

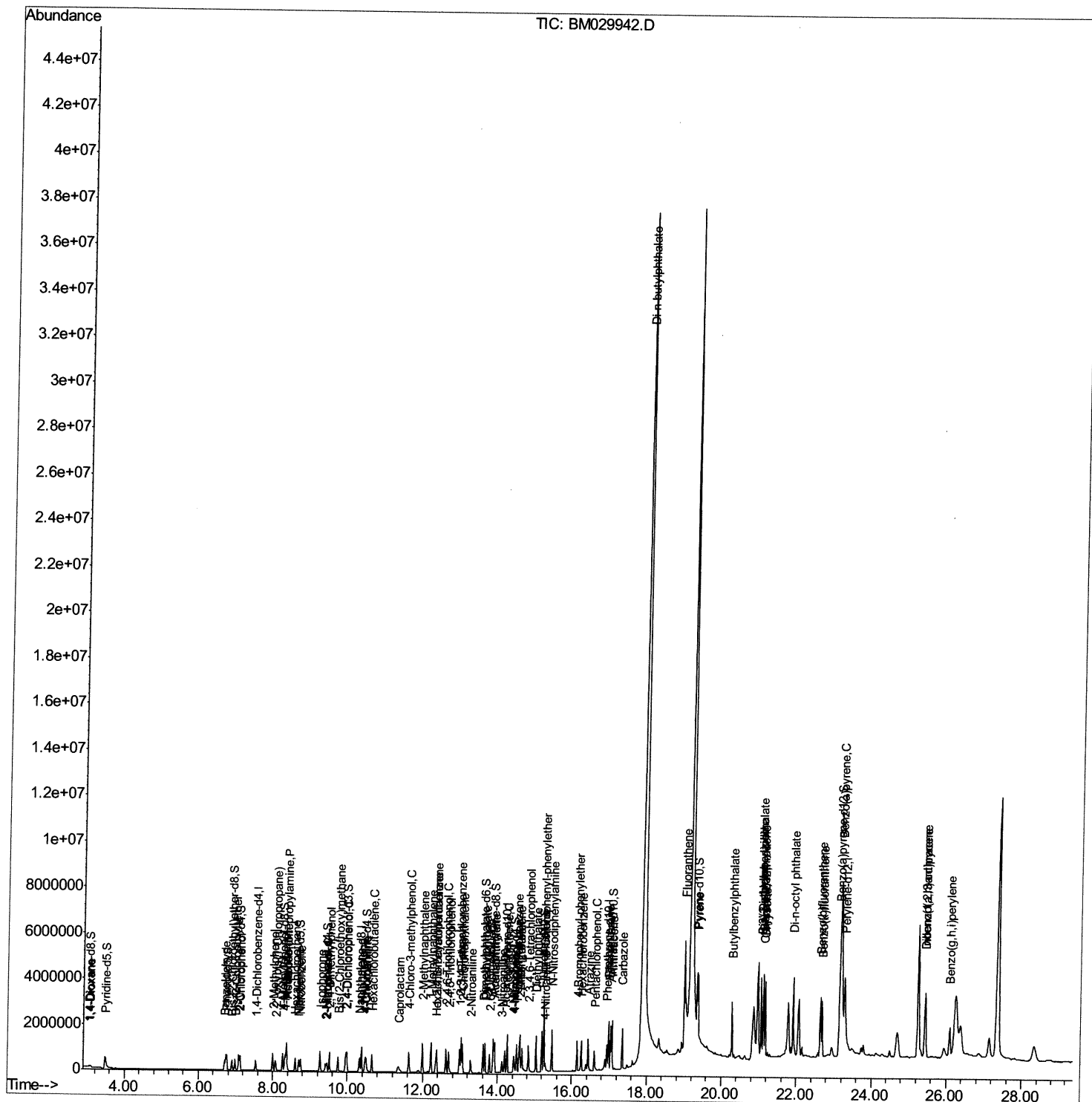
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Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
Data File : BM029942.D  
Acq On    : 11 May 2021   14:34  
Operator  : CG/JU  
Sample    : PB136055BS  
Misc      :  
ALS Vial  : 41      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**ClientSampleId :**  
SLCS055

## Manual Integrations

mohammad  
5/12/2021 2:06:33 PM

Quant Time: May 11 15:13:18 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue May 11 06:44:51 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

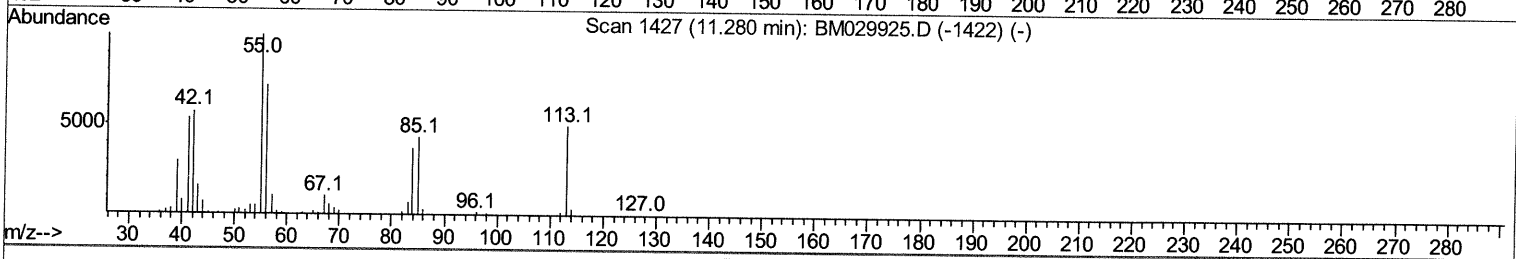
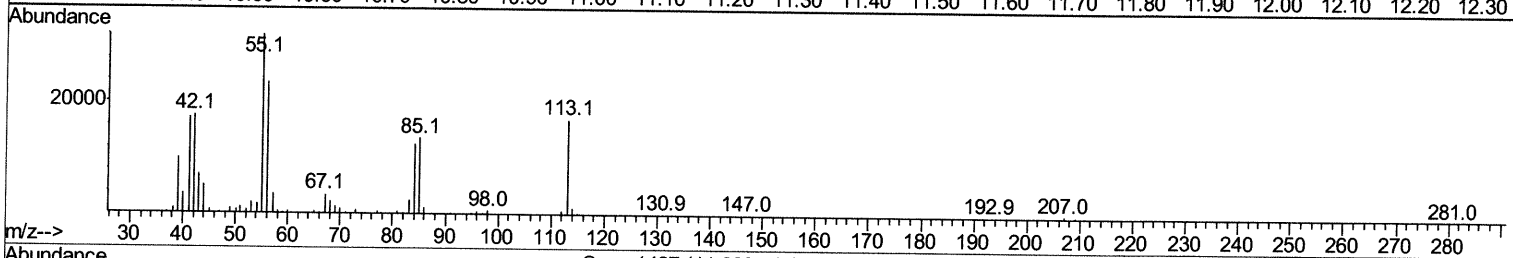
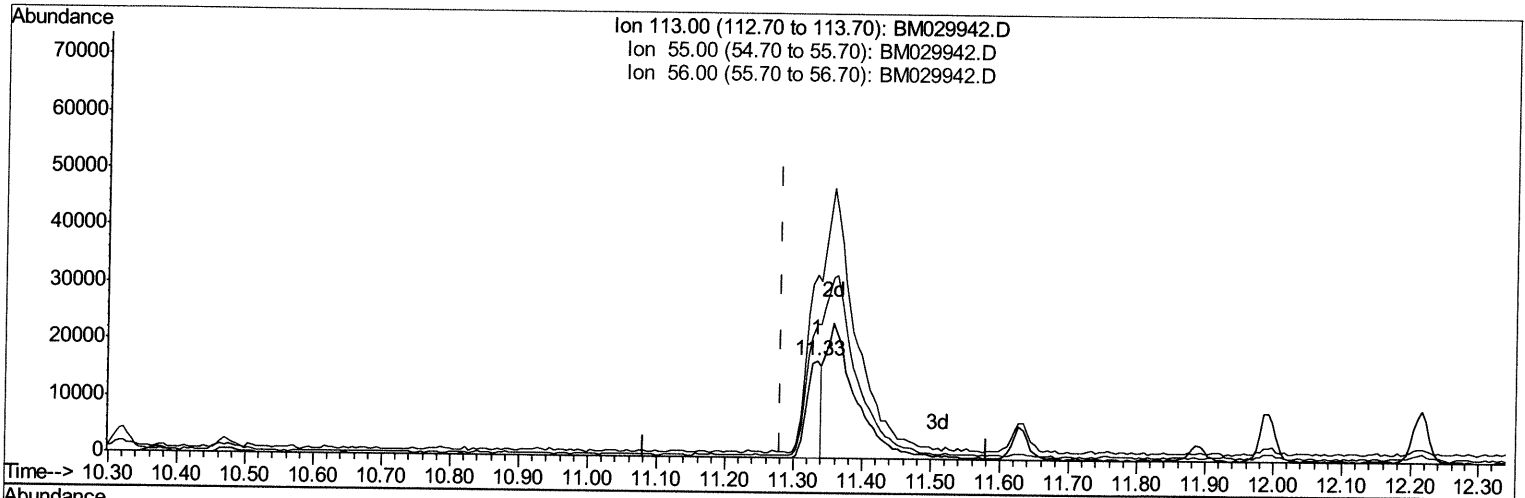
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 SLCS055

**Manual Integrations**  
**APPROVED**

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: May 11 15:05:21 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



TIC: BM029942.D

(34) Caprolactam

11.333min (+0.053) 9.63ng/ul

response 25795

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 173.60 | 187.86 |
| 56.00  | 131.80 | 137.54 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

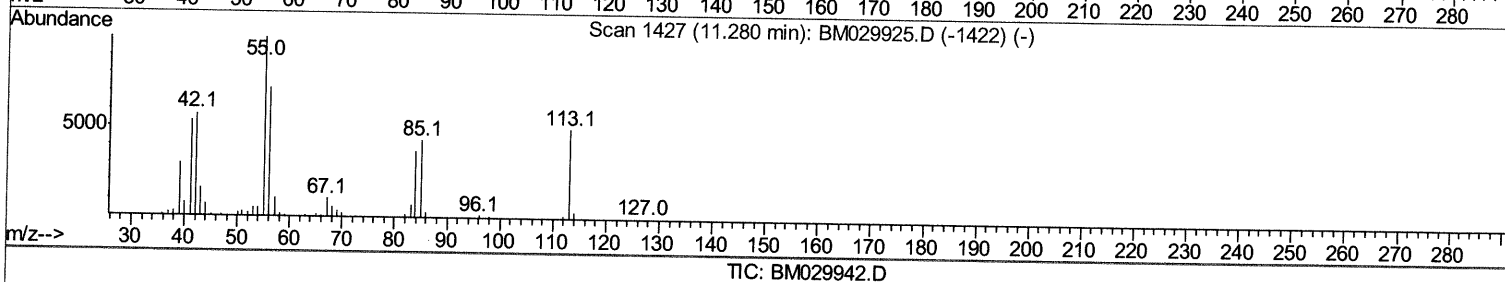
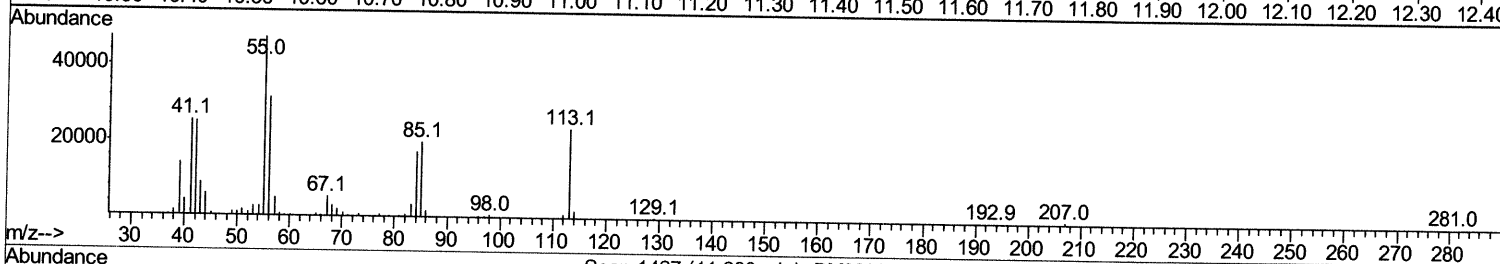
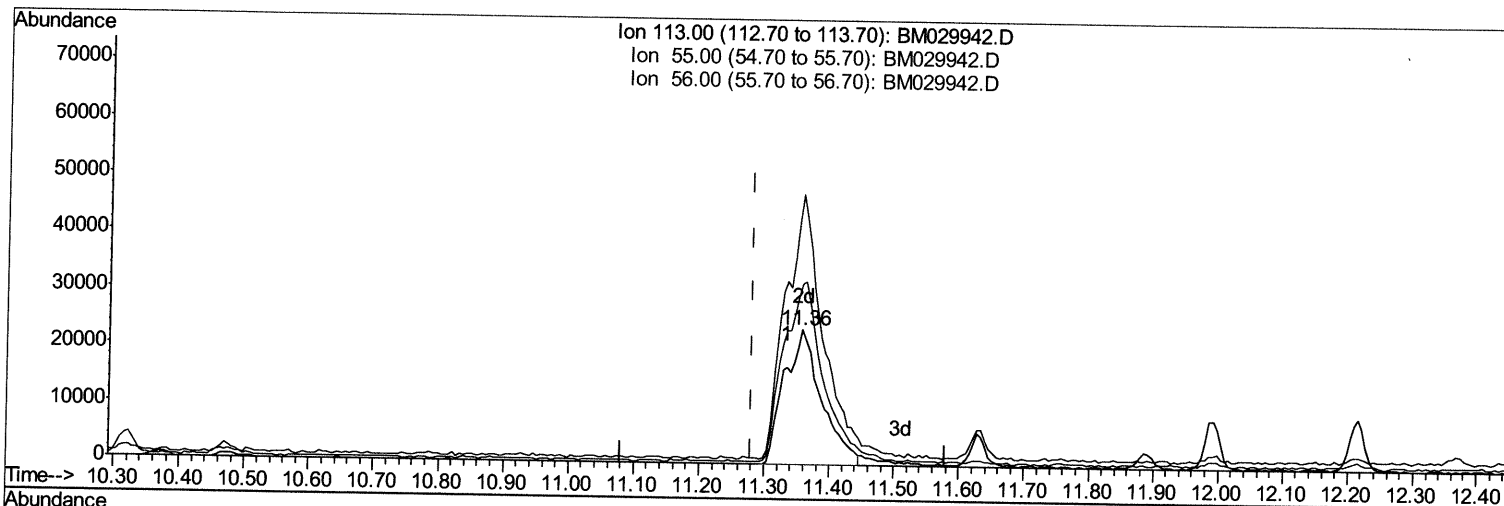
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
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 Operator : CG/JU  
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Instrument :  
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Manual Integrations  
 APPROVED

mohammad  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



TIC: BM029942.D

(34) Caprolactam

11.357min (+0.077) 35.40ng/ul m 05/13/21 JU

response 94829

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 173.60 | 200.34 |
| 56.00  | 131.80 | 133.67 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

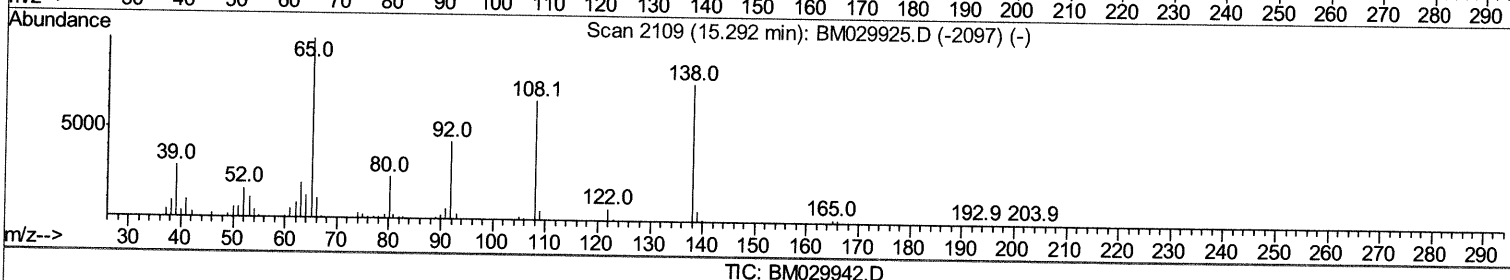
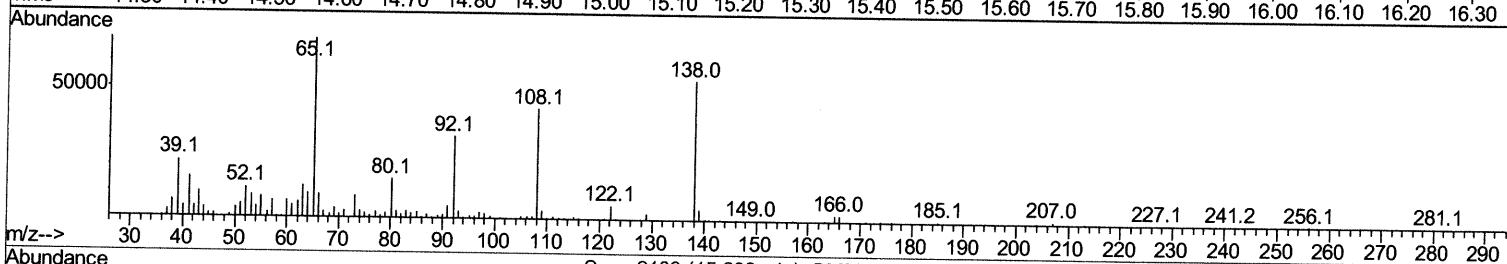
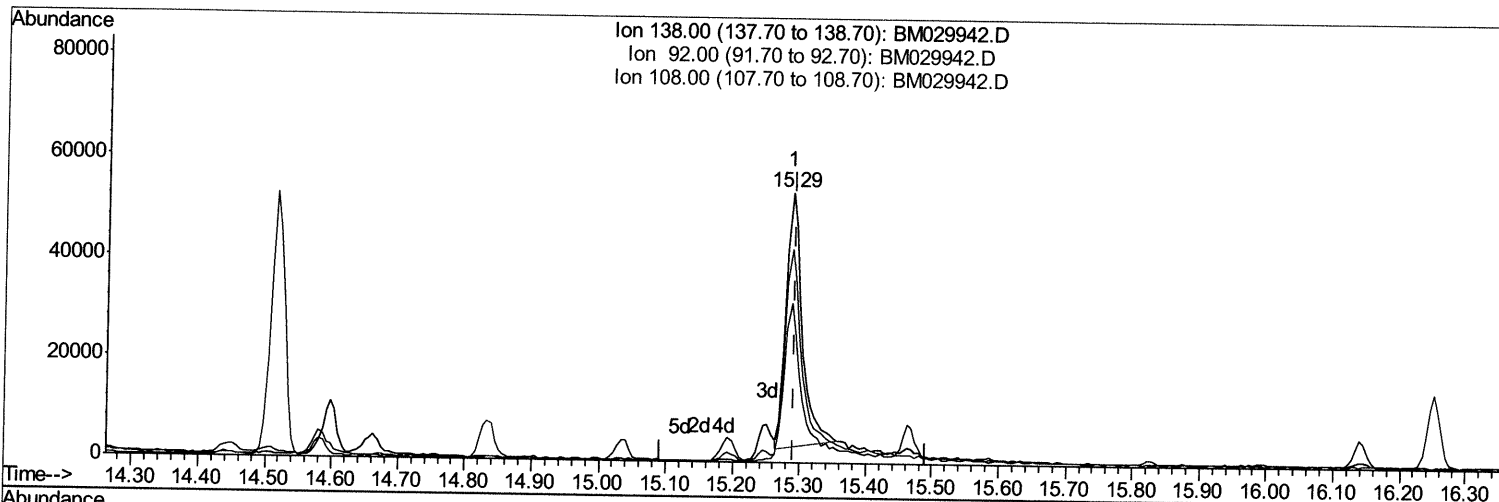
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
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Manual Integrations  
 APPROVED

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: Mav 11 15:05:21 2021  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



TIC: BM029942.D

(63) 4-Nitroaniline

15.286min (-0.006) 17.69ng/ul

response 84560

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 59.02 |
| 108.00 | 81.60 | 79.10 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

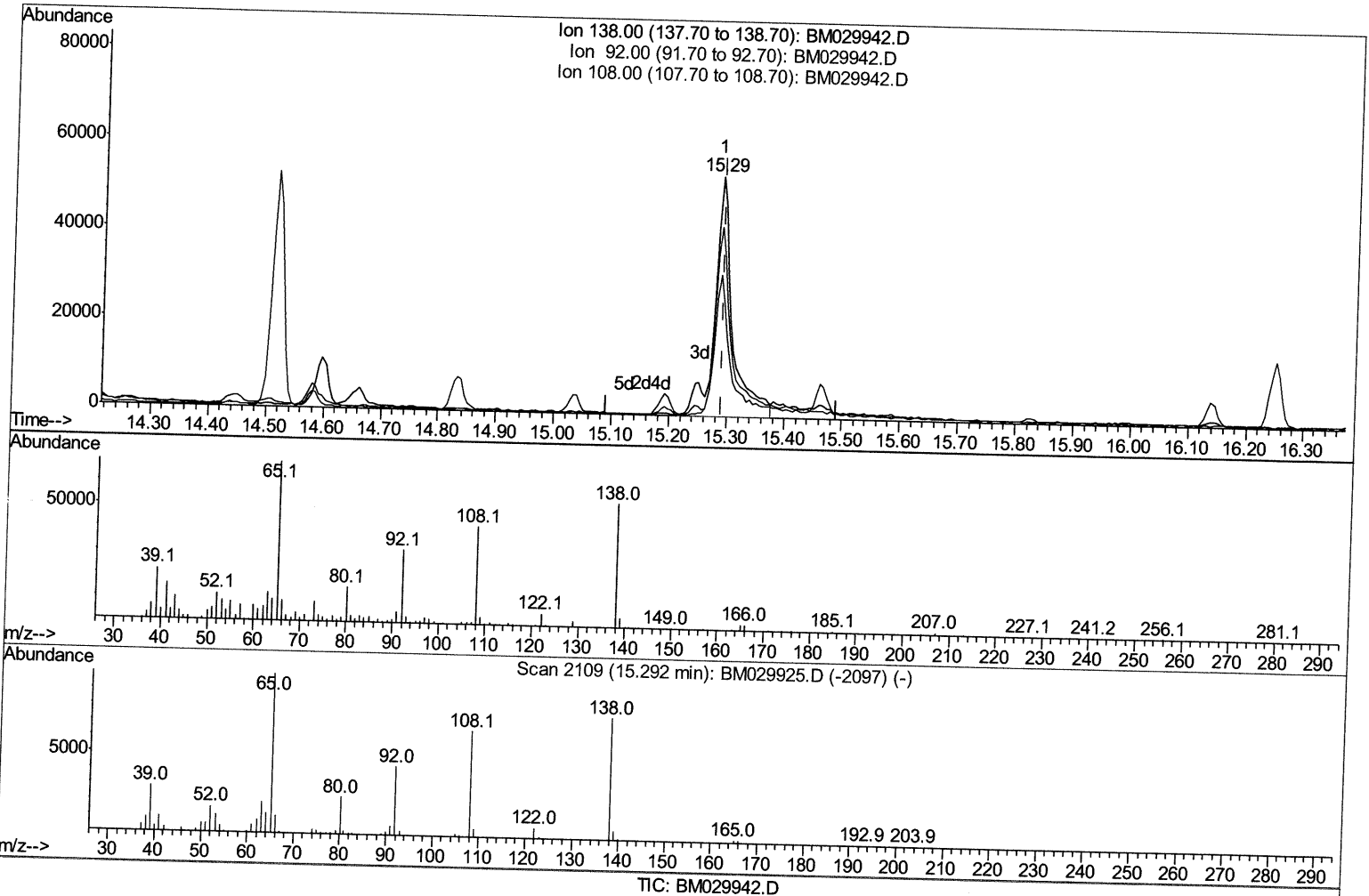
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS055

Manual Integrations  
 APPROVED

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: May 11 15:05:21 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



(63) 4-Nitroaniline

15.286min (-0.006) 25.06ng/ul m 05/13/2021

response 119813

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 59.02 |
| 108.00 | 81.60 | 79.10 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

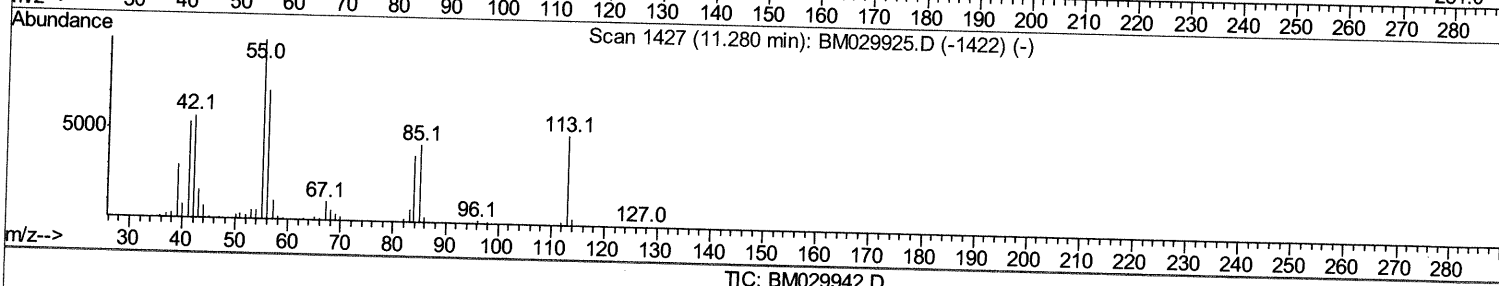
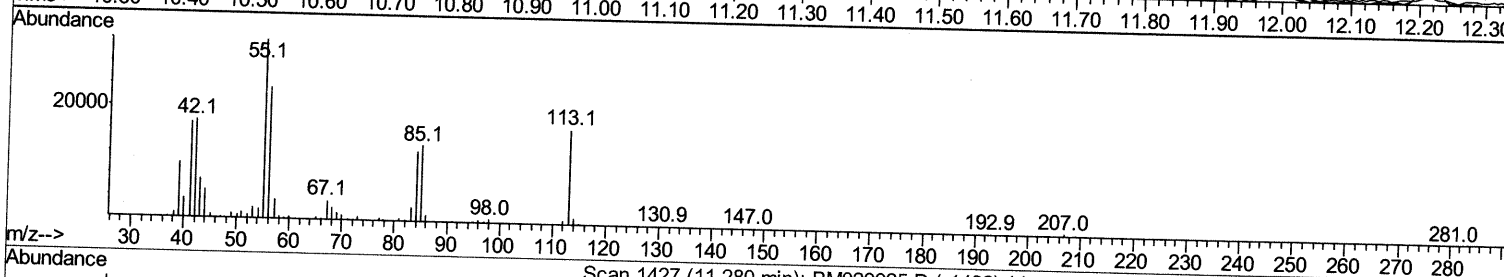
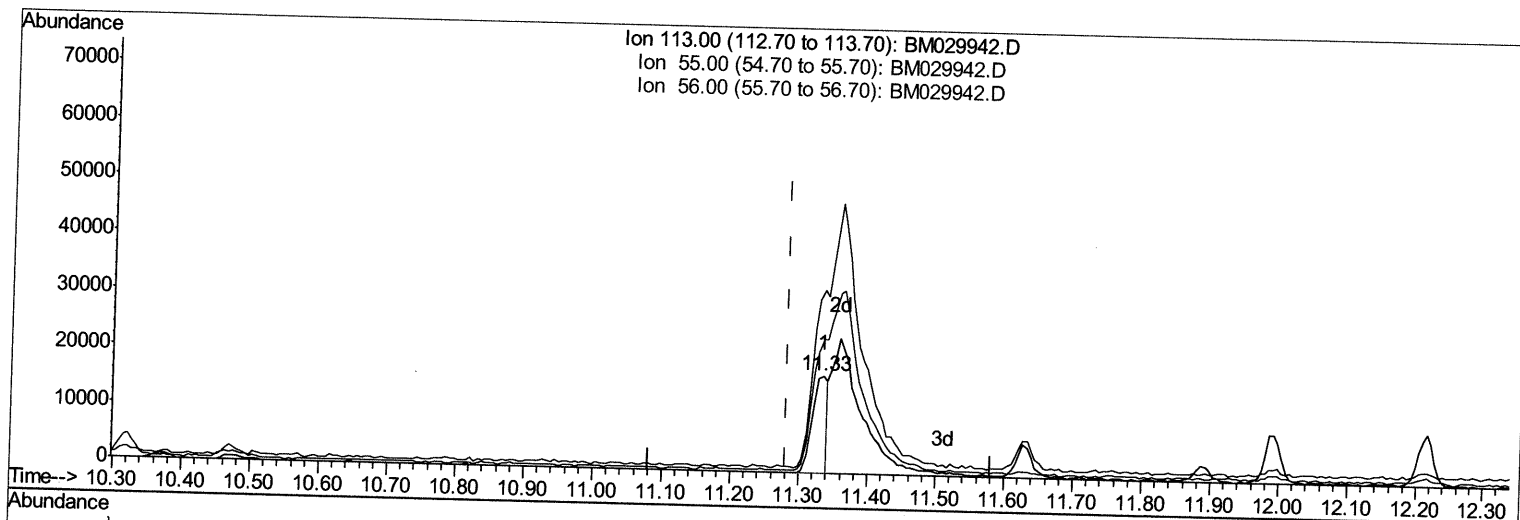
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 SLCS055

**Manual Integrations**  
**APPROVED**

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: May 11 15:09:54 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



TIC: BM029942.D

(34) Caprolactam

11.333min (+0.053) 9.63ng/ul

response 25795

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 173.60 | 187.86 |
| 56.00  | 131.80 | 137.54 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

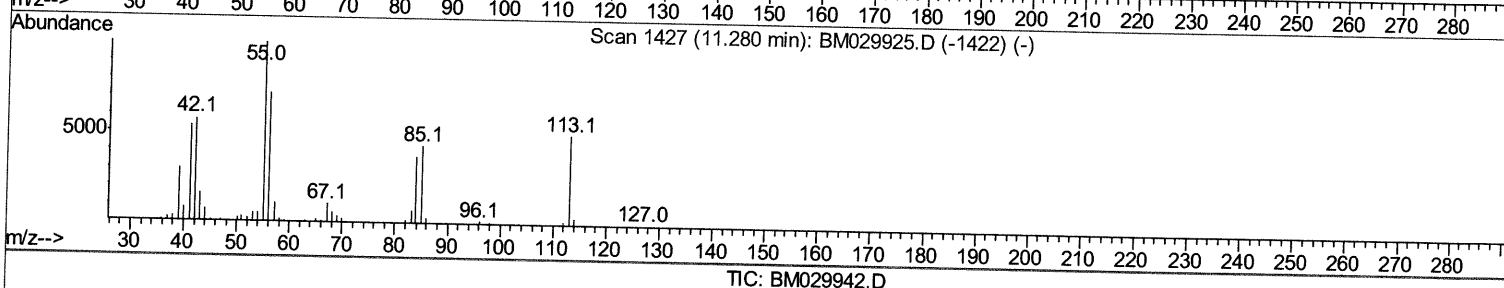
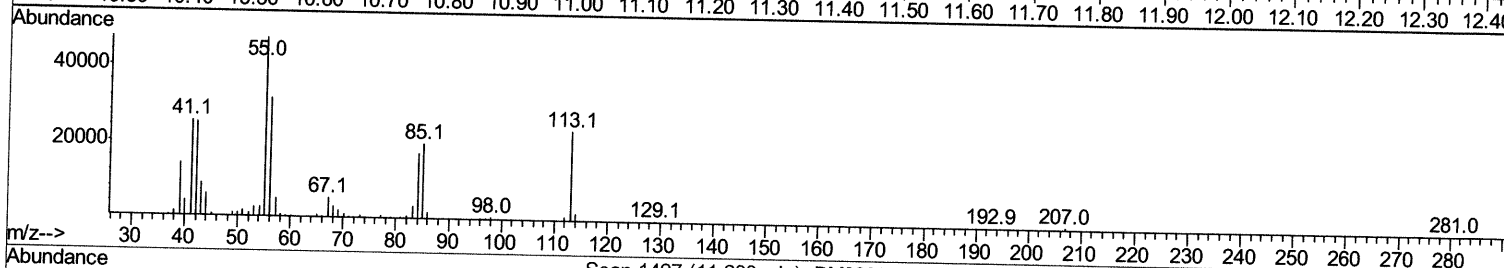
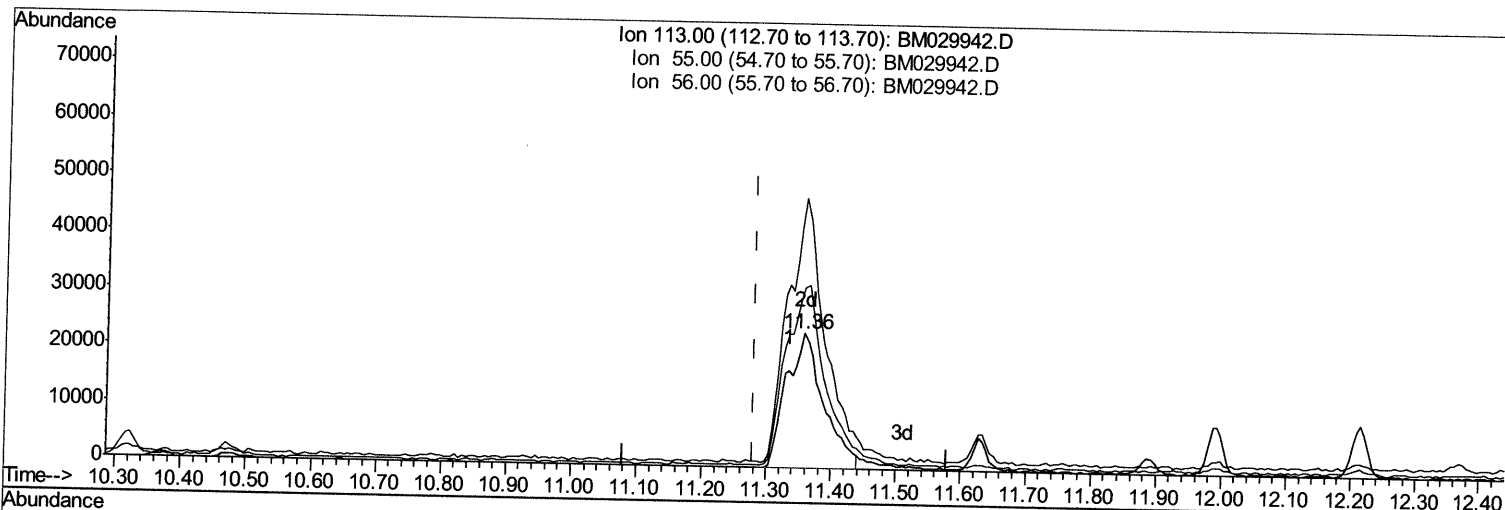
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 SLCS055

**Manual Integrations**  
**APPROVED**

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: May 11 15:09:54 2021  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



(34) Caprolactam

11.357min (+0.077) 35.18ng/ul m 05/11/2021

response 94233

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 173.60 | 200.34 |
| 56.00  | 131.80 | 133.67 |
| 0.00   | 0.00   | 0.00   |

## Quantitation Report (Qedit)

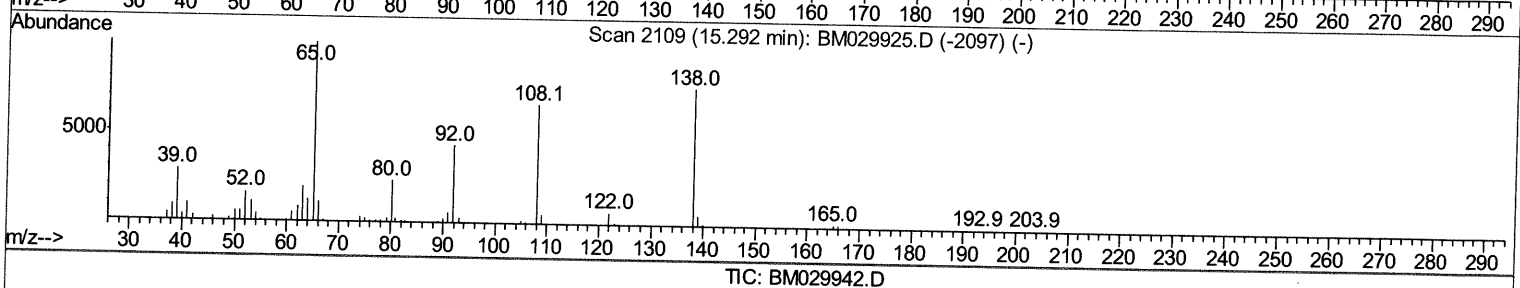
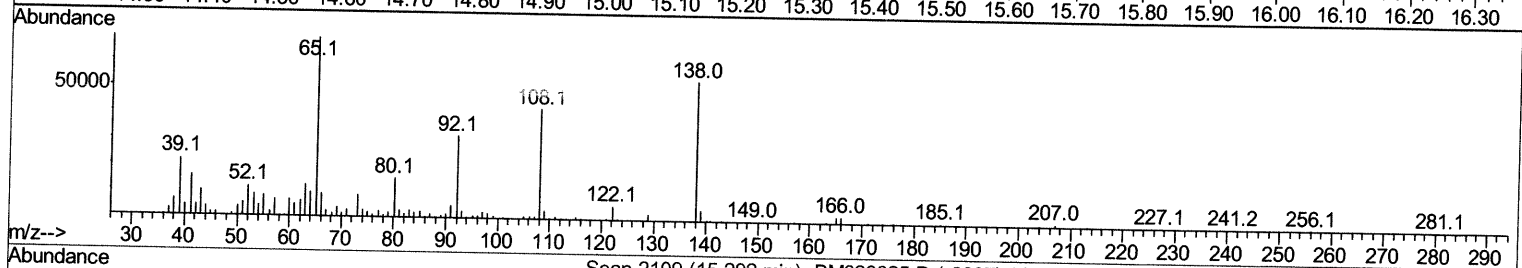
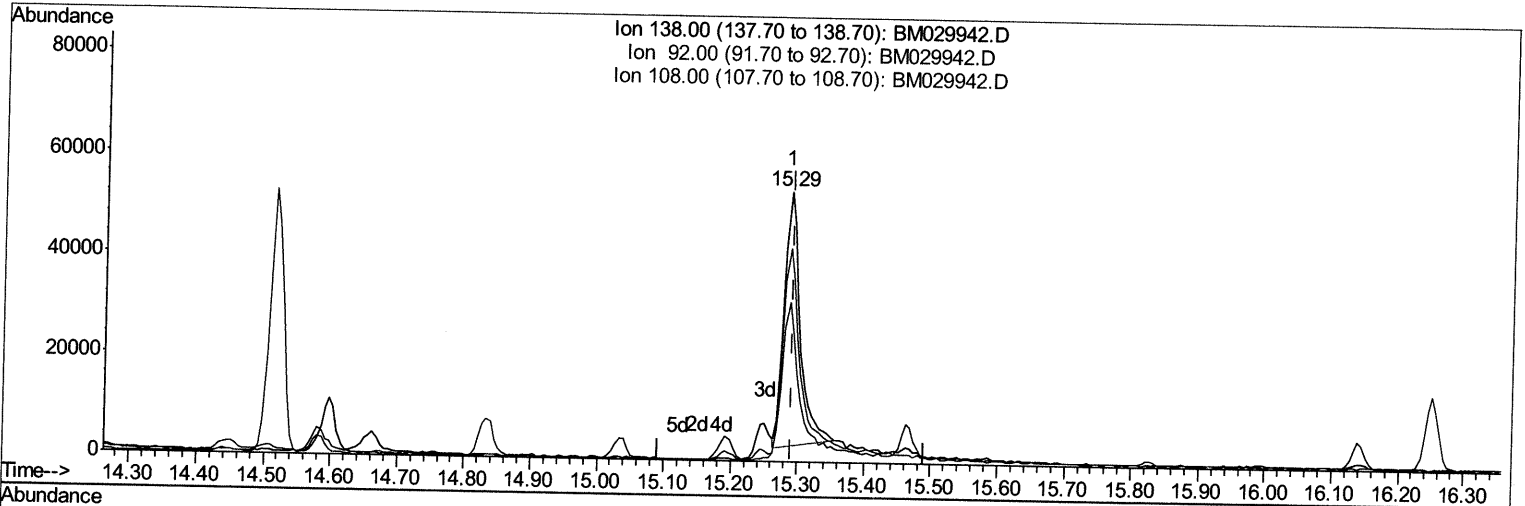
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
Data File : BM029942.D  
Acq On : 11 May 2021 14:34  
Operator : CG/JU  
Sample : PB136055BS  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS055

Manual Integrations  
APPROVED

mohammad  
5/12/2021 2:06:33 PM

Quant Time: May 11 15:09:54 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue May 11 06:44:51 2021  
Response via : Initial Calibration



(63) 4-Nitroaniline

15.286min (-0.006) 17.69ng/ul

response 84560

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 59.02 |
| 108.00 | 81.60 | 79.10 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

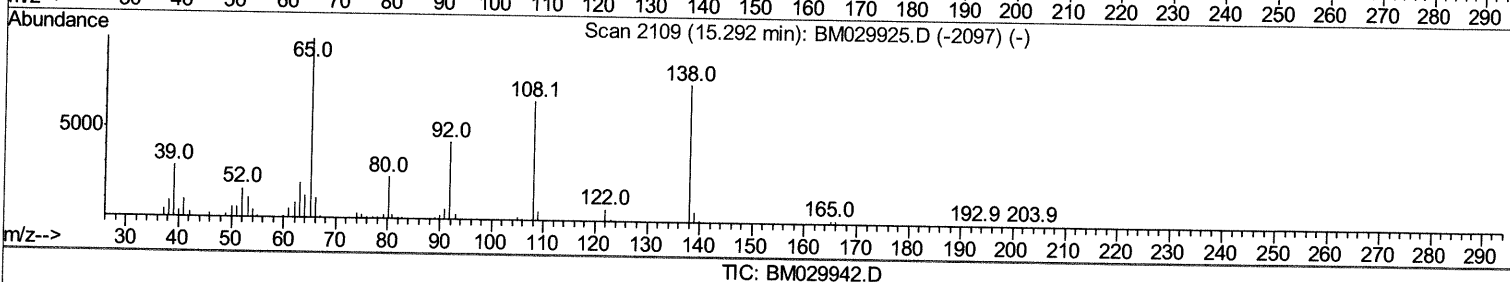
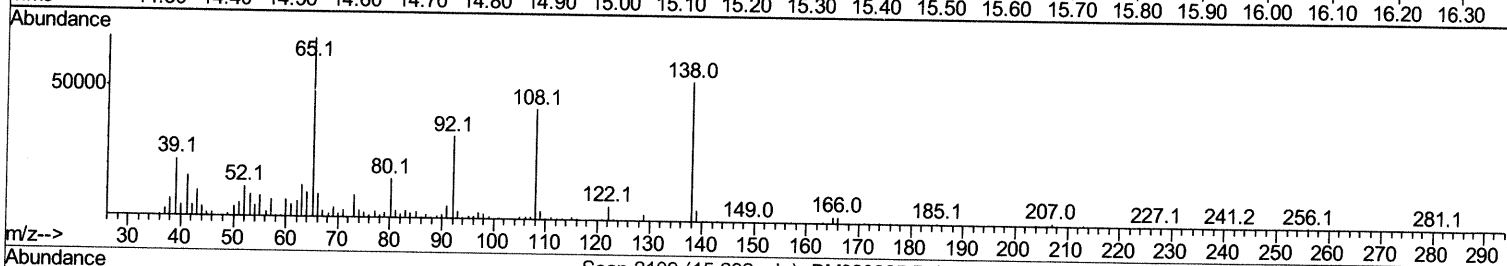
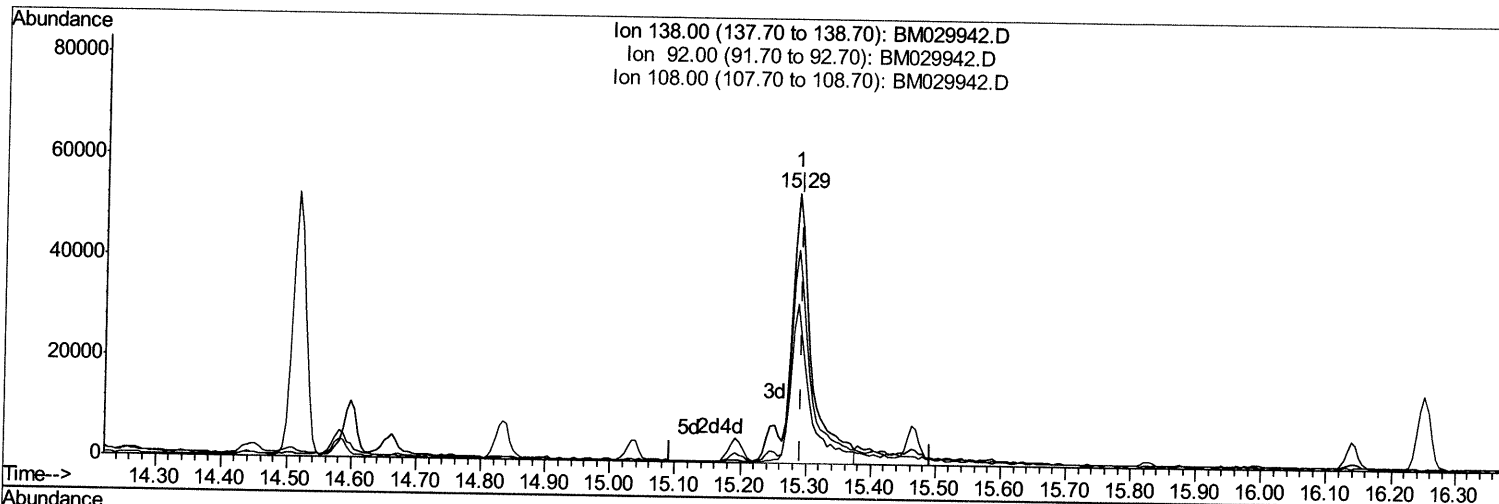
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
 Data File : BM029942.D  
 Acq On : 11 May 2021 14:34  
 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
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Manual Integrations  
 APPROVED

mohammad  
 5/12/2021 2:06:33 PM

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 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration



TIC: BM029942.D

(63) 4-Nitroaniline

15.286min (-0.006) 25.03ng/ul m 05/13/21JU

response 119673

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 59.02 |
| 108.00 | 81.60 | 79.10 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.55  | 152  | 115296   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.32 | 136  | 492792   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.20 | 164  | 308154   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.94 | 188  | 615803   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.13 | 240  | 636737   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.29 | 264  | 667531   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.06  | 96  | 19592   | 6.79  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.48  | 84  | 224988  | 31.13 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.73  | 99  | 340305  | 31.74 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.89  | 67  | 206782  | 31.10 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.08  | 132 | 279190  | 33.74 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.26  | 113 | 272166  | 30.69 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.70  | 128 | 107141  | 31.55 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.42  | 143 | 96979   | 25.30 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 9.95  | 165 | 288935  | 37.59 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.47 | 131 | 317333  | 27.13 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.62 | 166 | 783494  | 32.80 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 13.89 | 160 | 978004  | 32.31 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.43 | 143 | 110540  | 28.41 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.19 | 176 | 672666  | 33.17 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.34 | 200 | 1803    | 0.45  | ng/ul | 0.01 |
| 73) Anthracene-d10             | 17.04 | 188 | 1063716 | 34.30 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.34 | 212 | 1214244 | 37.17 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.16 | 264 | 1363554 | 36.03 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        | Ovalue |           |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.09  | 88  | 41847  | 13.394 | ng/uL# 86 |
| 5) Pyridine                    | 3.49  | 79  | 233527 | 31.497 | ng/ul 98  |
| 6) Benzaldehyde                | 6.70  | 77  | 160645 | 29.025 | ng/ul 97  |
| 8) Phenol                      | 6.76  | 94  | 349927 | 31.966 | ng/ul 99  |
| 10) Bis(2-Chloroethyl)ether    | 6.99  | 93  | 256931 | 29.684 | ng/ul 99  |
| 12) 2-Chlorophenol             | 7.12  | 128 | 280245 | 33.207 | ng/ul 99  |
| 13) 2-Methylphenol             | 8.00  | 108 | 247058 | 29.788 | ng/ul 98  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.08  | 45  | 430342 | 33.007 | ng/ul 95  |
| 16) Acetophenone               | 8.37  | 105 | 401313 | 28.202 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.36  | 70  | 205431 | 28.216 | ng/ul 96  |
| 18) 4-Methylphenol             | 8.33  | 108 | 278572 | 30.297 | ng/ul 98  |
| 19) Hexachloroethane           | 8.60  | 117 | 109992 | 30.848 | ng/ul 91  |
| 22) Nitrobenzene               | 8.75  | 77  | 295134 | 33.352 | ng/ul 98  |
| 23) Isophorone                 | 9.27  | 82  | 594102 | 31.741 | ng/ul 98  |
| 25) 2-Nitrophenol              | 9.45  | 139 | 111008 | 26.676 | ng/ul 96  |
| 26) 2,4-Dimethylphenol         | 9.52  | 107 | 284389 | 30.287 | ng/ul 98  |
| 27) Bis(2-Chloroethoxy)methane | 9.75  | 93  | 351972 | 32.106 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 9.98  | 162 | 274732 | 36.398 | ng/ul 99  |
| 30) Naphthalene                | 10.37 | 128 | 828275 | 30.281 | ng/ul 100 |
| 32) 4-Chloroaniline            | 10.49 | 127 | 303194 | 25.162 | ng/ul 99  |
| 33) Hexachlorobutadiene        | 10.65 | 225 | 148154 | 29.275 | ng/ul 97  |
| 34) Caprolactam                | 11.36 | 113 | 94233m | 35.179 | ng/ul >   |

05/13/21 JU

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 Operator : CG/JU  
 Sample : PB136055BS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS055

Manual Integrations  
 APPROVED

mohammad  
 5/12/2021 2:06:33 PM

Quant Time: May 11 15:13:18 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 06:44:51 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min)   |
|--------------------------------|-------|------|----------|--------|--------|------------|
| 35) 4-Chloro-3-methylphenol    | 11.63 | 107  | 298924   | 35.787 | ng/ul  | 99         |
| 36) 2-Methyl-naphthalene       | 11.99 | 142  | 587191   | 29.594 | ng/ul  | 98         |
| 37) 1-Methyl-naphthalene       | 12.22 | 142  | 571258   | 29.051 | ng/ul  | 99         |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.37 | 216  | 288931   | 30.343 | ng/ul  | 98         |
| 40) Hexachlorocyclopentadiene  | 12.35 | 237  | 118946   | 24.047 | ng/ul  | 99         |
| 41) 2,4,6-Trichlorophenol      | 12.62 | 196  | 226884   | 39.072 | ng/ul  | 99         |
| 42) 2,4,5-Trichlorophenol      | 12.70 | 196  | 242097   | 39.394 | ng/ul  | 98         |
| 43) 1,1'-Biphenyl              | 13.02 | 154  | 754013   | 31.477 | ng/ul  | 97         |
| 44) 2-Chloronaphthalene        | 13.06 | 162  | 581669   | 31.641 | ng/ul  | 98         |
| 45) 2-Nitroaniline             | 13.28 | 65   | 155392   | 33.443 | ng/ul  | 95         |
| 47) Dimethylphthalate          | 13.66 | 163  | 768641   | 32.138 | ng/ul  | 99         |
| 48) 2,6-Dinitrotoluene         | 13.79 | 165  | 134241   | 31.266 | ng/ul  | 99         |
| 50) Acenaphthylene             | 13.92 | 152  | 890006   | 30.402 | ng/ul  | 99         |
| 51) 3-Nitroaniline             | 14.12 | 138  | 127308   | 26.395 | ng/ul  | 97         |
| 52) Acenaphthene               | 14.26 | 153  | 651005   | 31.465 | ng/ul  | 99         |
| 55) 4-Nitrophenol              | 14.44 | 109  | 85602    | 28.618 | ng/ul  | 91         |
| 56) Dibenzofuran               | 14.60 | 168  | 889940   | 31.405 | ng/ul  | 99         |
| 57) 2,4-Dinitrotoluene         | 14.58 | 165  | 206615   | 32.521 | ng/ul  | 94         |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.83 | 232  | 213970   | 36.563 | ng/ul  | 98         |
| 59) Diethylphthalate           | 15.03 | 149  | 832391   | 33.369 | ng/ul  | 99         |
| 61) Fluorene                   | 15.25 | 166  | 768571   | 31.726 | ng/ul  | 100        |
| 62) 4-Chlorophenyl-phenylether | 15.25 | 204  | 370580   | 31.131 | ng/ul  | 97         |
| 63) 4-Nitroaniline             | 15.29 | 138  | 119673m> | 25.031 | ng/ul> | 05/13/21JU |
| 67) N-Nitrosodiphenylamine     | 15.46 | 169  | 665409   | 33.263 | ng/ul  | 99         |
| 68) 4-Bromophenyl-phenylether  | 16.14 | 248  | 229082   | 33.421 | ng/ul  | 97         |
| 69) Hexachlorobenzene          | 16.25 | 284  | 264198   | 32.458 | ng/ul  | 95         |
| 70) Atrazine                   | 16.44 | 200  | 231391   | 31.351 | ng/ul  | 99         |
| 71) Pentachlorophenol          | 16.60 | 266  | 168040   | 36.029 | ng/ul  | 100        |
| 72) Phenanthrene               | 16.99 | 178  | 1219273  | 33.579 | ng/ul  | 100        |
| 74) Anthracene                 | 17.07 | 178  | 1256944  | 33.860 | ng/ul  | 99         |
| 75) 1,2,3,4-Tetrachlorobenzene | 12.99 | 216  | 302293   | 31.033 | ng/uL  | 97         |
| 76) Pentachlorobenzene         | 14.52 | 250  | 284370   | 30.612 | ng/uL  | 98         |
| 77) Carbazole                  | 17.35 | 167  | 1132947  | 33.310 | ng/ul  | 99         |
| 78) Di-n-butylphthalate        | 17.92 | 149  | 1778658  | 44.058 | ng/ul  | 99         |
| 80) Fluoranthene               | 19.01 | 202  | 1498913  | 37.705 | ng/ul# | 96         |
| 82) Pvrene                     | 19.37 | 202  | 1532922  | 36.641 | ng/ul  | 100        |
| 83) Butylbenzylphthalate       | 20.29 | 149  | 702496   | 47.250 | ng/ul  | 97         |
| 84) 3,3'-Dichlorobenzidine     | 21.06 | 252  | 66543    | 4.594  | ng/ul# | 95         |
| 85) Benzo(a)anthracene         | 21.12 | 228  | 1471779  | 35.428 | ng/ul  | 99         |
| 86) Bis(2-ethylhexyl)phthalate | 21.06 | 149  | 1076676  | 42.000 | ng/ul  | 99         |
| 87) Chrysene                   | 21.17 | 228  | 1431592  | 34.599 | ng/ul  | 99         |
| 89) Di-n-octyl phthalate       | 21.92 | 149  | 1877582  | 36.756 | ng/ul  | 100        |
| 90) Benzo(b)fluoranthene       | 22.64 | 252  | 1538395  | 35.529 | ng/ul  | 99         |
| 91) Benzo(k)fluoranthene       | 22.69 | 252  | 1494610  | 32.833 | ng/ul  | 98         |
| 93) Benzo(a)pyrene             | 23.20 | 252  | 1476750  | 37.552 | ng/ul  | 98         |
| 94) Indeno(1,2,3-cd)pyrene     | 25.44 | 276  | 1796774  | 37.750 | ng/ul  | 98         |
| 95) Dibenzo(a,h)anthracene     | 25.45 | 278  | 1505295  | 37.310 | ng/ul  | 99         |
| 96) Benzo(a,h,i)perylene       | 26.10 | 276  | 1456120  | 37.758 | ng/ul  | 99         |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\  
Data File : BM029942.D  
Acq On : 11 May 2021 14:34  
Operator : CG/JU  
Sample : PB136055BS  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS055

Manual Integrations  
APPROVED

mohammad  
5/12/2021 2:06:33 PM

Quant Time: May 11 15:13:18 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue May 11 06:44:51 2021  
Response via : Initial Calibration

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