

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\  
 Data File : BF088213.D  
 Acq On : 21 Jun 2016 11:52  
 Operator : UM/SJ  
 Sample : H3580-15  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 P8-B7(0-12)C

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-BF060716.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         | 1.655       | 5             | 6           | 14           | rVB      | 203953         | 323079        | 8.70%           | 1.615%        |
| 2         | 1.770       | 14            | 16          | 53           | rVB      | 1729301        | 3713168       | 100.00%         | 18.567%       |
| 3         | 4.821       | 281           | 283         | 291          | rBV      | 62913          | 107334        | 2.89%           | 0.537%        |
| 4         | 5.267       | 319           | 322         | 324          | rBV      | 1204885        | 1670712       | 44.99%          | 8.354%        |
| 5         | 6.330       | 412           | 415         | 417          | rBV      | 1386892        | 1800628       | 48.49%          | 9.004%        |
| 6         | 6.444       | 422           | 425         | 427          | rBV      | 1437509        | 1774486       | 47.79%          | 8.873%        |
| 7         | 6.662       | 443           | 444         | 447          | rBV      | 342569         | 465383        | 12.53%          | 2.327%        |
| 8         | 6.822       | 456           | 458         | 461          | rBV      | 1531609        | 1340976       | 36.11%          | 6.705%        |
| 9         | 7.233       | 492           | 494         | 497          | rBV      | 827304         | 976344        | 26.29%          | 4.882%        |
| 10        | 7.953       | 555           | 557         | 559          | rBV      | 796934         | 643722        | 17.34%          | 3.219%        |
| 11        | 9.027       | 649           | 651         | 654          | rBV      | 1801553        | 1894529       | 51.02%          | 9.473%        |
| 12        | 9.428       | 684           | 686         | 688          | rBV      | 619025         | 504054        | 13.57%          | 2.520%        |
| 13        | 9.702       | 707           | 710         | 712          | rBV      | 736309         | 658945        | 17.75%          | 3.295%        |
| 14        | 10.136      | 747           | 748         | 750          | rVB      | 46099          | 46512         | 1.25%           | 0.233%        |
| 15        | 10.491      | 776           | 779         | 781          | rBV      | 723917         | 820152        | 22.09%          | 4.101%        |
| 16        | 10.616      | 788           | 790         | 794          | rBV      | 41754          | 72068         | 1.94%           | 0.360%        |
| 17        | 11.176      | 837           | 839         | 841          | rBV      | 636439         | 523538        | 14.10%          | 2.618%        |
| 18        | 12.754      | 975           | 977         | 980          | rBV      | 1135978        | 1335425       | 35.96%          | 6.677%        |
| 19        | 13.702      | 1058          | 1060        | 1066         | rVV3     | 32416          | 82989         | 2.23%           | 0.415%        |
| 20        | 13.805      | 1066          | 1069        | 1071         | rVV      | 558866         | 560844        | 15.10%          | 2.804%        |
| 21        | 13.897      | 1075          | 1077        | 1081         | rBV3     | 20397          | 38908         | 1.05%           | 0.195%        |
| 22        | 14.639      | 1140          | 1142        | 1144         | rBV      | 183158         | 167582        | 4.51%           | 0.838%        |
| 23        | 15.177      | 1187          | 1189        | 1194         | rBV      | 405165         | 477806        | 12.87%          | 2.389%        |

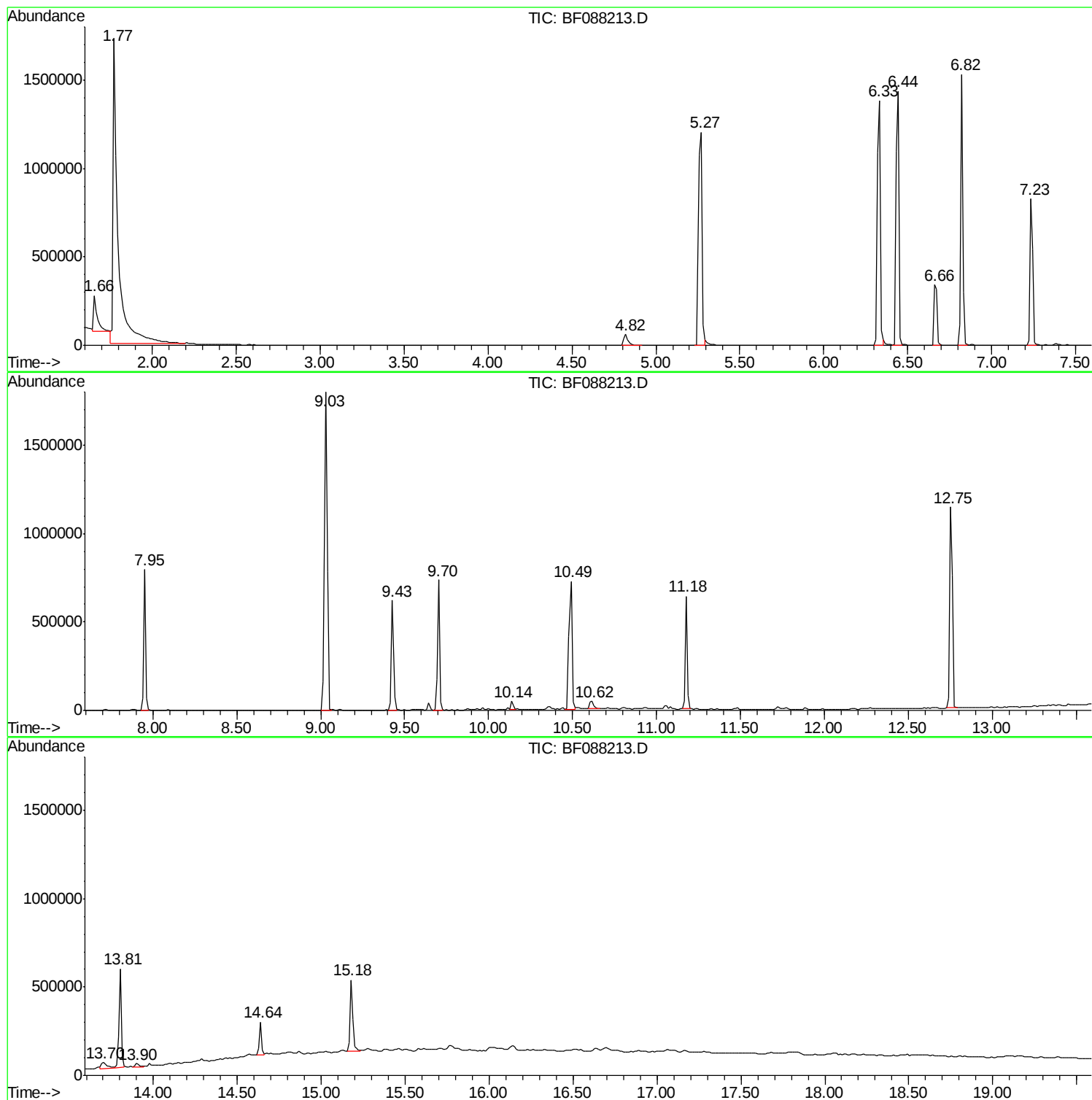
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ClientSampleId :  
P8-B7(0-12)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.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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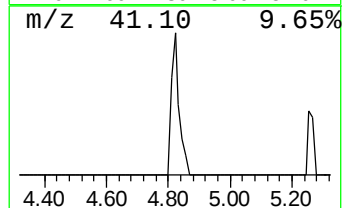
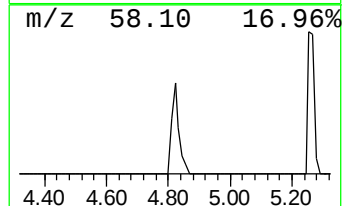
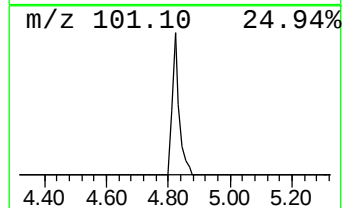
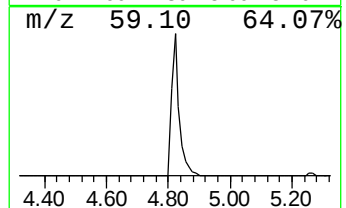
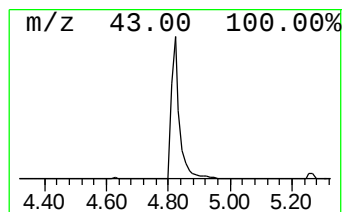
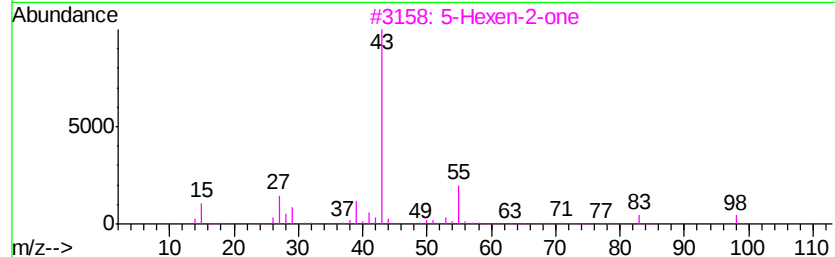
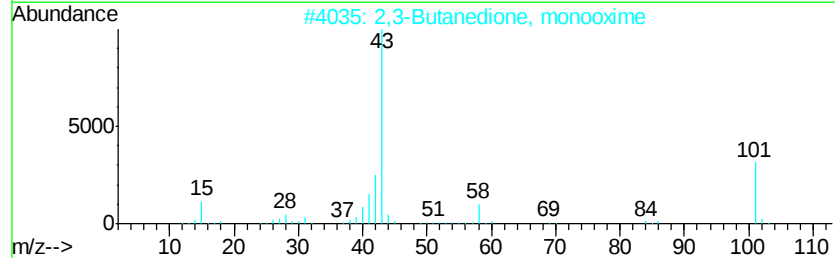
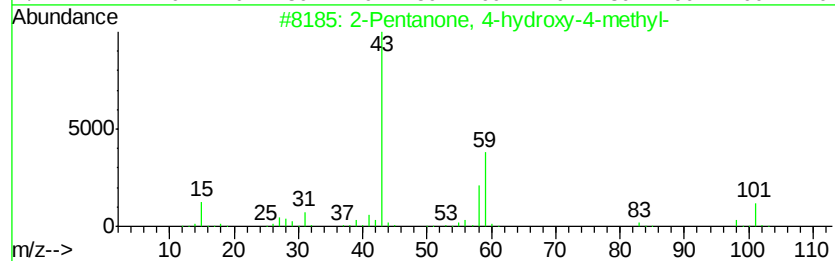
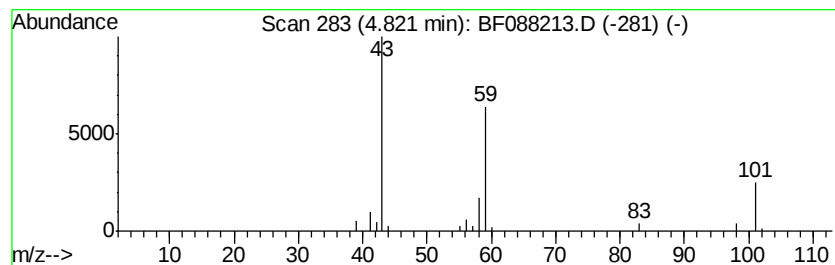
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.82 | 4.61 ng | 107334 | 1,4-Dichlorobenzene-d4 | 6.67 |

| Hit# | of                                 | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-   | 116 | C6H12O2      | 000123-42-2 | 56      |      |      |
| 2    | 2,3-Butanedione, monooxime         | 101 | C4H7NO2      | 000057-71-6 | 9       |      |      |
| 3    | 5-Hexen-2-one                      | 98  | C6H10O       | 000109-49-9 | 9       |      |      |
| 4    | Morpholine, 4-methyl-              | 101 | C5H11NO      | 000109-02-4 | 9       |      |      |
| 5    | (+)-4-Amino-4,5-dihydro-2(3H)-f... | 101 | C4H7NO2      | 016504-58-8 | 9       |      |      |



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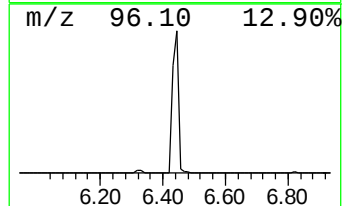
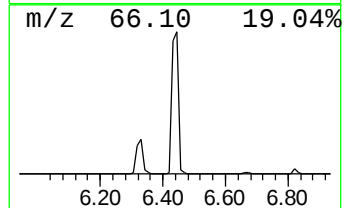
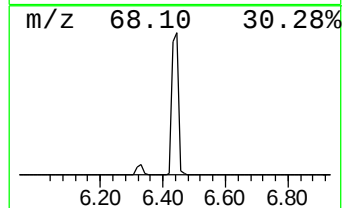
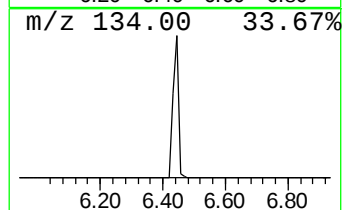
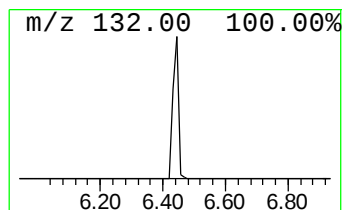
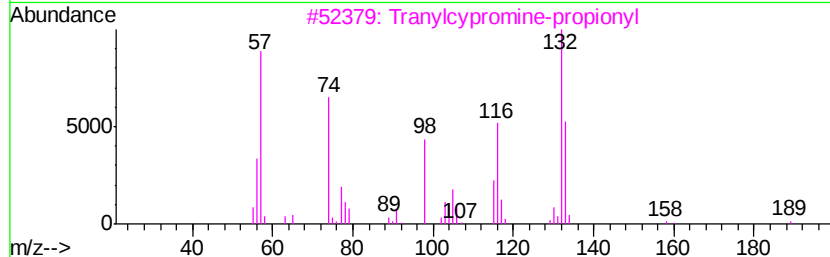
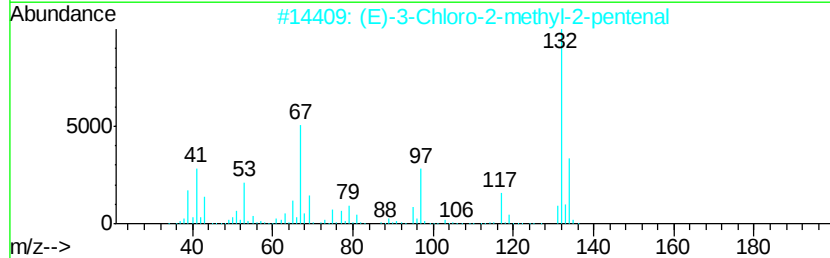
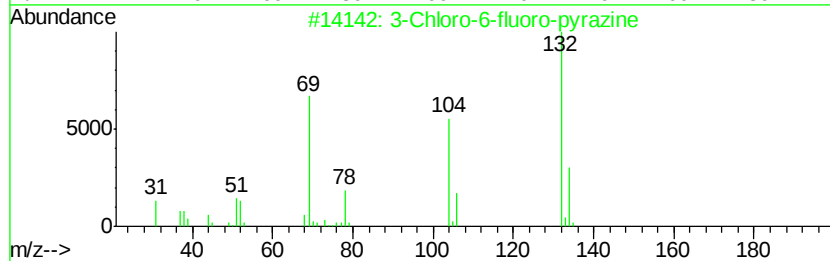
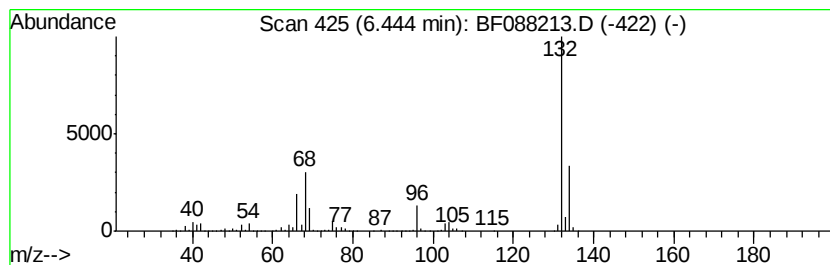
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 unknown6.44 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.44 | 76.26 ng | 1774490 | 1,4-Dichlorobenzene-d4 | 6.67 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Tranvlcypropine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | 1H-Indazole, 7-methyl-           | 132 | C8H8N2    | 1000316-00-7 | 10   |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



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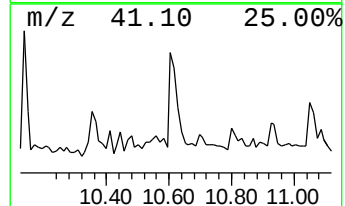
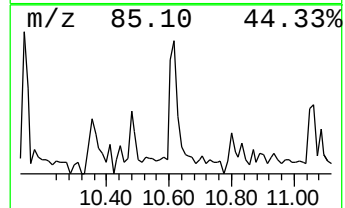
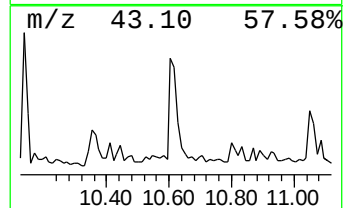
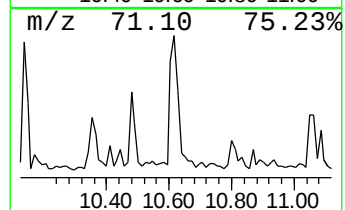
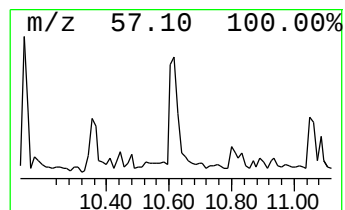
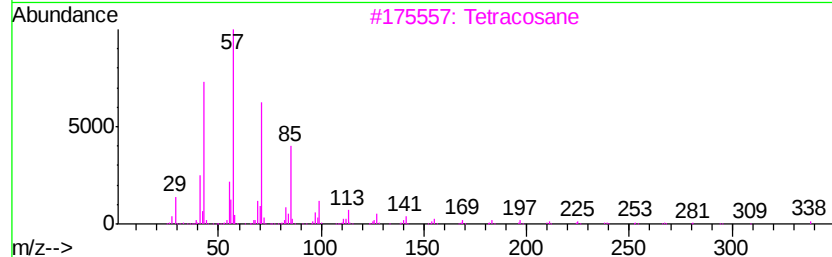
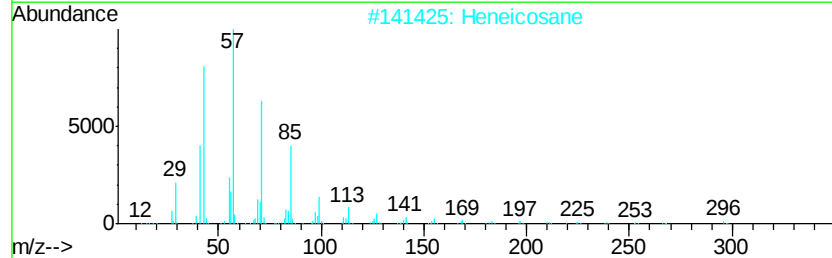
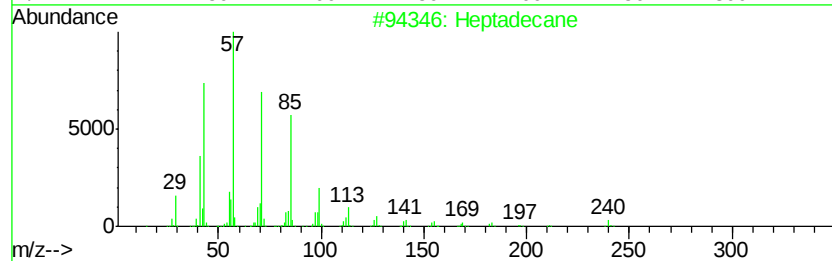
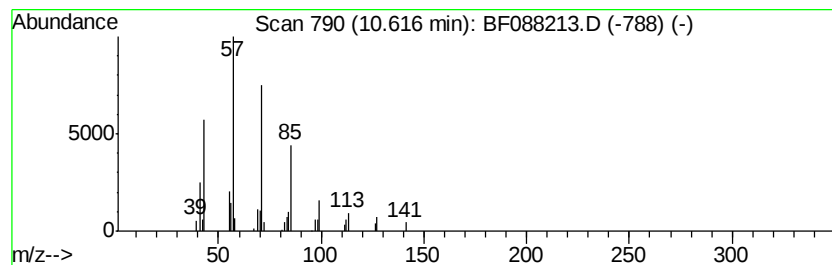
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BNA\_F  
ClientSampleId :  
P8-B7(0-12)C

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Peak Number 6 Heptadecane Concentration Rank 8

| R.T.      | EstConc      | Area  | Relative to ISTD |             | R.T.  |
|-----------|--------------|-------|------------------|-------------|-------|
| 10.62     | 2.75 ng      | 72068 | Phenanthrene-d10 |             | 11.18 |
| Hit# of 5 | Tentative ID | MW    | MolForm          | CAS#        | Qual  |
| 1         | Heptadecane  | 240   | C17H36           | 000629-78-7 | 91    |
| 2         | Heneicosane  | 296   | C21H44           | 000629-94-7 | 91    |
| 3         | Tetracosane  | 338   | C24H50           | 000646-31-1 | 91    |
| 4         | Hexadecane   | 226   | C16H34           | 000544-76-3 | 90    |
| 5         | Octadecane   | 254   | C18H38           | 000593-45-3 | 90    |



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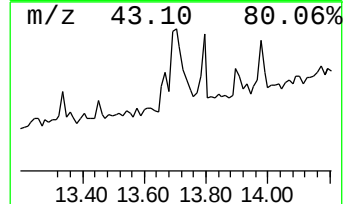
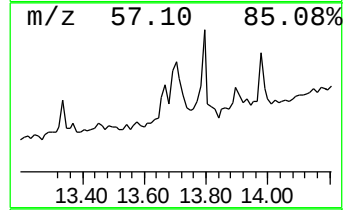
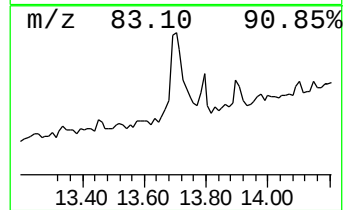
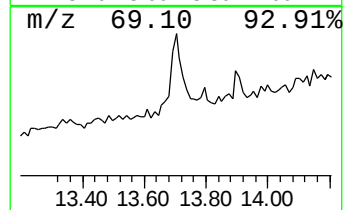
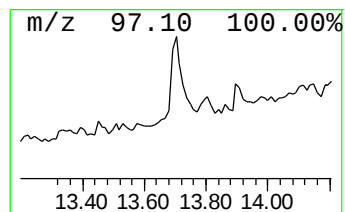
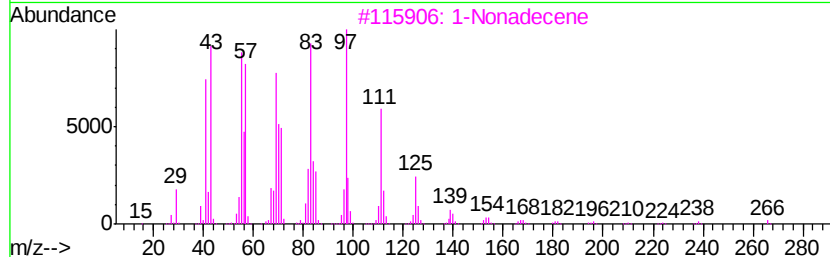
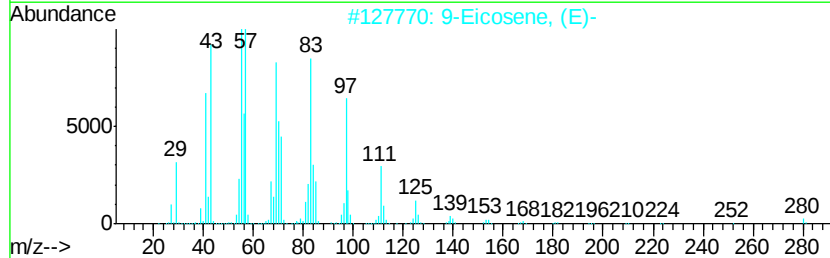
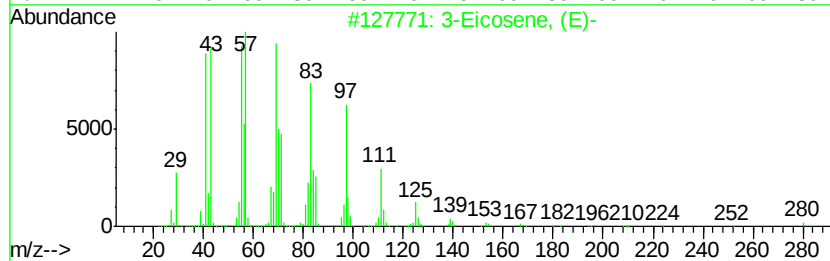
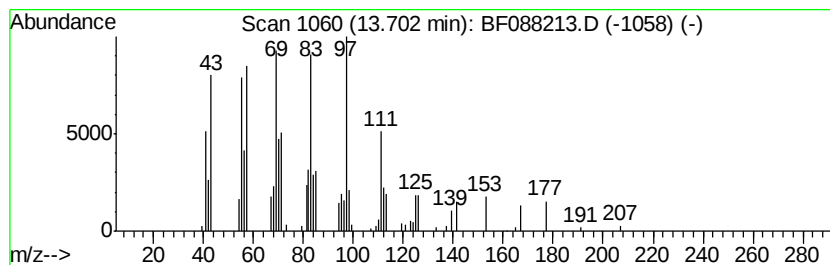
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\*\*\*\*\*  
Peak Number 7 3-Eicosene, (E)- Concentration Rank 7

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|-------|------------------|-------------|------|
| 13.70     | 2.96 ng                             | 82989 | Chrysene-d12     | 13.81       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#        | Qual |
| 1         | 3-Eicosene, (E)-                    | 280   | C20H40           | 074685-33-9 | 87   |
| 2         | 9-Eicosene, (E)-                    | 280   | C20H40           | 074685-29-3 | 87   |
| 3         | 1-Nonadecene                        | 266   | C19H38           | 018435-45-5 | 87   |
| 4         | Trichloroacetic acid, pentadecyl... | 372   | C17H31Cl3O2      | 074339-53-0 | 87   |
| 5         | Trifluoroacetoxy hexadecane         | 338   | C18H33F3O2       | 006222-03-3 | 87   |



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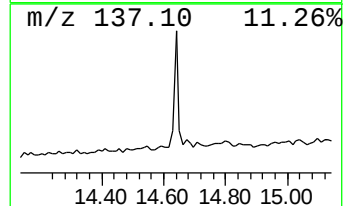
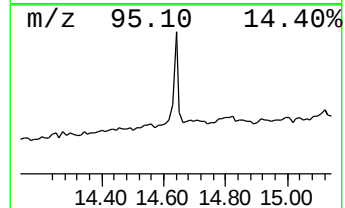
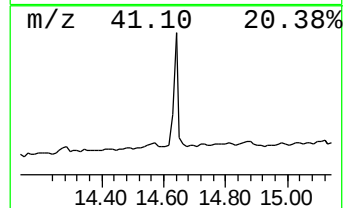
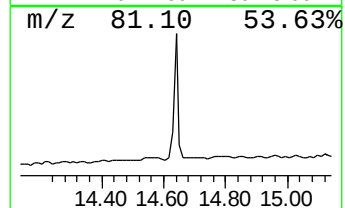
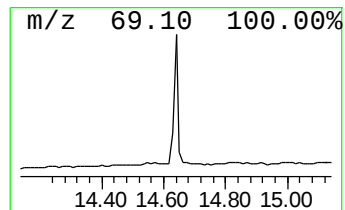
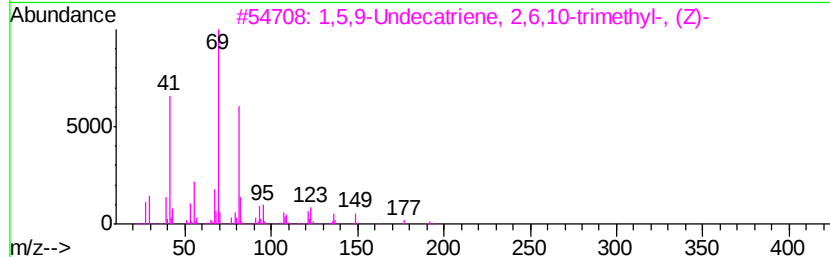
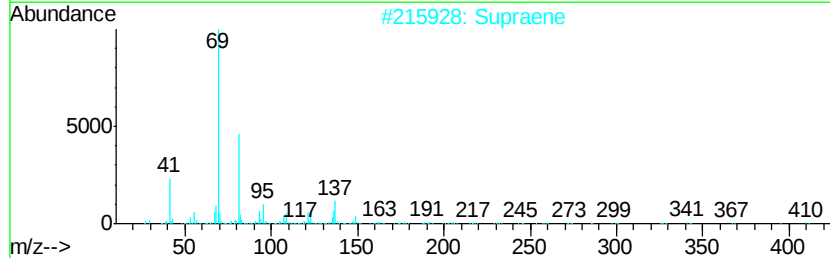
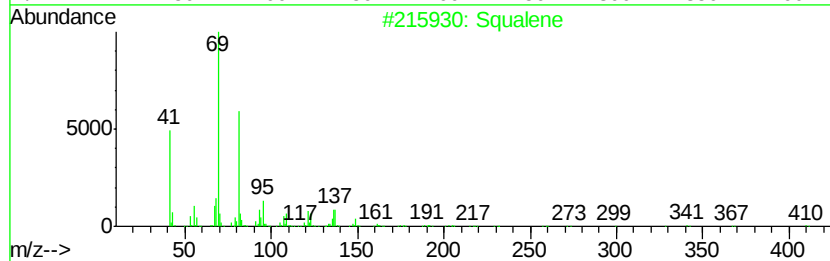
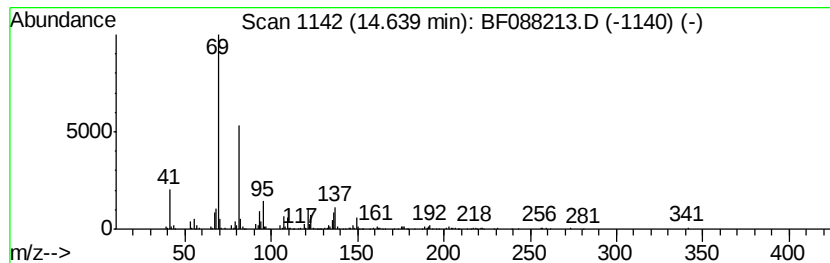
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Peak Number 8 Squalene Concentration Rank 5

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 14.64   | 7.01 ng                             | 167582       | Perylene-d12     | 15.18     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Squalene                            | 410 C30H50   | 000111-02-4      | 91        |
| 2       | Supraene                            | 410 C30H50   | 007683-64-9      | 91        |
| 3       | 1,5,9-Undecatriene, 2,6,10-trime... | 192 C14H24   | 062951-96-6      | 87        |
| 4       | 4,9,13,17-Tetramethyl-4,8,12,16-... | 316 C22H360  | 056882-09-8      | 83        |
| 5       | 4-Oxo-.beta.-damascone              | 206 C13H1802 | 053398-08-6      | 72        |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062016\  
Data File : BF088213.D  
Acq On : 21 Jun 2016 11:52  
Operator : UM/SJ  
Sample : H3580-15  
Misc :  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
P8-B7(0-12)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.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 |
| 2-Pentanone, 4-hy... | 4.82  | 4.6     | ng    | 107334   | 1                     | 6.67  | 465383 | 20.0 |
| unknown6.44          | 6.44  | 76.3    | ng    | 1774490  | 1                     | 6.67  | 465383 | 20.0 |
| Heptadecane          | 10.62 | 2.8     | ng    | 72068    | 4                     | 11.18 | 523538 | 20.0 |
| 3-Eicosene, (E)-     | 13.70 | 3.0     | ng    | 82989    | 5                     | 13.81 | 560844 | 20.0 |
| Squalene             | 14.64 | 7.0     | ng    | 167582   | 6                     | 15.18 | 477806 | 20.0 |