

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF070316\  
 Data File : BF088661.D  
 Acq On : 3 Jul 2016 21:52  
 Operator : UM/SJ  
 Sample : H3817-05  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BS-1C-6.5-7FT

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\_F\METHODS\8270-BF063016.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         | 1.644       | 3             | 5           | 8            | rVB      | 4841214        | 6886871       | 100.00%         | 19.325%       |
| 2         | 1.850       | 21            | 23          | 36           | rVV      | 148194         | 219693        | 3.19%           | 0.616%        |
| 3         | 2.032       | 36            | 39          | 42           | rVB      | 61463          | 86861         | 1.26%           | 0.244%        |
| 4         | 3.027       | 124           | 126         | 143          | rBV      | 77291          | 150136        | 2.18%           | 0.421%        |
| 5         | 4.958       | 292           | 295         | 301          | rBV      | 183540         | 270376        | 3.93%           | 0.759%        |
| 6         | 5.381       | 329           | 332         | 335          | rBV      | 2677662        | 2950637       | 42.84%          | 8.280%        |
| 7         | 6.422       | 420           | 423         | 425          | rBV      | 2829954        | 3245729       | 47.13%          | 9.108%        |
| 8         | 6.547       | 431           | 434         | 437          | rBV      | 2507290        | 3240269       | 47.05%          | 9.092%        |
| 9         | 6.787       | 452           | 455         | 456          | rBV      | 586724         | 705951        | 10.25%          | 1.981%        |
| 10        | 6.936       | 466           | 468         | 471          | rVB      | 2077163        | 2373790       | 34.47%          | 6.661%        |
| 11        | 7.347       | 501           | 504         | 506          | rBV      | 1889610        | 1961475       | 28.48%          | 5.504%        |
| 12        | 8.067       | 564           | 567         | 569          | rBV      | 1083000        | 942669        | 13.69%          | 2.645%        |
| 13        | 9.142       | 658           | 661         | 664          | rBB      | 2516376        | 2967055       | 43.08%          | 8.326%        |
| 14        | 9.542       | 693           | 696         | 698          | rBV      | 295764         | 374574        | 5.44%           | 1.051%        |
| 15        | 9.816       | 717           | 720         | 722          | rBV      | 1176169        | 1026935       | 14.91%          | 2.882%        |
| 16        | 10.605      | 786           | 789         | 791          | rBV      | 1927338        | 2145501       | 31.15%          | 6.020%        |
| 17        | 11.176      | 837           | 839         | 841          | rBV      | 107718         | 100451        | 1.46%           | 0.282%        |
| 18        | 11.291      | 847           | 849         | 851          | rVB      | 1191063        | 1022523       | 14.85%          | 2.869%        |
| 19        | 11.359      | 851           | 855         | 859          | rVB4     | 29358          | 70232         | 1.02%           | 0.197%        |
| 20        | 12.879      | 985           | 988         | 990          | rBV      | 2725637        | 2562338       | 37.21%          | 7.190%        |
| 21        | 13.554      | 1045          | 1047        | 1049         | rBV      | 160747         | 136301        | 1.98%           | 0.382%        |
| 22        | 13.759      | 1062          | 1065        | 1066         | rBV      | 108386         | 133731        | 1.94%           | 0.375%        |
| 23        | 13.817      | 1066          | 1070        | 1076         | rVV      | 83142          | 215586        | 3.13%           | 0.605%        |
| 24        | 13.919      | 1076          | 1079        | 1081         | rVB      | 888852         | 948043        | 13.77%          | 2.660%        |
| 25        | 14.468      | 1124          | 1127        | 1137         | rBV      | 25280          | 69414         | 1.01%           | 0.195%        |
| 26        | 14.765      | 1151          | 1153        | 1156         | rBV      | 90958          | 113080        | 1.64%           | 0.317%        |
| 27        | 15.302      | 1197          | 1200        | 1203         | rBV      | 480831         | 716978        | 10.41%          | 2.012%        |

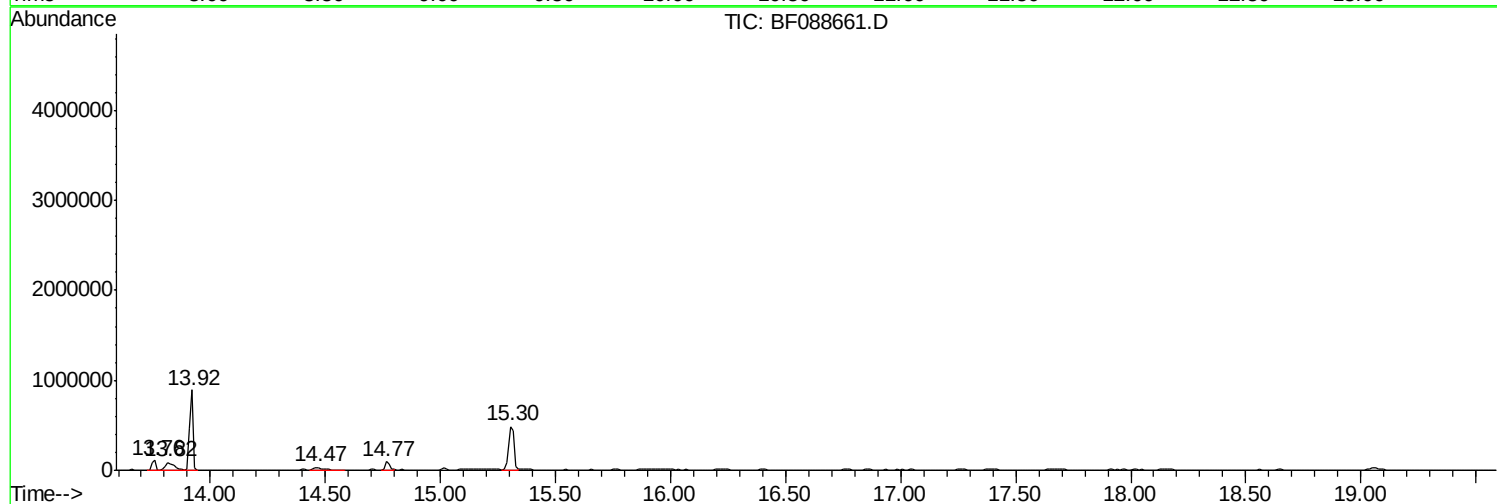
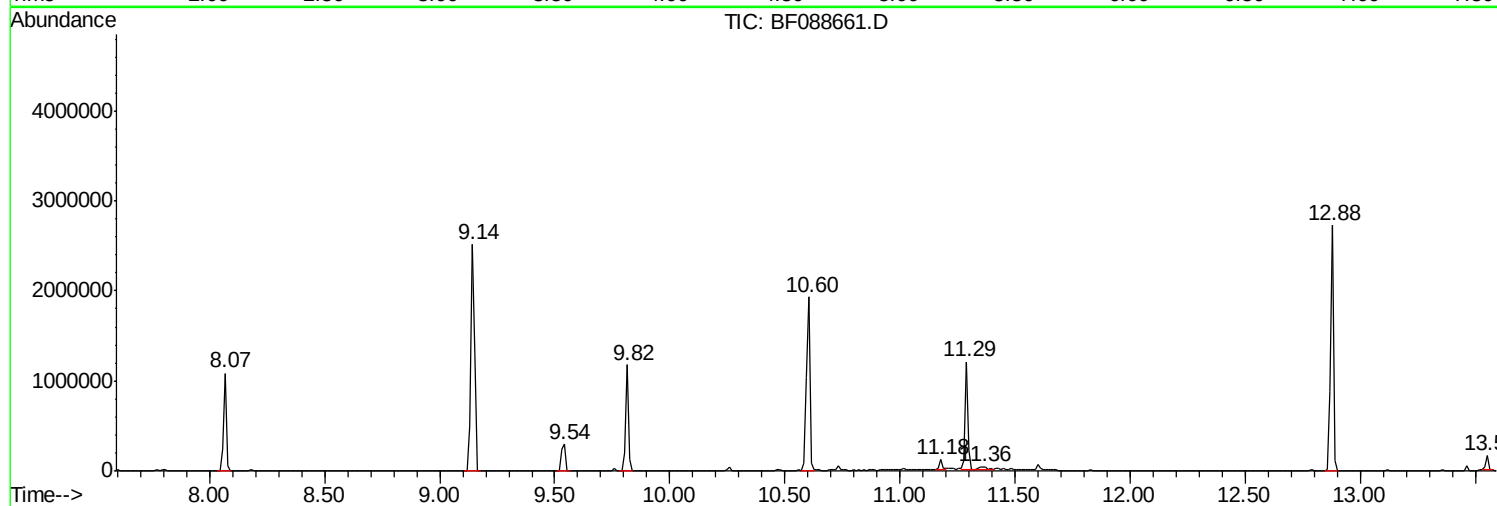
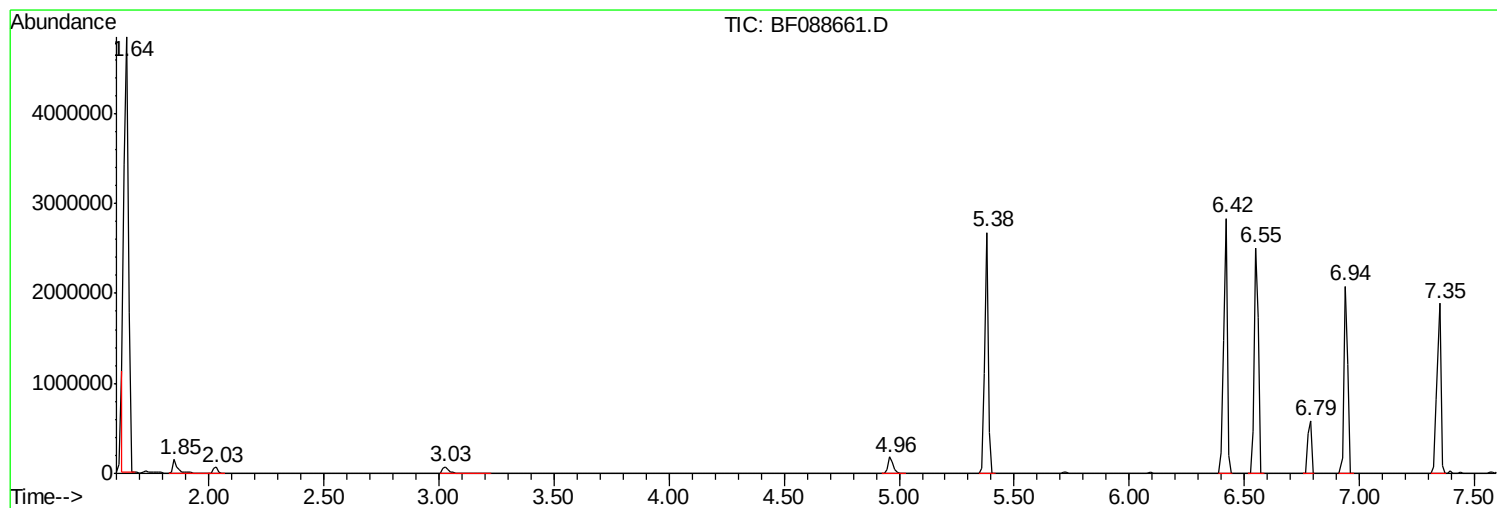
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BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P



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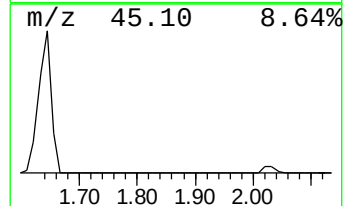
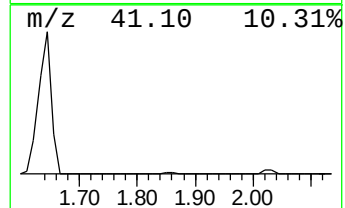
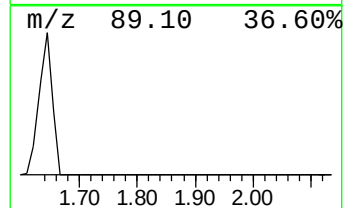
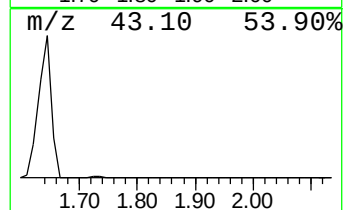
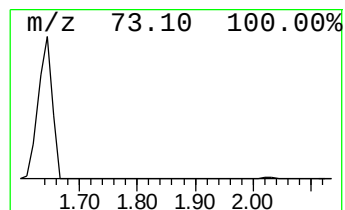
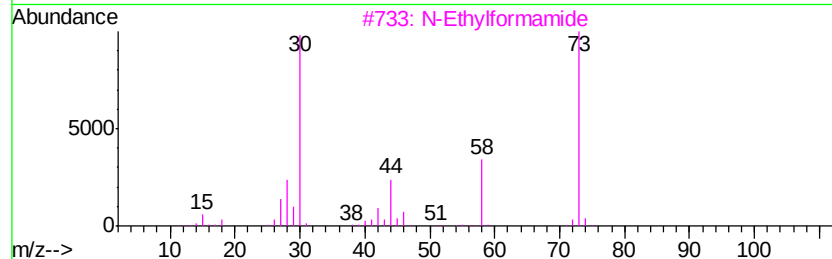
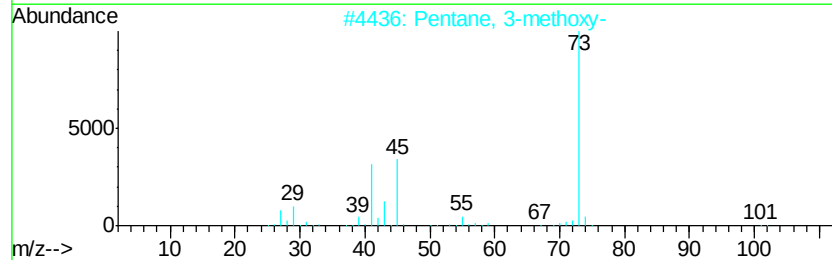
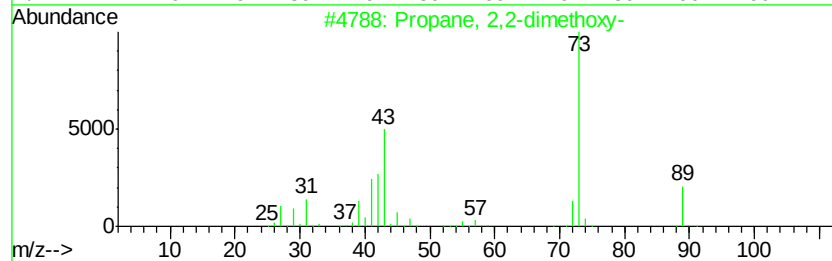
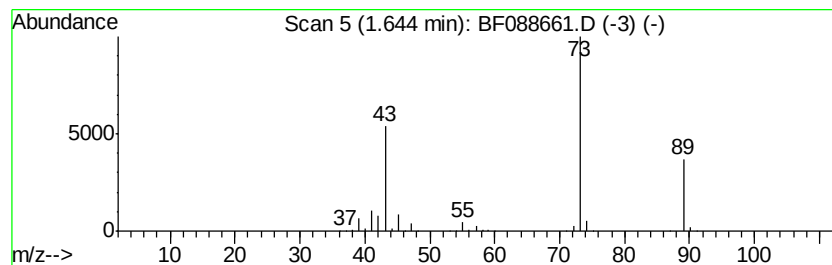
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown1.64 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 1.64 | 195.11 ng | 6886870 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy- | 104 | C5H12O2 | 000077-76-9 | 38   |
| 2    |    | Pentane, 3-methoxy-     | 102 | C6H14O  | 036839-67-5 | 12   |
| 3    |    | N-Ethylformamide        | 73  | C3H7NO  | 000627-45-2 | 9    |
| 4    |    | 1-Propanol, 2-methyl-   | 74  | C4H10O  | 000078-83-1 | 9    |
| 5    |    | 1,3-Diaminoguanidine    | 89  | CH7N5   | 004364-78-7 | 9    |



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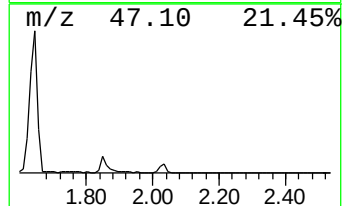
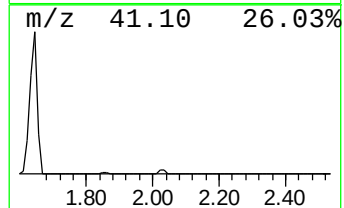
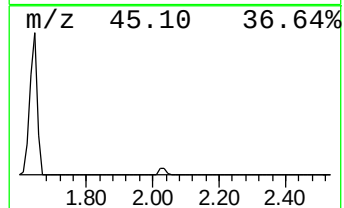
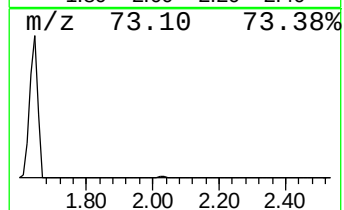
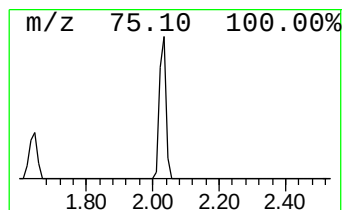
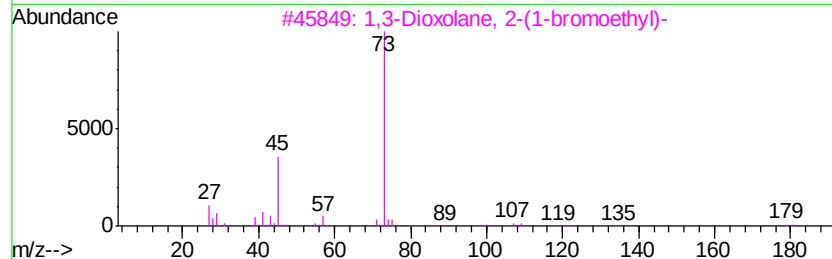
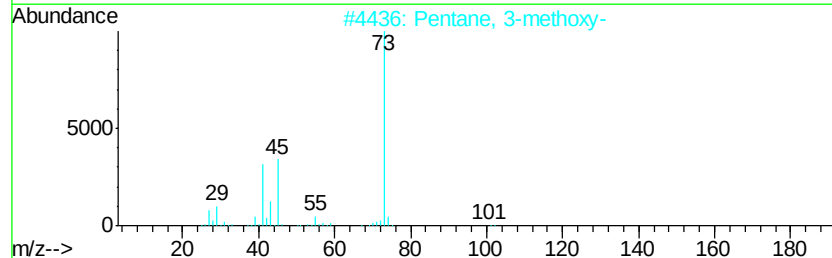
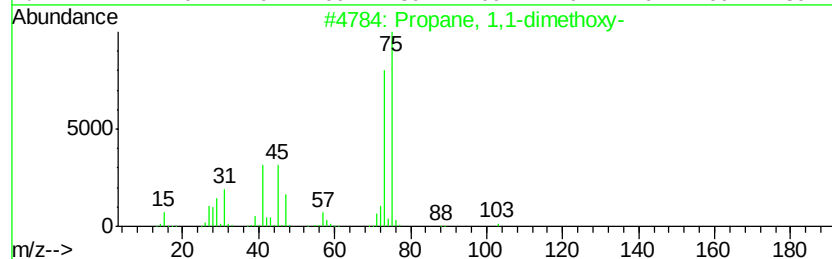
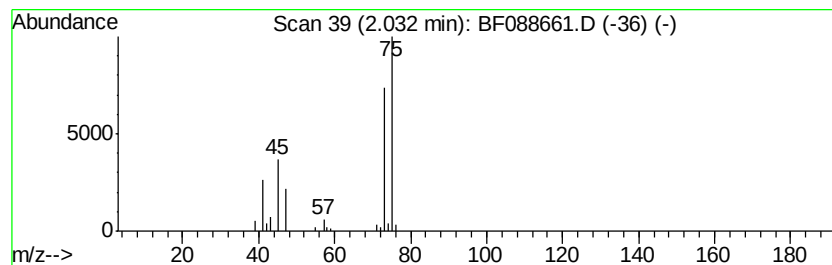
Instrument :  
BNA\_F  
ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 11

| R.T.    | EstConc                          | Area         | Relative to ISTD       | R.T.      |
|---------|----------------------------------|--------------|------------------------|-----------|
| 2.03    | 2.46 ng                          | 86861        | 1,4-Dichlorobenzene-d4 | 6.79      |
| Hit# of | 5                                | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | Propane, 1,1-dimethoxy-          | 104 C5H12O2  | 004744-10-9            | 78        |
| 2       | Pentane, 3-methoxy-              | 102 C6H14O   | 036839-67-5            | 14        |
| 3       | 1,3-Dioxolane, 2-(1-bromoethyl)- | 180 C5H9BrO2 | 005267-73-2            | 10        |
| 4       | 1,3-Dioxolane, 2-(chloromethyl)- | 122 C4H7ClO2 | 002568-30-1            | 9         |
| 5       | Methane, isothiocyanato-         | 73 C2H3NS    | 000556-61-6            | 7         |



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Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

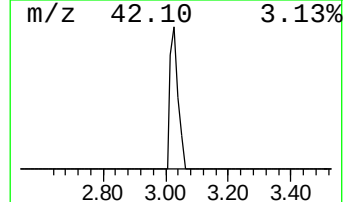
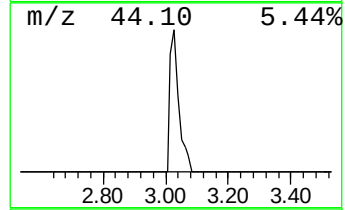
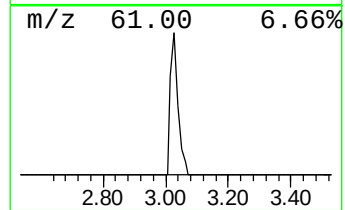
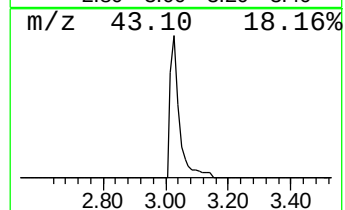
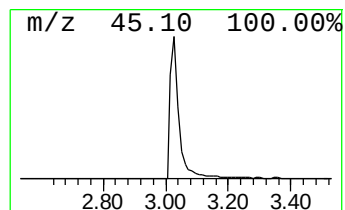
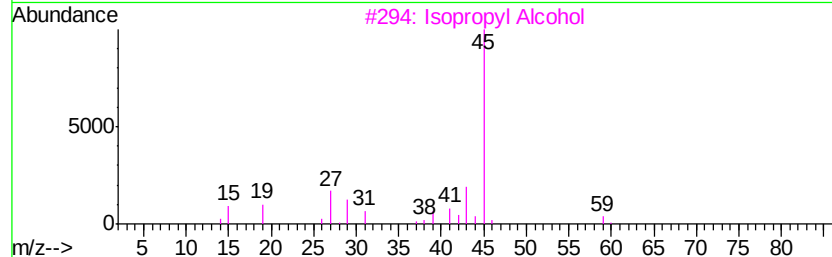
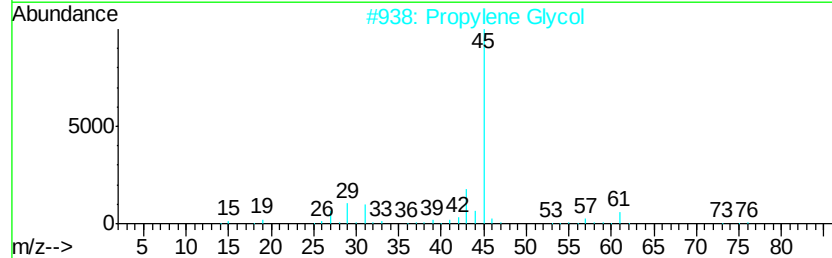
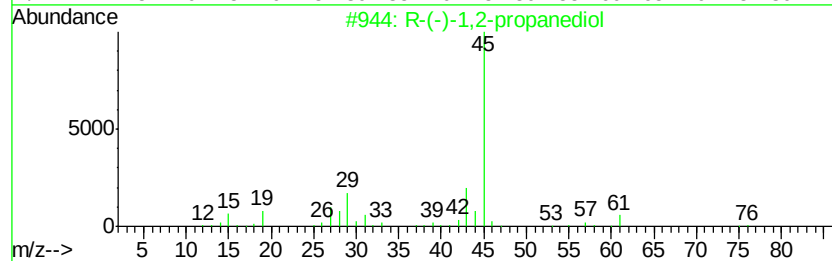
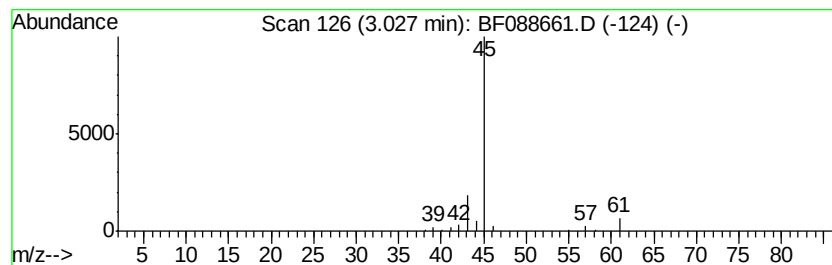
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 R-(-)-1,2-propanediol Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 3.03 | 4.25 ng | 150136 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | R-(-)-1,2-propanediol               | 76  | C3H8O2  | 004254-14-2 | 64   |
| 2    |    | Propylene Glycol                    | 76  | C3H8O2  | 000057-55-6 | 9    |
| 3    |    | Isopropyl Alcohol                   | 60  | C3H8O   | 000067-63-0 | 9    |
| 4    |    | Propanoic acid, 2-hydroxy-, meth... | 104 | C4H8O3  | 002155-30-8 | 9    |
| 5    |    | (S)-(+)-1,2-Propanediol             | 76  | C3H8O2  | 004254-15-3 | 9    |



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BNA\_F  
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BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

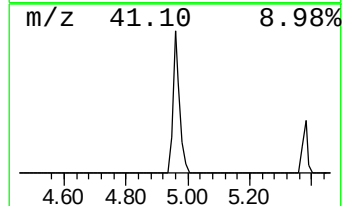
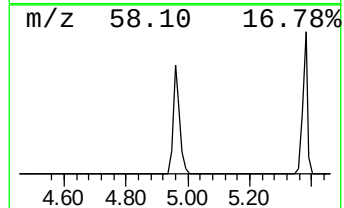
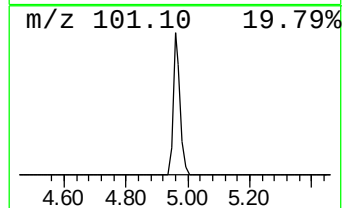
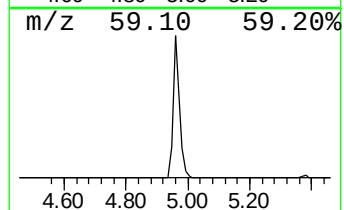
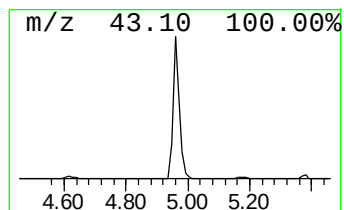
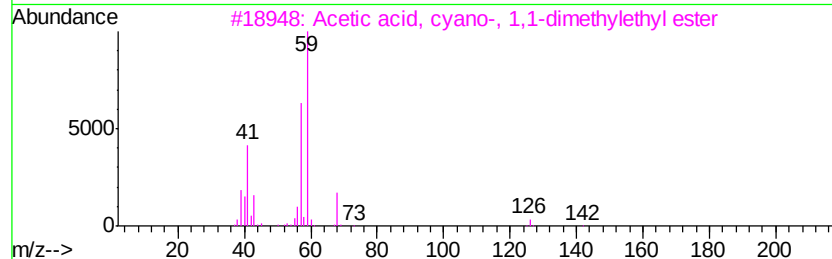
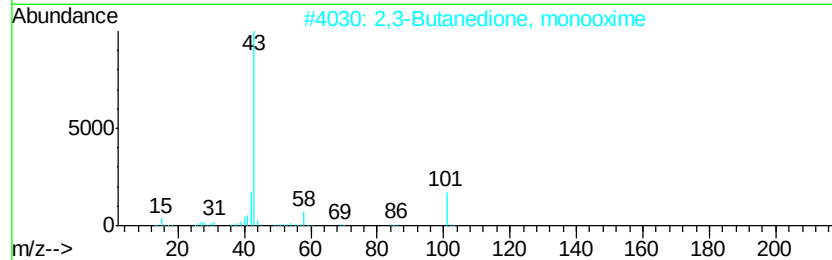
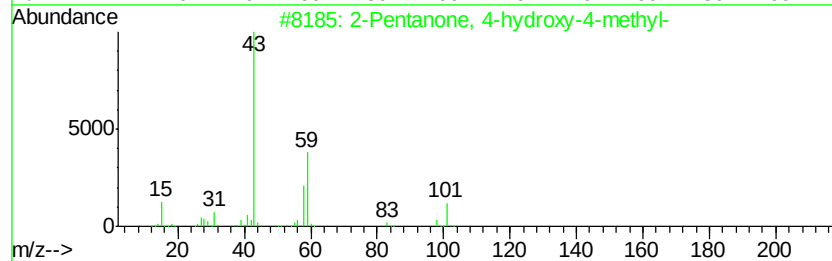
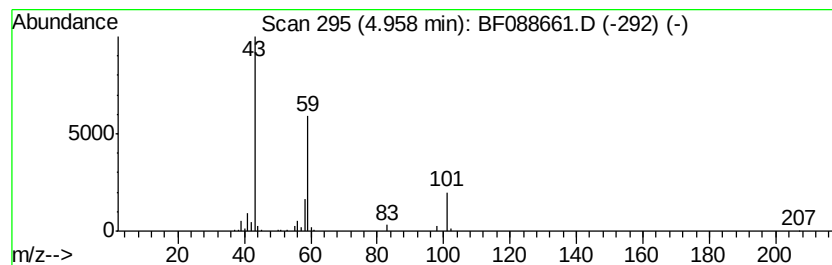
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.96 | 7.66 ng | 270376 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of                                   | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|--------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2      | 000123-42-2 | 53      |      |      |
| 2    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2      | 000057-71-6 | 25      |      |      |
| 3    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2     | 001116-98-9 | 23      |      |      |
| 4    | Butane, 1-ethoxy-                    | 102 | C6H14O       | 000628-81-9 | 9       |      |      |
| 5    | 5-Hexen-2-one                        | 98  | C6H10O       | 000109-49-9 | 9       |      |      |



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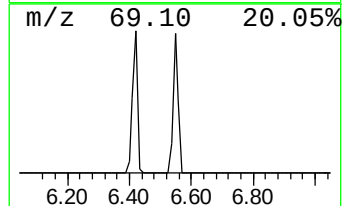
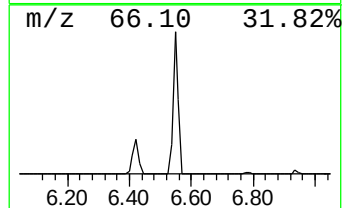
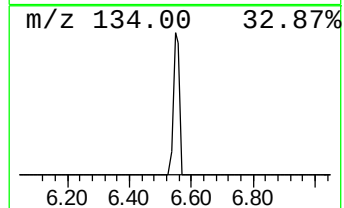
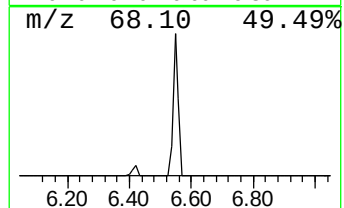
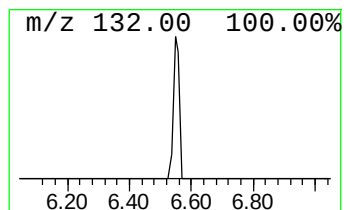
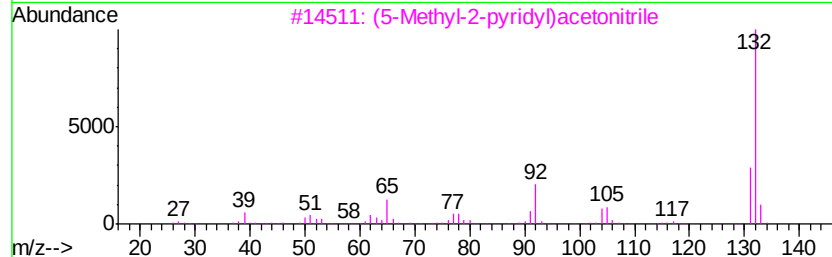
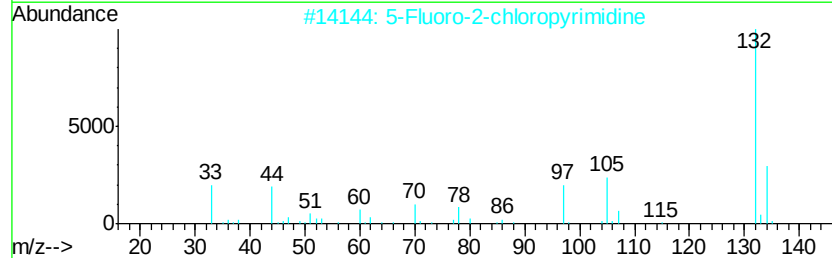
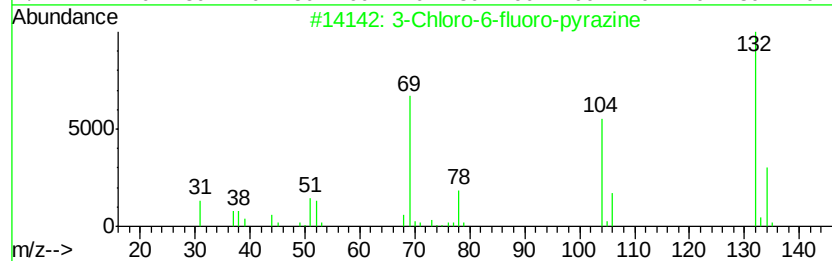
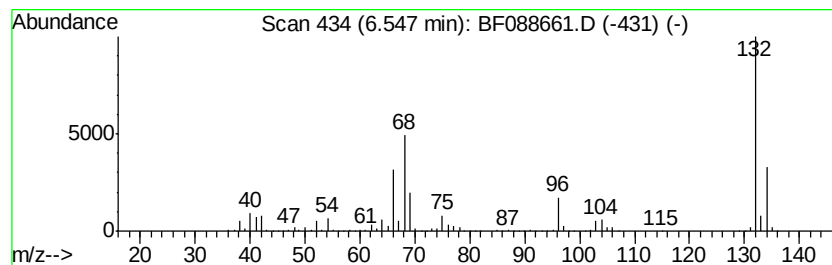
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Peak Number 6 unknown6.55 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.55 | 91.80 ng | 3240270 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of                               | Tentative ID | MW        | MolForm      | CAS# | Qual |
|------|----------------------------------|--------------|-----------|--------------|------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine       | 132          | C4H2ClFN2 | 1000146-10-7 | 27   |      |
| 2    | 5-Fluoro-2-chloropyrimidine      | 132          | C4H2ClFN2 | 062802-42-0  | 12   |      |
| 3    | (5-Methyl-2-pyridyl)acetonitrile | 132          | C8H8N2    | 1000241-93-9 | 10   |      |
| 4    | Cyclopropanamine, 1-phenyl-      | 133          | C9H11N    | 1000351-26-2 | 10   |      |
| 5    | 4-Methylpyrrolo[1,2-a]pyrazine   | 132          | C8H8N2    | 064608-60-2  | 9    |      |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\  
Data File : BF088661.D  
Acq On : 3 Jul 2016 21:52  
Operator : UM/SJ  
Sample : H3817-05  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

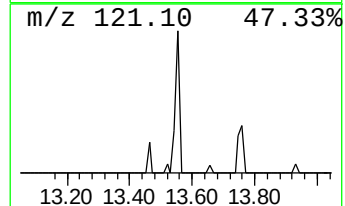
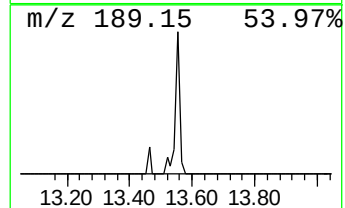
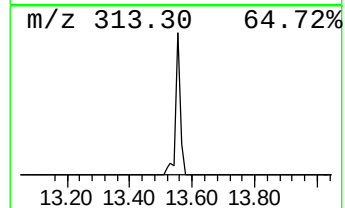
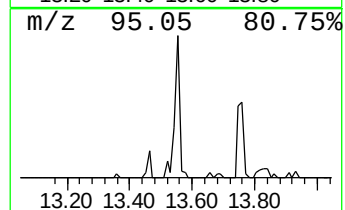
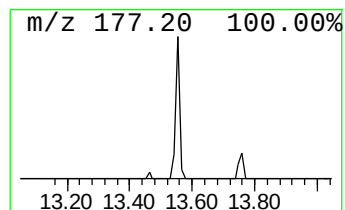
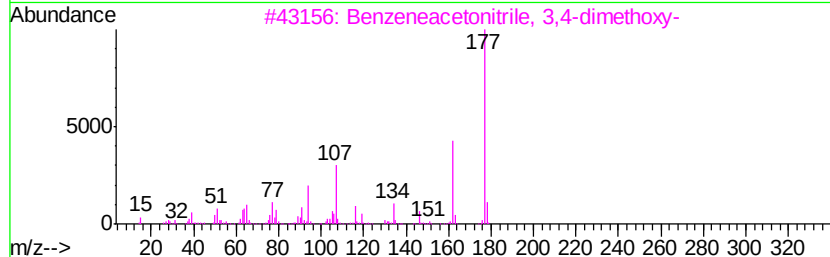
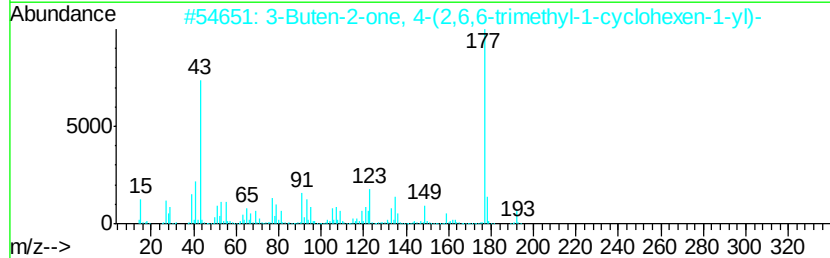
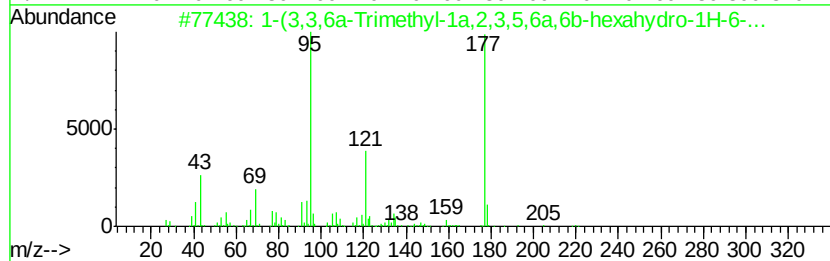
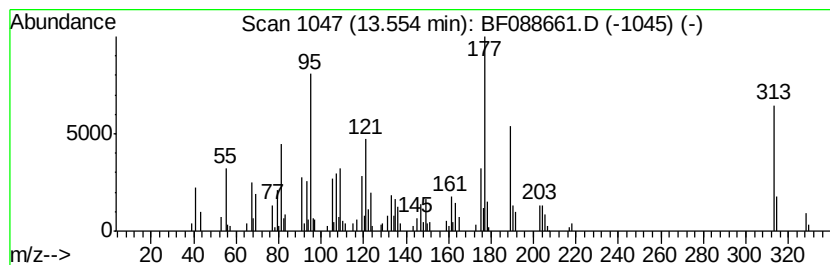
Instrument :  
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ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown13.55 Concentration Rank 9

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.55     | 2.88 ng                             | 136301 | Chrysene-d12     | 13.92        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-(3,3,6a-Trimethyl-1a,2,3,5,6a,... | 220    | C14H20O2         | 1000190-30-6 | 32   |
| 2         | 3-Buten-2-one, 4-(2,6,6-trimethy... | 192    | C13H20O          | 014901-07-6  | 22   |
| 3         | Benzeneacetonitrile, 3,4-dimethoxy- | 177    | C10H11NO2        | 000093-17-4  | 14   |
| 4         | Bicyclo[2.2.1]heptane-2-carboxam... | 328    | C19H24N2OS       | 1000270-21-2 | 11   |
| 5         | 5-Aminomethylene-6-hydroxy-4-met... | 177    | C8H7N3O2         | 1000223-27-4 | 10   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\  
Data File : BF088661.D  
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Operator : UM/SJ  
Sample : H3817-05  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

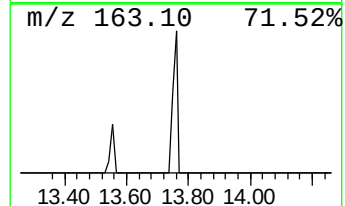
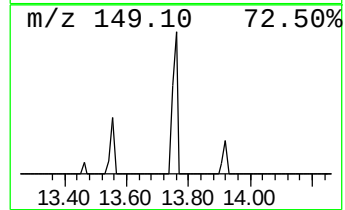
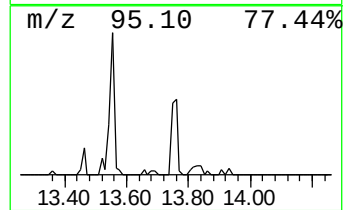
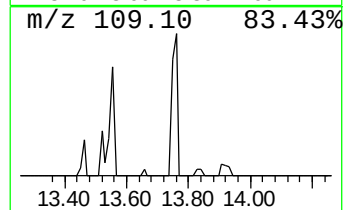
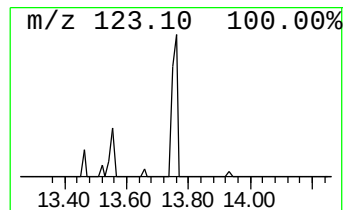
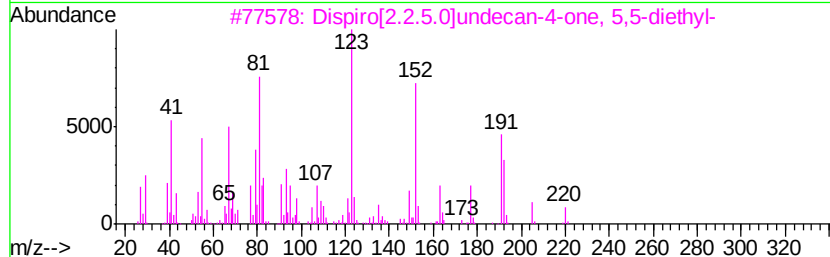
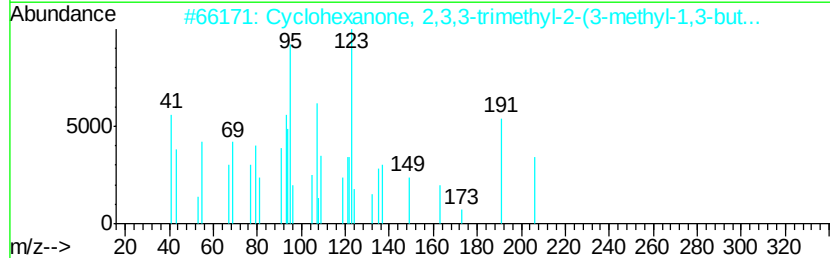
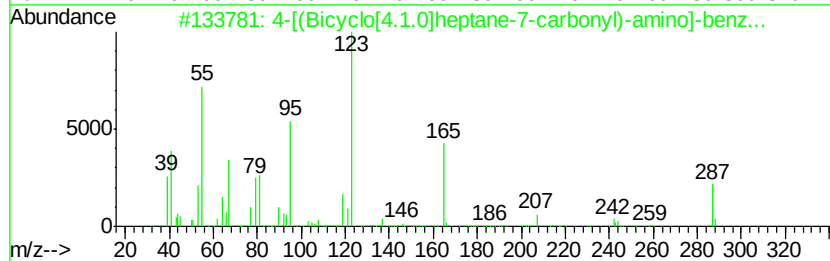
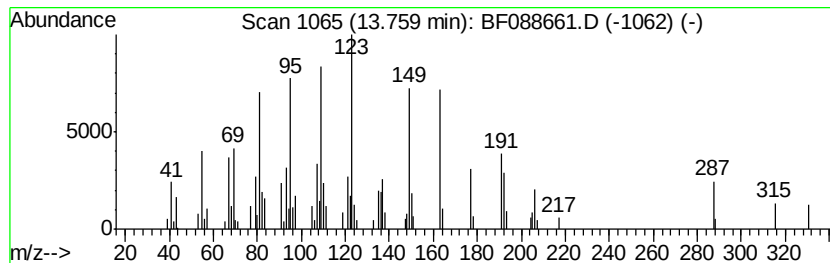
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ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown13.76 Concentration Rank 10

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.76     | 2.82 ng                             | 133731 | Chrysene-d12     | 13.92        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 4-[(Bicyclo[4.1.0]heptane-7-carb... | 287    | C17H21NO3        | 1000297-42-7 | 38   |
| 2         | Cyclohexanone, 2,3,3-trimethyl-2... | 206    | C14H22O          | 069296-90-8  | 38   |
| 3         | Dispiro[2.2.5.0]undecan-4-one, 5... | 220    | C15H24O          | 1000152-13-8 | 27   |
| 4         | Cyclohexene, 3-methyl-6-(1-methy... | 138    | C10H18           | 005256-65-5  | 25   |
| 5         | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138    | C10H18           | 020536-40-7  | 25   |



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Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

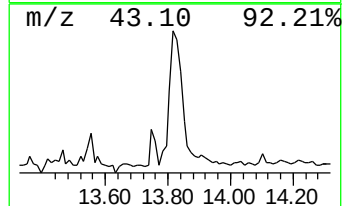
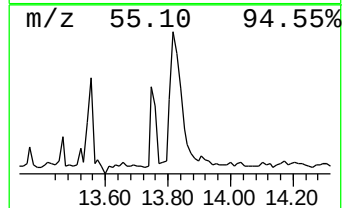
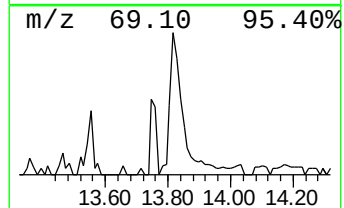
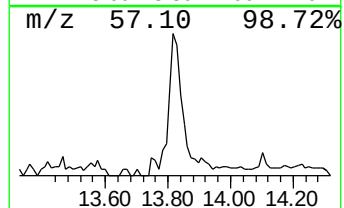
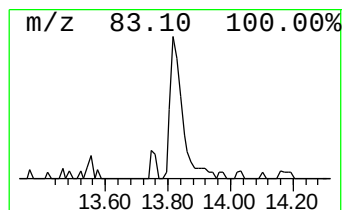
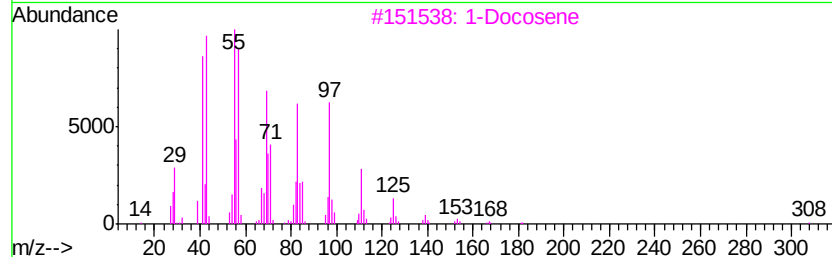
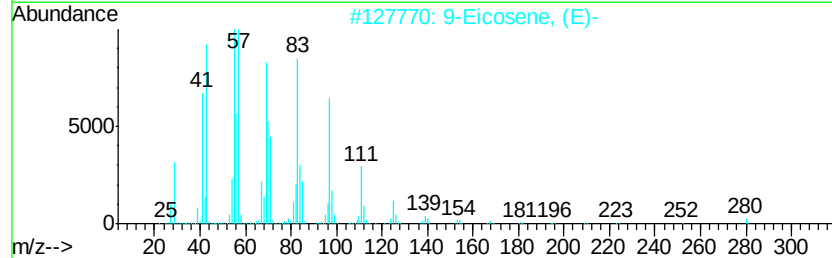
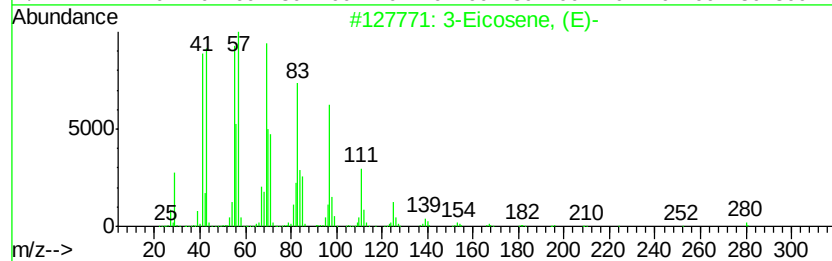
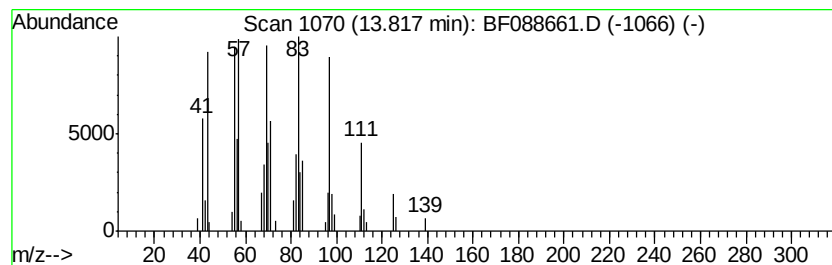
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 3-Eicosene, (E)- Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.82 | 4.55 ng | 215586 | Chrysene-d12     | 13.92 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 3-Eicosene, (E)-                    | 280 | C20H40      | 074685-33-9 | 90   |
| 2    |    |   | 9-Eicosene, (E)-                    | 280 | C20H40      | 074685-29-3 | 87   |
| 3    |    |   | 1-Docosene                          | 308 | C22H44      | 001599-67-3 | 81   |
| 4    |    |   | Trichloroacetic acid, pentadecyl... | 372 | C17H31Cl3O2 | 074339-53-0 | 76   |
| 5    |    |   | 1-Hexadecanol                       | 242 | C16H34O     | 036653-82-4 | 68   |



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Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

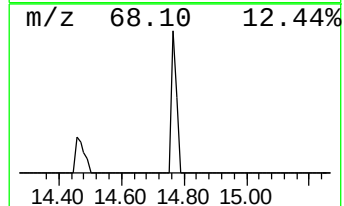
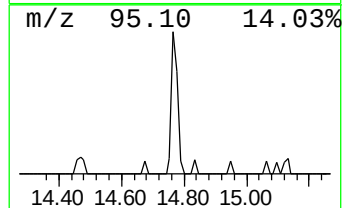
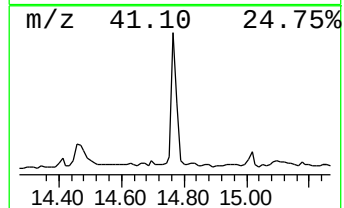
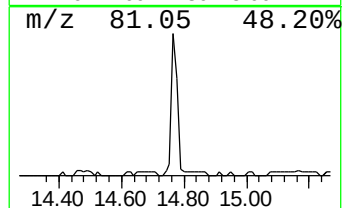
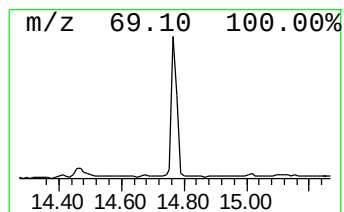
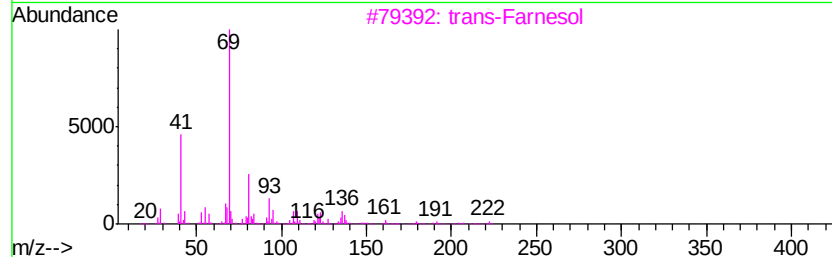
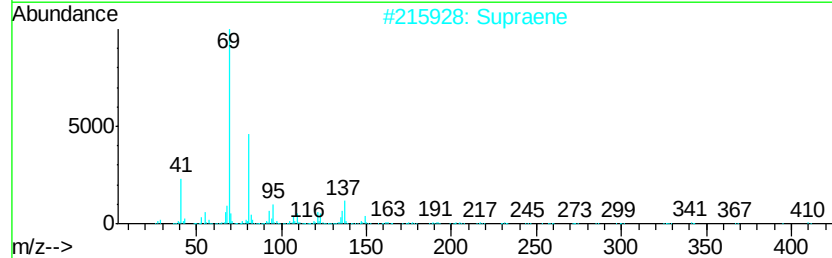
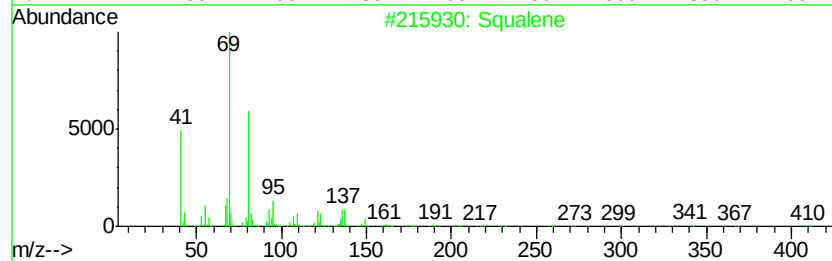
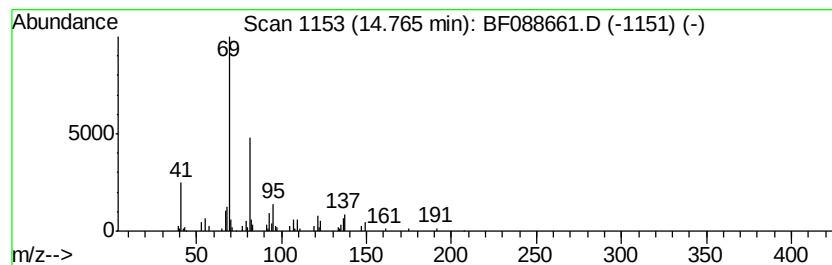
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Squalene Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.77 | 3.15 ng | 113080 | Perylene-d12     | 15.31 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Squalene                            | 410 | C30H50  | 000111-02-4 | 91   |
| 2    |    | Supraene                            | 410 | C30H50  | 007683-64-9 | 91   |
| 3    |    | trans-Farnesol                      | 222 | C15H26O | 000106-28-5 | 83   |
| 4    |    | 7,11-Dimethyldodeca-2,6,10-trien... | 208 | C14H24O | 125529-10-4 | 72   |
| 5    |    | 3,7,11-Tridecatrienitrile, 4,8...   | 231 | C16H25N | 006006-01-5 | 56   |



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Sample : H3817-05  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BS-1C-6.5-7FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| unknown1.64          | 1.64  | 195.1   | ng    | 6886870  | 1                     | 6.79  | 705951 | 20.0 |
| Propane, 1,1-dime... | 2.03  | 2.5     | ng    | 86861    | 1                     | 6.79  | 705951 | 20.0 |
| R-(-)-1,2-propane... | 3.03  | 4.3     | ng    | 150136   | 1                     | 6.79  | 705951 | 20.0 |
| 2-Pentanone, 4-hy... | 4.96  | 7.7     | ng    | 270376   | 1                     | 6.79  | 705951 | 20.0 |
| unknown6.55          | 6.55  | 91.8    | ng    | 3240270  | 1                     | 6.79  | 705951 | 20.0 |
| unknown13.55         | 13.55 | 2.9     | ng    | 136301   | 5                     | 13.92 | 948043 | 20.0 |
| unknown13.76         | 13.76 | 2.8     | ng    | 133731   | 5                     | 13.92 | 948043 | 20.0 |
| 3-Eicosene, (E)-     | 13.82 | 4.5     | ng    | 215586   | 5                     | 13.92 | 948043 | 20.0 |
| Squalene             | 14.77 | 3.1     | ng    | 113080   | 6                     | 15.31 | 716978 | 20.0 |