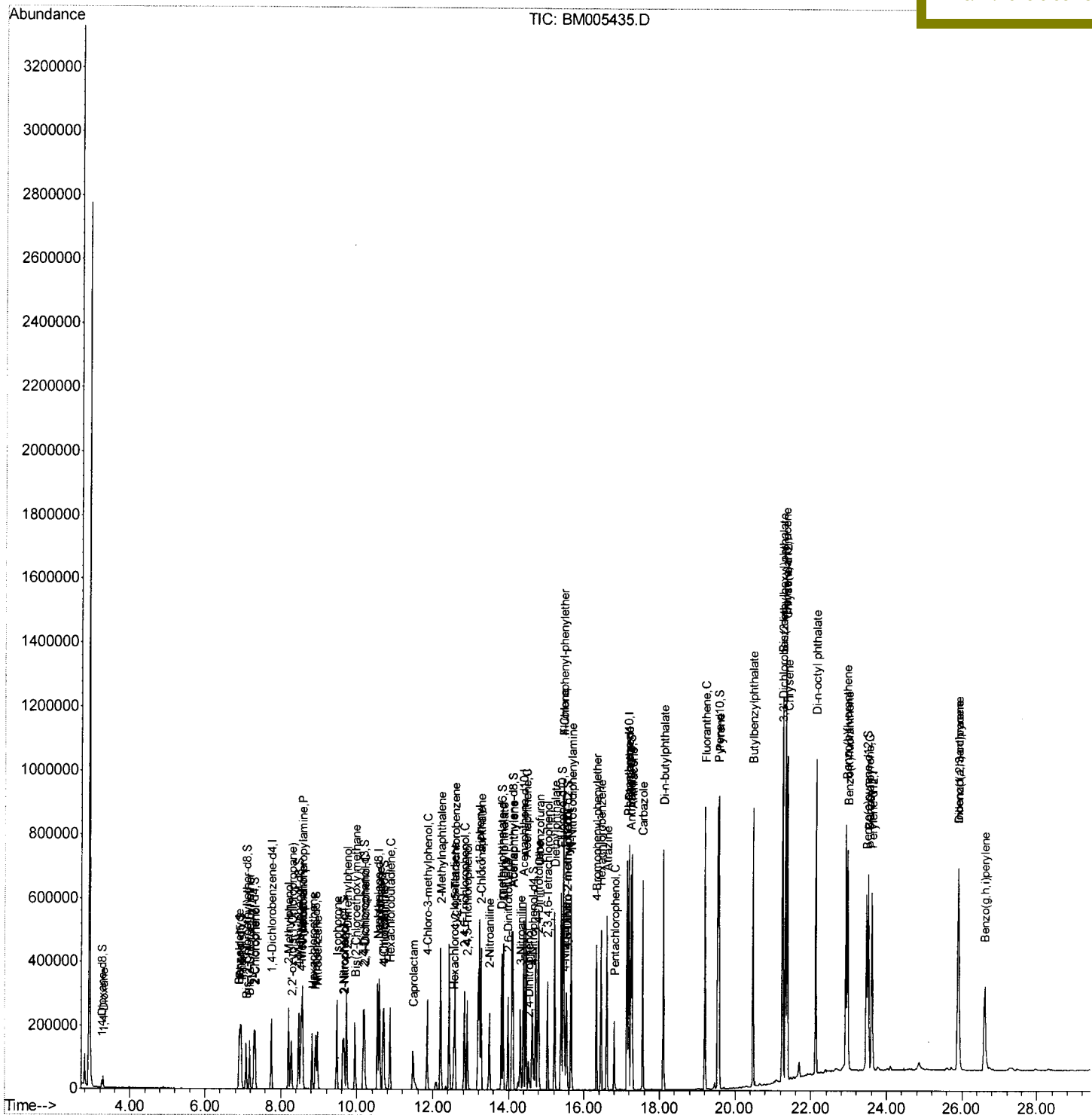


Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleID :  
 SSTD02034

Quant Time: May 14 00:47:48 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration

Manual Integrations  
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 5/14/2016 9:58:48 AM



Quantitation Report (Qedit)

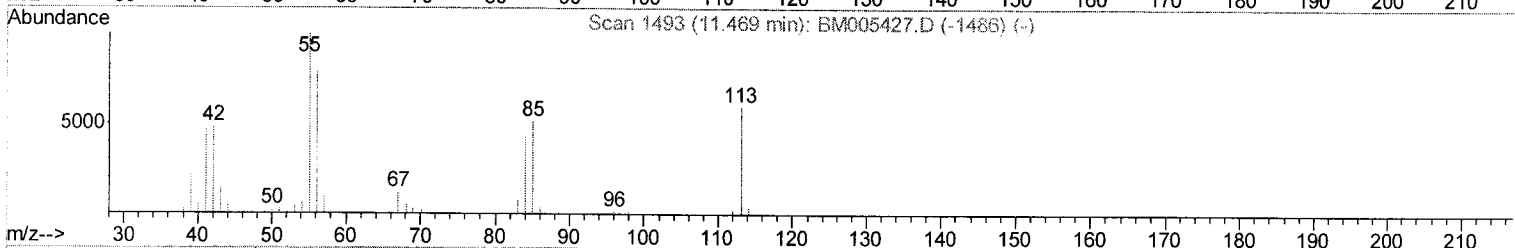
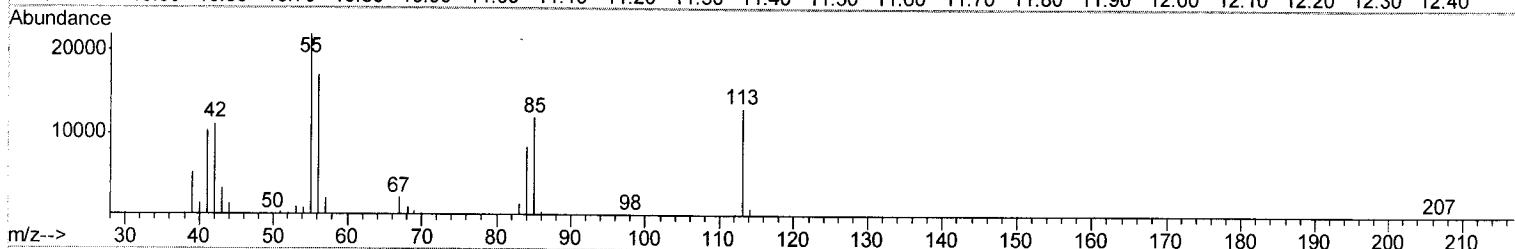
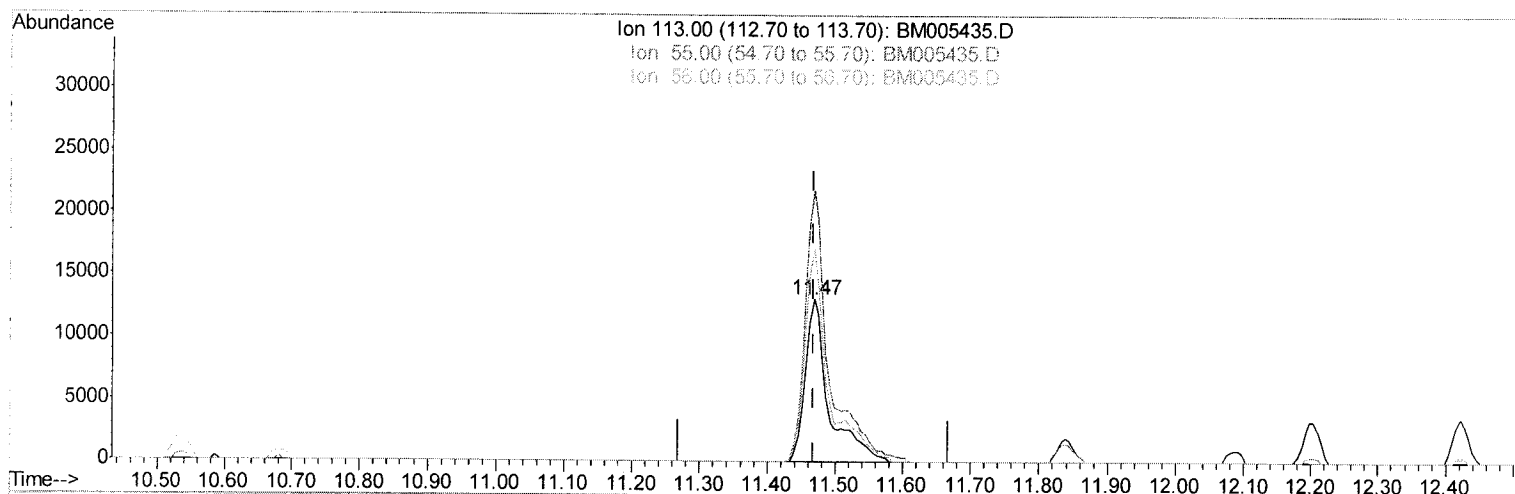
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Manual Integrations  
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 5/14/2016 9:58:48 AM

Quant Time: May 14 00:21:13 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration



TIC: BM005435.D

(32) Caprolactam

11.469min (+0.000) 20.08ng/ul m *U.M*

response 31987

*05/17/16*

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 168.20 | 167.14 |
| 56.00  | 120.80 | 130.40 |
| 0.00   | 0.00   | 0.00   |

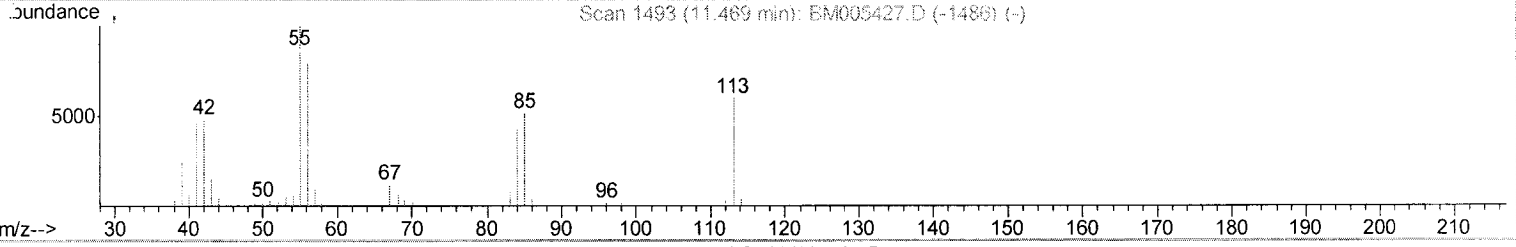
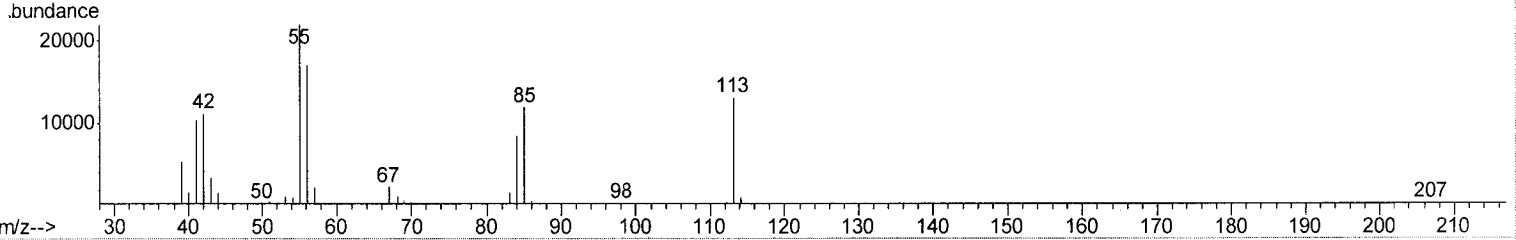
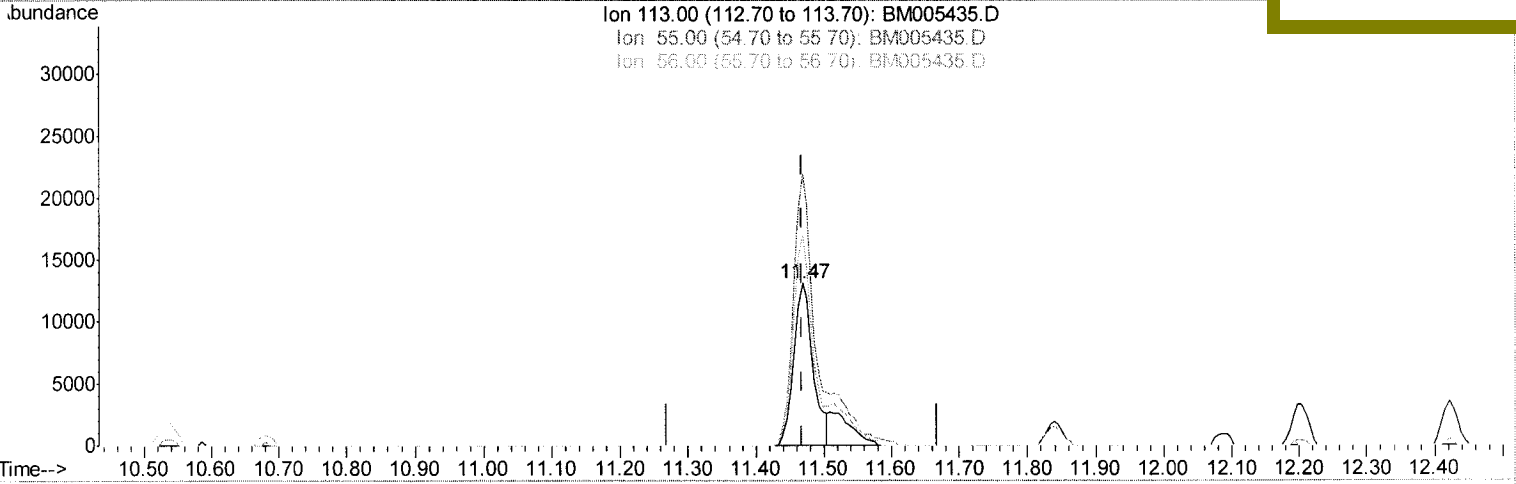
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Quant Time: May 14 00:21:13 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED  
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 5/14/2016 9:58:48 AM



TIC: BM005435.D

(32) Caprolactam

11.469min (+0.000) 16.17ng/ul

response 25766

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 168.20 | 167.14 |
| 56.00  | 120.80 | 130.40 |
| 0.00   | 0.00   | 0.00   |

Quantitation Report (Qedit)

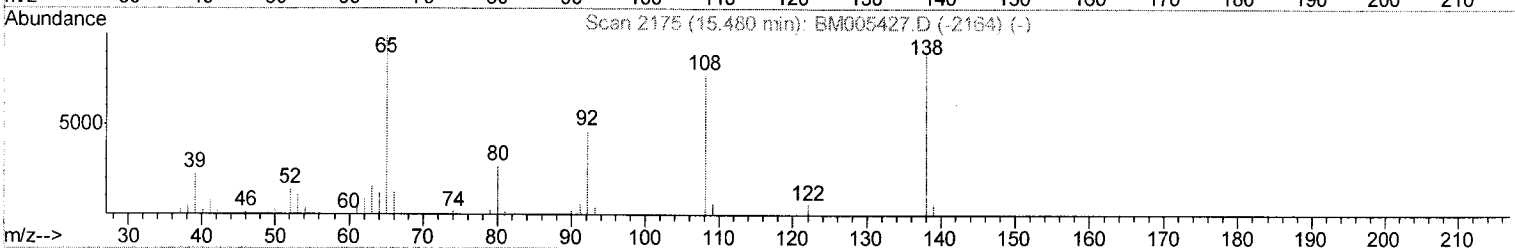
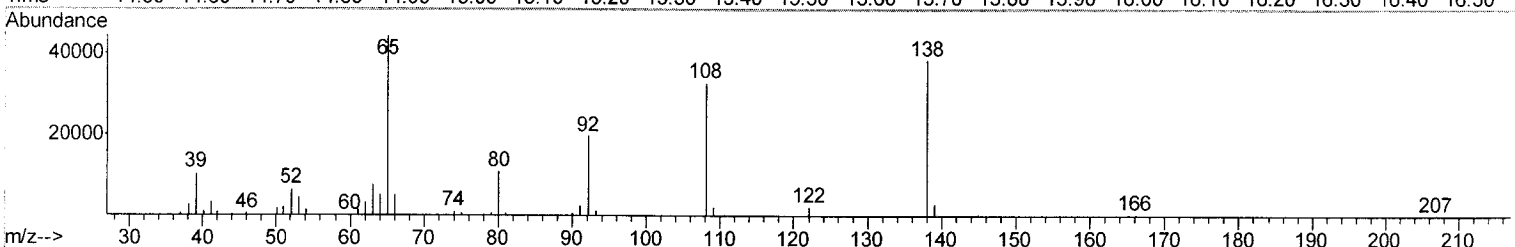
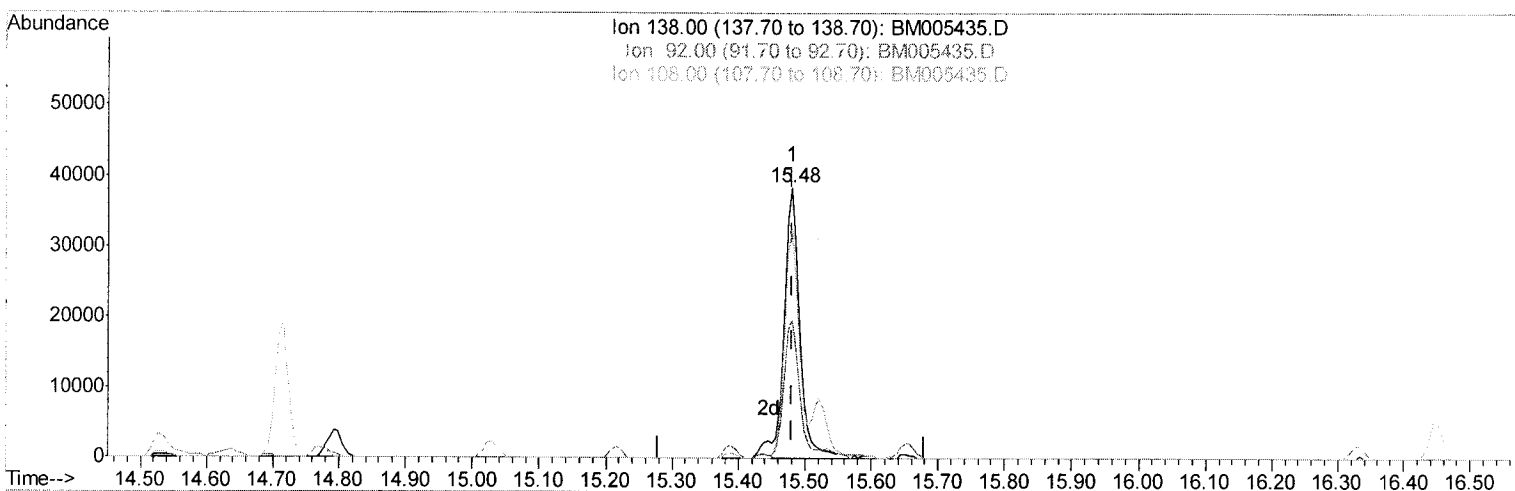
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Manual Integrations  
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 5/14/2016 9:58:48 AM

Quant Time: May 14 00:21:13 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration



TIC: BM005435.D

(60) 4-Nitroaniline

15.480min (+0.000) 21.00ng/ul m U.M

response 65233

*05/17/16*

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 55.40 | 51.50 |
| 108.00 | 82.90 | 85.22 |
| 0.00   | 0.00  | 0.00  |

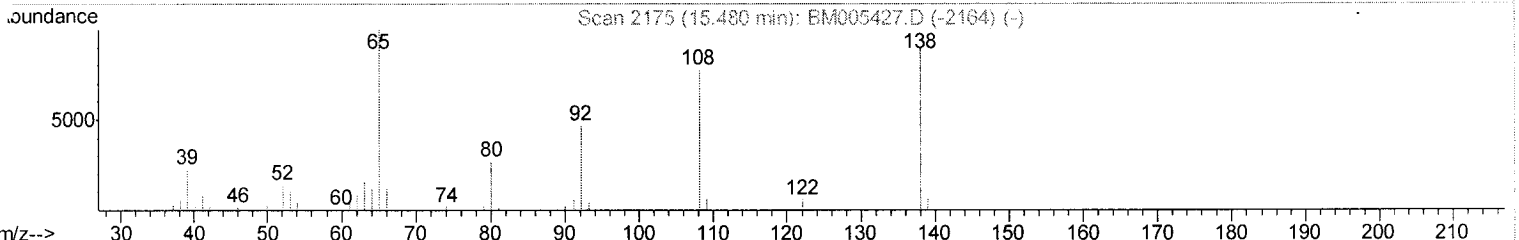
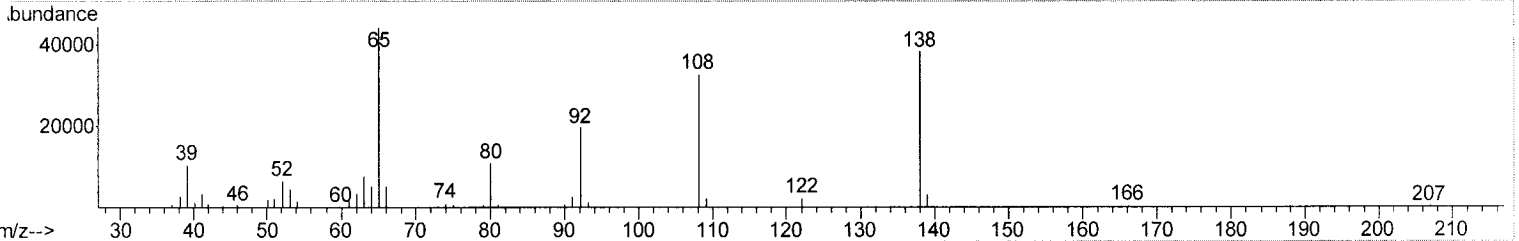
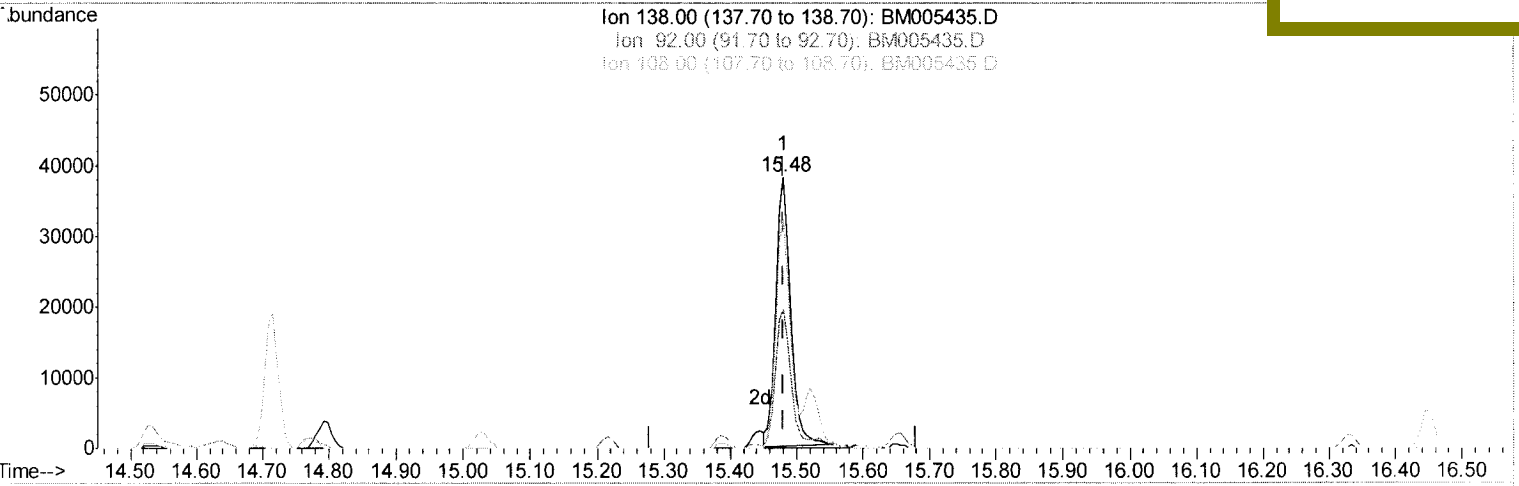
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleID :  
 SSTD02034

Quant Time: May 14 00:21:13 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED  
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 5/14/2016 9:58:48 AM



TIC: BM005435.D

(60) 4-Nitroaniline

15.480min (+0.000) 19.11ng/ul

response 59387

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 55.40 | 51.50 |
| 108.00 | 82.90 | 85.22 |
| 0.00   | 0.00  | 0.00  |

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleID :  
 SSTD02034

Manual Integrations  
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 5/14/2016 9:58:48 AM

Quant Time: May 14 00:47:48 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.57 | 216  | 112506   | 21.24 | ng/ul  | 98       |
| 37) Hexachlorocyclopentadiene  | 12.55 | 237  | 42927    | 15.87 | ng/ul  | 98       |
| 38) 2,4,6-Trichlorophenol      | 12.82 | 196  | 74688    | 21.17 | ng/ul  | 95       |
| 39) 2,4,5-Trichlorophenol      | 12.90 | 196  | 82983    | 21.20 | ng/ul  | 97       |
| 40) 1,1'-Biphenyl              | 13.22 | 154  | 285955   | 20.73 | ng/ul  | 99       |
| 41) 2-Chloronaphthalene        | 13.26 | 162  | 216607   | 20.77 | ng/ul  | 99       |
| 42) 2-Nitroaniline             | 13.48 | 65   | 70564    | 21.98 | ng/ul  | 95       |
| 44) Dimethylphthalate          | 13.85 | 163  | 289065   | 20.73 | ng/ul  | 100      |
| 45) 2,6-Dinitrotoluene         | 13.98 | 165  | 59180    | 21.52 | ng/ul# | 88       |
| 47) Acenaphthylene             | 14.12 | 152  | 360748   | 20.83 | ng/ul  | 100      |
| 48) 3-Nitroaniline             | 14.31 | 138  | 63989    | 21.83 | ng/ul  | 93       |
| 49) Acenaphthene               | 14.46 | 153  | 236335   | 20.71 | ng/ul  | 99       |
| 50) 2,4-Dinitrophenol          | 14.53 | 184  | 23933    | 15.13 | ng/ul  | 99       |
| 52) 4-Nitrophenol              | 14.63 | 109  | 39795    | 18.79 | ng/ul  | 97       |
| 53) Dibenzofuran               | 14.79 | 168  | 342310   | 20.67 | ng/ul  | 100      |
| 54) 2,4-Dinitrotoluene         | 14.77 | 165  | 88569    | 21.60 | ng/ul  | 99       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.03 | 232  | 67080    | 20.16 | ng/ul# | 96       |
| 56) Diethylphthalate           | 15.22 | 149  | 296405   | 21.04 | ng/ul  | 99       |
| 58) Fluorene                   | 15.44 | 166  | 274397   | 21.20 | ng/ul  | 100      |
| 59) 4-Chlorophenyl-phenylether | 15.44 | 204  | 138348   | 21.57 | ng/ul  | 98       |
| 60) 4-Nitroaniline             | 15.48 | 138  | 65233m   | 21.00 | ng/ul  |          |
| 63) 4,6-Dinitro-2-methylphenol | 15.54 | 198  | 48563    | 19.69 | ng/ul# | 93       |
| 64) N-Nitrosodiphenylamine     | 15.66 | 169  | 242600   | 21.23 | ng/ul  | 99       |
| 65) 4-Bromophenyl-phenylether  | 16.33 | 248  | 88368    | 22.09 | ng/ul  | 96       |
| 66) Hexachlorobenzene          | 16.45 | 284  | 98745    | 22.00 | ng/ul  | 96       |
| 67) Atrazine                   | 16.60 | 200  | 96006    | 23.02 | ng/ul  | 100      |
| 68) Pentachlorophenol          | 16.80 | 266  | 43436    | 17.42 | ng/ul  | 99       |
| 69) Phenanthrene               | 17.19 | 178  | 461069   | 21.01 | ng/ul  | 99       |
| 71) Anthracene                 | 17.27 | 178  | 464292   | 21.22 | ng/ul  | 99       |
| 72) Carbazole                  | 17.55 | 167  | 433371   | 22.52 | ng/ul  | 99       |
| 73) Di-n-butylphthalate        | 18.10 | 149  | 520162   | 22.16 | ng/ul  | 100      |
| 74) Fluoranthene               | 19.21 | 202  | 566127   | 23.29 | ng/ul  | 96       |
| 77) Pyrene                     | 19.57 | 202  | 576547   | 21.15 | ng/ul  | 97       |
| 78) Butylbenzylphthalate       | 20.47 | 149  | 253618   | 22.41 | ng/ul  | 97       |
| 79) 3,3'-Dichlorobenzidine     | 21.26 | 252  | 187330   | 22.17 | ng/ul  | 98       |
| 80) Benzo(a)anthracene         | 21.33 | 228  | 565341   | 20.91 | ng/ul  | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.24 | 149  | 359721   | 22.88 | ng/ul# | 97       |
| 82) Chrysene                   | 21.38 | 228  | 537694   | 20.96 | ng/ul  | 100      |
| 84) Di-n-octyl phthalate       | 22.13 | 149  | 619924   | 25.55 | ng/ul  | 99       |
| 85) Benzo(b)fluoranthene       | 22.93 | 252  | 533943   | 21.99 | ng/ul  | 98       |
| 86) Benzo(k)fluoranthene       | 22.97 | 252  | 498813   | 21.82 | ng/ul  | 99       |
| 88) Benzo(a)pyrene             | 23.51 | 252  | 487738   | 21.24 | ng/ul  | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.90 | 276  | 447971   | 17.88 | ng/ul  | 97       |
| 90) Dibenzo(a,h)anthracene     | 25.91 | 278  | 376136   | 17.94 | ng/ul  | 98       |
| 91) Benzo(g,h,i)perylene       | 26.61 | 276  | 363228   | 17.12 | ng/ul  | 98       |

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 05/17/16

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

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051316\  
 Data File : BM005435.D  
 Acq On : 13 May 2016 17:05  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Quant Time: May 14 00:47:48 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 14 00:15:57 2016  
 Response via : Initial Calibration

Manual Integrations  
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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.75  | 152  | 59500    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.54 | 136  | 277888   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.39 | 164  | 174095   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.14 | 188  | 417314   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.34 | 240  | 470014   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.61 | 264  | 415334   | 20.00 | ng/ul | 0.00     |

System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.25  | 96  | 10424  | 8.24  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.93  | 99  | 108211 | 20.05 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl) ether-d | 7.09  | 67  | 63812  | 20.73 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.29  | 132 | 83431  | 20.47 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.46  | 113 | 87757  | 19.67 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.91  | 128 | 41680  | 21.01 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.63  | 143 | 48702  | 21.68 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.17 | 165 | 88800  | 21.27 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.68 | 131 | 113736 | 22.63 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.80 | 166 | 291682 | 20.90 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.09 | 160 | 342929 | 20.95 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.62 | 143 | 45520  | 17.87 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.39 | 176 | 247746 | 20.56 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.52 | 200 | 46141  | 19.66 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.24 | 188 | 392585 | 21.28 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.54 | 212 | 457296 | 21.08 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.47 | 264 | 390208 | 21.22 | ng/ul | 0.00 |

Target Compounds

| Target Compounds               | R.T.  | QIon | Response | Conc  | Units  | Qvalue |
|--------------------------------|-------|------|----------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.29  | 88   | 18702    | 8.11  | ng/uL# | 93     |
| 4) Benzaldehyde                | 6.90  | 77   | 68199    | 26.08 | ng/ul  | 97     |
| 6) Phenol                      | 6.96  | 94   | 113435   | 20.34 | ng/ul  | 99     |
| 8) Bis(2-Chloroethyl) ether    | 7.17  | 93   | 85379    | 20.33 | ng/ul  | 97     |
| 10) 2-Chlorophenol             | 7.32  | 128  | 85207    | 20.44 | ng/ul  | 98     |
| 11) 2-Methylphenol             | 8.20  | 108  | 85497    | 19.83 | ng/ul  | 100    |
| 12) 2,2'-oxybis(1-Chloropropan | 8.28  | 45   | 120976   | 21.24 | ng/ul  | 98     |
| 14) Acetophenone               | 8.57  | 105  | 139141   | 21.06 | ng/ul  | 96     |
| 15) N-Nitroso-di-n-propylamine | 8.55  | 70   | 71706    | 20.77 | ng/ul  | 100    |
| 16) 4-Methylphenol             | 8.53  | 108  | 93794    | 19.61 | ng/ul  | 100    |
| 17) Hexachloroethane           | 8.82  | 117  | 32996    | 21.03 | ng/ul  | 94     |
| 20) Nitrobenzene               | 8.95  | 77   | 104295   | 21.00 | ng/ul  | 96     |
| 21) Isophorone                 | 9.47  | 82   | 201145   | 20.86 | ng/ul  | 99     |
| 23) 2-Nitrophenol              | 9.66  | 139  | 51674    | 21.35 | ng/ul  | 96     |
| 24) 2,4-Dimethylphenol         | 9.72  | 107  | 106957   | 20.83 | ng/ul  | 97     |
| 25) Bis(2-Chloroethoxy)methane | 9.95  | 93   | 120688   | 20.09 | ng/ul  | 98     |
| 27) 2,4-Dichlorophenol         | 10.19 | 162  | 88930    | 20.80 | ng/ul  | 100    |
| 28) Naphthalene                | 10.59 | 128  | 287252   | 20.54 | ng/ul  | 99     |
| 30) 4-Chloroaniline            | 10.70 | 127  | 113782   | 22.20 | ng/ul  | 99     |
| 31) Hexachlorobutadiene        | 10.86 | 225  | 56755    | 22.33 | ng/ul  | 98     |
| 32) Caprolactam                | 11.47 | 113  | 31987m   | 20.08 | ng/ul  | 98     |
| 33) 4-Chloro-3-methylphenol    | 11.84 | 107  | 107902   | 21.05 | ng/ul  | 98     |
| 34) 2-Methylnaphthalene        | 12.20 | 142  | 215843   | 20.64 | ng/ul  | 98     |

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