

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
 Data File : BG016163.D  
 Acq On : 27 Mar 2015 5:38  
 Operator : TP/IZ  
 Sample : G1653-14  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 FD9

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.791       | 13            | 17          | 34           | rVB      | 679729         | 790661        | 30.95%          | 4.697%        |
| 2         | 2.932       | 36            | 41          | 50           | rVV      | 76214          | 121299        | 4.75%           | 0.721%        |
| 3         | 3.008       | 50            | 54          | 59           | rVB      | 65612          | 71737         | 2.81%           | 0.426%        |
| 4         | 4.917       | 374           | 379         | 388          | rVB      | 87734          | 118584        | 4.64%           | 0.705%        |
| 5         | 5.381       | 451           | 458         | 471          | rVB      | 720464         | 1006873       | 39.41%          | 5.982%        |
| 6         | 6.962       | 720           | 727         | 738          | rBV      | 745323         | 1126577       | 44.10%          | 6.693%        |
| 7         | 7.326       | 781           | 789         | 799          | rBV      | 735567         | 1132187       | 44.32%          | 6.726%        |
| 8         | 7.790       | 862           | 868         | 875          | rBV      | 182346         | 278887        | 10.92%          | 1.657%        |
| 9         | 8.113       | 915           | 923         | 929          | rBV      | 506767         | 825692        | 32.32%          | 4.906%        |
| 10        | 8.947       | 1057          | 1065        | 1072         | rBV      | 395617         | 664054        | 25.99%          | 3.945%        |
| 11        | 10.345      | 1296          | 1303        | 1311         | rBV      | 25894          | 39674         | 1.55%           | 0.236%        |
| 12        | 10.574      | 1335          | 1342        | 1351         | rVB      | 272268         | 436227        | 17.08%          | 2.592%        |
| 13        | 13.041      | 1754          | 1762        | 1770         | rVB      | 1124909        | 1709226       | 66.91%          | 10.155%       |
| 14        | 13.870      | 1898          | 1903        | 1910         | rVB      | 26098          | 37239         | 1.46%           | 0.221%        |
| 15        | 14.363      | 1982          | 1987        | 1990         | rBV      | 47766          | 66769         | 2.61%           | 0.397%        |
| 16        | 14.416      | 1990          | 1996        | 2006         | rVB2     | 427350         | 653553        | 25.58%          | 3.883%        |
| 17        | 15.896      | 2242          | 2248        | 2256         | rBV      | 924435         | 1323095       | 51.79%          | 7.861%        |
| 18        | 16.061      | 2270          | 2276        | 2281         | rBV      | 21640          | 30686         | 1.20%           | 0.182%        |
| 19        | 16.167      | 2288          | 2294        | 2302         | rVV      | 150360         | 236495        | 9.26%           | 1.405%        |
| 20        | 16.237      | 2302          | 2306        | 2310         | rVV      | 23959          | 35215         | 1.38%           | 0.209%        |
| 21        | 16.296      | 2312          | 2316        | 2322         | rVV2     | 22532          | 41818         | 1.64%           | 0.248%        |
| 22        | 16.355      | 2322          | 2326        | 2331         | rVV2     | 23140          | 29083         | 1.14%           | 0.173%        |
| 23        | 16.554      | 2356          | 2360        | 2367         | rVB4     | 18985          | 35707         | 1.40%           | 0.212%        |
| 24        | 16.625      | 2367          | 2372        | 2375         | rBV2     | 25039          | 34344         | 1.34%           | 0.204%        |
| 25        | 17.148      | 2454          | 2461        | 2471         | rBV2     | 558723         | 809967        | 31.70%          | 4.812%        |
| 26        | 17.371      | 2495          | 2499        | 2505         | rBV2     | 22315          | 30362         | 1.19%           | 0.180%        |
| 27        | 17.694      | 2549          | 2554        | 2565         | rBV      | 72218          | 114211        | 4.47%           | 0.679%        |
| 28        | 18.046      | 2609          | 2614        | 2623         | rBV2     | 79561          | 131280        | 5.14%           | 0.780%        |
| 29        | 19.321      | 2827          | 2831        | 2843         | rBV4     | 35970          | 61810         | 2.42%           | 0.367%        |
| 30        | 19.774      | 2902          | 2908        | 2915         | rBV      | 2099446        | 2554699       | 100.00%         | 15.178%       |
| 31        | 20.296      | 2992          | 2997        | 3007         | rBV      | 41325          | 62923         | 2.46%           | 0.374%        |
| 32        | 21.031      | 3117          | 3122        | 3131         | rBV      | 252507         | 322949        | 12.64%          | 1.919%        |
| 33        | 21.324      | 3166          | 3172        | 3181         | rBV      | 678048         | 901243        | 35.28%          | 5.354%        |
| 34        | 22.511      | 3369          | 3374        | 3379         | rBV      | 27220          | 39438         | 1.54%           | 0.234%        |

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Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 23.604 | 3553 | 3560 | 3570 | rVB2 | 452349 | 875160 | 34.26% | 5.199% |
| 36 | 24.244 | 3664 | 3669 | 3672 | rBV5 | 14514  | 26460  | 1.04%  | 0.157% |
| 37 | 26.165 | 3991 | 3996 | 4003 | rVB8 | 23218  | 55775  | 2.18%  | 0.331% |

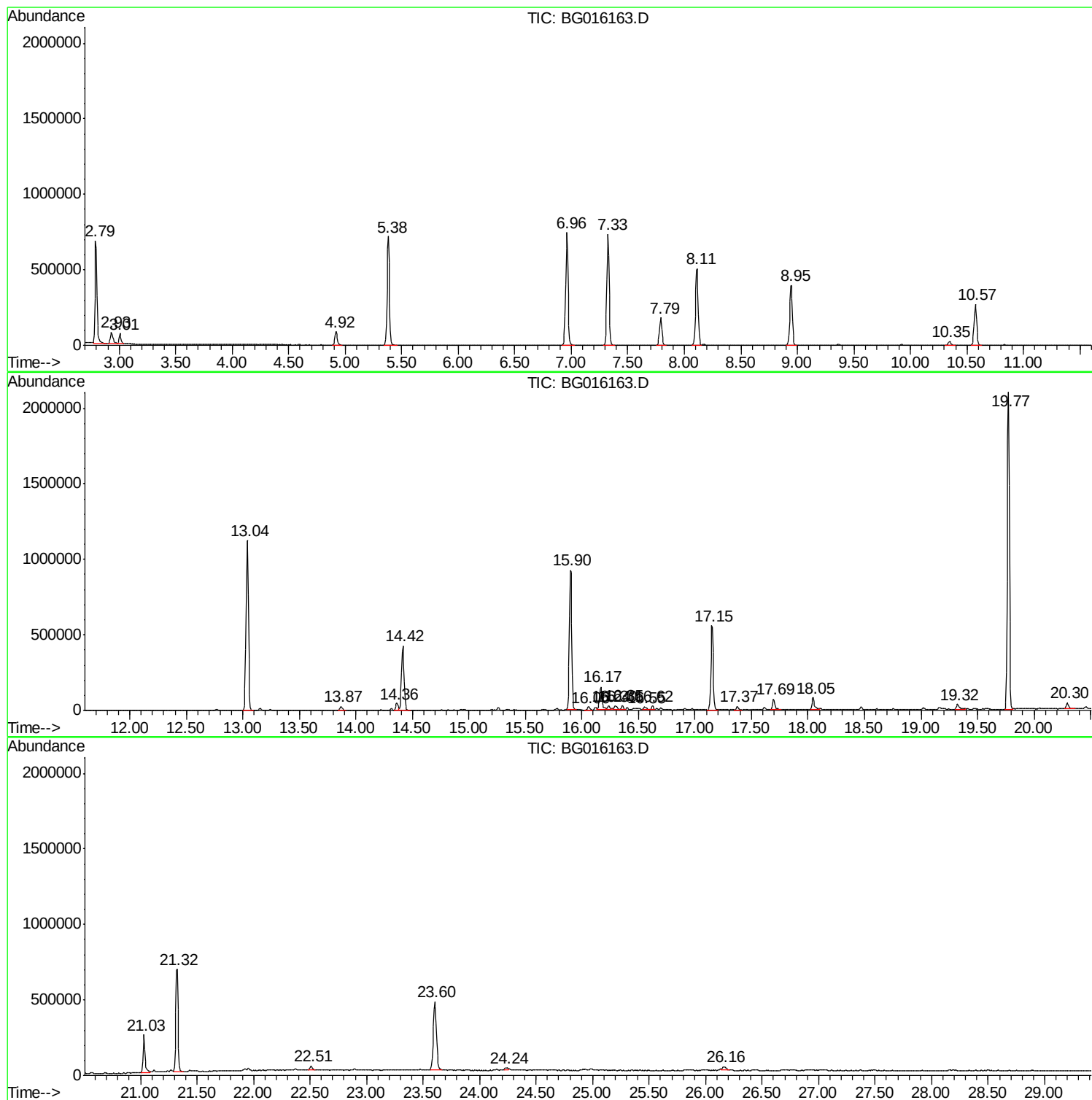
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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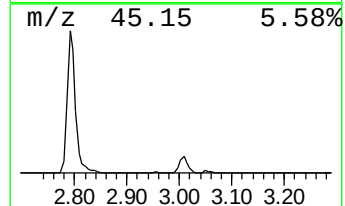
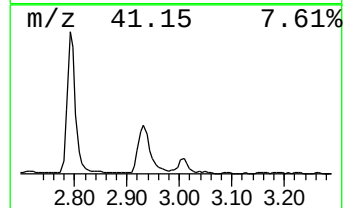
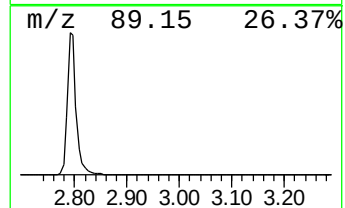
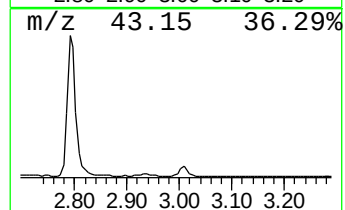
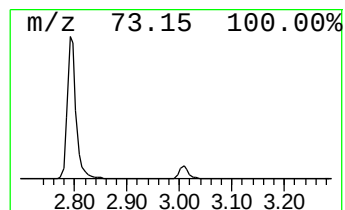
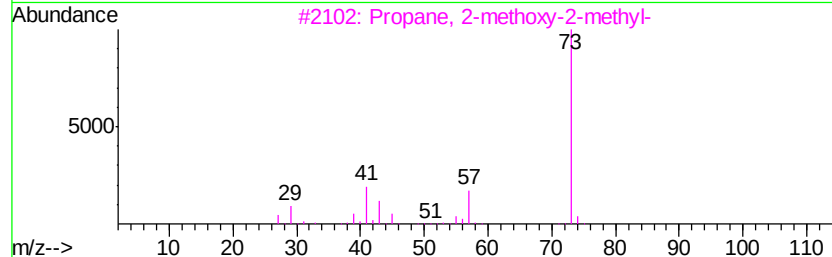
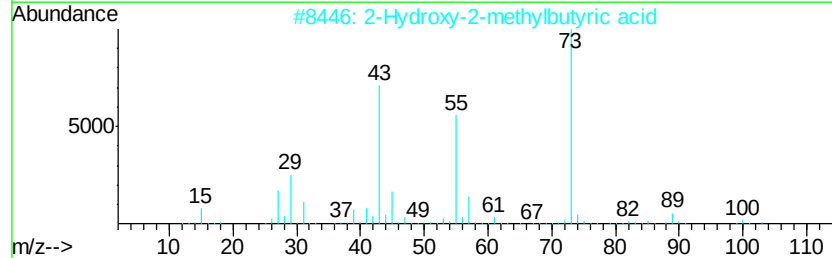
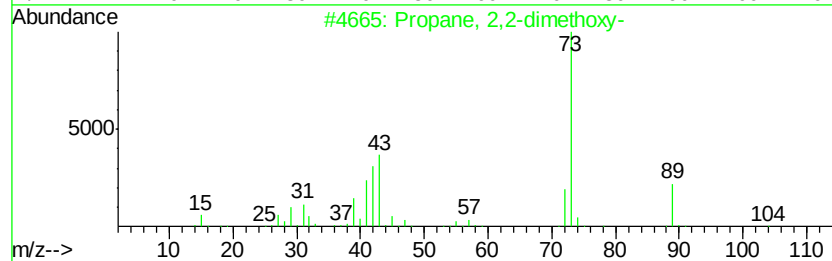
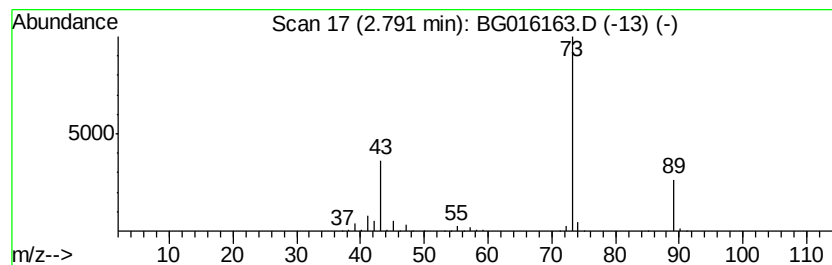
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown2.79 Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 2.79 | 56.70 ng | 790661 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy-        | 104 | C5H12O2 | 000077-76-9 | 45   |
| 2    |    | 2-Hydroxy-2-methylbutyric acid | 118 | C5H10O3 | 003739-30-8 | 9    |
| 3    |    | Propane, 2-methoxy-2-methyl-   | 88  | C5H12O  | 001634-04-4 | 9    |
| 4    |    | Methane, isothiocyanato-       | 73  | C2H3NS  | 000556-61-6 | 7    |
| 5    |    | N-Ethylformamide               | 73  | C3H7NO  | 000627-45-2 | 7    |



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BNA\_G  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

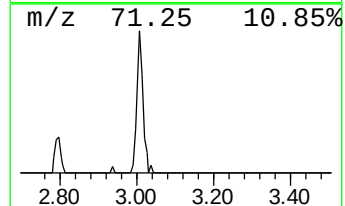
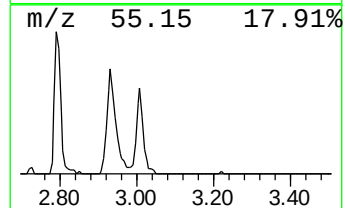
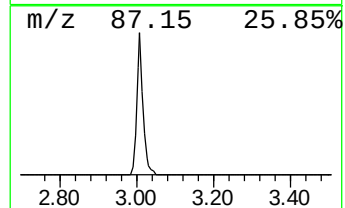
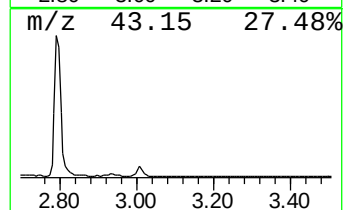
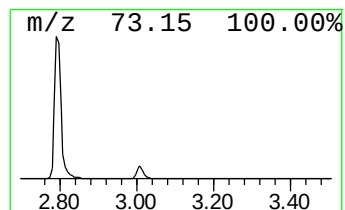
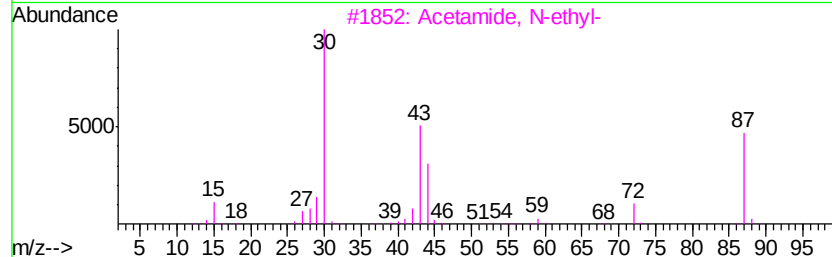
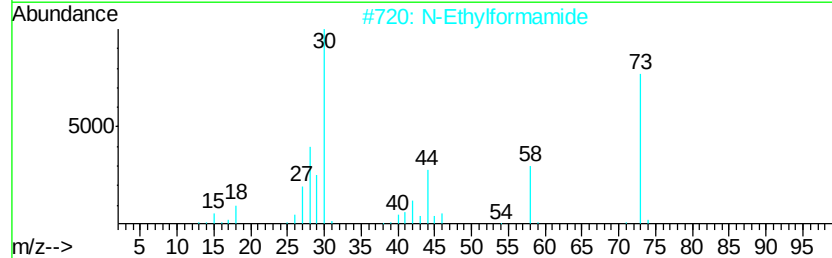
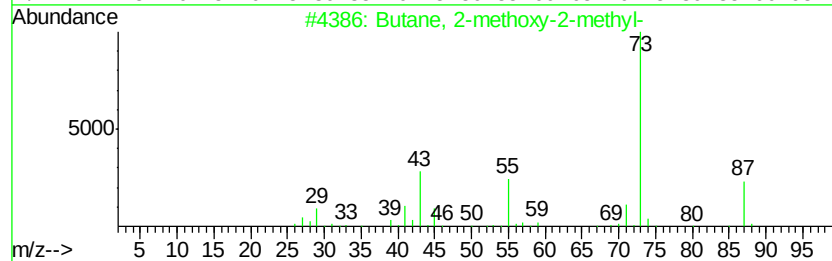
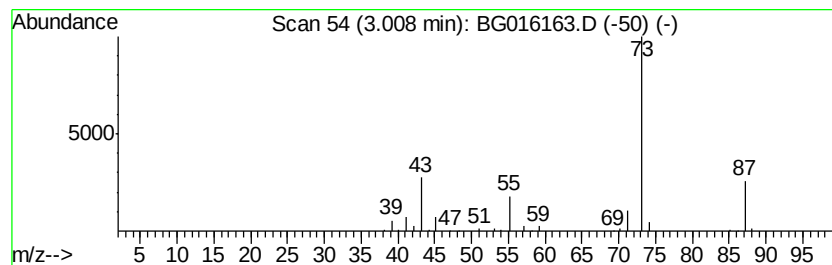
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 8

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 3.01 | 5.14 ng | 71737 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2       |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 3       |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |
| 4       |   | 3-Pentanol, 2,4-dimethyl-   | 116 | C7H16O  | 000600-36-2 | 9    |
| 5       |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 9    |



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Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

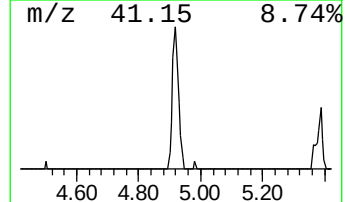
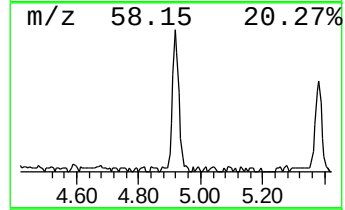
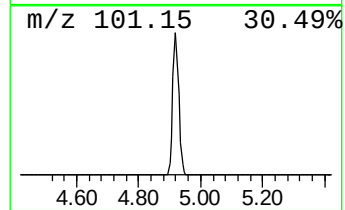
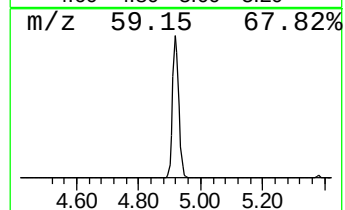
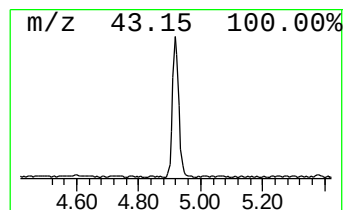
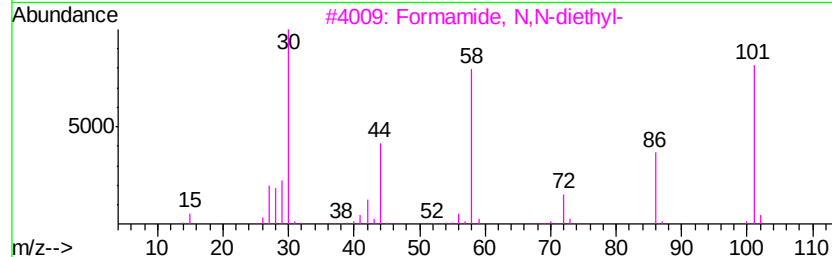
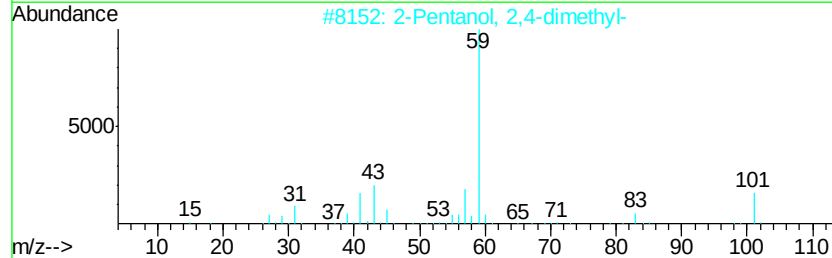
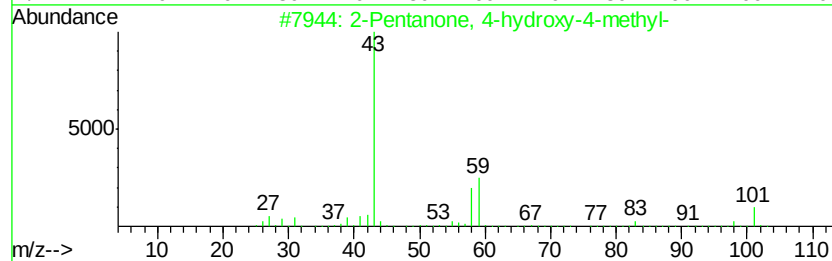
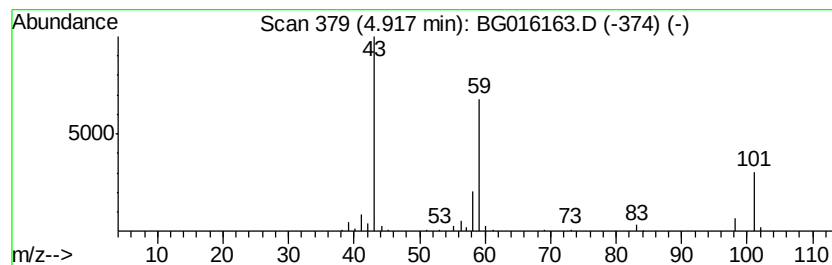
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.92 | 8.50 ng | 118584 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|---------|---|----------------------------------|-----|---------|-------------|------|
| 1       |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2       |   | 2-Pentanol, 2,4-dimethyl-        | 116 | C7H16O  | 000625-06-9 | 17   |
| 3       |   | Formamide, N,N-diethyl-          | 101 | C5H11NO | 000617-84-5 | 16   |
| 4       |   | Propanamide, N-ethyl-            | 101 | C5H11NO | 005129-72-6 | 14   |
| 5       |   | Morpholine, 4-methyl-            | 101 | C5H11NO | 000109-02-4 | 12   |



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ALS Vial : 25 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

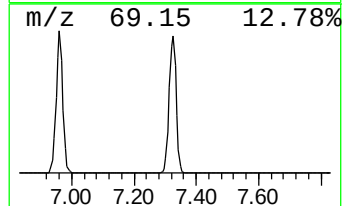
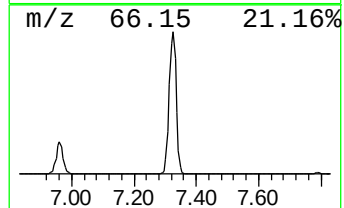
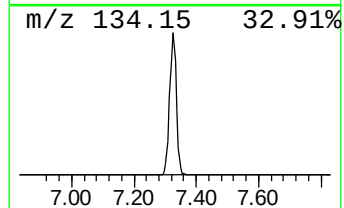
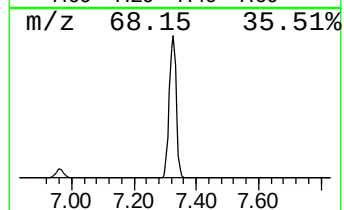
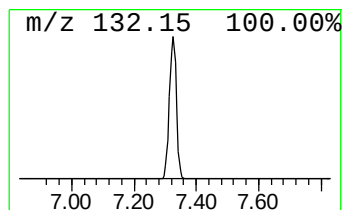
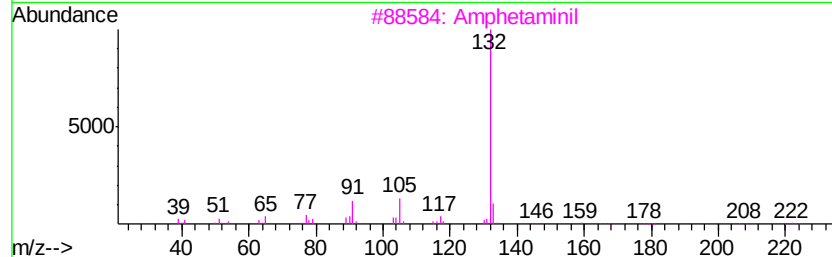
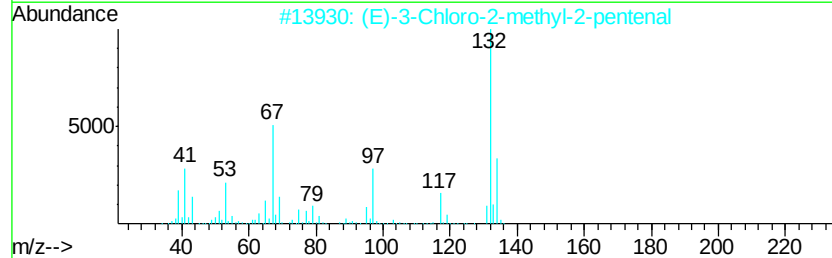
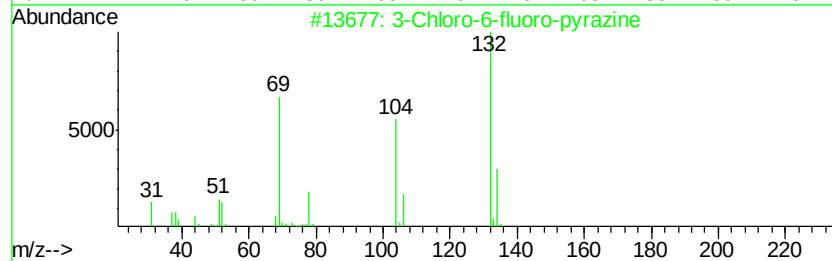
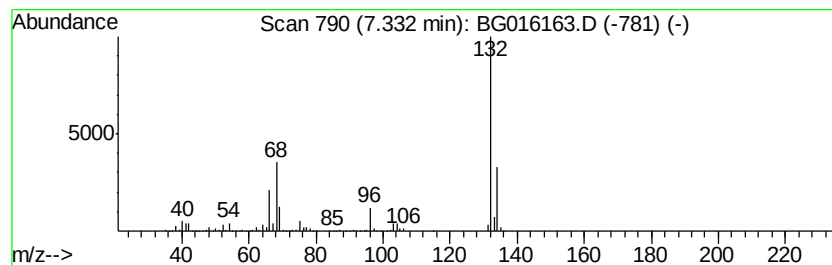
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown7.33 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.33 | 81.19 ng | 1132190 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 35   |
| 3    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 12   |
| 4    |    | Tranlylcypromine-propionyl       | 189 | C12H15NO  | 1000123-86-3 | 9    |
| 5    |    | Xenon                            | 132 | Xe        | 007440-63-3  | 9    |



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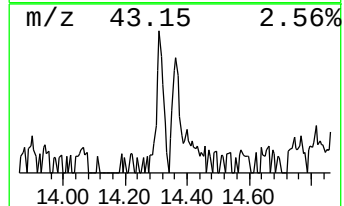
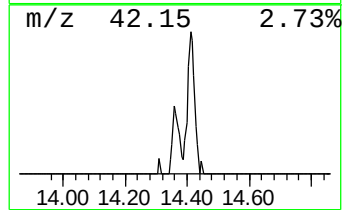
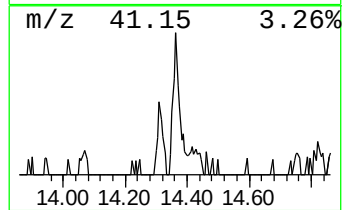
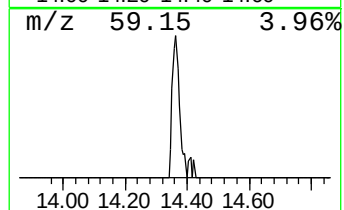
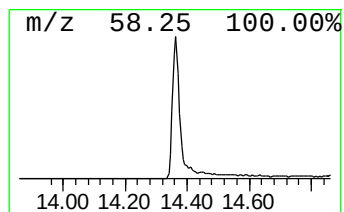
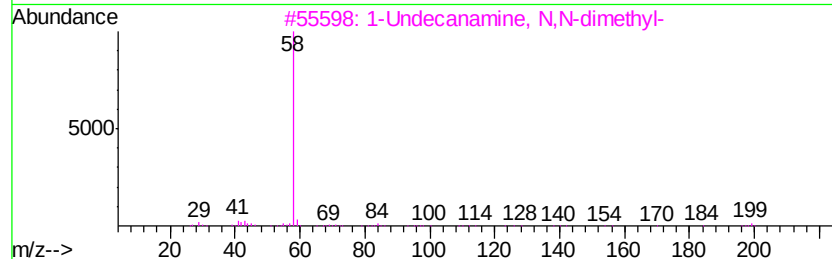
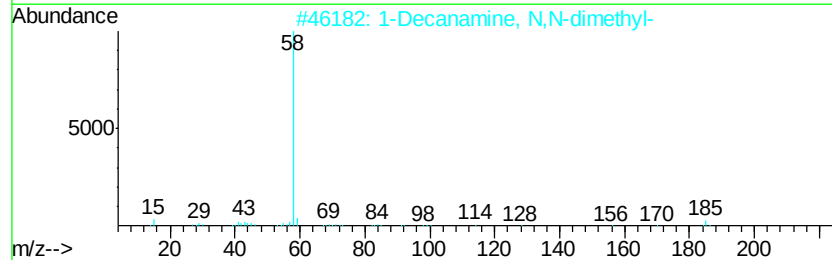
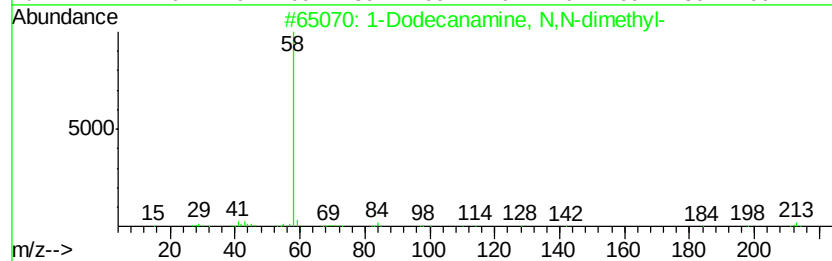
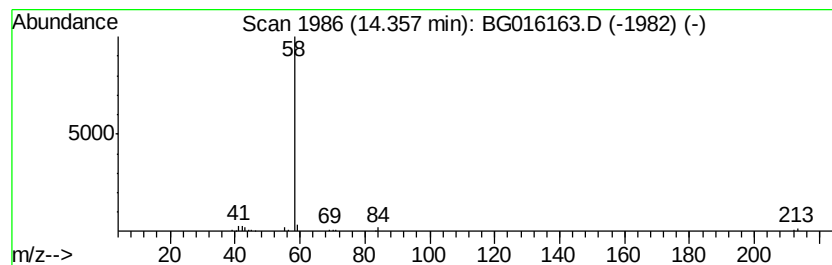
Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 1-Dodecanamine, N,N-dimethyl- Concentration Rank 11

| R.T.    | EstConc | Area                           | Relative to ISTD | R.T.           |
|---------|---------|--------------------------------|------------------|----------------|
| 14.36   | 2.04 ng | 66769                          | Acenaphthene-d10 | 14.42          |
| Hit# of | 5       | Tentative ID                   | MW MolForm       | CAS# Qual      |
| 1       |         | 1-Dodecanamine, N,N-dimethyl-  | 213 C14H31N      | 000112-18-5 91 |
| 2       |         | 1-Decanamine, N,N-dimethyl-    | 185 C12H27N      | 001120-24-7 78 |
| 3       |         | 1-Undecanamine, N,N-dimethyl-  | 199 C13H29N      | 017373-28-3 72 |
| 4       |         | 1-Tridecanamine, N,N-dimethyl- | 227 C15H33N      | 017373-29-4 72 |
| 5       |         | Tetradonium Bromide            | 335 C17H38BrN    | 001119-97-7 72 |





Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016163.D  
Acq On : 27 Mar 2015 5:38  
Operator : TP/IZ  
Sample : G1653-14  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

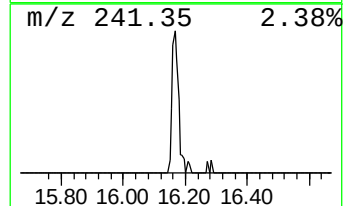
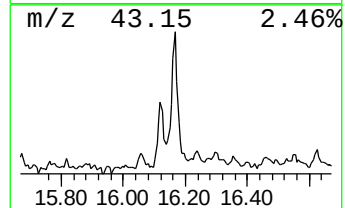
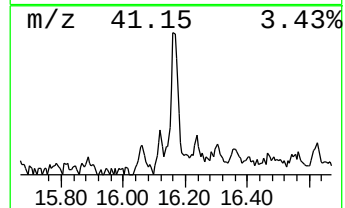
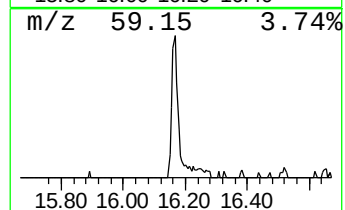
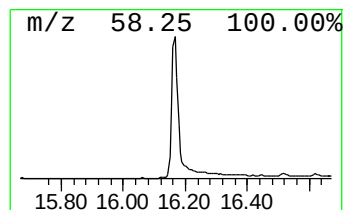
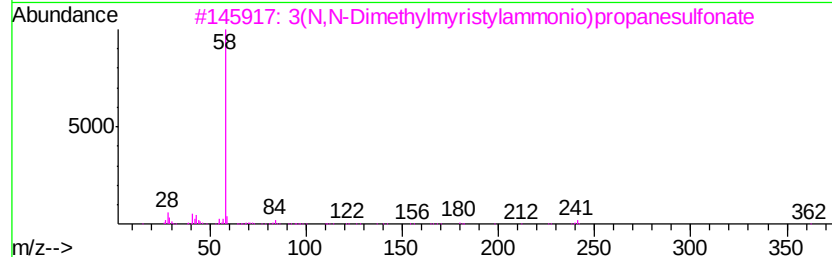
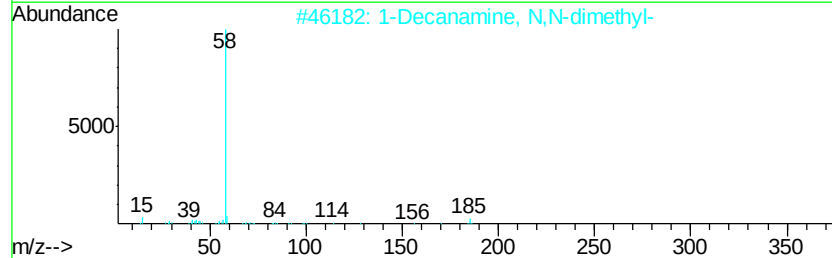
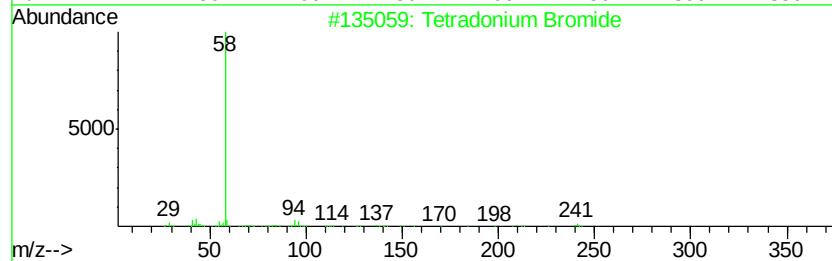
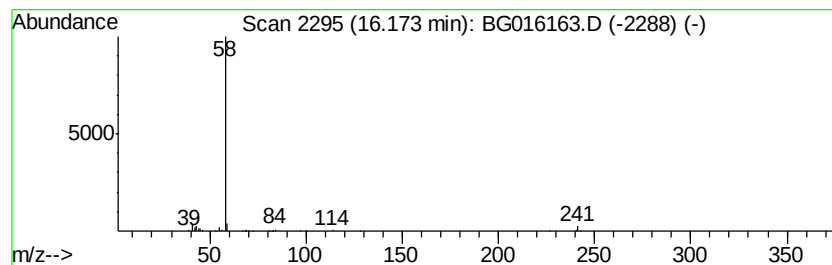
Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Tetradonium Bromide Concentration Rank 7

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 16.17     | 5.84 ng                             | 236495 | Phenanthrene-d10 | 17.15       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Tetradonium Bromide                 | 335    | C17H38BrN        | 001119-97-7 | 78   |
| 2         | 1-Decanamine, N,N-dimethyl-         | 185    | C12H27N          | 001120-24-7 | 72   |
| 3         | 3(N,N-Dimethylmyristylammonio)pr... | 363    | C19H41N03S       | 014933-09-6 | 64   |
| 4         | 1-Dodecanamine, N,N-dimethyl-       | 213    | C14H31N          | 000112-18-5 | 56   |
| 5         | 1-Hexadecanamine, N,N-dimethyl-     | 269    | C18H39N          | 000112-69-6 | 56   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016163.D  
Acq On : 27 Mar 2015 5:38  
Operator : TP/IZ  
Sample : G1653-14  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

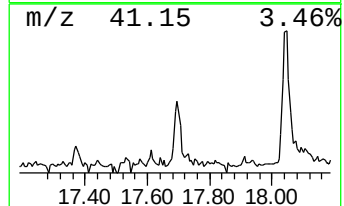
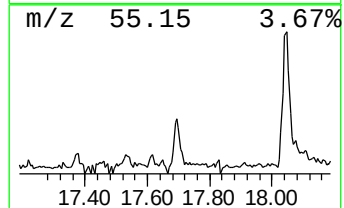
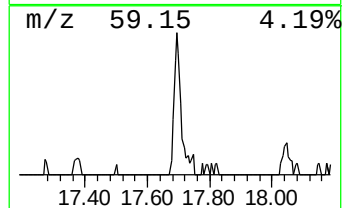
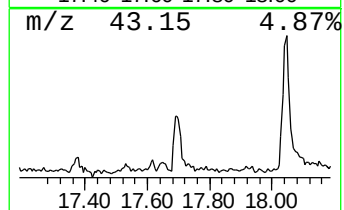
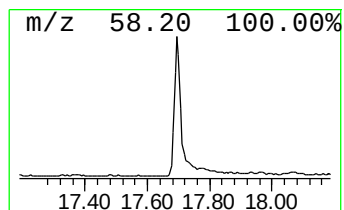
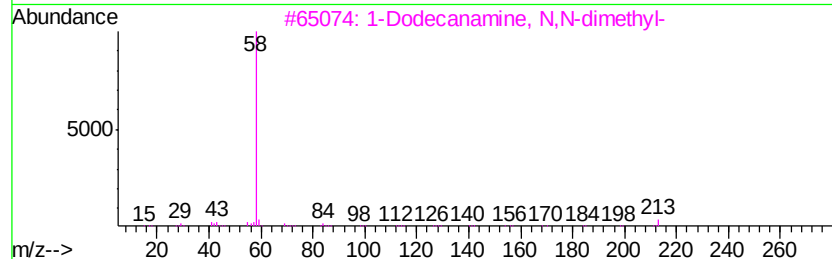
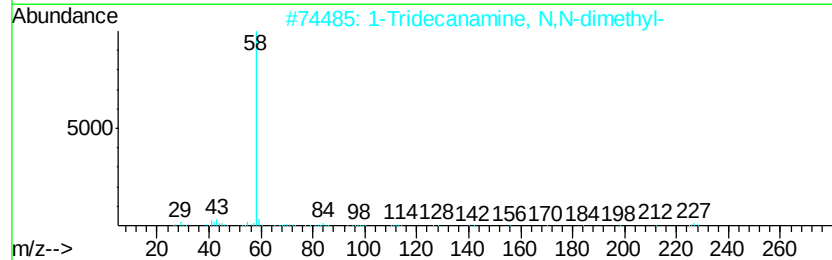
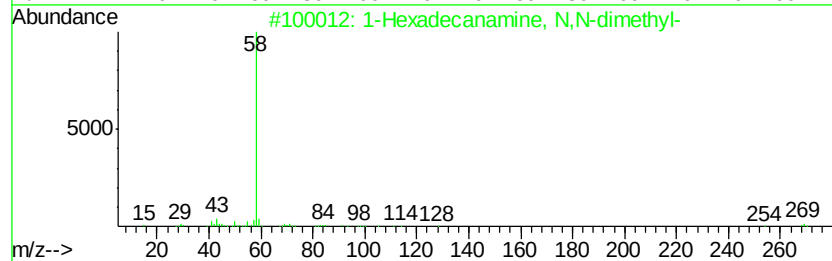
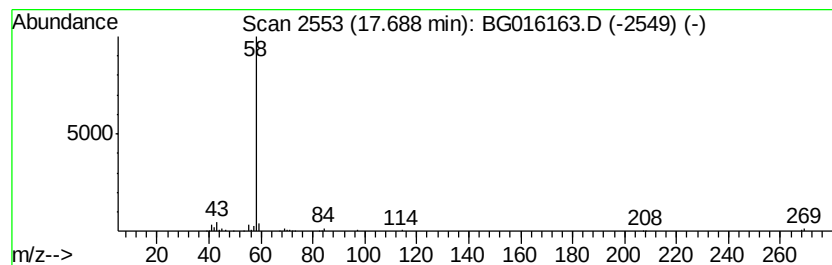
Instrument :  
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ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 1-Hexadecanamine, N,N-dimet... Concentration Rank 10

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 17.69     | 2.82 ng                             | 114211 | Phenanthrene-d10 | 17.15       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 1-Hexadecanamine, N,N-dimethyl-     | 269    | C18H39N          | 000112-69-6 | 91   |
| 2         | 1-Tridecanamine, N,N-dimethyl-      | 227    | C15H33N          | 017373-29-4 | 72   |
| 3         | 1-Dodecanamine, N,N-dimethyl-       | 213    | C14H31N          | 000112-18-5 | 64   |
| 4         | 1-Octadecanamine, N,N-dimethyl-     | 297    | C20H43N          | 000124-28-7 | 64   |
| 5         | 3(N,N-Dimethylmyristylammonio)pr... | 363    | C19H41N03S       | 014933-09-6 | 64   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016163.D  
Acq On : 27 Mar 2015 5:38  
Operator : TP/IZ  
Sample : G1653-14  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

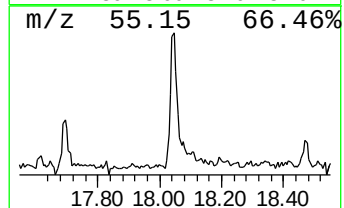
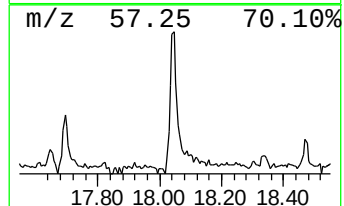
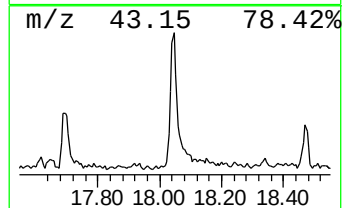
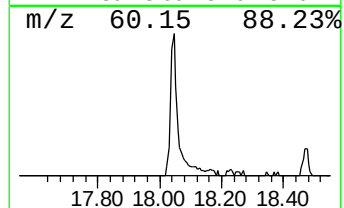
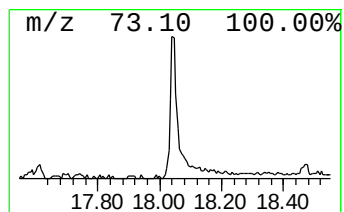
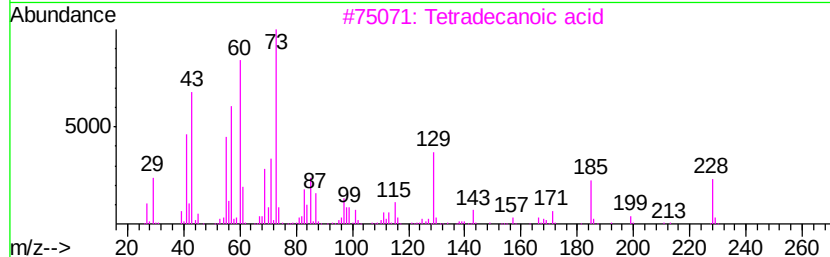
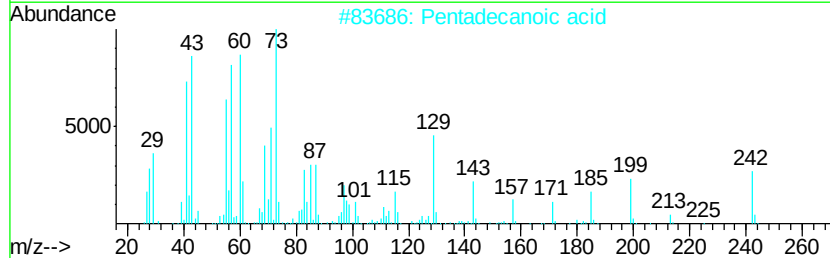
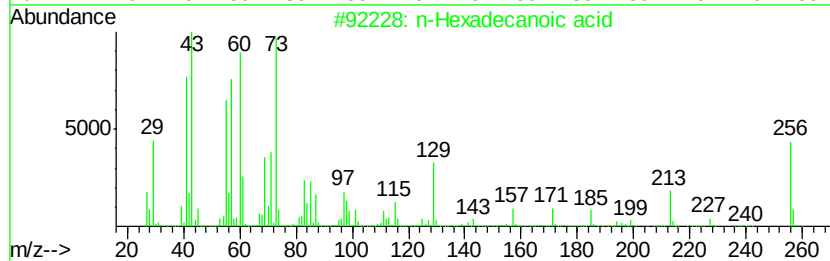
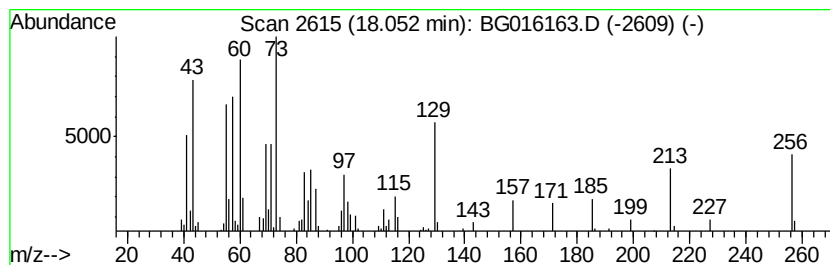
Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 n-Hexadecanoic acid Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD    |     | R.T.     |             |      |
|-------|---------|--------|---------------------|-----|----------|-------------|------|
| 18.05 | 3.24 ng | 131280 | Phenanthrene-d10    |     | 17.15    |             |      |
| Hit#  | of      | 5      | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
| 1     |         |        | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2     |         |        | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 81   |
| 3     |         |        | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 72   |
| 4     |         |        | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 70   |
| 5     |         |        | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 62   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016163.D  
Acq On : 27 Mar 2015 5:38  
Operator : TP/IZ  
Sample : G1653-14  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

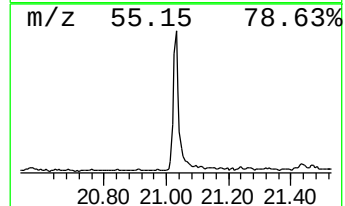
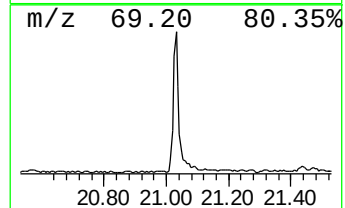
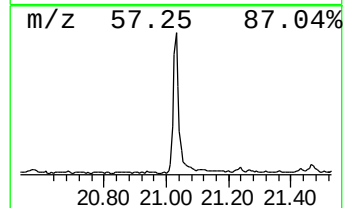
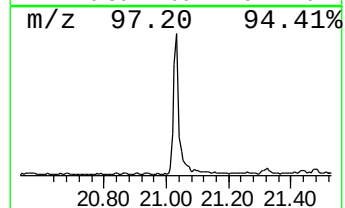
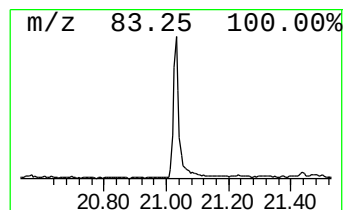
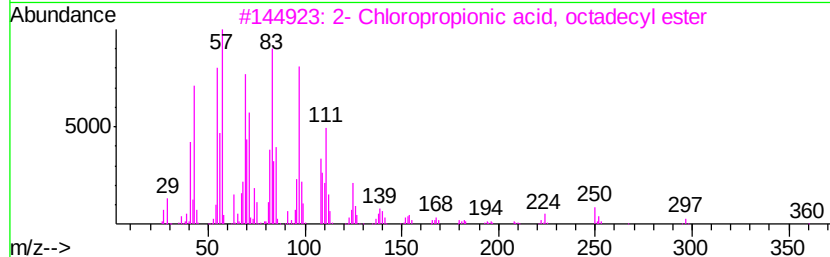
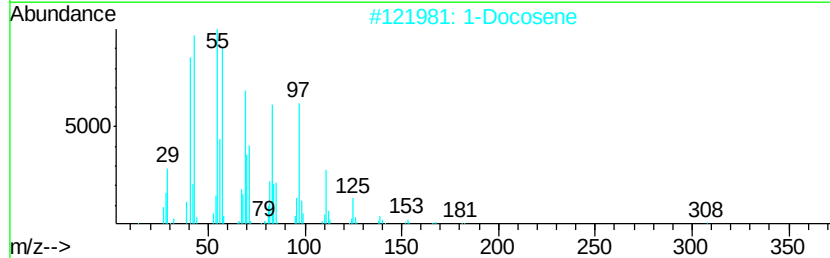
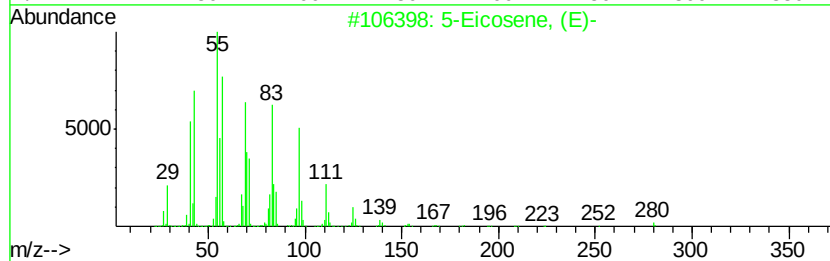
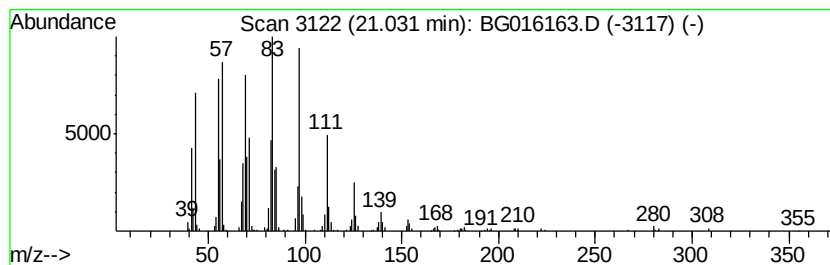
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 5-Eicosene, (E)- Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.03 | 7.17 ng | 322949 | Chrysene-d12     | 21.32 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 5-Eicosene, (E)-                    | 280 | C20H40     | 074685-30-6 | 98   |
| 2    |    | 1-Docosene                          | 308 | C22H44     | 001599-67-3 | 93   |
| 3    |    | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2 | 088104-31-8 | 93   |
| 4    |    | 1-Heneicosyl formate                | 340 | C22H44O2   | 077899-03-7 | 91   |
| 5    |    | Cyclopentadecane                    | 210 | C15H30     | 000295-48-7 | 90   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016163.D  
Acq On : 27 Mar 2015 5:38  
Operator : TP/IZ  
Sample : G1653-14  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| unknown2.79          | 2.79  | 56.7    | ng    | 790661   | 1                     | 7.79  | 278887 | 20.0 |
| Butane, 2-methoxy... | 3.01  | 5.1     | ng    | 71737    | 1                     | 7.79  | 278887 | 20.0 |
| 2-Pentanone, 4-hy... | 4.92  | 8.5     | ng    | 118584   | 1                     | 7.79  | 278887 | 20.0 |
| unknown7.33          | 7.33  | 81.2    | ng    | 1132190  | 1                     | 7.79  | 278887 | 20.0 |
| 1-Dodecanamine, N... | 14.36 | 2.0     | ng    | 66769    | 3                     | 14.42 | 653553 | 20.0 |
| Tetradonium Bromide  | 16.17 | 5.8     | ng    | 236495   | 4                     | 17.15 | 809967 | 20.0 |
| 1-Hexadecanamine,... | 17.69 | 2.8     | ng    | 114211   | 4                     | 17.15 | 809967 | 20.0 |
| n-Hexadecanoic acid  | 18.05 | 3.2     | ng    | 131280   | 4                     | 17.15 | 809967 | 20.0 |
| 5-Eicosene, (E)-     | 21.03 | 7.2     | ng    | 322949   | 5                     | 21.32 | 901243 | 20.0 |