

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
 Data File : BF082289.D  
 Acq On : 14 Oct 2015 20:07  
 Operator : UM/IZ  
 Sample : G4016-06  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CS-6

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-BF101015.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         | 3.291       | 97            | 101         | 112          | rVB      | 59044          | 123228        | 2.05%           | 0.429%        |
| 2         | 4.343       | 188           | 193         | 196          | rBV      | 2492053        | 6014012       | 100.00%         | 20.925%       |
| 3         | 5.017       | 247           | 252         | 255          | rBV      | 1710375        | 3860736       | 64.20%          | 13.433%       |
| 4         | 6.446       | 375           | 377         | 387          | rBV      | 363937         | 409993        | 6.82%           | 1.427%        |
| 5         | 6.583       | 387           | 389         | 393          | rVB      | 116896         | 102207        | 1.70%           | 0.356%        |
| 6         | 6.823       | 408           | 410         | 412          | rBV      | 374721         | 395683        | 6.58%           | 1.377%        |
| 7         | 6.972       | 421           | 423         | 426          | rBV      | 1122143        | 1148919       | 19.10%          | 3.997%        |
| 8         | 7.143       | 437           | 438         | 441          | rBV3     | 99591          | 118367        | 1.97%           | 0.412%        |
| 9         | 7.395       | 457           | 460         | 462          | rBV      | 717462         | 880741        | 14.64%          | 3.064%        |
| 10        | 8.103       | 520           | 522         | 525          | rBV      | 392624         | 558478        | 9.29%           | 1.943%        |
| 11        | 9.189       | 614           | 617         | 619          | rBV      | 2132011        | 1910208       | 31.76%          | 6.646%        |
| 12        | 9.589       | 649           | 652         | 654          | rBV      | 328025         | 298940        | 4.97%           | 1.040%        |
| 13        | 9.863       | 674           | 676         | 679          | rBV      | 642701         | 697315        | 11.59%          | 2.426%        |
| 14        | 10.298      | 710           | 714         | 716          | rBV2     | 34345          | 61283         | 1.02%           | 0.213%        |
| 15        | 10.766      | 753           | 755         | 760          | rVB      | 53957          | 79141         | 1.32%           | 0.275%        |
| 16        | 11.349      | 804           | 806         | 809          | rBV      | 462562         | 677509        | 11.27%          | 2.357%        |
| 17        | 11.406      | 809           | 811         | 815          | rVB      | 53120          | 77581         | 1.29%           | 0.270%        |
| 18        | 11.601      | 823           | 828         | 829          | rBV3     | 39530          | 77669         | 1.29%           | 0.270%        |
| 19        | 11.646      | 830           | 832         | 833          | rBV      | 86307          | 68018         | 1.13%           | 0.237%        |
| 20        | 11.806      | 844           | 846         | 849          | rBV2     | 33419          | 89537         | 1.49%           | 0.312%        |
| 21        | 11.966      | 857           | 860         | 861          | rBV3     | 41124          | 78456         | 1.30%           | 0.273%        |
| 22        | 12.058      | 866           | 868         | 869          | rVV      | 96986          | 85531         | 1.42%           | 0.298%        |
| 23        | 12.344      | 890           | 893         | 895          | rBV      | 108728         | 232518        | 3.87%           | 0.809%        |
| 24        | 12.447      | 901           | 902         | 904          | rBV      | 111440         | 107532        | 1.79%           | 0.374%        |
| 25        | 12.549      | 909           | 911         | 913          | rBV2     | 92067          | 137542        | 2.29%           | 0.479%        |
| 26        | 12.687      | 919           | 923         | 925          | rBV2     | 112306         | 282906        | 4.70%           | 0.984%        |
| 27        | 12.801      | 931           | 933         | 934          | rBV      | 125219         | 153938        | 2.56%           | 0.536%        |
| 28        | 12.949      | 943           | 946         | 949          | rBV      | 1592613        | 2153803       | 35.81%          | 7.494%        |
| 29        | 13.224      | 968           | 970         | 973          | rBV2     | 181101         | 395282        | 6.57%           | 1.375%        |
| 30        | 13.315      | 976           | 978         | 979          | rVV      | 240157         | 292201        | 4.86%           | 1.017%        |
| 31        | 13.395      | 983           | 985         | 987          | rBV2     | 184656         | 364569        | 6.06%           | 1.268%        |
| 32        | 13.875      | 1023          | 1027        | 1033         | rBV2     | 370146         | 1267921       | 21.08%          | 4.412%        |
| 33        | 14.035      | 1039          | 1041        | 1044         | rBV      | 417812         | 488603        | 8.12%           | 1.700%        |
| 34        | 14.332      | 1065          | 1067        | 1071         | rBV2     | 301185         | 556738        | 9.26%           | 1.937%        |

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Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

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\_F\METHODS\8270-BF101015.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 15.693 | 1184 | 1186 | 1192 | rVB2 | 362071 | 763550  | 12.70% | 2.657% |
| 36 | 16.241 | 1232 | 1234 | 1243 | rVB  | 686588 | 1929173 | 32.08% | 6.712% |
| 37 | 16.664 | 1268 | 1271 | 1278 | rVB  | 806012 | 1801156 | 29.95% | 6.267% |

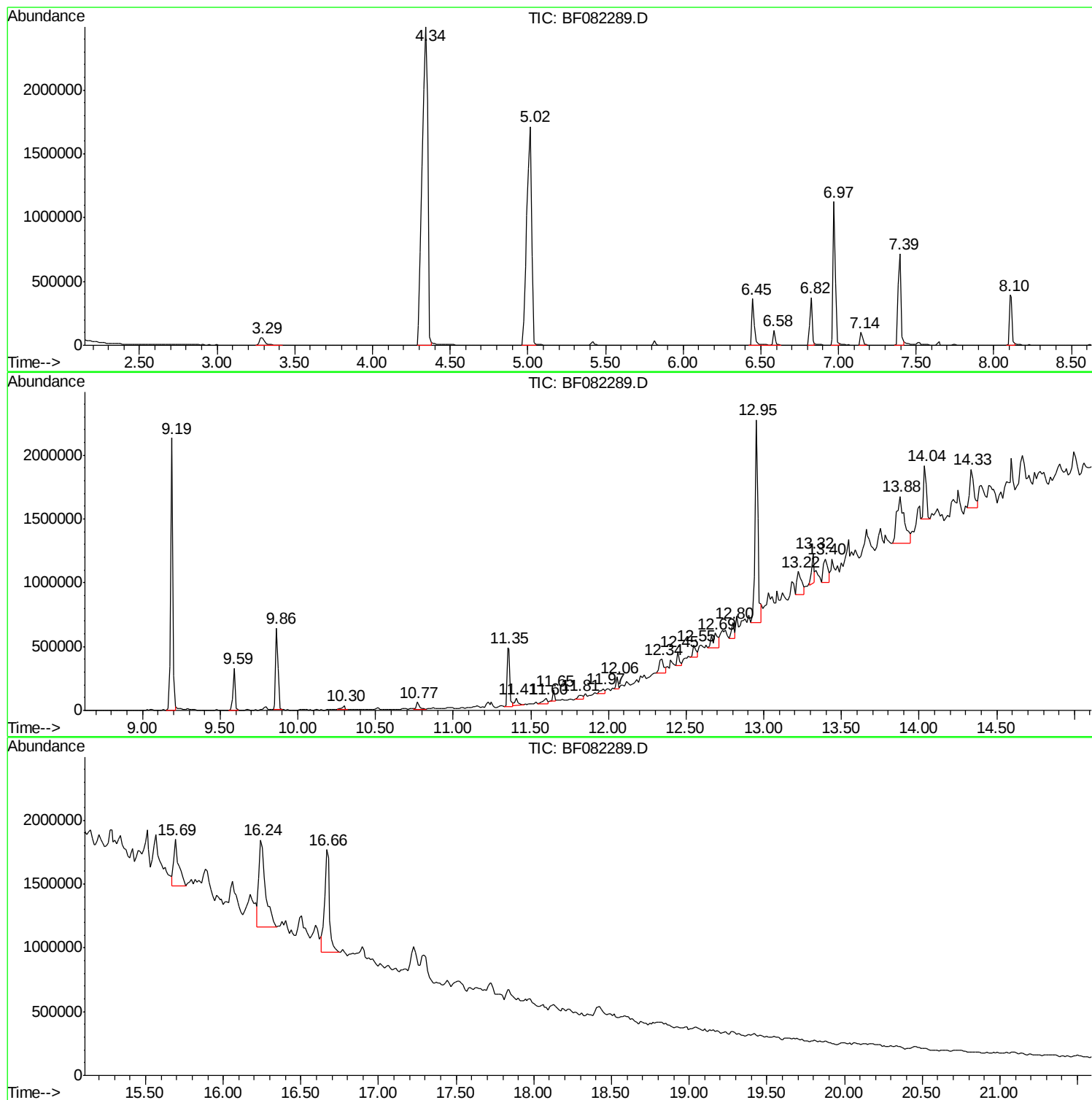
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : G4016-06  
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Instrument :  
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CS-6

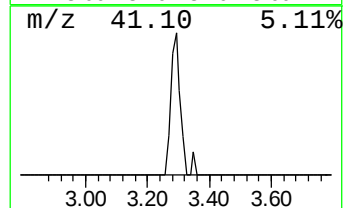
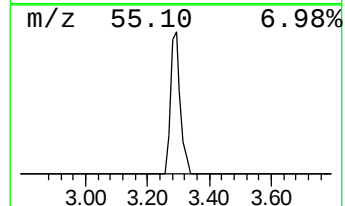
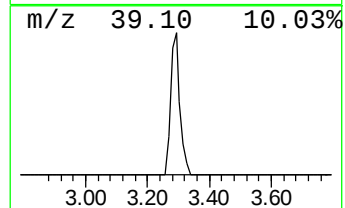
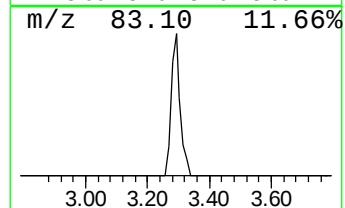
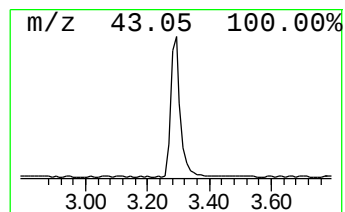
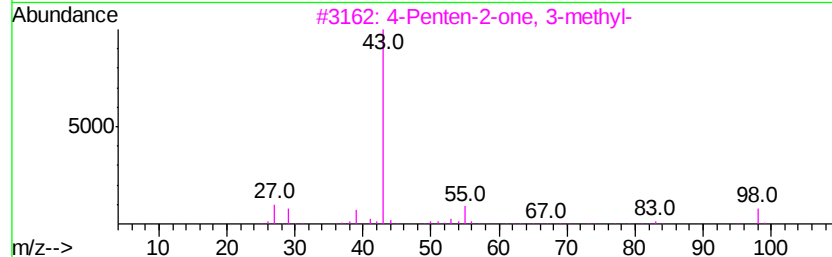
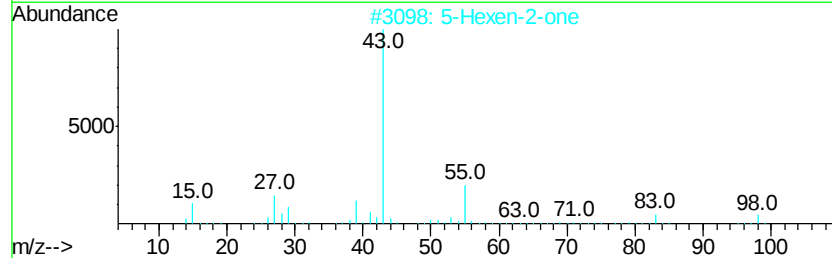
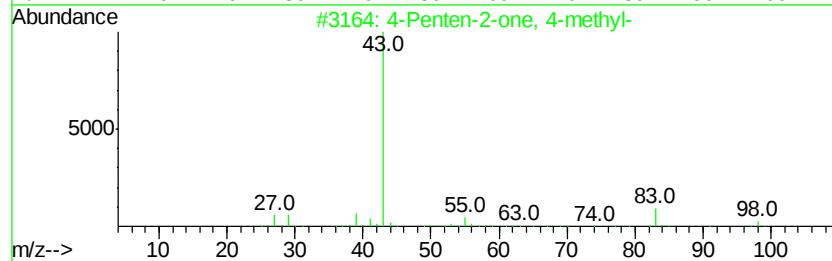
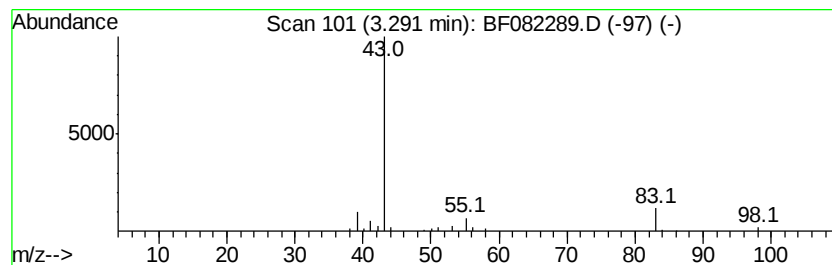
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 15

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 3.29 | 6.23 ng | 123228 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of | 5 | Tentative ID                      | MW | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|----|---------|-------------|------|
| 1    |    |   | 4-Penten-2-one, 4-methyl-         | 98 | C6H10O  | 003744-02-3 | 78   |
| 2    |    |   | 5-Hexen-2-one                     | 98 | C6H10O  | 000109-49-9 | 50   |
| 3    |    |   | 4-Penten-2-one, 3-methyl-         | 98 | C6H10O  | 000758-87-2 | 25   |
| 4    |    |   | 2-Pentanone, 3-methyl-            | 98 | C6H10O  | 004359-77-7 | 17   |
| 5    |    |   | Methyl 1-methylcyclopropyl ketone | 98 | C6H10O  | 001567-75-5 | 10   |



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BNA\_F  
ClientSampleId :  
CS-6

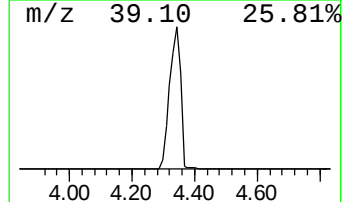
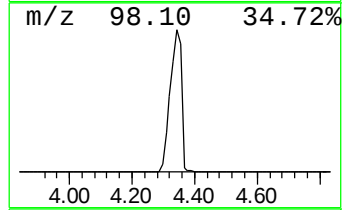
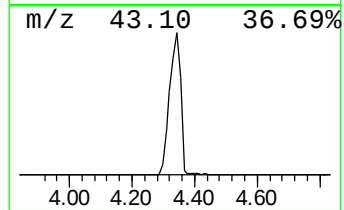
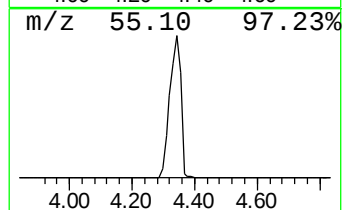
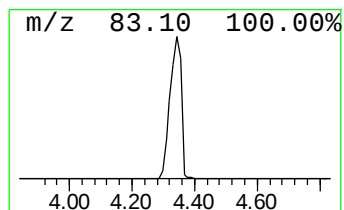
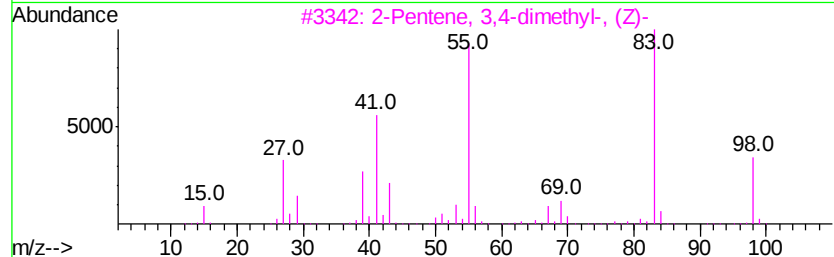
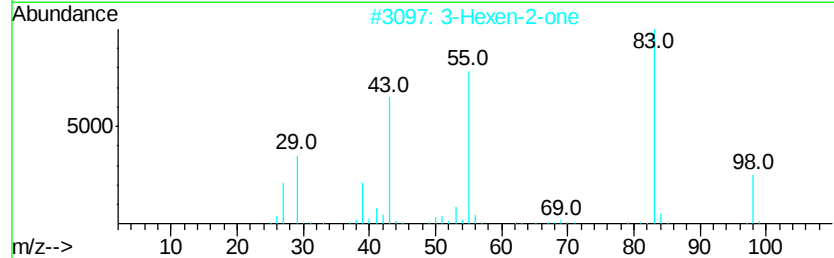
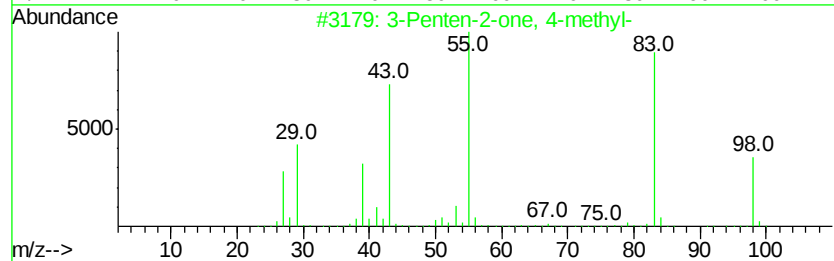
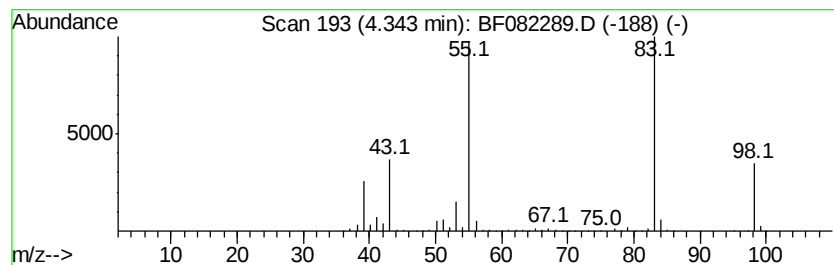
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 4.34 | 303.98 ng | 6014010 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|--------------------------------|----|---------|-------------|------|
| 1    | 5  | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 91   |
| 2    |    | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 91   |
| 3    |    | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 86   |
| 4    |    | 2-Pentene, 4,4-dimethyl-, (E)- | 98 | C7H14   | 000690-08-4 | 86   |
| 5    |    | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 86   |



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Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

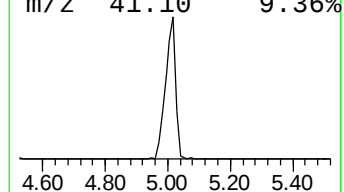
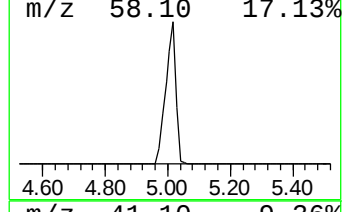
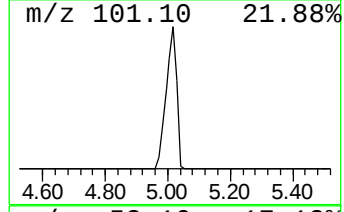
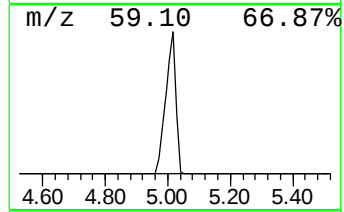
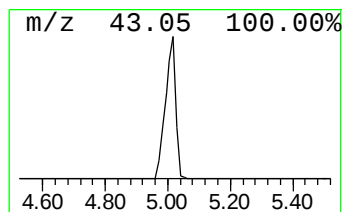
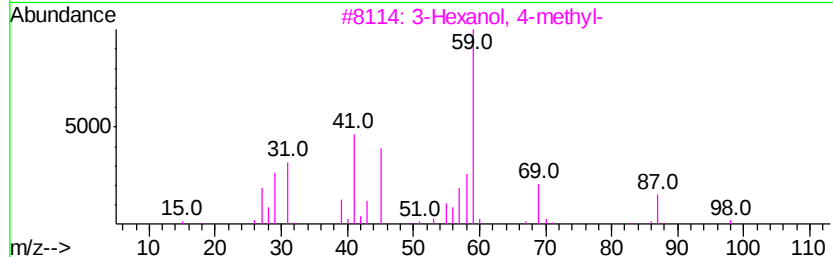
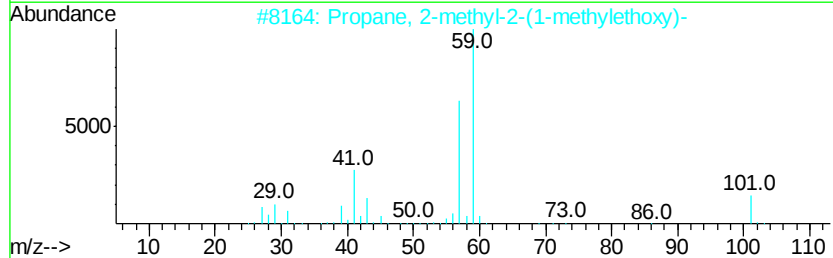
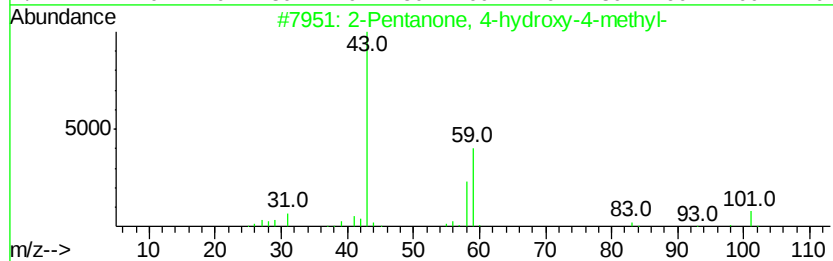
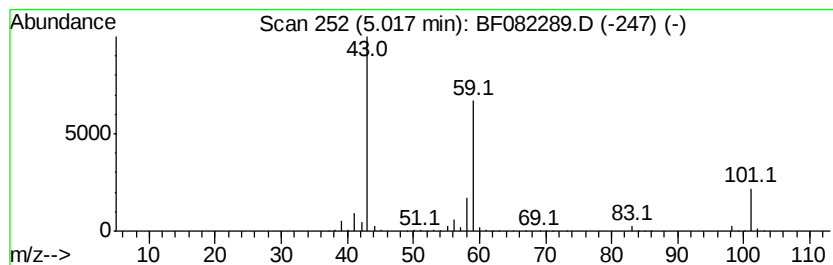
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 5.02 | 195.14 ng | 3860740 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of | Tentative ID                          | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-      | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    | Propane, 2-methyl-2-(1-methylethoxy)- | 116 | C7H16O   | 017348-59-3 | 53   |
| 3    |    | 3-Hexanol, 4-methyl-                  | 116 | C7H16O   | 000615-29-2 | 28   |
| 4    |    | Acetic acid, cyano-, 1,1-dimethyl-    | 141 | C7H11NO2 | 001116-98-9 | 25   |
| 5    |    | Butane, 1-ethoxy-                     | 102 | C6H14O   | 000628-81-9 | 9    |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

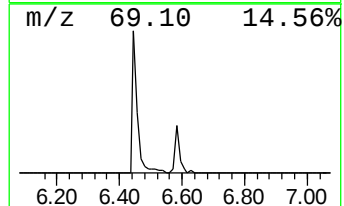
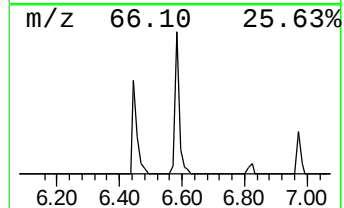
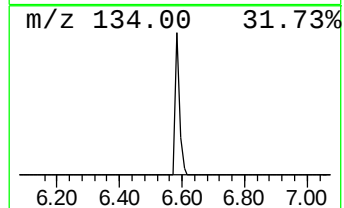
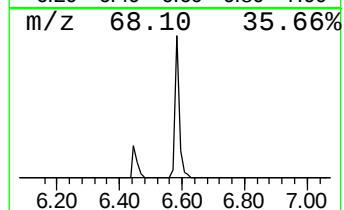
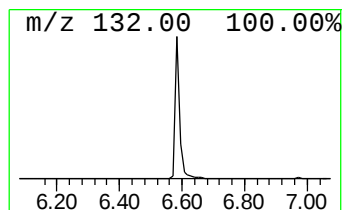
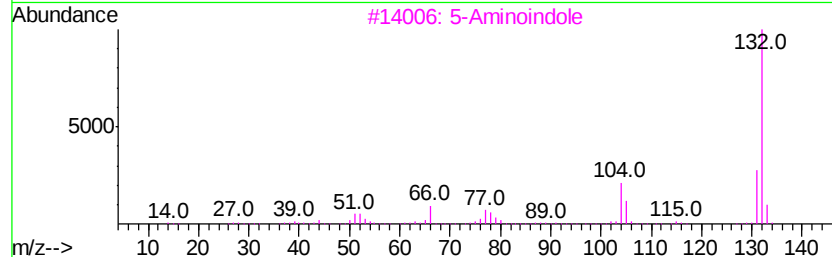
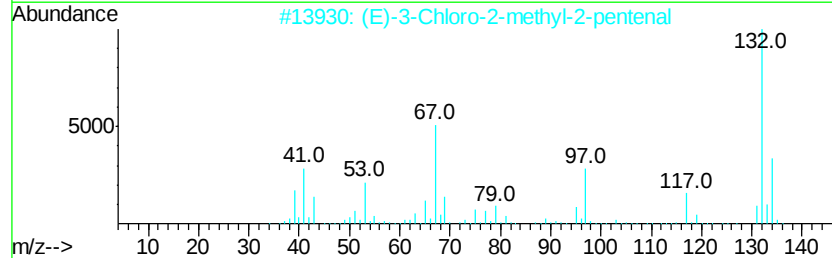
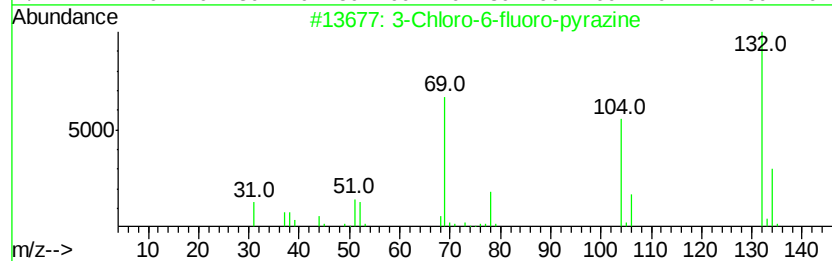
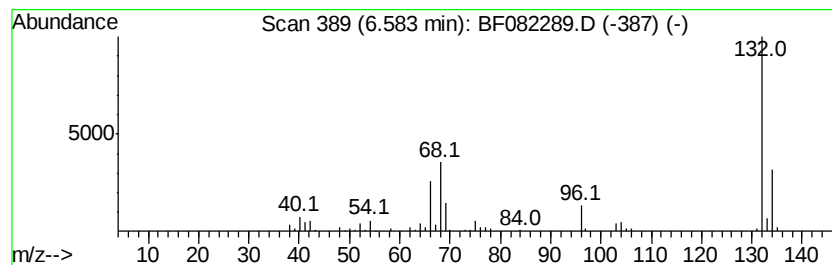
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown6.58 Concentration Rank 17

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.58 | 5.17 ng | 102207 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|---|----------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    |   | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 25   |
| 3    |    |   | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 10   |
| 4    |    |   | 4-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-60-2  | 9    |
| 5    |    |   | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

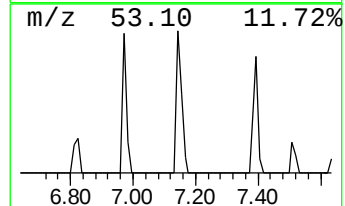
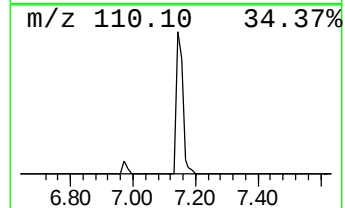
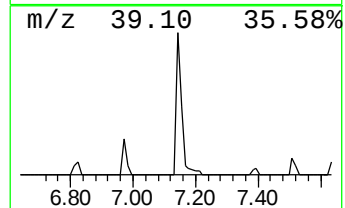
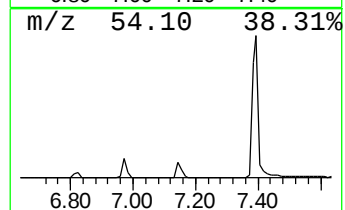
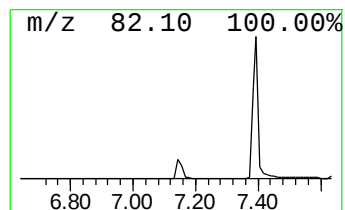
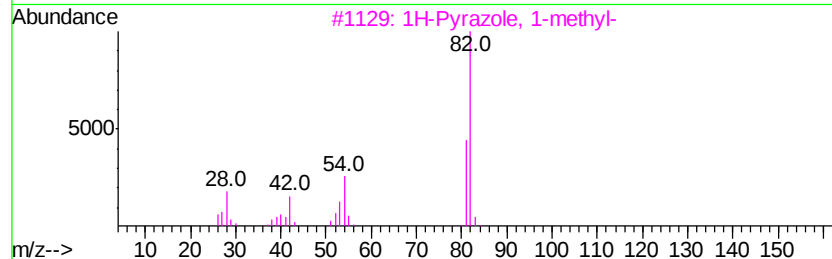
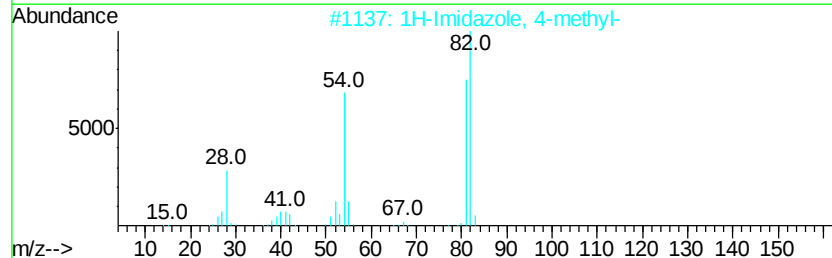
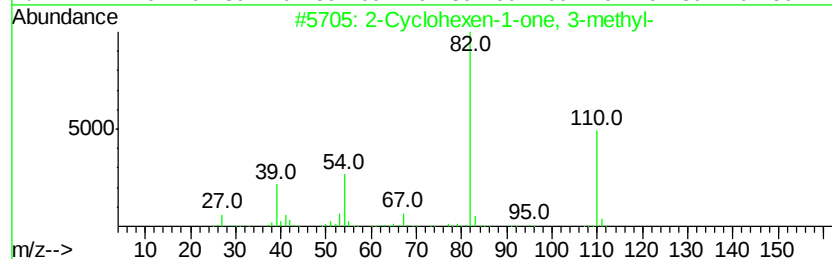
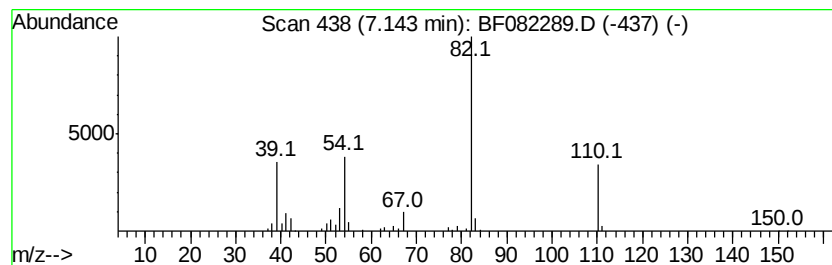
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 2-Cyclohexen-1-one, 3-methyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.14 | 5.98 ng | 118367 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Cyclohexen-1-one, 3-methyl-       | 110 | C7H10O   | 001193-18-6 | 91   |
| 2    |    |   | 1H-Imidazole, 4-methyl-             | 82  | C4H6N2   | 000822-36-6 | 59   |
| 3    |    |   | 1H-Pyrazole, 1-methyl-              | 82  | C4H6N2   | 000930-36-9 | 50   |
| 4    |    |   | 1H-Pyrazole, 3-methyl-              | 82  | C4H6N2   | 001453-58-3 | 49   |
| 5    |    |   | 2-Cyclohexen-1-one, 6-(1-hydroxy... | 168 | C10H16O2 | 087791-00-2 | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

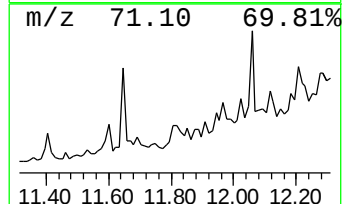
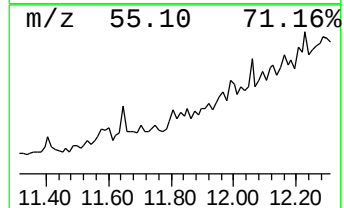
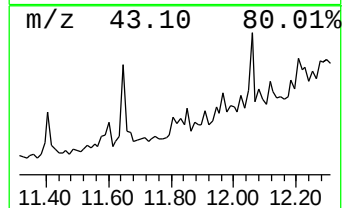
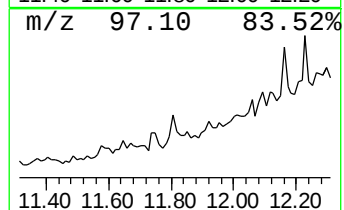
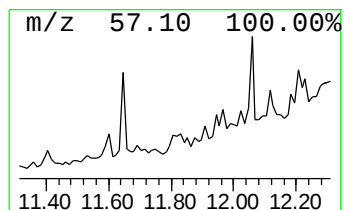
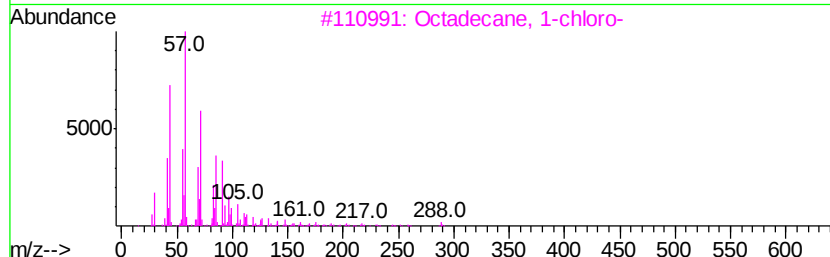
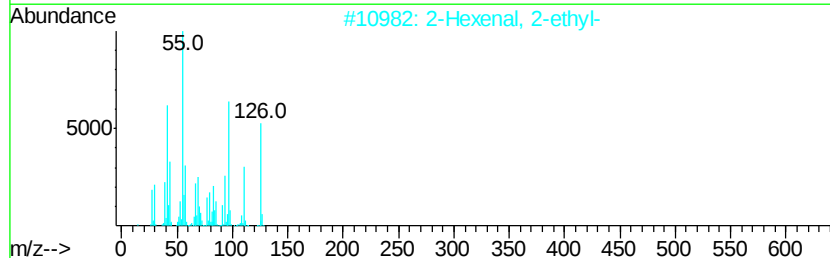
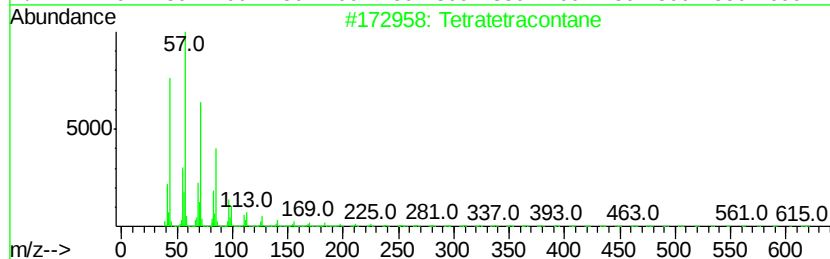
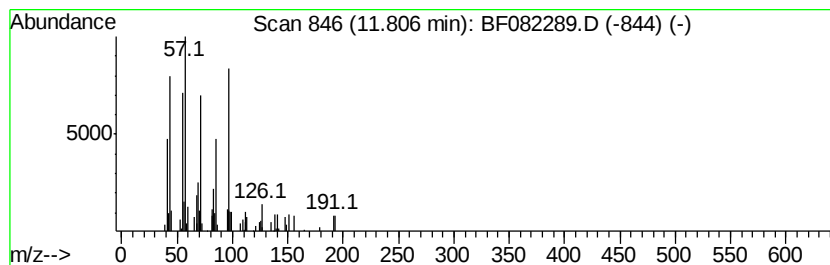
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown11.81 Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.81 | 2.64 ng | 89537 | Phenanthrene-d10 | 11.36 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Tetratetracontane                   | 619 | C44H90   | 007098-22-8 | 43   |
| 2    |    |   | 2-Hexenal, 2-ethyl-                 | 126 | C8H14O   | 000645-62-5 | 43   |
| 3    |    |   | Octadecane, 1-chloro-               | 288 | C18H37Cl | 003386-33-2 | 43   |
| 4    |    |   | Tritetracontane                     | 605 | C43H88   | 007098-21-7 | 43   |
| 5    |    |   | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18    | 004923-78-8 | 41   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

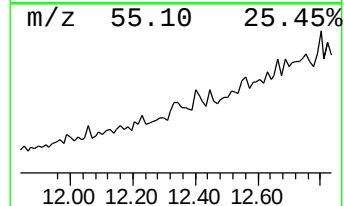
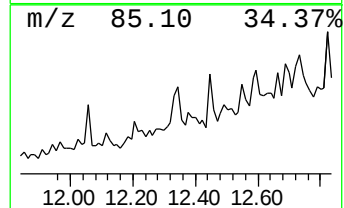
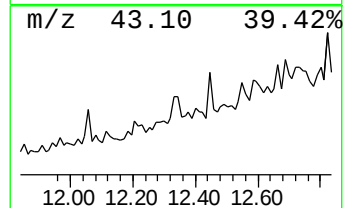
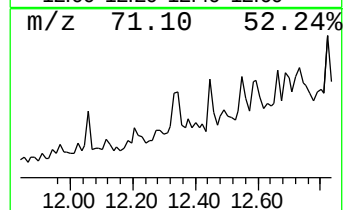
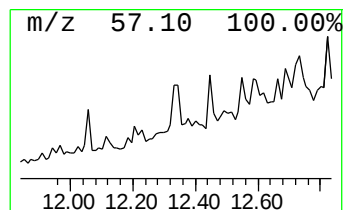
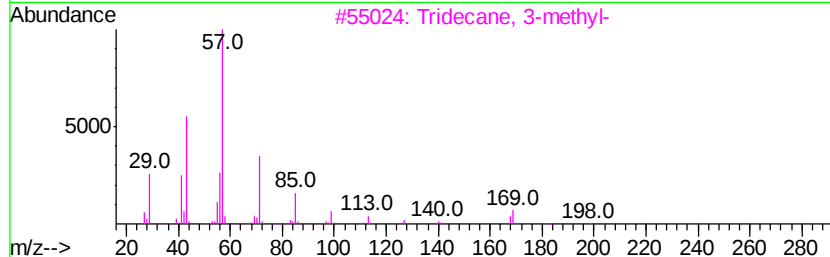
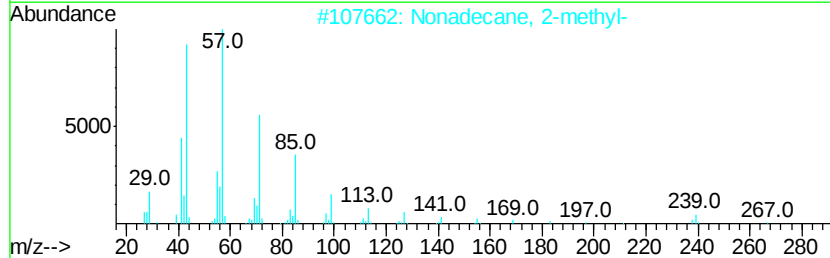
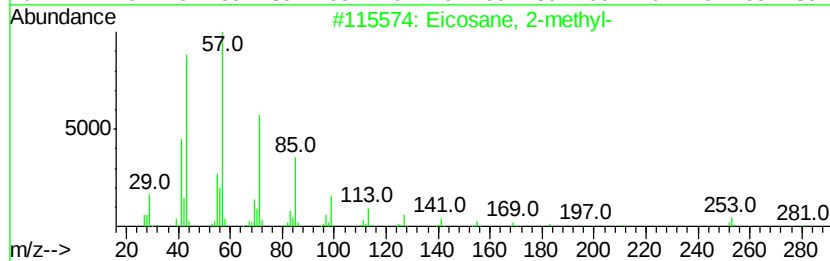
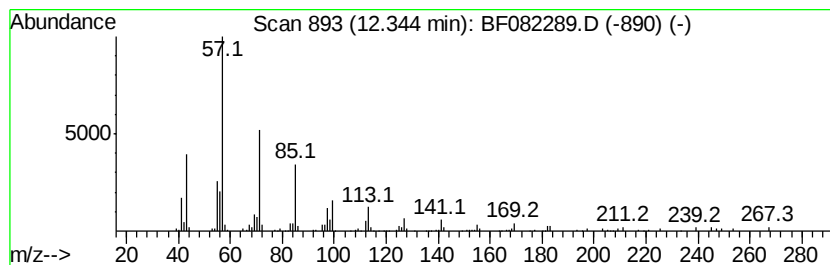
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Eicosane, 2-methyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.34 | 6.86 ng | 232518 | Phenanthrene-d10 | 11.36 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane, 2-methyl-   | 296 | C21H44  | 001560-84-5 | 86   |
| 2    |    | Nonadecane, 2-methyl- | 282 | C20H42  | 001560-86-7 | 86   |
| 3    |    | Tridecane, 3-methyl-  | 198 | C14H30  | 006418-41-3 | 86   |
| 4    |    | Octacosane            | 394 | C28H58  | 000630-02-4 | 83   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

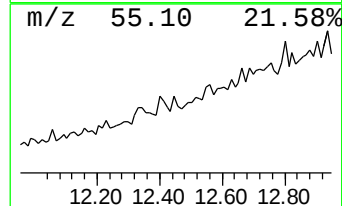
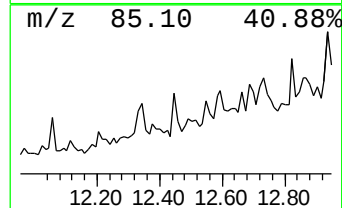
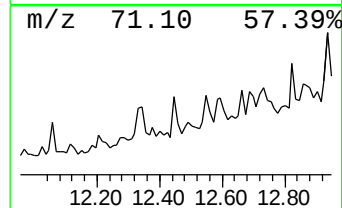
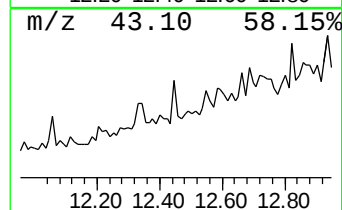
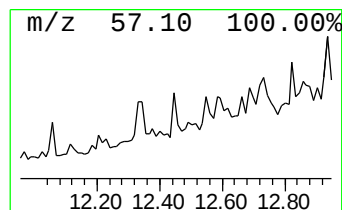
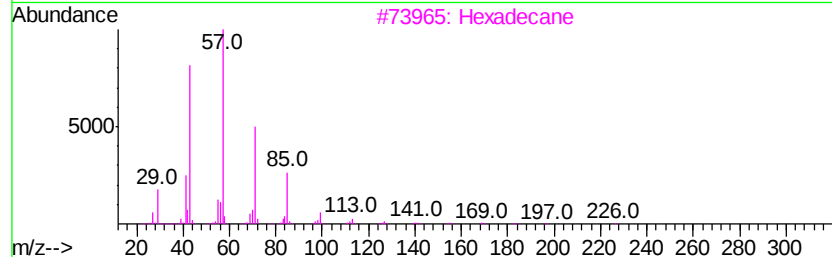
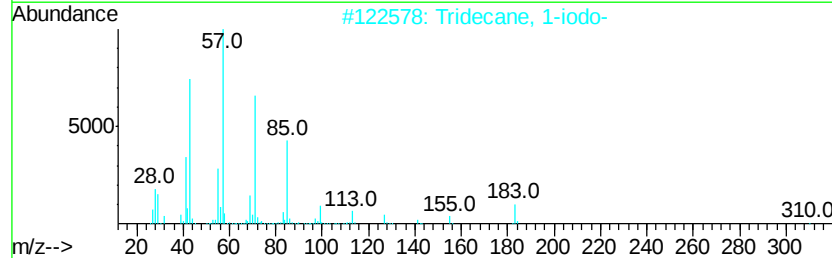
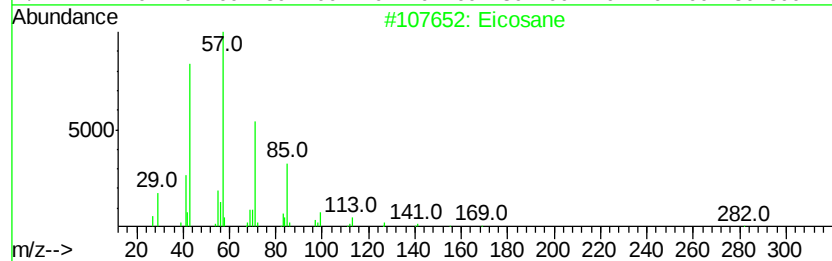
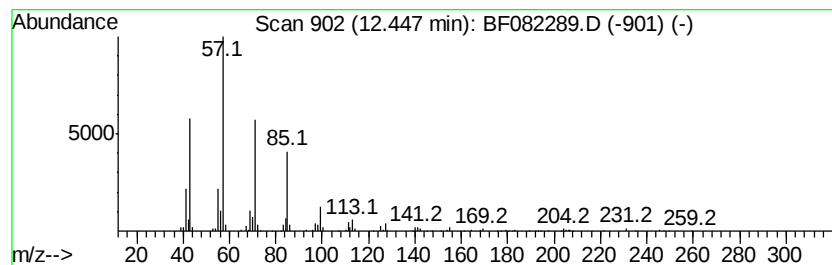
Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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 9 Eicosane Concentration Rank 19

| R.T.    | EstConc            | Area         | Relative to ISTD | R.T.           |
|---------|--------------------|--------------|------------------|----------------|
| 12.45   | 3.17 ng            | 107532       | Phenanthrene-d10 | 11.36          |
| Hit# of | 5                  | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Eicosane           |              | 282 C20H42       | 000112-95-8 86 |
| 2       | Tridecane, 1-iodo- |              | 310 C13H27I      | 035599-77-0 86 |
| 3       | Hexadecane         |              | 226 C16H34       | 000544-76-3 86 |
| 4       | Pentadecane        |              | 212 C15H32       | 000629-62-9 83 |
| 5       | Heneicosane        |              | 296 C21H44       | 000629-94-7 83 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

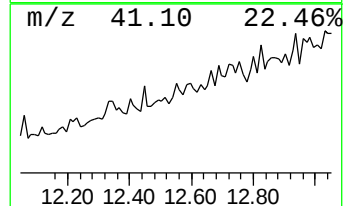
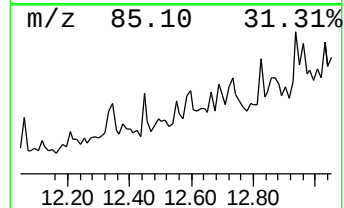
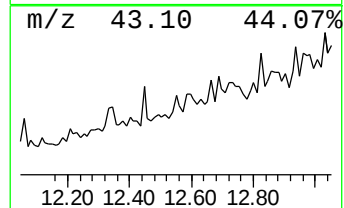
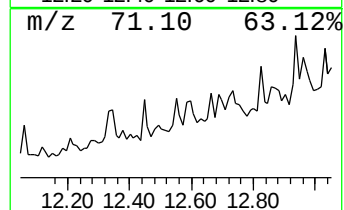
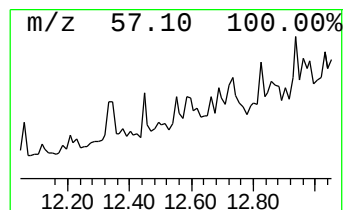
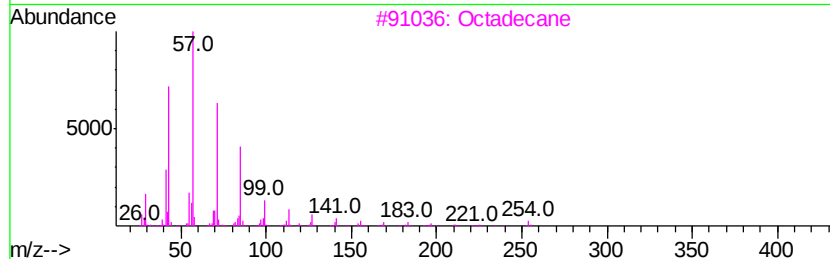
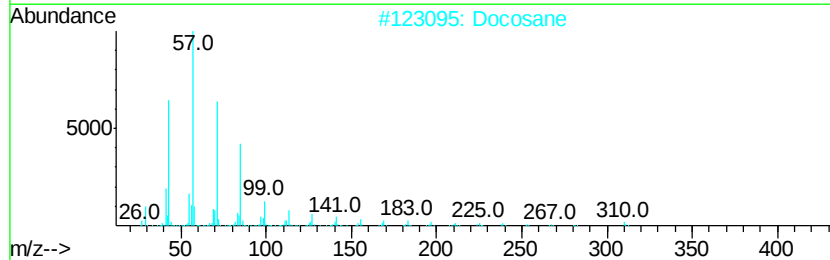
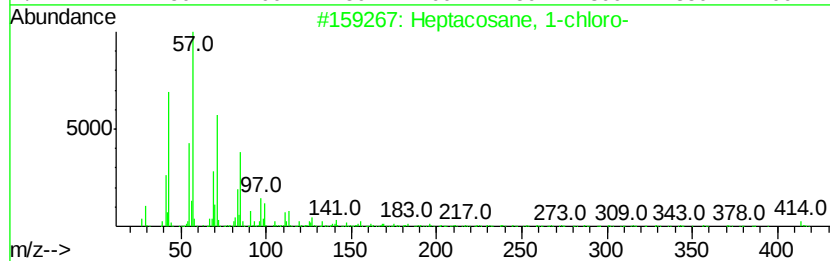
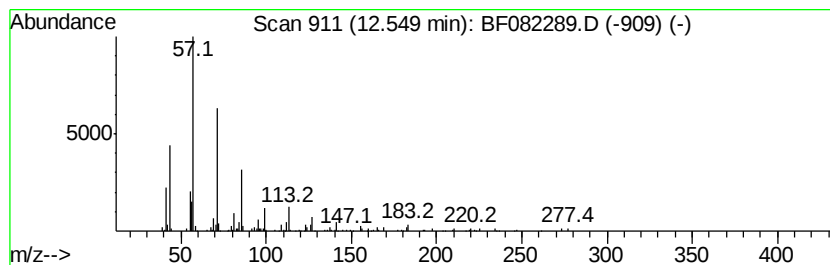
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Heptacosane, 1-chloro- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.55 | 4.06 ng | 137542 | Phenanthrene-d10 | 11.36 |

| Hit# of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|---------|---|------------------------|-----|----------|-------------|------|
| 1       |   | Heptacosane, 1-chloro- | 414 | C27H55Cl | 062016-79-9 | 72   |
| 2       |   | Docosane               | 310 | C22H46   | 000629-97-0 | 72   |
| 3       |   | Octadecane             | 254 | C18H38   | 000593-45-3 | 72   |
| 4       |   | Hexadecane             | 226 | C16H34   | 000544-76-3 | 72   |
| 5       |   | Hexatriacontane        | 507 | C36H74   | 000630-06-8 | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

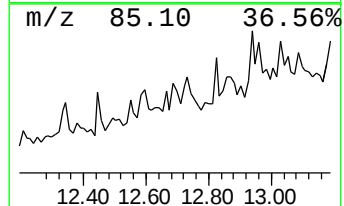
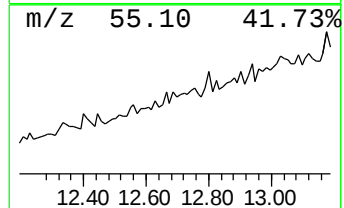
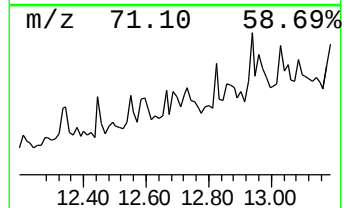
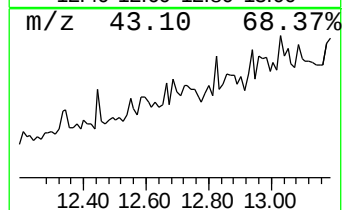
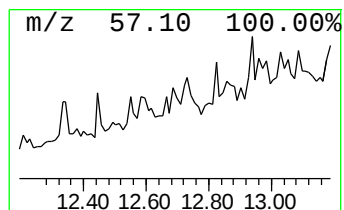
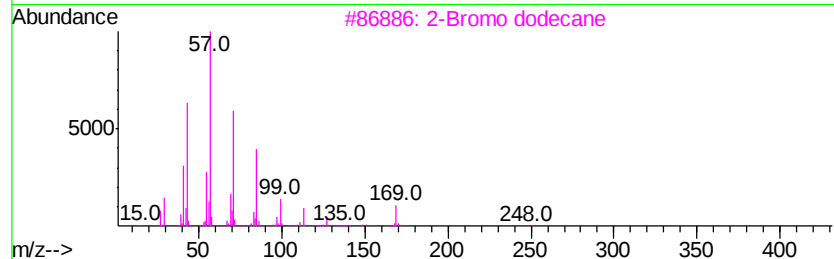
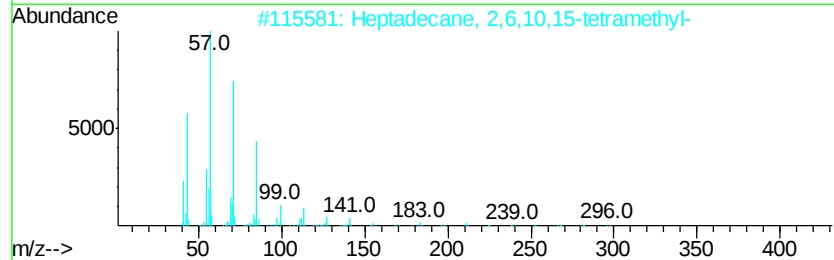
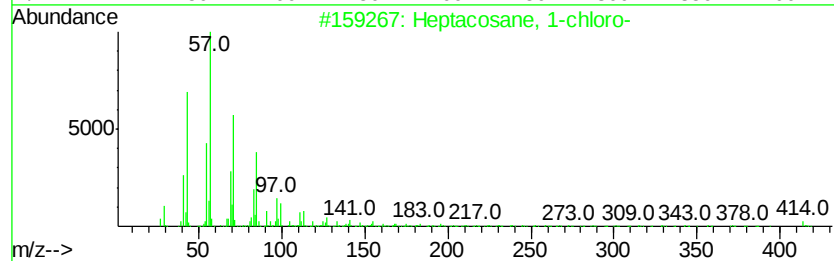
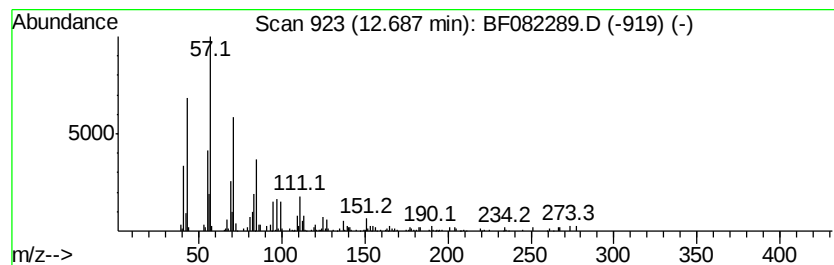
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Heptadecane, 2,6,10,15-tetr... Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.69 | 8.35 ng | 282906 | Phenanthrene-d10 | 11.36 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Heptacosane, 1-chloro-              | 414 | C27H55Cl | 062016-79-9 | 72   |
| 2    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 64   |
| 3    |    |   | 2-Bromo dodecane                    | 248 | C12H25Br | 013187-99-0 | 64   |
| 4    |    |   | Hexadecane, 1-bromo-                | 304 | C16H33Br | 000112-82-3 | 64   |
| 5    |    |   | Heptacosane                         | 380 | C27H56   | 000593-49-7 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

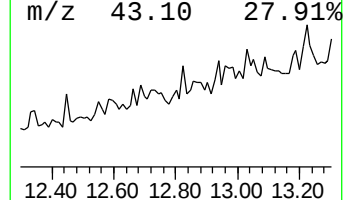
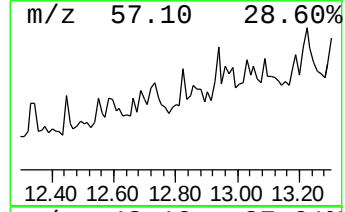
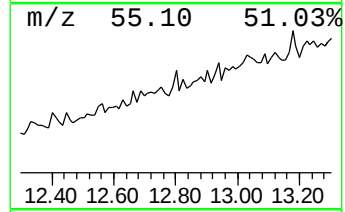
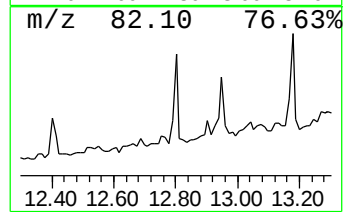
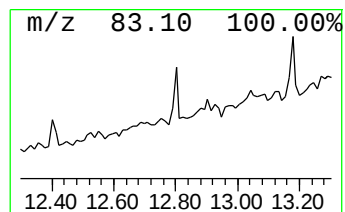
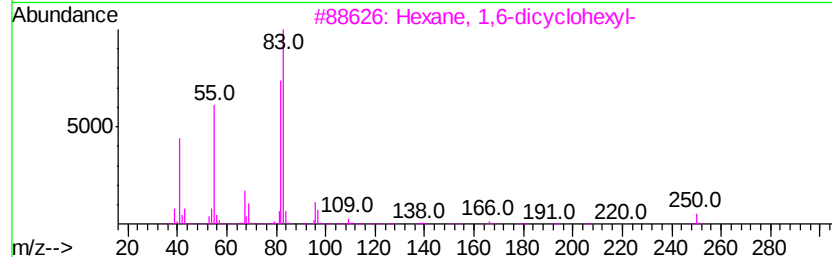
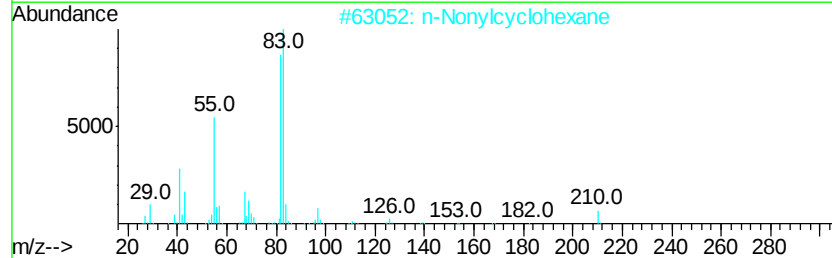
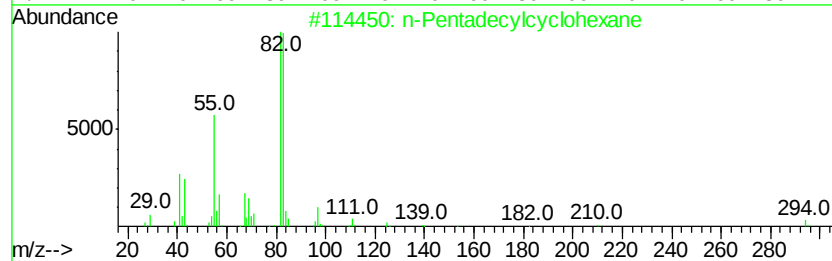
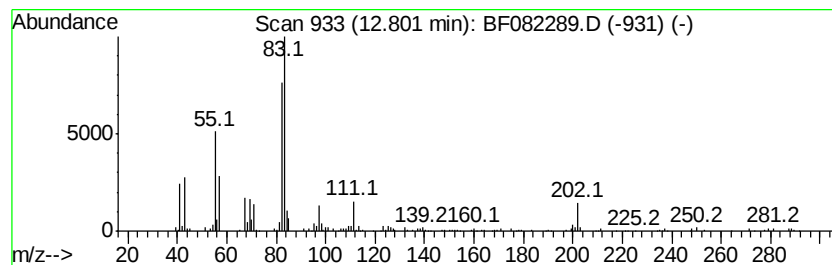
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BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 n-Pentadecylcyclohexane Concentration Rank 14

| R.T.      | EstConc                   | Area   | Relative to ISTD |             | R.T.  |
|-----------|---------------------------|--------|------------------|-------------|-------|
| 12.80     | 6.30 ng                   | 153938 | Chrysene-d12     |             | 14.04 |
| Hit# of 5 | Tentative ID              | MW     | MolForm          | CAS#        | Qual  |
| 1         | n-Pentadecylcyclohexane   | 294    | C21H42           | 006006-95-7 | 80    |
| 2         | n-Nonylcyclohexane        | 210    | C15H30           | 002883-02-5 | 80    |
| 3         | Hexane, 1,6-dicyclohexyl- | 250    | C18H34           | 001610-23-7 | 74    |
| 4         | Dodecylcyclohexane        | 252    | C18H36           | 001795-17-1 | 72    |
| 5         | Cyclohexane, eicosyl-     | 364    | C26H52           | 004443-55-4 | 72    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

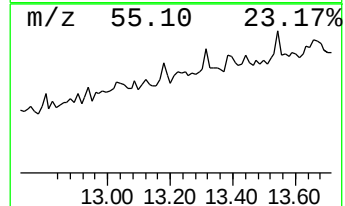
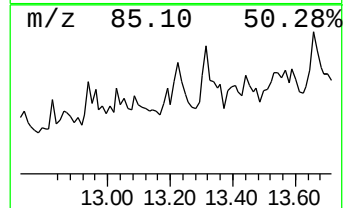
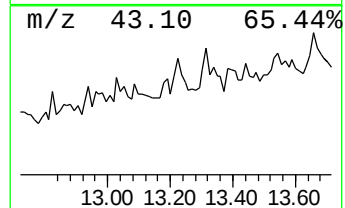
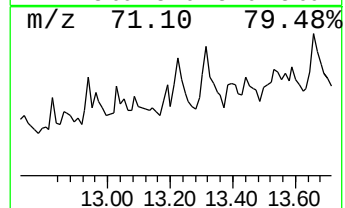
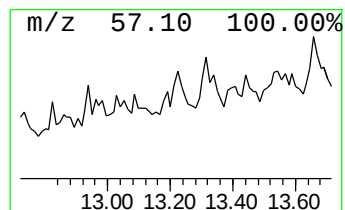
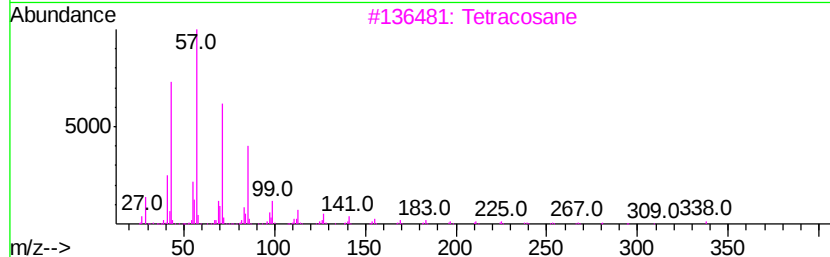
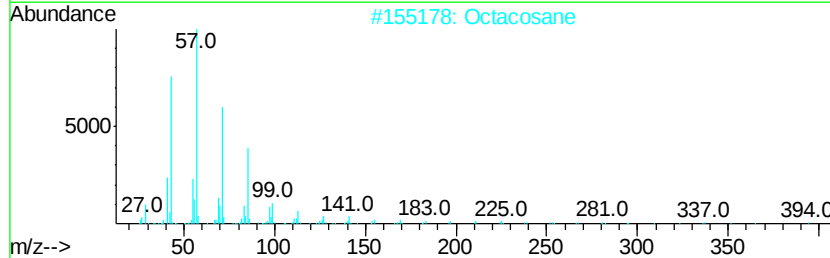
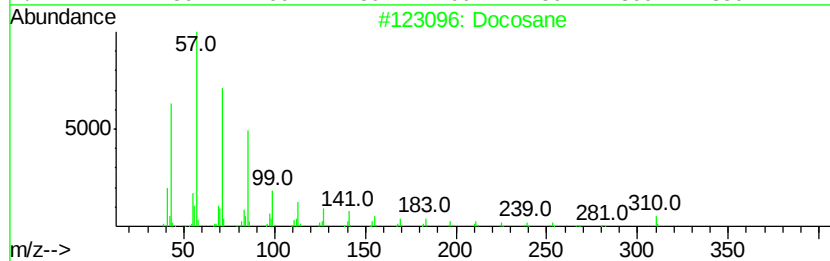
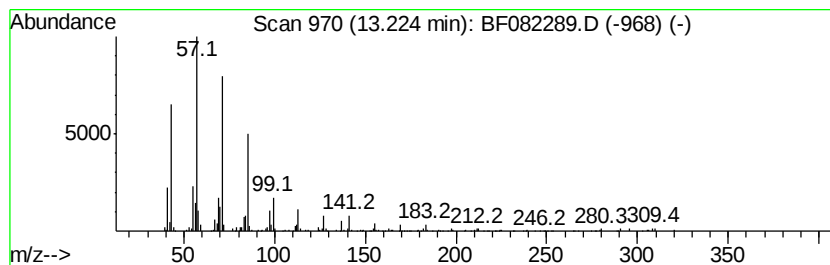
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Docosane Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.22 | 16.18 ng | 395282 | Chrysene-d12     | 14.04 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Docosane               | 310 | C22H46  | 000629-97-0 | 91   |
| 2    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |
| 3    |    | Tetracosane            | 338 | C24H50  | 000646-31-1 | 91   |
| 4    |    | Tetratetracontane      | 619 | C44H90  | 007098-22-8 | 91   |
| 5    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 90   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

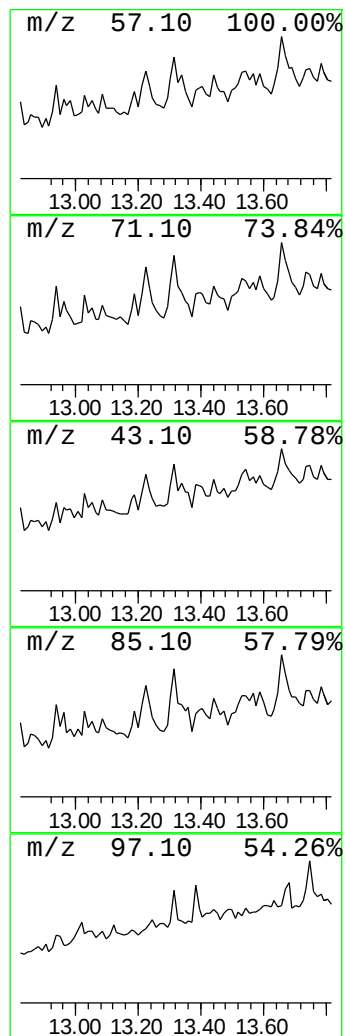
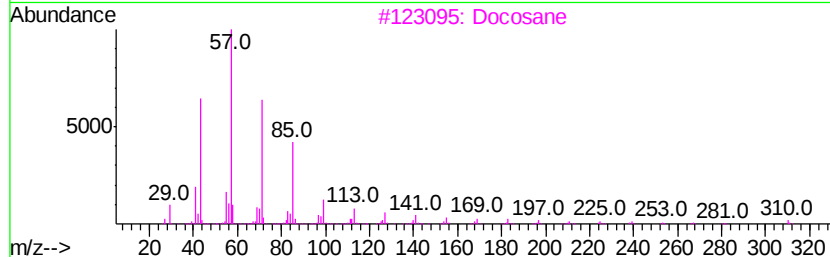
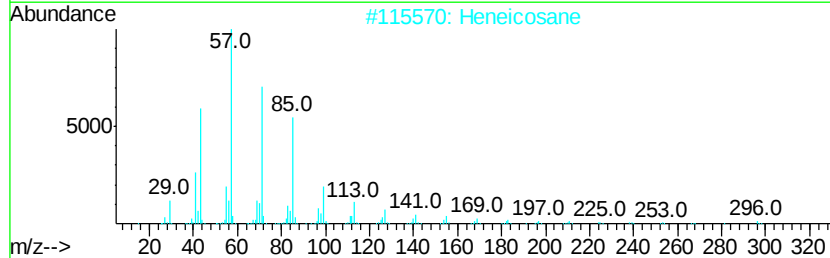
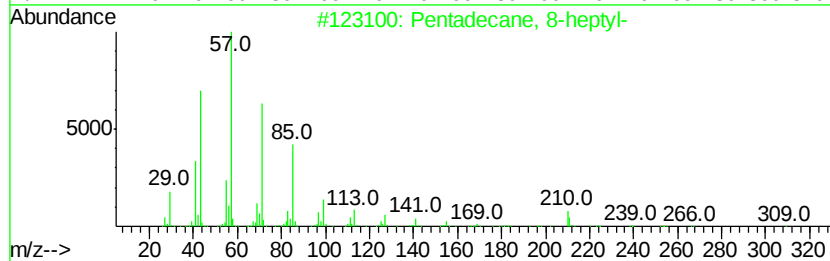
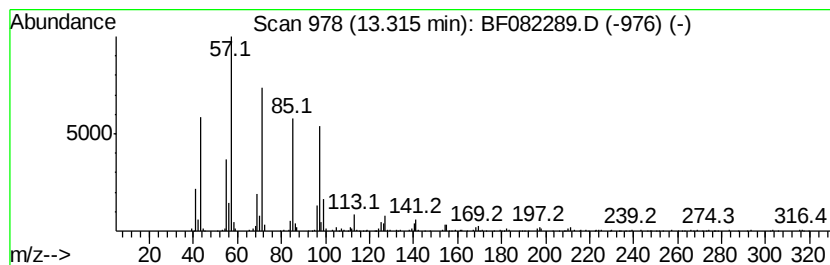
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Pentadecane, 8-heptyl- Concentration Rank 11

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.32 | 11.96 ng | 292201 | Chrysene-d12     | 14.04 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 64   |
| 2    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 64   |
| 3    |    | Docosane               | 310 | C22H46  | 000629-97-0 | 64   |
| 4    |    | Tetradecane            | 198 | C14H30  | 000629-59-4 | 64   |
| 5    |    | Octadecane             | 254 | C18H38  | 000593-45-3 | 62   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
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Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

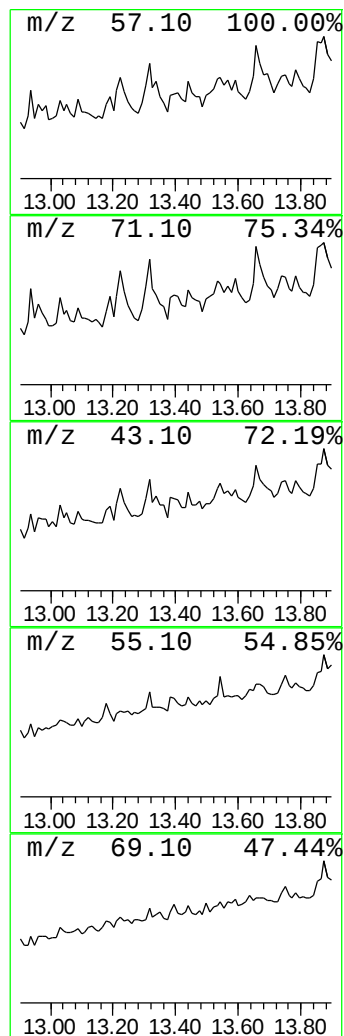
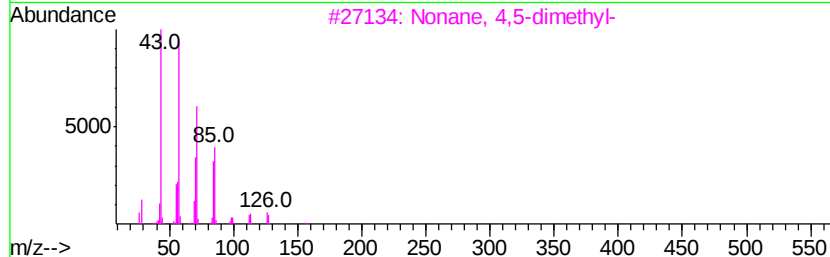
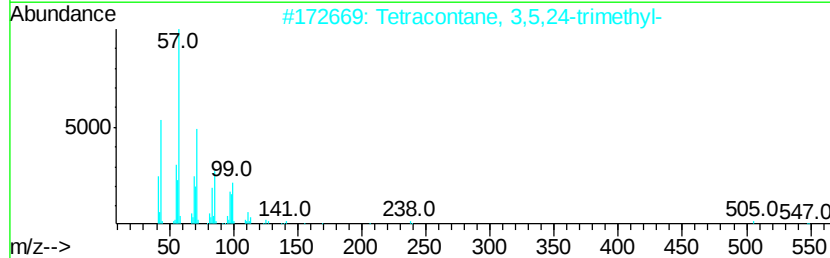
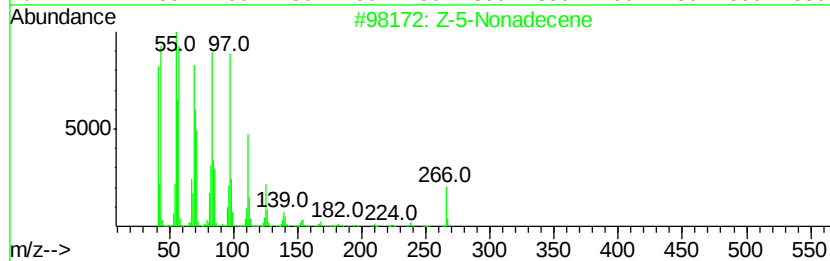
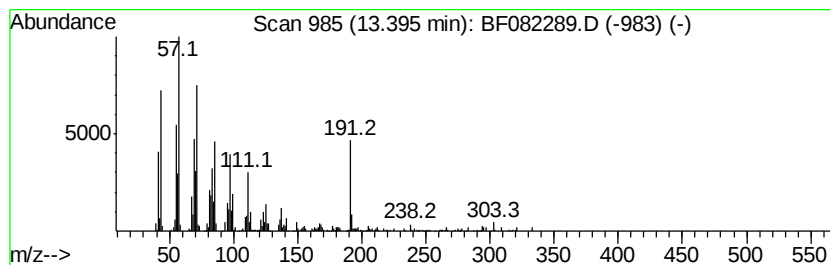
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ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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 15 Z-5-Nonadecene Concentration Rank 10

| R.T.      | EstConc                         | Area   | Relative to ISTD | R.T.         |      |
|-----------|---------------------------------|--------|------------------|--------------|------|
| 13.40     | 14.92 ng                        | 364569 | Chrysene-d12     | 14.04        |      |
| Hit# of 5 | Tentative ID                    | MW     | MolForm          | CAS#         | Qual |
| 1         | Z-5-Nonadecene                  | 266    | C19H38           | 1000131-11-8 | 60   |
| 2         | Tetracontane, 3,5,24-trimethyl- | 605    | C43H88           | 055162-61-3  | 52   |
| 3         | Nonane, 4,5-dimethyl-           | 156    | C11H24           | 017302-23-7  | 46   |
| 4         | Pentadecane                     | 212    | C15H32           | 000629-62-9  | 46   |
| 5         | Tetradecane, 1-bromo-           | 276    | C14H29Br         | 000112-71-0  | 45   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

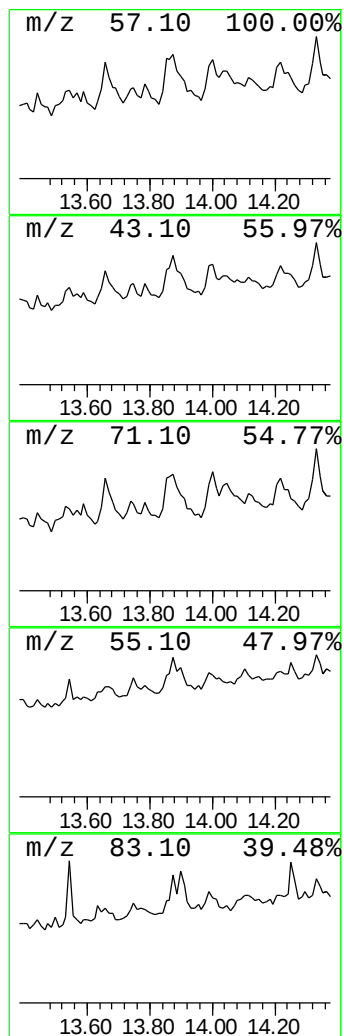
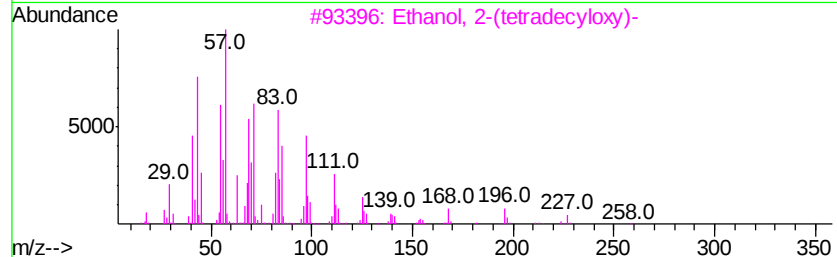
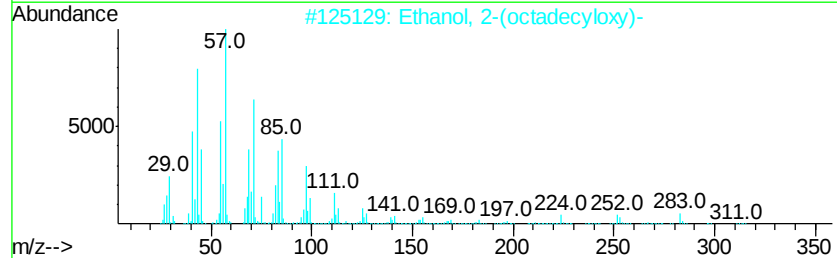
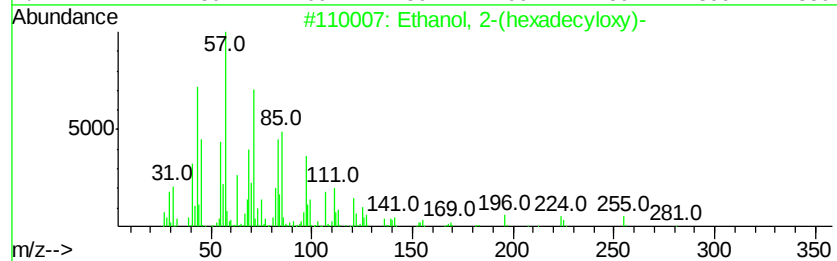
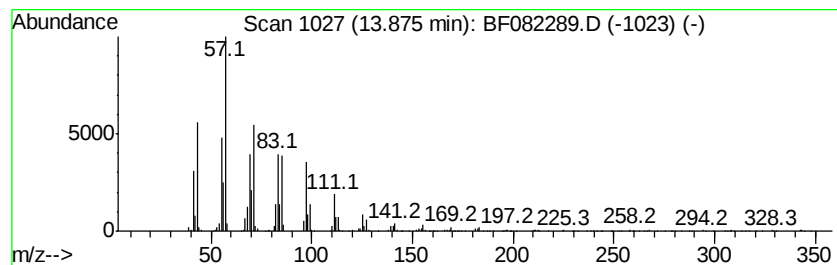
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Ethanol, 2-(hexadecyloxy)- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.88 | 51.90 ng | 1267920 | Chrysene-d12     | 14.04 |

| Hit# | of | Tentative ID                | MW  | MolForm  | CAS#         | Qual |
|------|----|-----------------------------|-----|----------|--------------|------|
| 1    | 5  | Ethanol, 2-(hexadecyloxy)-  | 286 | C18H38O2 | 002136-71-2  | 86   |
| 2    |    | Ethanol, 2-(octadecyloxy)-  | 314 | C20H42O2 | 002136-72-3  | 80   |
| 3    |    | Ethanol, 2-(tetradecyloxy)- | 258 | C16H34O2 | 002136-70-1  | 64   |
| 4    |    | E-15-Heptadecenal           | 252 | C17H32O  | 1000130-97-9 | 64   |
| 5    |    | 17-Pentatriacontene         | 491 | C35H70   | 006971-40-0  | 62   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

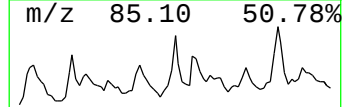
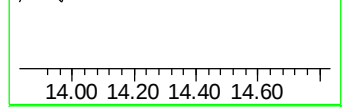
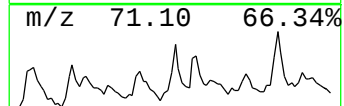
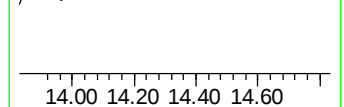
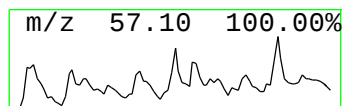
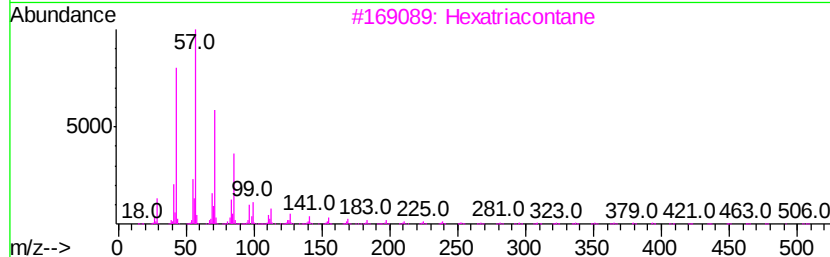
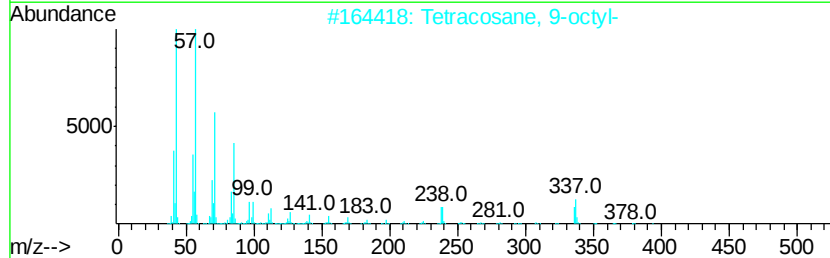
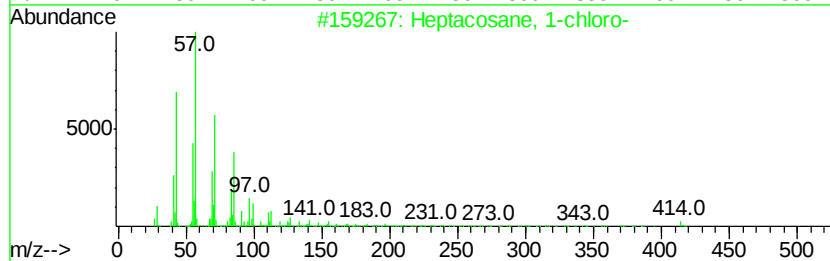
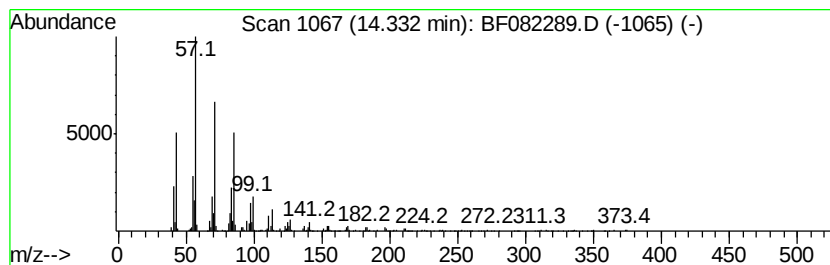
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Tetracosane, 9-octyl- Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.33 | 22.79 ng | 556738 | Chrysene-d12     | 14.04 |

| Hit# | of | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptacosane, 1-chloro- | 414 | C27H55Cl | 062016-79-9 | 91   |
| 2    |    | Tetracosane, 9-octyl-  | 451 | C32H66   | 055401-54-2 | 90   |
| 3    |    | Hexatriacontane        | 507 | C36H74   | 000630-06-8 | 86   |
| 4    |    | Heneicosane            | 296 | C21H44   | 000629-94-7 | 83   |
| 5    |    | Tetracosane            | 338 | C24H50   | 000646-31-1 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

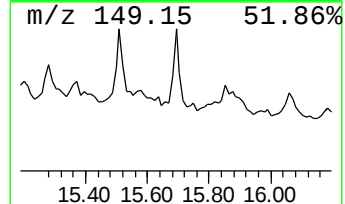
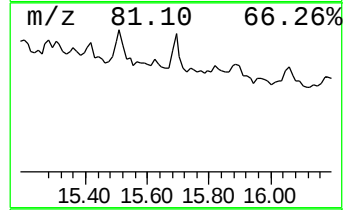
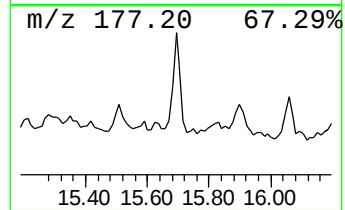
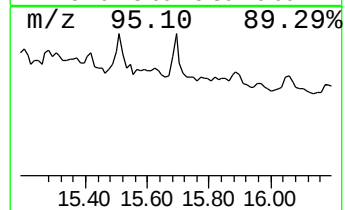
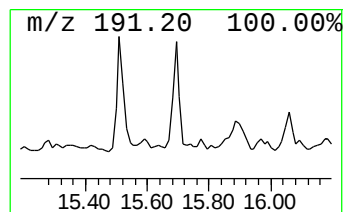
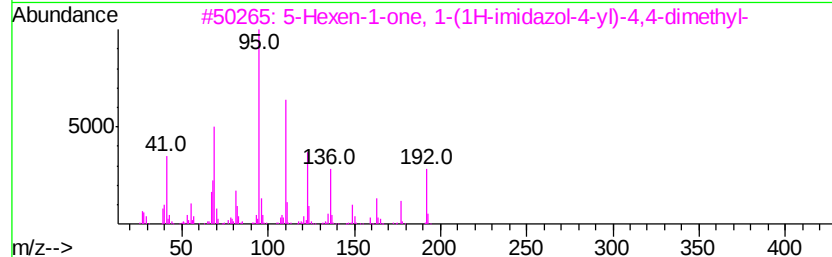
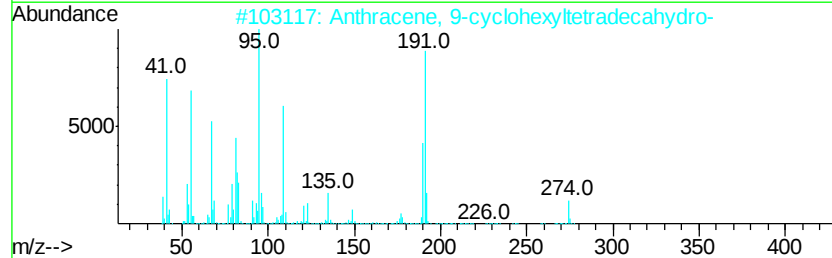
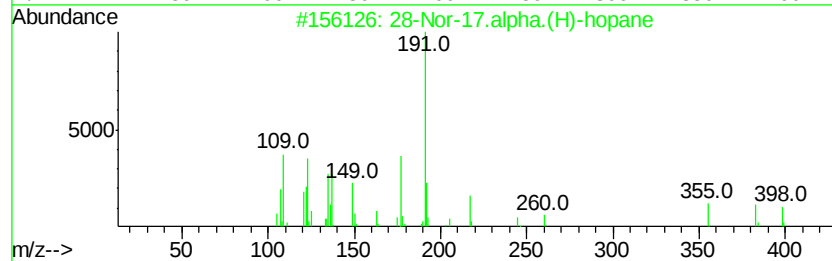
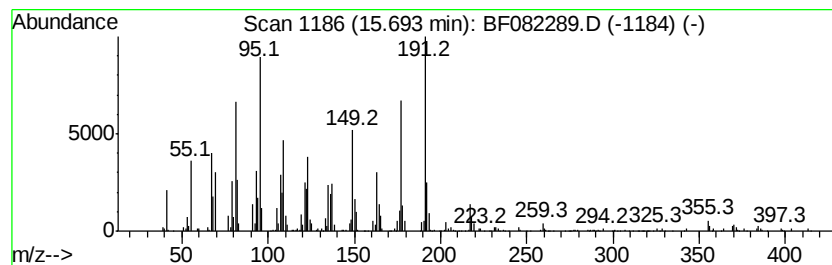
Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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 18 28-Nor-17.alpha.(H)-hopane Concentration Rank 8

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.    |              |      |
|---------|-------------------------------------|--------------|------------------|---------|--------------|------|
| 15.69   | 20.00 ng                            | 763550       | Perylene-d12     | 15.57   |              |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm | CAS#         | Qual |
| 1       | 28-Nor-17.alpha.(H)-hopane          | 398          | C29H50           |         | 053584-60-4  | 59   |
| 2       | Anthracene, 9-cyclohexyltetradec... | 274          | C20H34           |         | 055255-70-4  | 43   |
| 3       | 5-Hexen-1-one, 1-(1H-imidazol-4-... | 192          | C11H16N2O        |         | 069393-41-5  | 30   |
| 4       | Pyrido[3,4-d]pvridazin-1(2H)-one... | 191          | C8H9N5O          |         | 1000263-69-7 | 25   |
| 5       | 5-(p-Aminophenyl)-2-thiazolamine    | 191          | C9H9N3S          |         | 090349-87-4  | 14   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

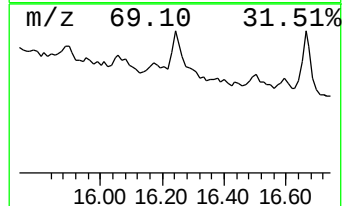
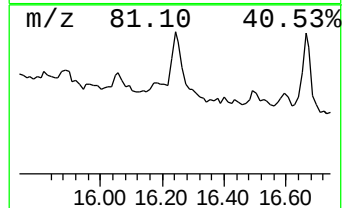
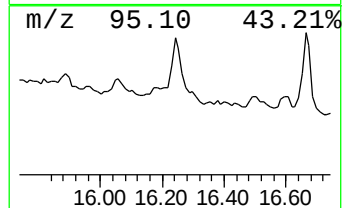
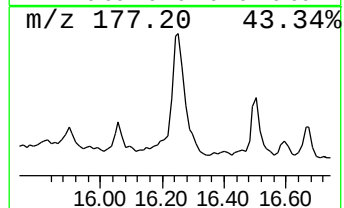
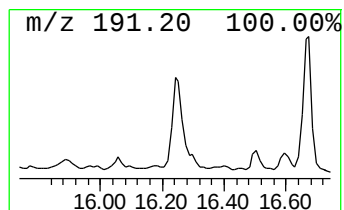
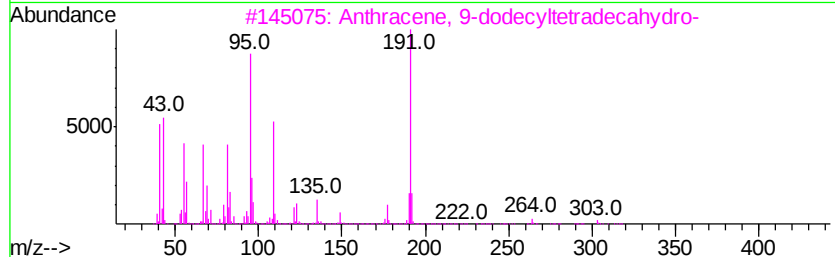
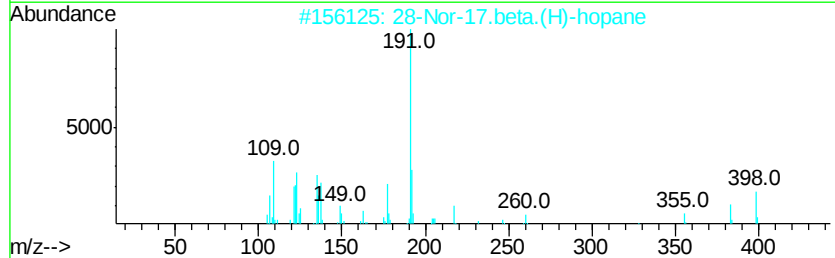
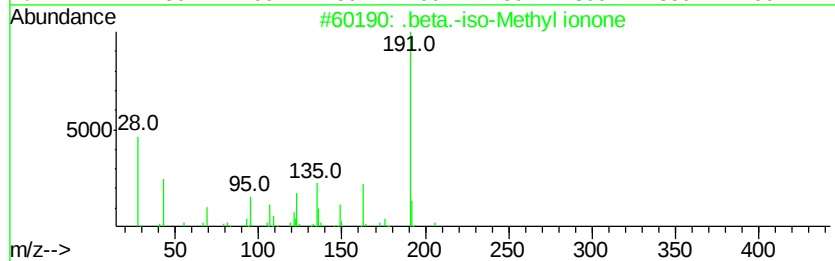
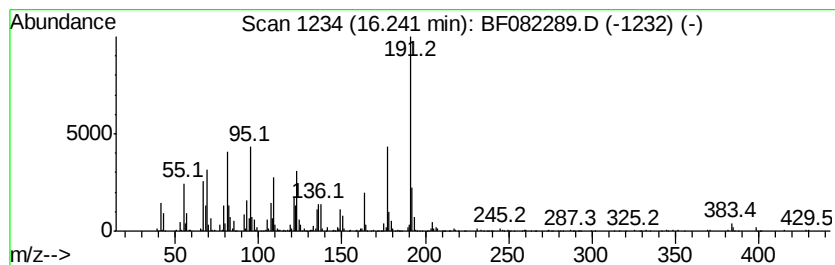
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown16.24 Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 16.24 | 50.53 ng | 1929170 | Perylene-d12     | 15.57 |

| Hit# | of | Tentative ID                                                       | MW  | MolForm      | CAS#         | Qual |
|------|----|--------------------------------------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | .beta.-iso-Methyl ionone                                           | 206 | C14H22O      | 1000285-40-2 | 43   |
| 2    |    | 28-Nor-17.beta.(H)-hopane                                          | 398 | C29H50       | 036728-72-0  | 38   |
| 3    |    | Anthracene, 9-dodecyltetradecahydro-                               | 360 | C26H48       | 055401-75-7  | 38   |
| 4    |    | 2,4,5,5,8a-Pentamethyl-6,7,8,8a-tetrahydronaphthalene              | 206 | C14H22O      | 1000195-40-9 | 32   |
| 5    |    | 2-(2-((2-Chloro-6-fluorobenzyl)sulfonyl)ethyl)-1-methylpyrrolidine | 334 | C17H16ClFN2S | 1000298-18-2 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
Data File : BF082289.D  
Acq On : 14 Oct 2015 20:07  
Operator : UM/IZ  
Sample : G4016-06  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

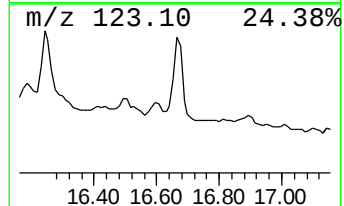
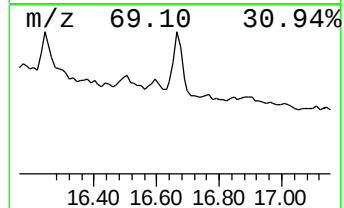
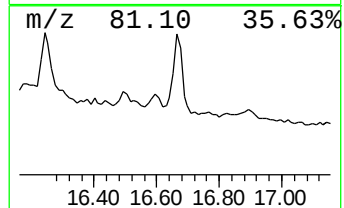
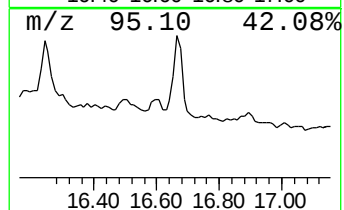
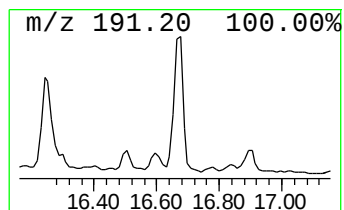
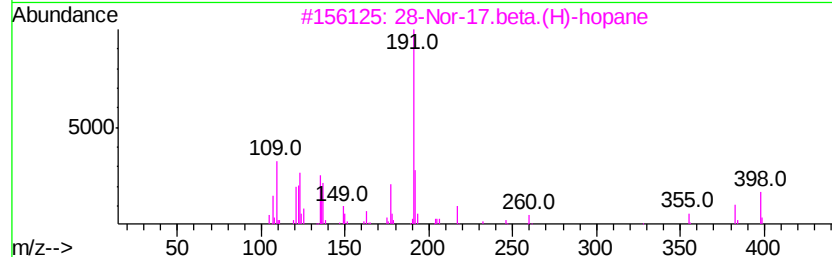
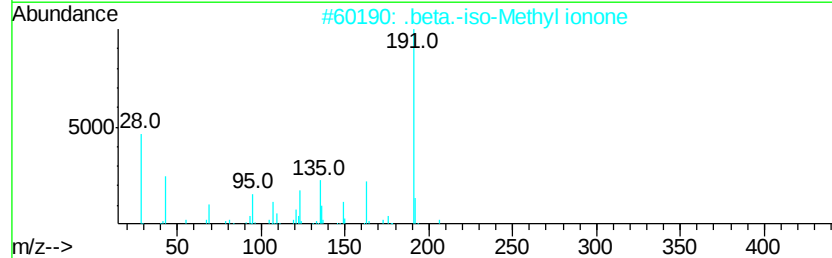
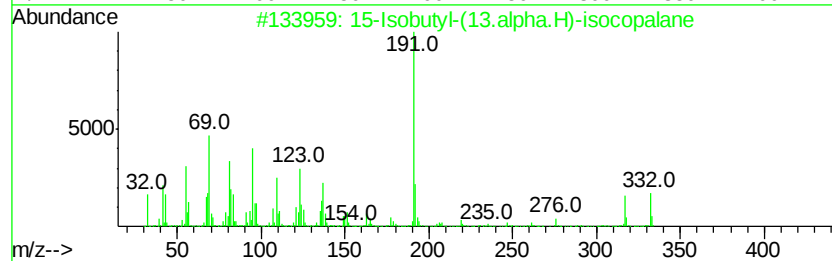
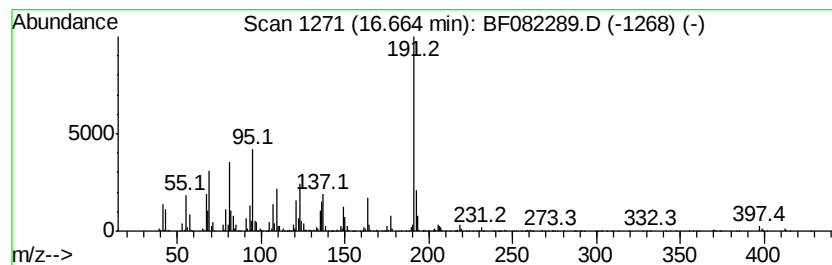
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 15-Isobutyl-(13.alpha.H)-is... Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 16.66 | 47.18 ng | 1801160 | Perylene-d12     | 15.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 15-Isobutyl-(13.alpha.H)-isocopa... | 332 | C24H44  | 087953-47-7  | 64   |
| 2    |    | .beta.-iso-Methyl ionone            | 206 | C14H22O | 1000285-40-2 | 58   |
| 3    |    | 28-Nor-17.beta.(H)-hopane           | 398 | C29H50  | 036728-72-0  | 56   |
| 4    |    | Anthracene, 9-dodecyltetradecahy... | 360 | C26H48  | 055401-75-7  | 53   |
| 5    |    | 1-Cyclohexene, 1,3,3-trimethyl-2... | 206 | C14H22O | 1000197-08-4 | 45   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF101515\  
 Data File : BF082289.D  
 Acq On : 14 Oct 2015 20:07  
 Operator : UM/IZ  
 Sample : G4016-06  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CS-6

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF101015.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 |
| 4-Penten-2-one, 4... | 3.29  | 6.2     | ng    | 123228   | 1                     | 6.82  | 395683 | 20.0 |
| 3-Penten-2-one, 4... | 4.34  | 304.0   | ng    | 6014010  | 1                     | 6.82  | 395683 | 20.0 |
| 2-Pentanone, 4-hy... | 5.02  | 195.1   | ng    | 3860740  | 1                     | 6.82  | 395683 | 20.0 |
| unknown6.58          | 6.58  | 5.2     | ng    | 102207   | 1                     | 6.82  | 395683 | 20.0 |
| 2-Cyclohexen-1-on... | 7.14  | 6.0     | ng    | 118367   | 1                     | 6.82  | 395683 | 20.0 |
| unknown11.81         | 11.81 | 2.6     | ng    | 89537    | 4                     | 11.36 | 677509 | 20.0 |
| Eicosane, 2-methyl-  | 12.34 | 6.9     | ng    | 232518   | 4                     | 11.36 | 677509 | 20.0 |
| Eicosane             | 12.45 | 3.2     | ng    | 107532   | 4                     | 11.36 | 677509 | 20.0 |
| Heptacosane, 1-ch... | 12.55 | 4.1     | ng    | 137542   | 4                     | 11.36 | 677509 | 20.0 |
| Heptadecane, 2,6,... | 12.69 | 8.3     | ng    | 282906   | 4                     | 11.36 | 677509 | 20.0 |
| n-Pentadecylcyclo... | 12.80 | 6.3     | ng    | 153938   | 5                     | 14.04 | 488603 | 20.0 |
| Docosane             | 13.22 | 16.2    | ng    | 395282   | 5                     | 14.04 | 488603 | 20.0 |
| Pentadecane, 8-he... | 13.32 | 12.0    | ng    | 292201   | 5                     | 14.04 | 488603 | 20.0 |
| Z-5-Nonadecene       | 13.40 | 14.9    | ng    | 364569   | 5                     | 14.04 | 488603 | 20.0 |
| Ethanol, 2-(hexad... | 13.88 | 51.9    | ng    | 1267920  | 5                     | 14.04 | 488603 | 20.0 |
| Tetracosane, 9-oc... | 14.33 | 22.8    | ng    | 556738   | 5                     | 14.04 | 488603 | 20.0 |
| 28-Nor-17.alpha.(... | 15.69 | 20.0    | ng    | 763550   | 6                     | 15.57 | 763550 | 20.0 |
| unknown16.24         | 16.24 | 50.5    | ng    | 1929170  | 6                     | 15.57 | 763550 | 20.0 |
| 15-Isobutyl-(13.a... | 16.66 | 47.2    | ng    | 1801160  | 6                     | 15.57 | 763550 | 20.0 |