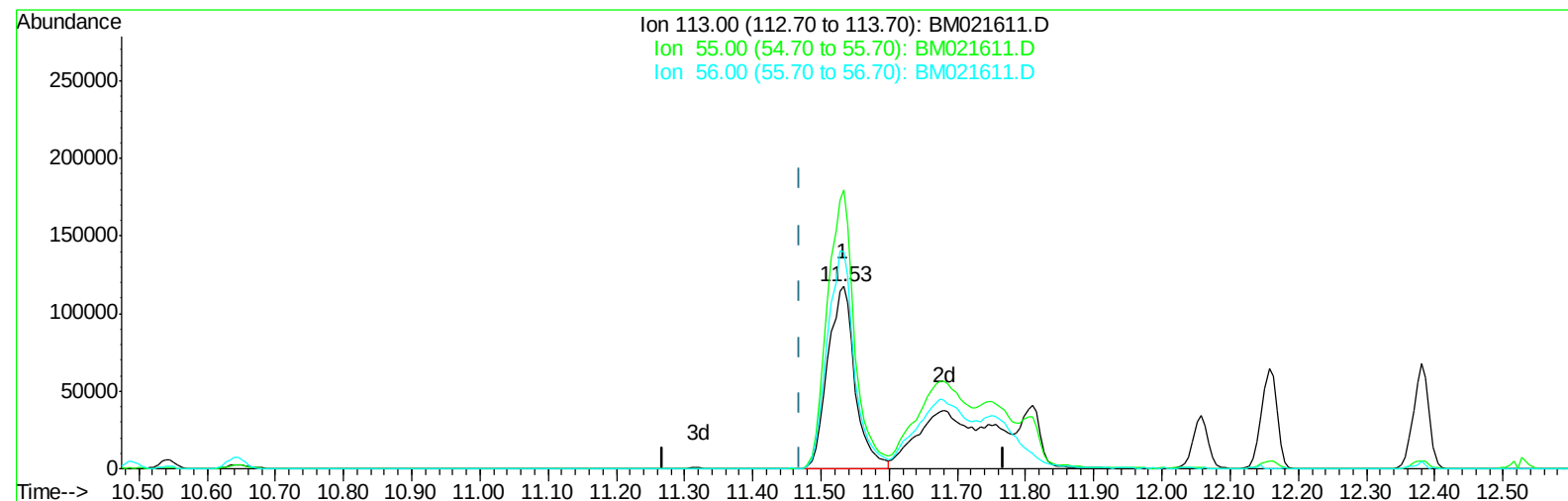
Quant Time: Jul 23 12:58:59 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM072319MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Jul 23 11:52:52 2019

Response via : Initial Calibration



TIC: BM021611.D

(32) Caprolactam

11.533min (+0.064) 107.70ng/ul

response 327262

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 162.00 | 153.16 |
| 56.00  | 119.40 | 118.93 |
| 0.00   | 0.00   | 0.00   |

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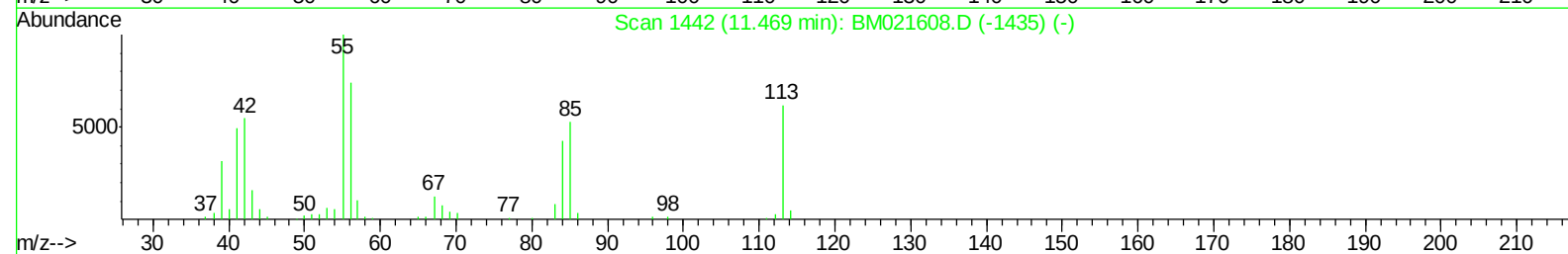
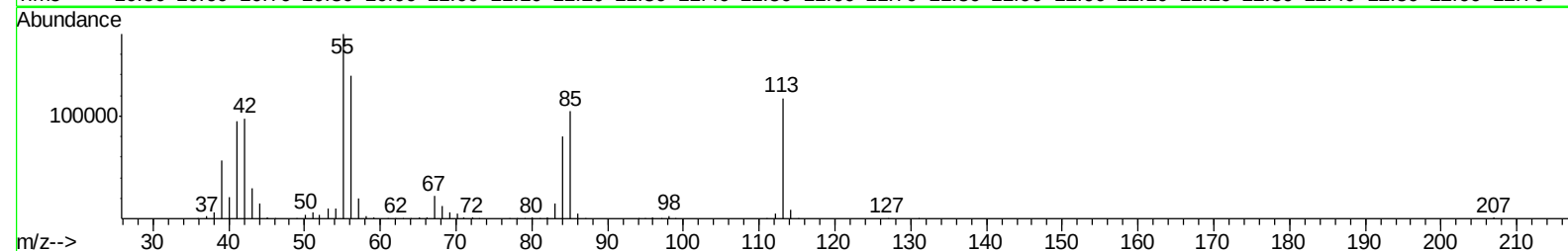
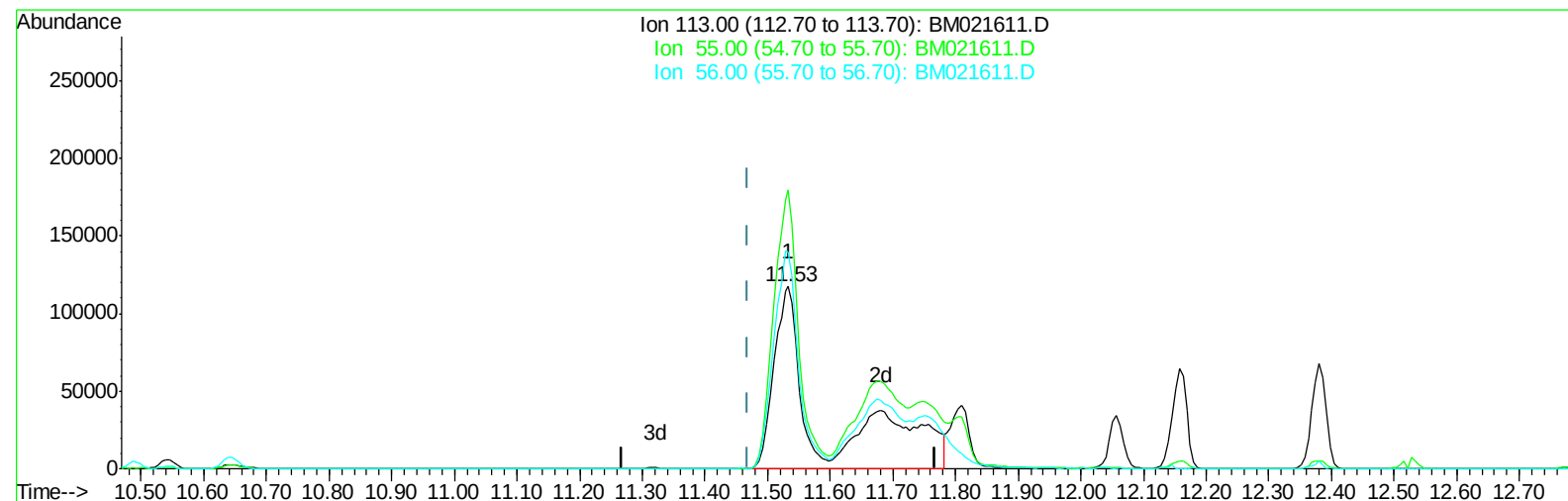
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TIC: BM021611.D

(32) Caprolactam

11.533min (+0.064) 198.93ng/ul m

response 604502

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 162.00 | 153.16 |
| 56.00  | 119.40 | 118.93 |
| 0.00   | 0.00   | 0.00   |

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Quant Time: Jul 23 13:05:38 2019  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 23 11:52:52 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.70  | 152  | 140744   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.49 | 136  | 685248   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.35 | 164  | 460809   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.11 | 188  | 1155845  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.30 | 240  | 1359510  | 20.00 | ng/ul | 0.01     |
| 85) Perylene-d12          | 23.55 | 264  | 1560729  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |        |       |      |
|--------------------------------|-------|-----|---------|--------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d      | 0.00   | ng/uL |      |
| 5) Phenol-d5                   | 6.89  | 99  | 2149022 | 196.37 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.06  | 67  | 1182192 | 182.19 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 0.00  | 132 | 0d      | 0.00   | ng/ul |      |
| 13) 4-Methylphenol-d8          | 8.43  | 113 | 1829366 | 209.87 | ng/ul | 0.02 |
| 19) Nitrobenzene-d5            | 0.00  | 128 | 0d      | 0.00   | ng/ul |      |
| 22) 2-Nitrophenol-d4           | 0.00  | 143 | 0d      | 0.00   | ng/ul |      |
| 26) 2,4-Dichlorophenol-d3      | 0.00  | 165 | 0d      | 0.00   | ng/ul |      |
| 29) 4-Chloroaniline-d4         | 10.65 | 131 | 1369734 | 119.01 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 0.00  | 166 | 0d      | 0.00   | ng/ul |      |
| 46) Acenaphthylene-d8          | 0.00  | 160 | 0d      | 0.00   | ng/ul |      |
| 51) 4-Nitrophenol-d4           | 14.61 | 143 | 1263120 | 197.50 | ng/ul | 0.04 |
| 57) Fluorene-d10               | 0.00  | 176 | 0d      | 0.00   | ng/ul |      |
| 62) 4,6-Dinitro-2-methylphenol | 15.51 | 200 | 1548575 | 199.30 | ng/ul | 0.02 |
| 70) Anthracene-d10             | 0.00  | 188 | 0d      | 0.00   | ng/ul |      |
| 78) Pyrene-d10                 | 0.00  | 212 | 0d      | 0.00   | ng/ul |      |
| 89) Benzo(a)pyrene-d12         | 0.00  | 264 | 0d      | 0.00   | ng/ul |      |

## Target Compounds

|                                |       |     |          |         |        | Qvalue |
|--------------------------------|-------|-----|----------|---------|--------|--------|
| 4) Benzaldehyde                | 6.87  | 77  | 669447   | 113.310 | ng/ul  | 97     |
| 6) Phenol                      | 6.92  | 94  | 2039073  | 194.668 | ng/ul  | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.15  | 93  | 1419811  | 175.313 | ng/ul  | 98     |
| 11) 2-Methylphenol             | 8.16  | 108 | 1611759  | 201.934 | ng/ul  | 100    |
| 12) 2,2'-oxybis(1-Chloropropan | 8.24  | 45  | 1835959  | 185.910 | ng/ul  | 99     |
| 14) Acetophenone               | 8.55  | 105 | 2578689  | 196.409 | ng/ul  | 96     |
| 16) 4-Methylphenol             | 8.50  | 108 | 1747789  | 206.537 | ng/ul  | 98     |
| 30) 4-Chloroaniline            | 10.67 | 127 | 1331191  | 121.007 | ng/ul  | 99     |
| 32) Caprolactam                | 11.53 | 113 | 604502m  | 198.929 | ng/ul  |        |
| 37) Hexachlorocyclopentadiene  | 12.49 | 237 | 1604918  | 176.037 | ng/ul  | 99     |
| 48) 3-Nitroaniline             | 14.30 | 138 | 1171130  | 190.241 | ng/ul  | 96     |
| 50) 2,4-Dinitrophenol          | 14.50 | 184 | 951476   | 231.188 | ng/ul  | 95     |
| 52) 4-Nitrophenol              | 14.63 | 109 | 938083   | 196.916 | ng/ul  | 93     |
| 60) 4-Nitroaniline             | 15.49 | 138 | 1317628  | 208.456 | ng/ul  | 95     |
| 63) 4,6-Dinitro-2-methylphenol | 15.53 | 198 | 1465915  | 189.643 | ng/ul# | 88     |
| 67) Atrazine                   | 16.60 | 200 | 1850574  | 145.681 | ng/ul  | 99     |
| 68) Pentachlorophenol          | 16.76 | 266 | 1729780  | 193.005 | ng/ul  | 99     |
| 74) Carbazole                  | 17.53 | 167 | 8999382  | 163.452 | ng/ul  | 99     |
| 76) Fluoranthene               | 19.17 | 202 | 12343693 | 159.502 | ng/ul  | 99     |
| 81) 3,3'-Dichlorobenzidine     | 21.23 | 252 | 4896928  | 162.000 | ng/ul  | 98     |
| 86) Di-n-octyl phthalate       | 22.09 | 149 | 12127103 | 143.838 | ng/ul  | 100    |

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

