

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
 Data File : BF084870.D  
 Acq On : 5 Feb 2016 20:06  
 Operator : SJ/IZ  
 Sample : H1357-02  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EFF-WW

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-BF020116.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         | 4.578       | 215           | 218         | 221          | rBV      | 893721         | 1059579       | 1.24%           | 0.293%        |
| 2         | 4.807       | 235           | 238         | 242          | rBV      | 1353759        | 1624257       | 1.89%           | 0.449%        |
| 3         | 4.967       | 248           | 252         | 255          | rBV      | 2156885        | 2573503       | 3.00%           | 0.712%        |
| 4         | 5.253       | 276           | 277         | 280          | rVB      | 1618008        | 1754120       | 2.05%           | 0.485%        |
| 5         | 6.316       | 367           | 370         | 374          | rVV      | 691471         | 1359886       | 1.59%           | 0.376%        |
| 6         | 6.430       | 378           | 380         | 382          | rVV2     | 3588622        | 3886509       | 4.53%           | 1.075%        |
| 7         | 6.533       | 382           | 389         | 392          | rVB      | 863502         | 3143904       | 3.67%           | 0.869%        |
| 8         | 6.659       | 398           | 400         | 402          | rBV      | 1331892        | 1116867       | 1.30%           | 0.309%        |
| 9         | 6.784       | 405           | 411         | 412          | rBV      | 9480347        | 24662767      | 28.77%          | 6.820%        |
| 10        | 6.819       | 412           | 414         | 416          | rVV      | 4373902        | 4283725       | 5.00%           | 1.185%        |
| 11        | 7.070       | 433           | 436         | 441          | rBV      | 2223264        | 2937623       | 3.43%           | 0.812%        |
| 12        | 7.242       | 447           | 451         | 452          | rBV      | 2590355        | 4490306       | 5.24%           | 1.242%        |
| 13        | 7.276       | 452           | 454         | 461          | rVV2     | 456887         | 1303018       | 1.52%           | 0.360%        |
| 14        | 7.824       | 461           | 502         | 504          | rVV      | 7113244        | 85717967      | 100.00%         | 23.705%       |
| 15        | 7.950       | 507           | 513         | 514          | rBV      | 3516111        | 6242100       | 7.28%           | 1.726%        |
| 16        | 8.042       | 514           | 521         | 524          | rVB2     | 4637867        | 19251562      | 22.46%          | 5.324%        |
| 17        | 8.396       | 549           | 552         | 559          | rBV2     | 569687         | 922191        | 1.08%           | 0.255%        |
| 18        | 9.002       | 593           | 605         | 606          | rBV      | 3339425        | 13189729      | 15.39%          | 3.648%        |
| 19        | 9.036       | 606           | 608         | 610          | rVB      | 5123844        | 5713678       | 6.67%           | 1.580%        |
| 20        | 9.093       | 610           | 613         | 614          | rBV      | 1227545        | 1585067       | 1.85%           | 0.438%        |
| 21        | 9.699       | 663           | 666         | 668          | rBV2     | 1790910        | 2256555       | 2.63%           | 0.624%        |
| 22        | 10.019      | 684           | 694         | 696          | rBV      | 4902385        | 15513484      | 18.10%          | 4.290%        |
| 23        | 10.488      | 732           | 735         | 737          | rVB      | 2536451        | 2488227       | 2.90%           | 0.688%        |
| 24        | 10.728      | 754           | 756         | 757          | rBV      | 2528669        | 2732431       | 3.19%           | 0.756%        |
| 25        | 10.762      | 757           | 759         | 760          | rVV      | 4694659        | 5983964       | 6.98%           | 1.655%        |
| 26        | 10.785      | 760           | 761         | 765          | rVB      | 3404932        | 3591251       | 4.19%           | 0.993%        |
| 27        | 10.888      | 767           | 770         | 772          | rVV      | 1405306        | 1537965       | 1.79%           | 0.425%        |
| 28        | 10.990      | 775           | 779         | 781          | rBV2     | 2196049        | 3861619       | 4.51%           | 1.068%        |
| 29        | 11.162      | 791           | 794         | 799          | rVB2     | 7273663        | 8408175       | 9.81%           | 2.325%        |
| 30        | 11.425      | 813           | 817         | 821          | rBV2     | 2255091        | 2739318       | 3.20%           | 0.758%        |
| 31        | 11.653      | 834           | 837         | 841          | rVV4     | 287453         | 958521        | 1.12%           | 0.265%        |
| 32        | 11.745      | 841           | 845         | 849          | rVV4     | 2036091        | 5992392       | 6.99%           | 1.657%        |
| 33        | 11.802      | 849           | 850         | 854          | rVV2     | 1237969        | 2781513       | 3.24%           | 0.769%        |
| 34        | 11.859      | 854           | 855         | 857          | rVV      | 764045         | 1249691       | 1.46%           | 0.346%        |

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 Max Peaks: 100  
 Peak Location: TOP

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 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 35 | 11.916 | 859  | 860  | 864  | rVV  | 1462168 | 2852285  | 3.33%  | 0.789% |
| 36 | 11.973 | 864  | 865  | 866  | rVV  | 973792  | 1042646  | 1.22%  | 0.288% |
| 37 | 12.019 | 866  | 869  | 873  | rVV2 | 7121055 | 15430701 | 18.00% | 4.267% |
| 38 | 12.076 | 873  | 874  | 878  | rVV  | 1570246 | 2611332  | 3.05%  | 0.722% |
| 39 | 12.133 | 878  | 879  | 881  | rVV  | 1166640 | 1243739  | 1.45%  | 0.344% |
| 40 | 12.202 | 881  | 885  | 887  | rVV3 | 568276  | 1391444  | 1.62%  | 0.385% |
| 41 | 12.248 | 887  | 889  | 891  | rVV  | 1602154 | 2609620  | 3.04%  | 0.722% |
| 42 | 12.293 | 891  | 893  | 894  | rVV  | 827652  | 1306269  | 1.52%  | 0.361% |
| 43 | 12.328 | 894  | 896  | 897  | rVV  | 666440  | 863701   | 1.01%  | 0.239% |
| 44 | 12.408 | 900  | 903  | 905  | rVV3 | 550957  | 1561102  | 1.82%  | 0.432% |
| 45 | 12.499 | 908  | 911  | 916  | rVV  | 2178307 | 4658058  | 5.43%  | 1.288% |
| 46 | 12.671 | 924  | 926  | 928  | rBV2 | 751729  | 1446959  | 1.69%  | 0.400% |
| 47 | 12.716 | 928  | 930  | 933  | rBV2 | 2158771 | 4055950  | 4.73%  | 1.122% |
| 48 | 12.762 | 933  | 934  | 936  | rVV  | 2936937 | 3949774  | 4.61%  | 1.092% |
| 49 | 12.796 | 936  | 937  | 941  | rVB2 | 1183598 | 1835206  | 2.14%  | 0.508% |
| 50 | 12.911 | 946  | 947  | 950  | rVV  | 1024011 | 2471237  | 2.88%  | 0.683% |
| 51 | 12.968 | 950  | 952  | 954  | rVV2 | 1109944 | 2583841  | 3.01%  | 0.715% |
| 52 | 13.014 | 954  | 956  | 960  | rVV2 | 1135311 | 3314009  | 3.87%  | 0.916% |
| 53 | 13.082 | 960  | 962  | 964  | rVV  | 2343049 | 3909613  | 4.56%  | 1.081% |
| 54 | 13.128 | 964  | 966  | 968  | rVV  | 1245446 | 2519305  | 2.94%  | 0.697% |
| 55 | 13.174 | 968  | 970  | 973  | rVV2 | 2333227 | 5221211  | 6.09%  | 1.444% |
| 56 | 13.231 | 973  | 975  | 980  | rVV  | 3633036 | 10048002 | 11.72% | 2.779% |
| 57 | 13.356 | 984  | 986  | 988  | rVV  | 5484022 | 5939879  | 6.93%  | 1.643% |
| 58 | 13.391 | 988  | 989  | 992  | rVV  | 595726  | 1009113  | 1.18%  | 0.279% |
| 59 | 13.436 | 992  | 993  | 996  | rVV  | 1793419 | 2825278  | 3.30%  | 0.781% |
| 60 | 13.528 | 996  | 1001 | 1005 | rVV  | 1875106 | 4510125  | 5.26%  | 1.247% |
| 61 | 13.585 | 1005 | 1006 | 1010 | rVV  | 569825  | 861629   | 1.01%  | 0.238% |
| 62 | 13.688 | 1012 | 1015 | 1018 | rVB  | 9597568 | 11142836 | 13.00% | 3.082% |
| 63 | 13.779 | 1022 | 1023 | 1027 | rBV2 | 1110334 | 2143826  | 2.50%  | 0.593% |
| 64 | 14.682 | 1100 | 1102 | 1107 | rVB  | 894854  | 1133431  | 1.32%  | 0.313% |
| 65 | 14.911 | 1120 | 1122 | 1131 | rVB  | 560587  | 860067   | 1.00%  | 0.238% |
| 66 | 15.174 | 1141 | 1145 | 1147 | rVV  | 725851  | 1059648  | 1.24%  | 0.293% |
| 67 | 17.585 | 1348 | 1356 | 1370 | rVV  | 1277417 | 6273834  | 7.32%  | 1.735% |

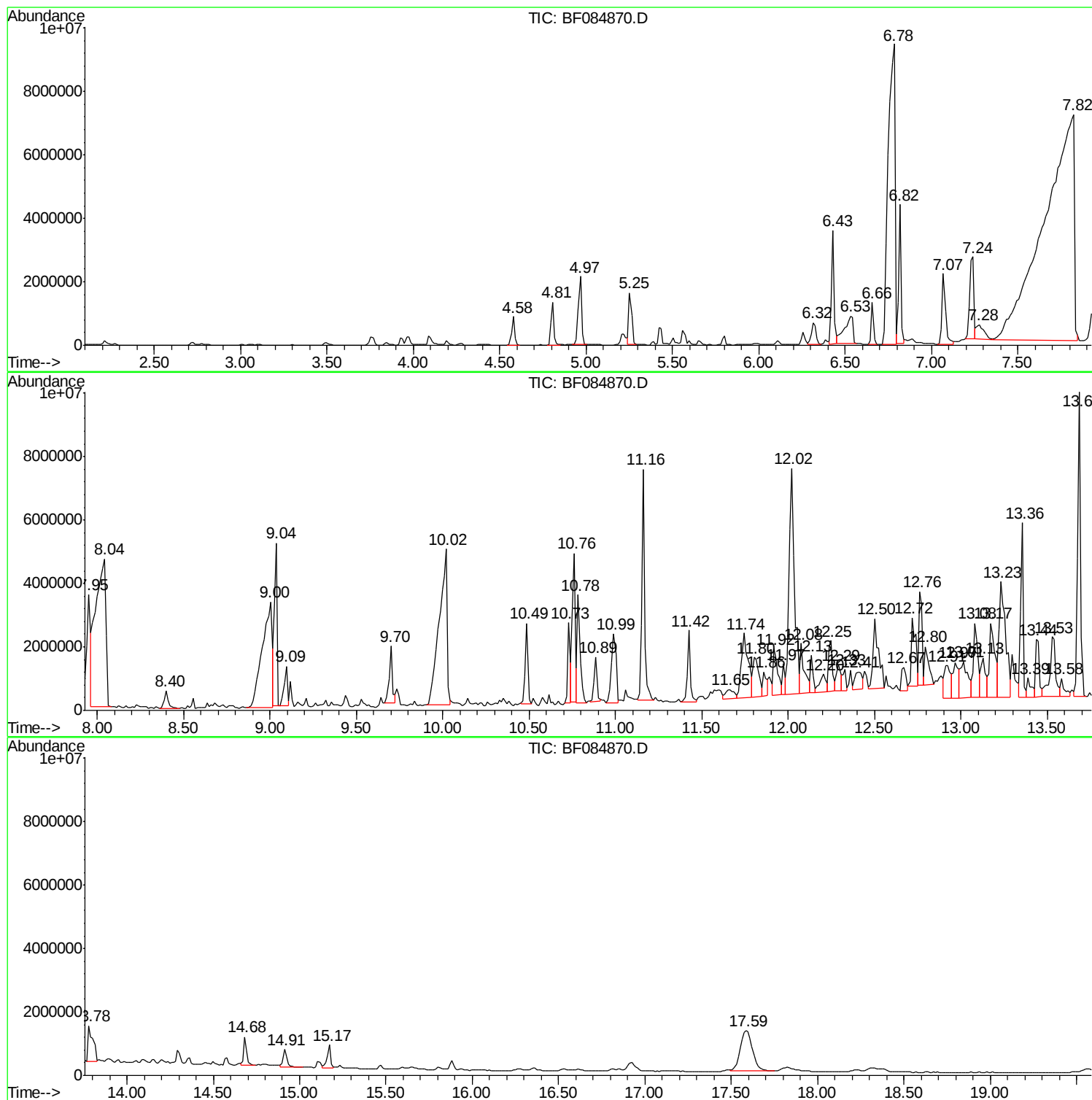
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Misc :  
ALS Vial : 18 Sample Multiplier: 1

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BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.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 : H1357-02  
Misc :  
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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

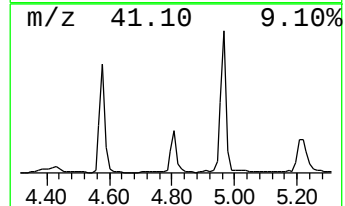
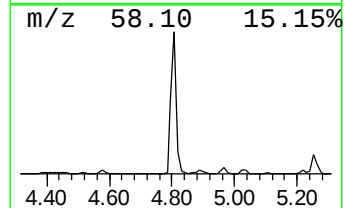
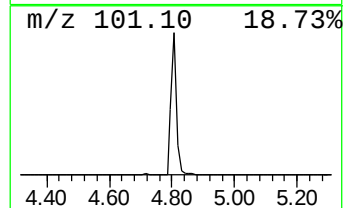
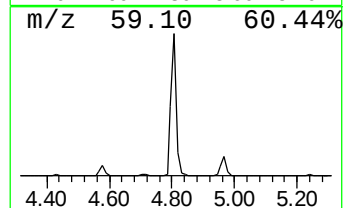
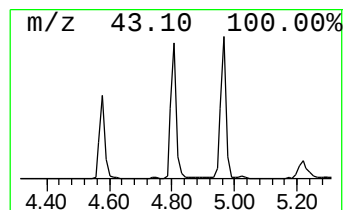
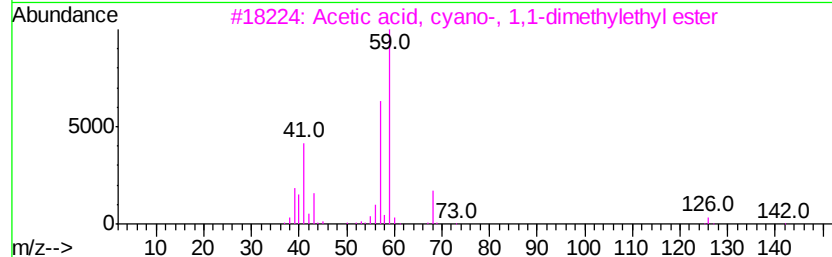
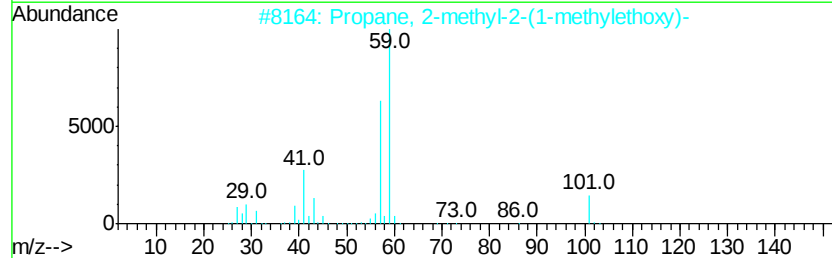
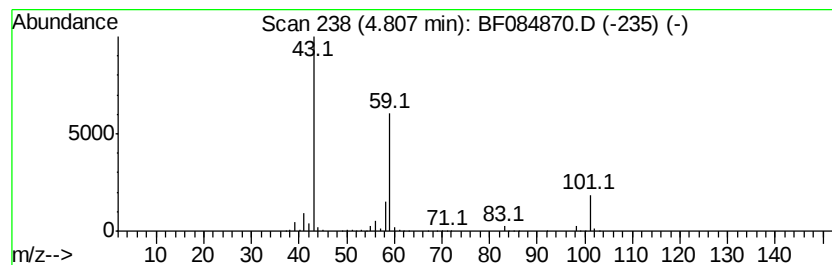
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 20

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.81 | 29.09 ng | 1624260 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-      | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2    |    |   | Propane, 2-methyl-2-(1-methylethoxy)- | 116 | C7H16O   | 017348-59-3 | 33   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethyl-    | 141 | C7H11NO2 | 001116-98-9 | 32   |
| 4    |    |   | Morpholine, 4-methyl-                 | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    |   | 2,3-Butanedione, monooxime            | 101 | C4H7NO2  | 000057-71-6 | 9    |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

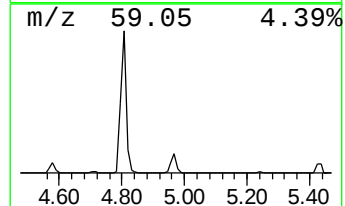
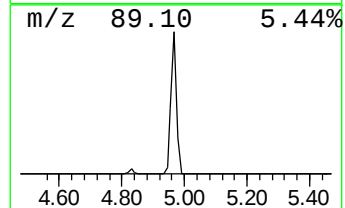
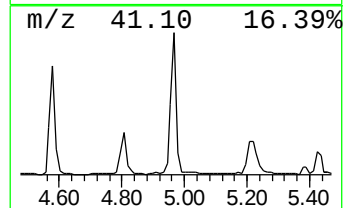
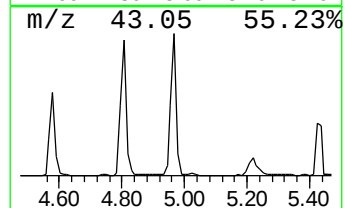
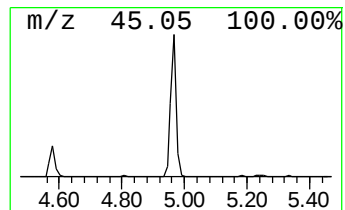
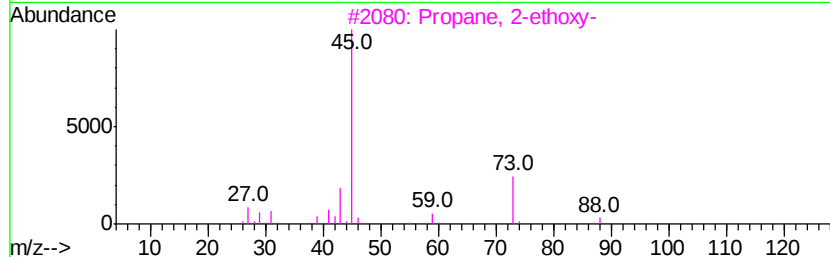
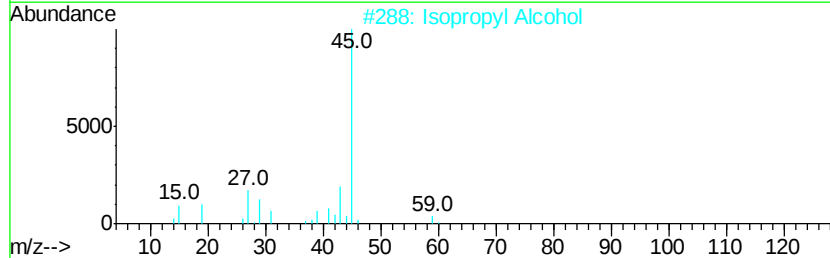
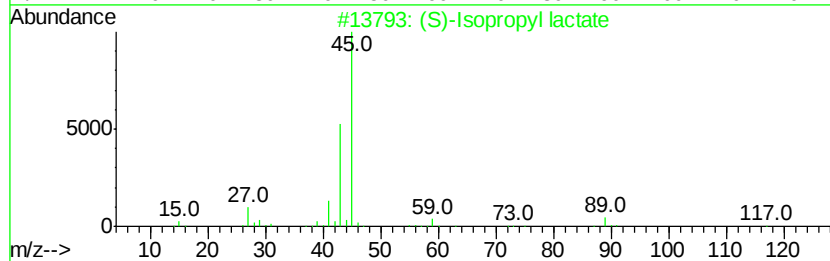
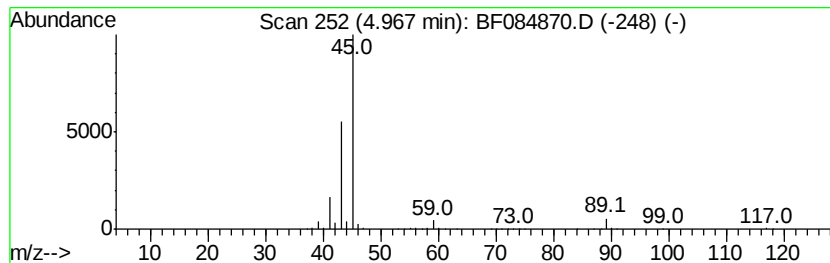
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (S)-Isopropyl lactate Concentration Rank 14

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.97 | 46.08 ng | 2573500 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# of | 5                           | Tentative ID | MW      | MolForm     | CAS# | Qual |
|---------|-----------------------------|--------------|---------|-------------|------|------|
| 1       | (S)-Isopropyl lactate       | 132          | C6H12O3 | 063697-00-7 | 78   |      |
| 2       | Isopropyl Alcohol           | 60           | C3H8O   | 000067-63-0 | 64   |      |
| 3       | Propane, 2-ethoxy-          | 88           | C5H12O  | 000625-54-7 | 45   |      |
| 4       | 2,3-Butanediol              | 90           | C4H10O2 | 000513-85-9 | 9    |      |
| 5       | (S)-2-Hydroxypropanoic acid | 90           | C3H6O3  | 000079-33-4 | 9    |      |



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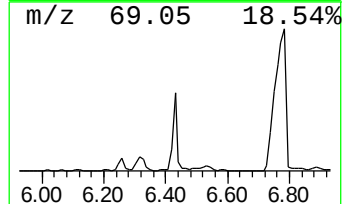
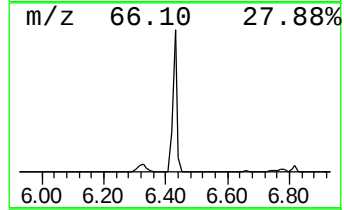
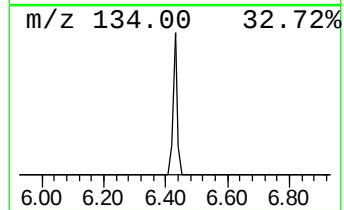
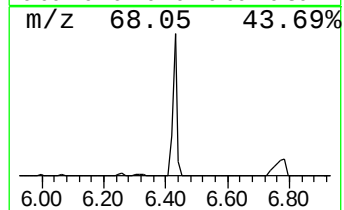
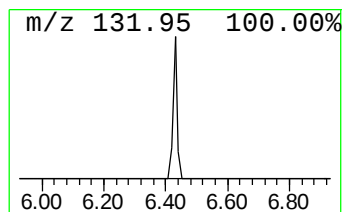
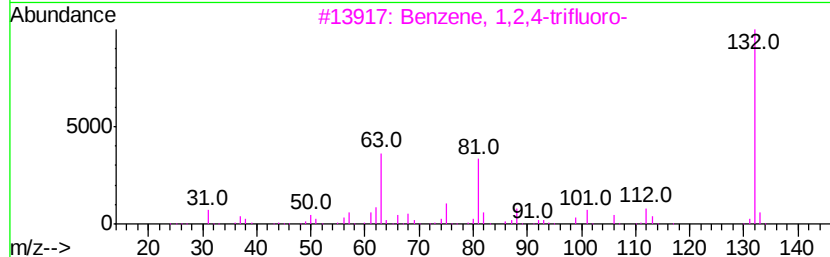
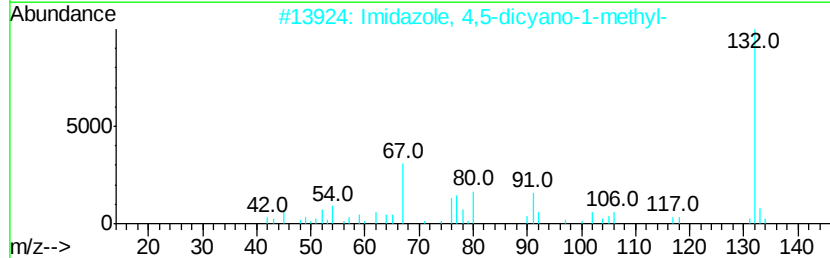
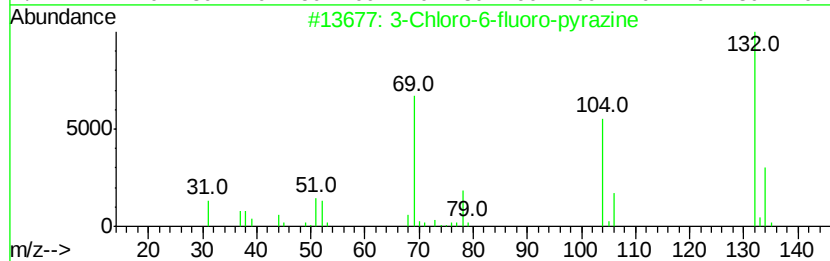
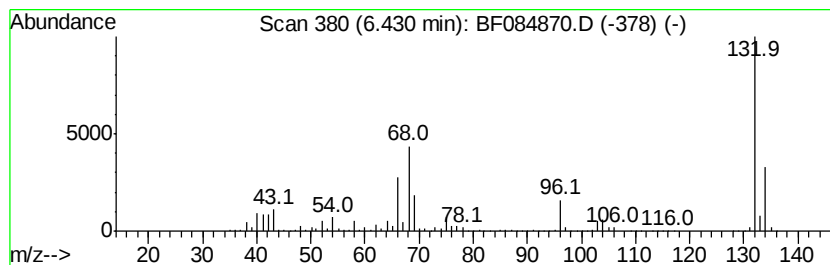
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown6.43 Concentration Rank 8

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.43 | 69.60 ng | 3886510 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | Imidazole, 4,5-dicyano-1-methyl- | 132 | C6H4N4    | 019485-35-9  | 9    |
| 3    |    | Benzene, 1,2,4-trifluoro-        | 132 | C6H3F3    | 000367-23-7  | 9    |
| 4    |    | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 9    |
| 5    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

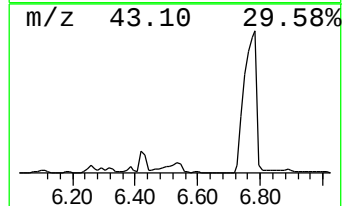
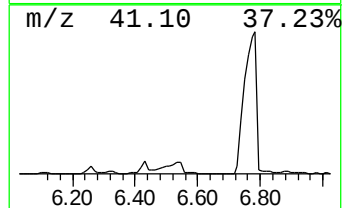
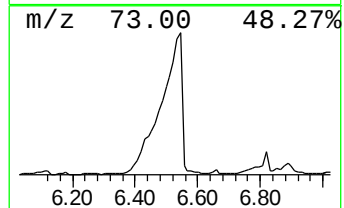
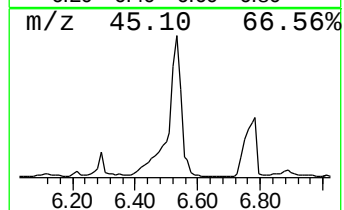
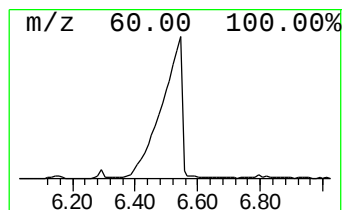
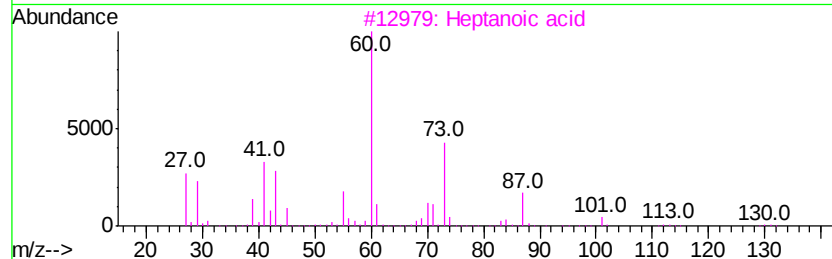
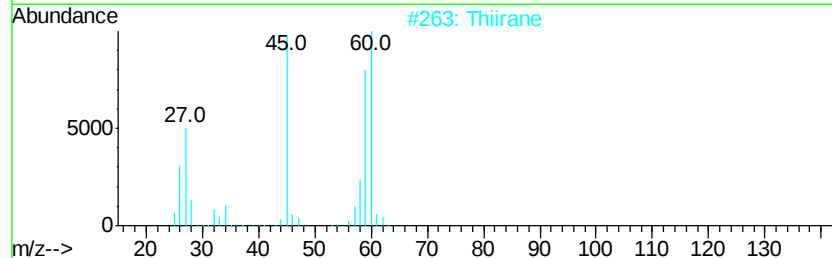
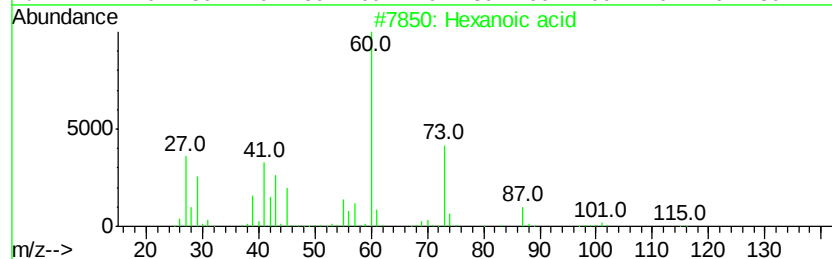
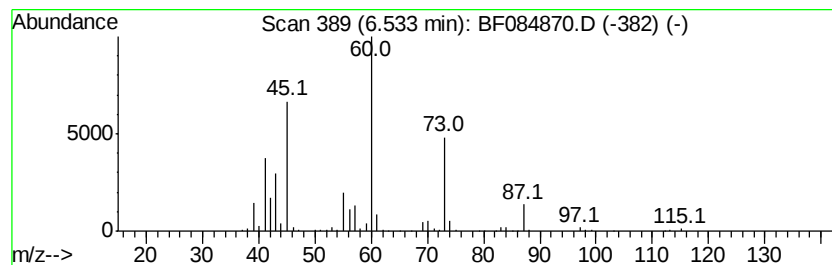
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Hexanoic acid Concentration Rank 10

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.53 | 56.30 ng | 3143900 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexanoic acid              | 116 | C6H12O2 | 000142-62-1 | 72   |
| 2    |    | Thiirane                   | 60  | C2H4S   | 000420-12-2 | 53   |
| 3    |    | Heptanoic acid             | 130 | C7H14O2 | 000111-14-8 | 50   |
| 4    |    | Propanedioic acid, propyl- | 146 | C6H10O4 | 000616-62-6 | 50   |
| 5    |    | Butanoic acid, 3-methyl-   | 102 | C5H10O2 | 000503-74-2 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

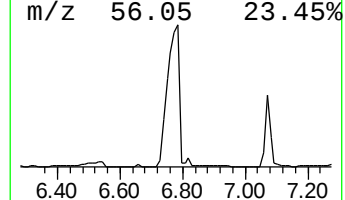
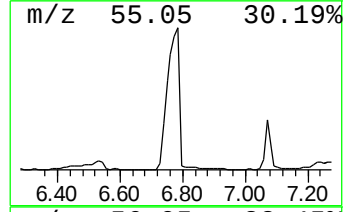
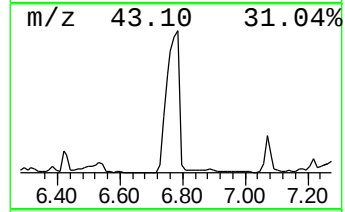
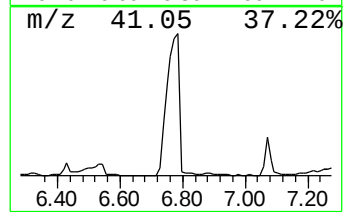
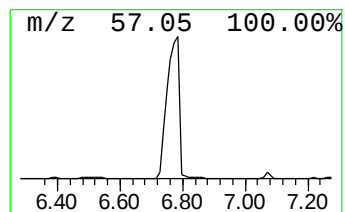
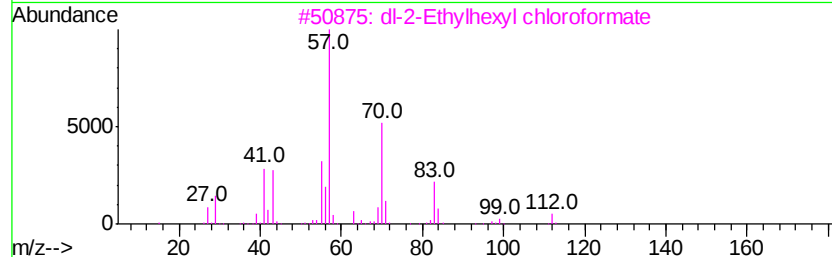
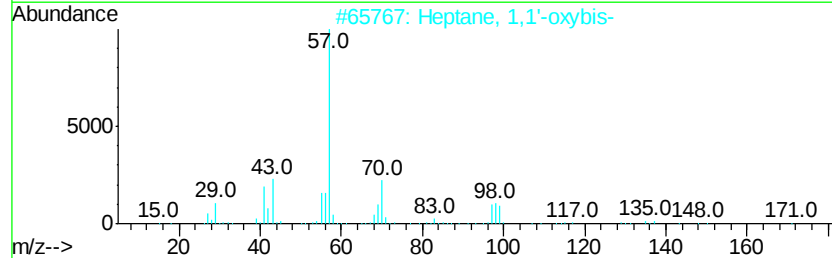
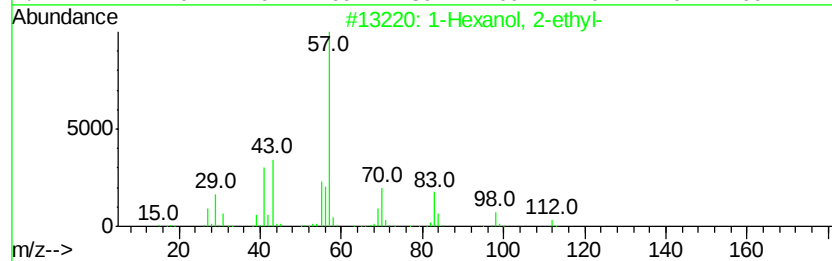
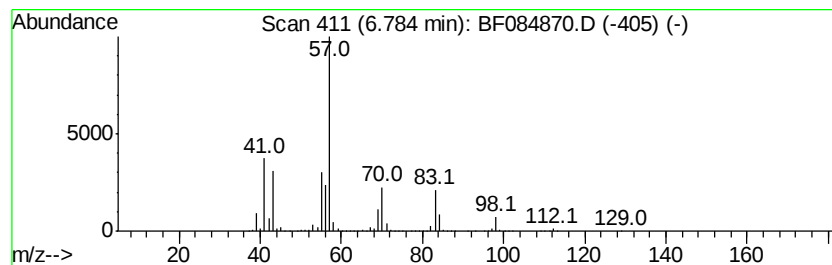
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 1

| R.T. | EstConc   | Area     | Relative to ISTD       | R.T. |
|------|-----------|----------|------------------------|------|
| 6.78 | 441.64 ng | 24662800 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                  | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1-Hexanol, 2-ethyl-           | 130 | C8H18O    | 000104-76-7 | 78   |
| 2    |    | Heptane, 1,1'-oxybis-         | 214 | C14H30O   | 000629-64-1 | 50   |
| 3    |    | dl-2-Ethylhexyl chloroformate | 192 | C9H17ClO2 | 024468-13-1 | 42   |
| 4    |    | Octane, 2,3-dimethyl-         | 142 | C10H22    | 007146-60-3 | 40   |
| 5    |    | 1-Pentanol, 2-ethyl-4-methyl- | 130 | C8H18O    | 000106-67-2 | 39   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

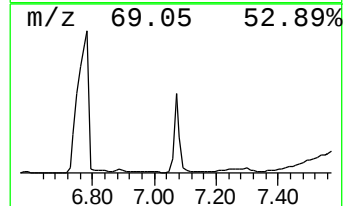
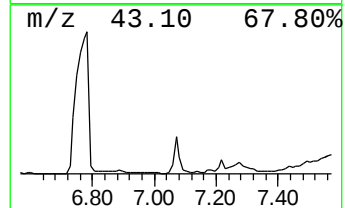
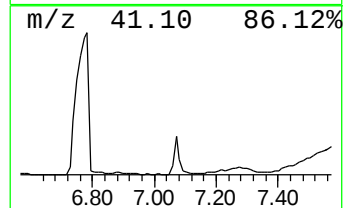
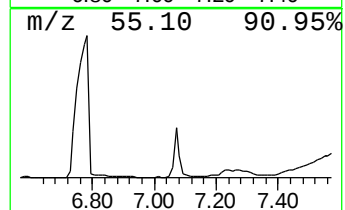
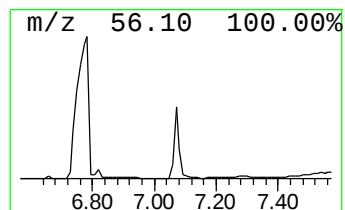
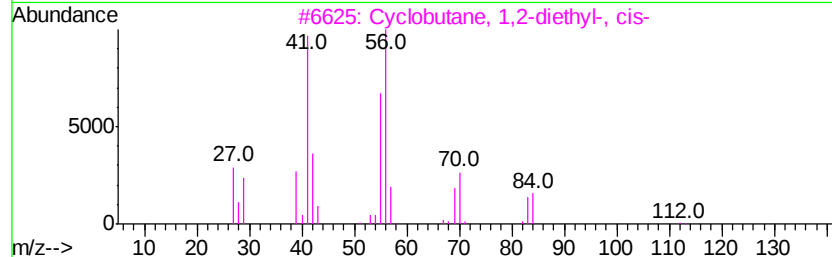
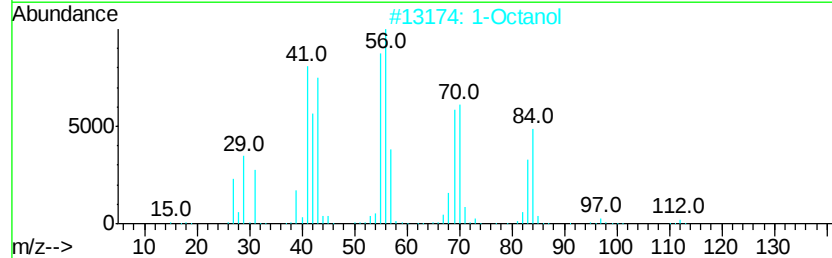
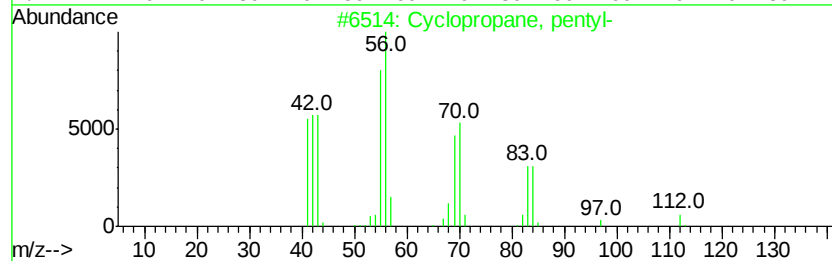
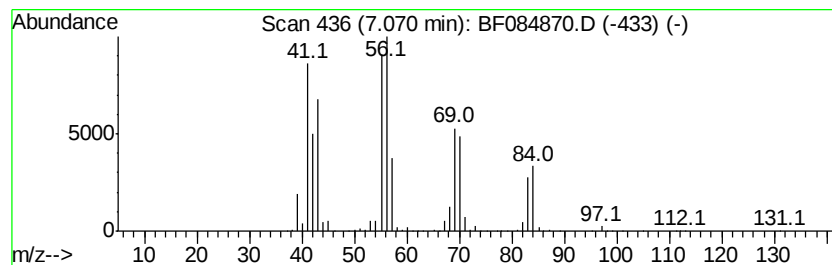
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Cyclopropane, pentyl- Concentration Rank 12

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.07 | 52.60 ng | 2937620 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopropane, pentyl-           | 112 | C8H16   | 002511-91-3 | 86   |
| 2    |    |   | 1-Octanol                       | 130 | C8H18O  | 000111-87-5 | 83   |
| 3    |    |   | Cyclobutane, 1,2-diethyl-, cis- | 112 | C8H16   | 061141-50-2 | 72   |
| 4    |    |   | 1-Heptene                       | 98  | C7H14   | 000592-76-7 | 64   |
| 5    |    |   | Cyclobutane, 1,2-diethyl-       | 112 | C8H16   | 061141-83-1 | 59   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

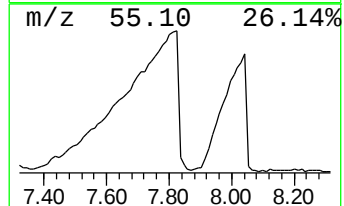
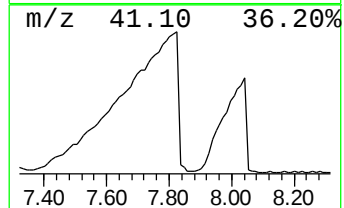
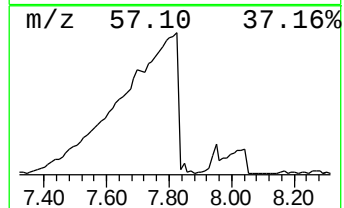
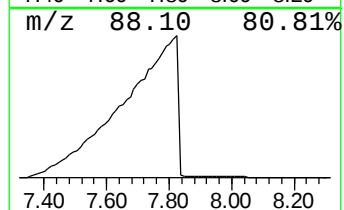
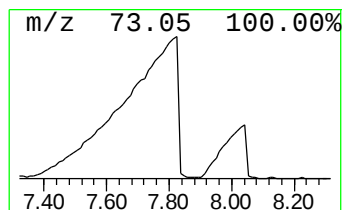
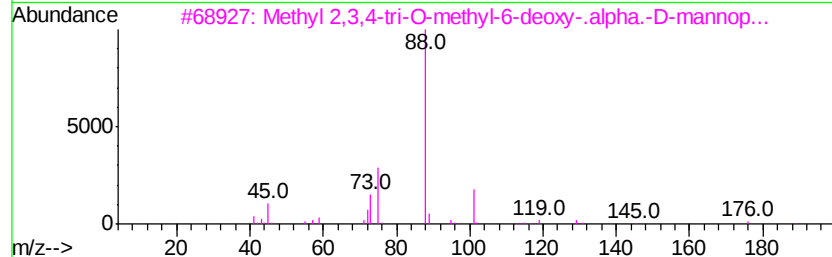
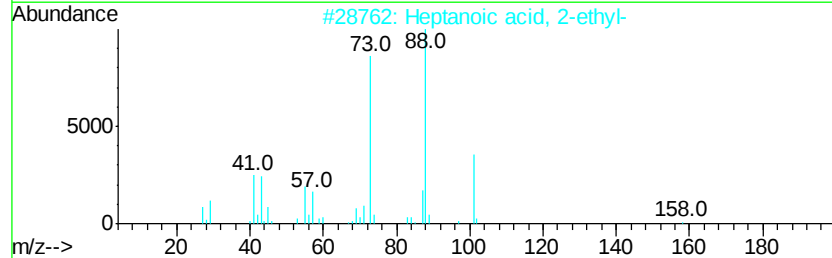
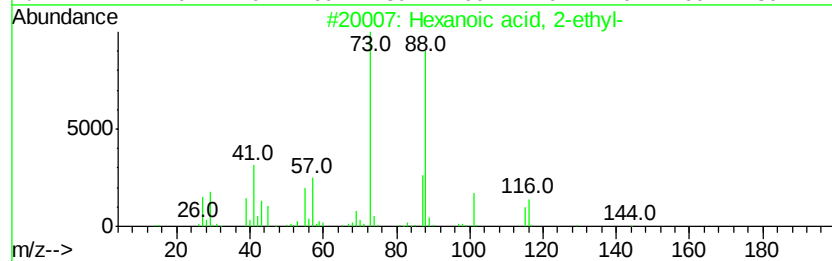
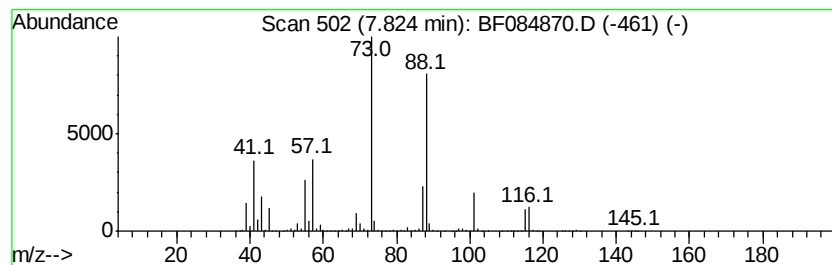
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Hexanoic acid, 2-ethyl- Concentration Rank 2

| R.T. | EstConc   | Area     | Relative to ISTD | R.T. |
|------|-----------|----------|------------------|------|
| 7.82 | 274.65 ng | 85718000 | Naphthalene-d8   | 7.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2  | 000149-57-5 | 90   |
| 2    |    | Heptanoic acid, 2-ethyl-            | 158 | C9H18O2  | 003274-29-1 | 53   |
| 3    |    | Methyl 2,3,4-tri-O-methyl-6-deox... | 220 | C10H20O5 | 072983-16-5 | 50   |
| 4    |    | Butanoic acid, 2-ethyl-, 1,2,3-p... | 386 | C21H38O6 | 056554-54-2 | 45   |
| 5    |    | Methyl L-(+)-.beta.-hydroxyisobu... | 118 | C5H10O3  | 080657-57-4 | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

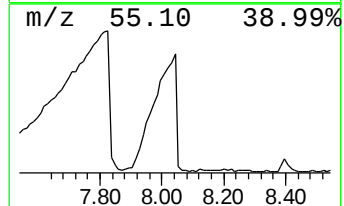
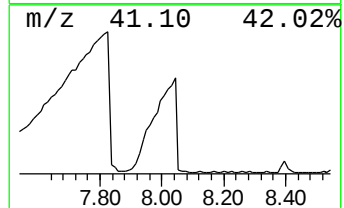
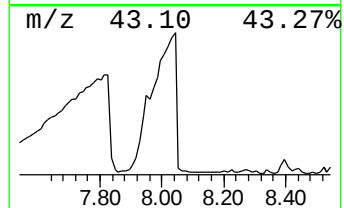
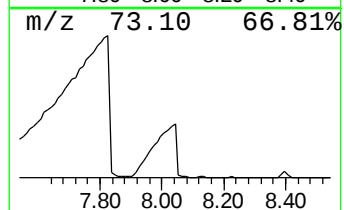
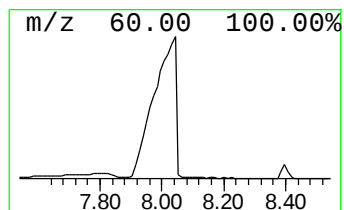
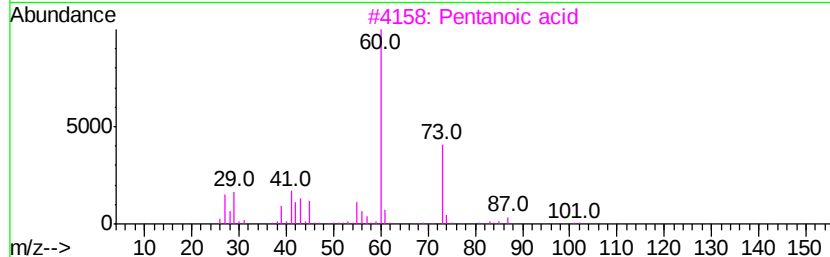
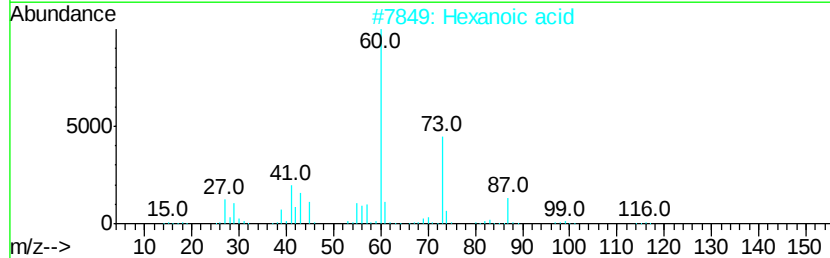
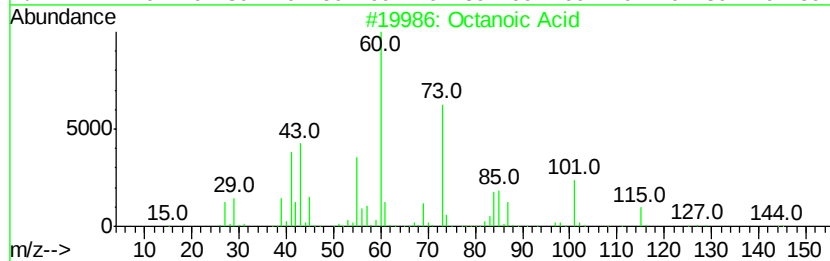
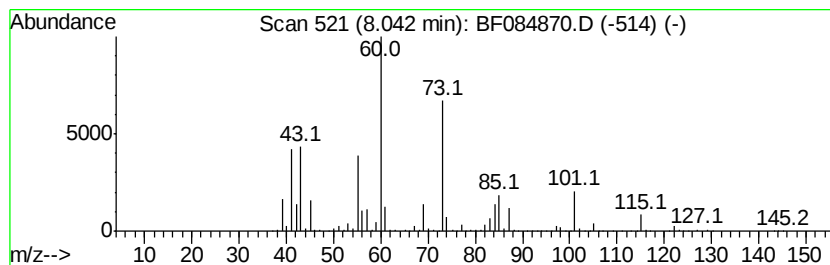
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Octanoic Acid Concentration Rank 9

| R.T. | EstConc  | Area     | Relative to ISTD | R.T. |
|------|----------|----------|------------------|------|
| 8.04 | 61.68 ng | 19251600 | Naphthalene-d8   | 7.95 |

| Hit# of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|-----------|----------------------------|-----|---------|-------------|------|
| 1         | Octanoic Acid              | 144 | C8H16O2 | 000124-07-2 | 97   |
| 2         | Hexanoic acid              | 116 | C6H12O2 | 000142-62-1 | 64   |
| 3         | Pentanoic acid             | 102 | C5H10O2 | 000109-52-4 | 52   |
| 4         | Butanoic acid, 3-methyl-   | 102 | C5H10O2 | 000503-74-2 | 50   |
| 5         | Propanedioic acid, propyl- | 146 | C6H10O4 | 000616-62-6 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

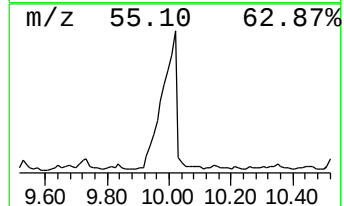
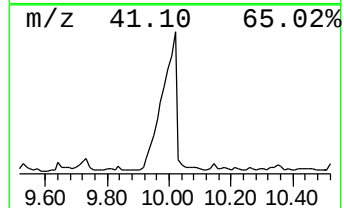
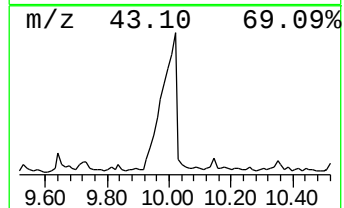
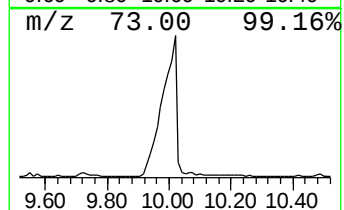
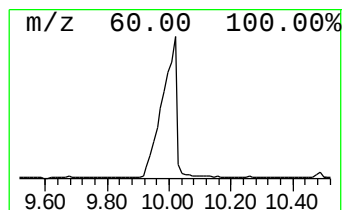
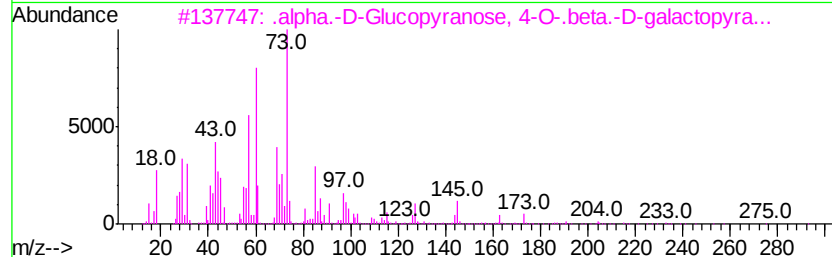
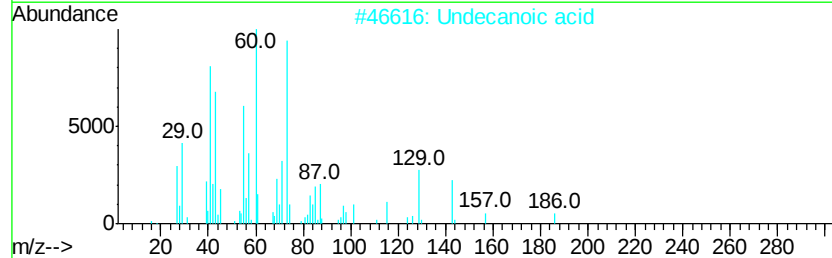
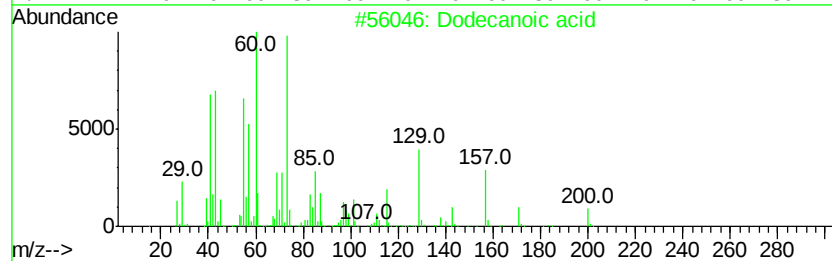
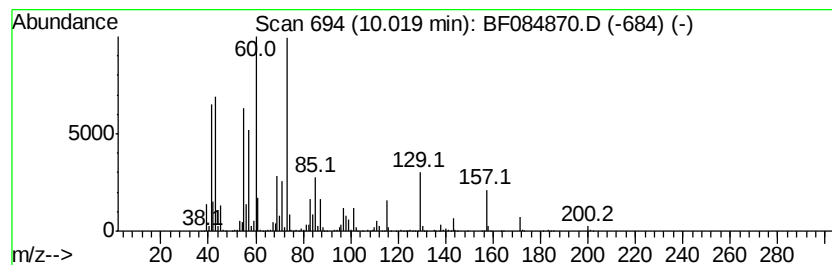
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Dodecanoic acid Concentration Rank 3

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T. |
|-------|-----------|----------|------------------|------|
| 10.02 | 137.50 ng | 15513500 | Acenaphthene-d10 | 9.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Dodecanoic acid                     | 200 | C12H24O2  | 000143-07-7 | 98   |
| 2    |    | Undecanoic acid                     | 186 | C11H22O2  | 000112-37-8 | 80   |
| 3    |    | .alpha.-D-Glucopyranose, 4-O-.be... | 342 | C12H22O11 | 014641-93-1 | 47   |
| 4    |    | Ethanone, 1-(4,5-dihydro-2-thiaz... | 129 | C5H7NOS   | 029926-41-8 | 43   |
| 5    |    | Hexanoic acid                       | 116 | C6H12O2   | 000142-62-1 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

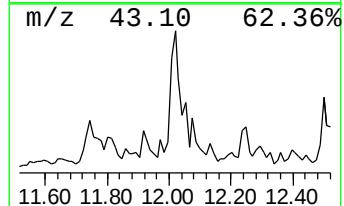
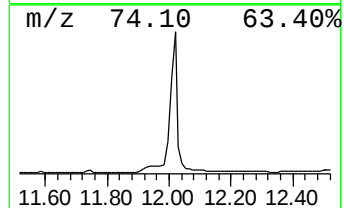
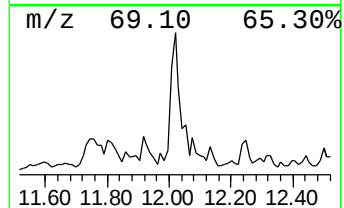
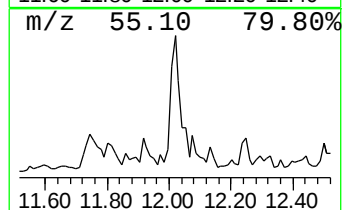
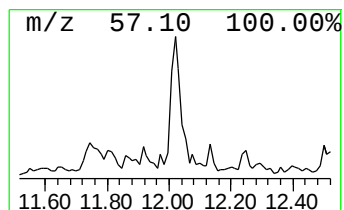
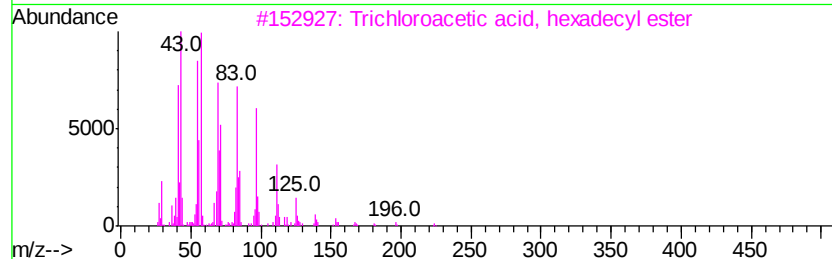
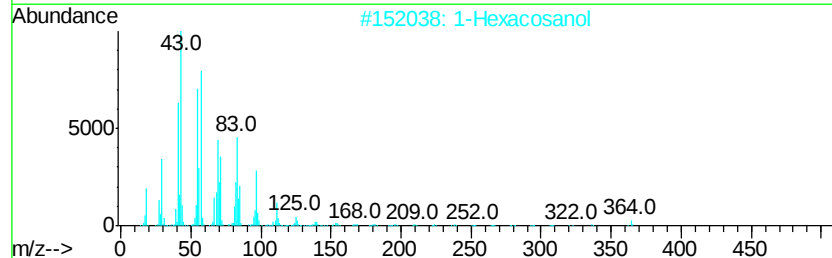
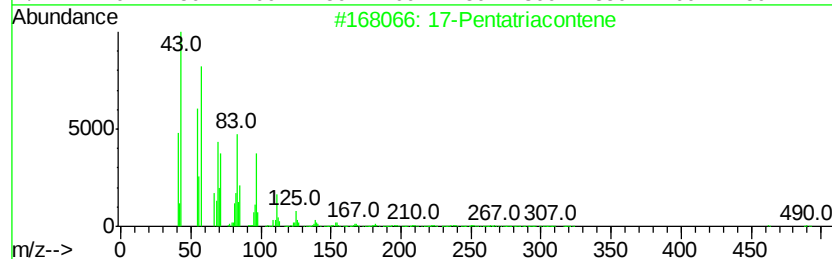
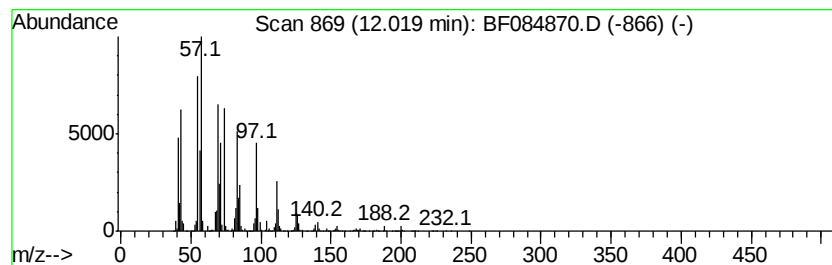
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 17-Pentatriacontene Concentration Rank 17

| R.T.  | EstConc  | Area     | Relative to ISTD | R.T.  |
|-------|----------|----------|------------------|-------|
| 12.02 | 36.70 ng | 15430700 | Phenanthrene-d10 | 11.17 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0  | 74   |
| 2    |    |   | 1-Hexacosanol                       | 382 | C26H54O     | 000506-52-5  | 72   |
| 3    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 68   |
| 4    |    |   | Pentafluoropropionic acid, octad... | 416 | C21H37F5O2  | 1000280-07-7 | 68   |
| 5    |    |   | Acetic acid, 3,7,11,15-tetrameth... | 340 | C22H44O2    | 1000193-63-0 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

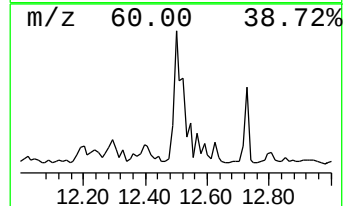
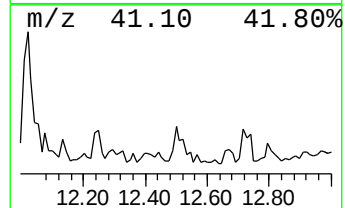
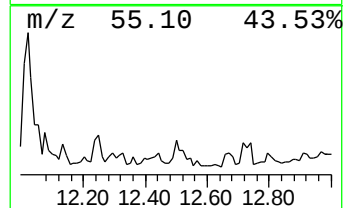
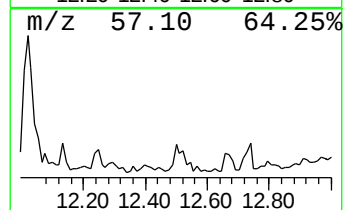
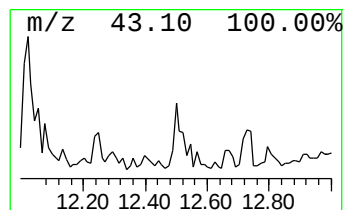
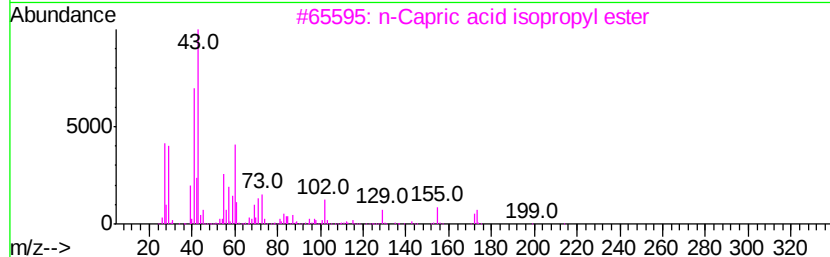
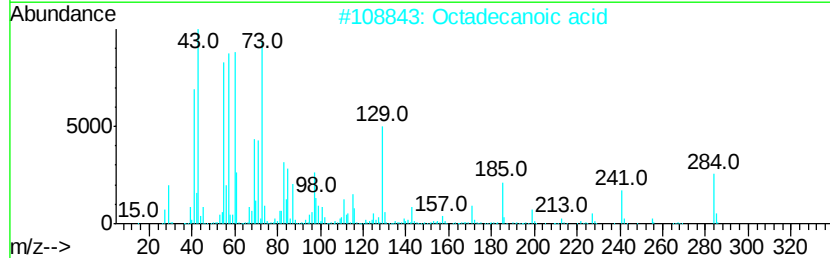
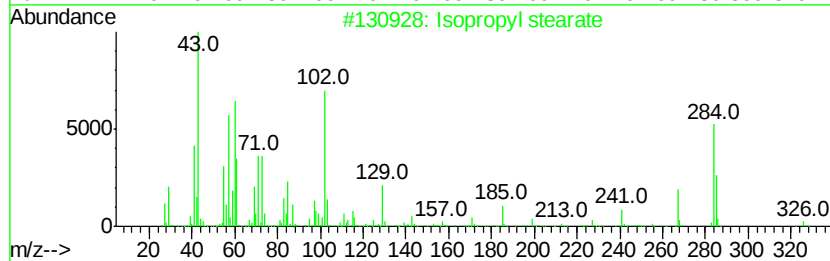
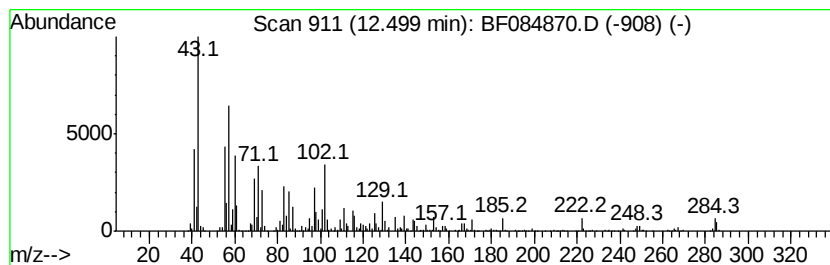
Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.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 12 Isopropyl stearate Concentration Rank 15

| R.T.      | EstConc                         | Area    | Relative to ISTD | R.T.        |      |
|-----------|---------------------------------|---------|------------------|-------------|------|
| 12.50     | 43.46 ng                        | 4658060 | Chrysene-d12     | 13.80       |      |
| Hit# of 5 | Tentative ID                    | MW      | MolForm          | CAS#        | Qual |
| 1         | Isopropyl stearate              | 326     | C21H42O2         | 000112-10-7 | 50   |
| 2         | Octadecanoic acid               | 284     | C18H36O2         | 000057-11-4 | 41   |
| 3         | n-Capric acid isopropyl ester   | 214     | C13H26O2         | 002311-59-3 | 38   |
| 4         | n-Octanoic acid isopropyl ester | 186     | C11H22O2         | 005458-59-3 | 27   |
| 5         | Heptanoic acid, propyl ester    | 172     | C10H20O2         | 007778-87-2 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

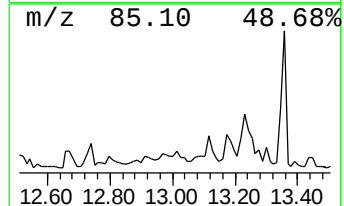
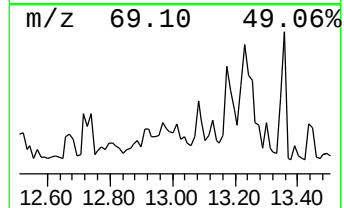
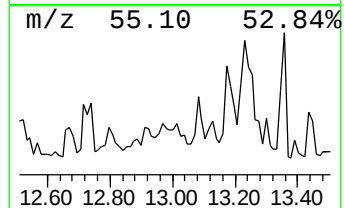
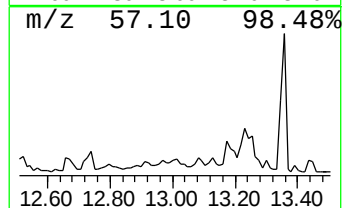
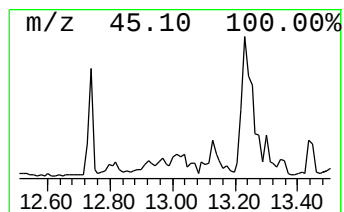
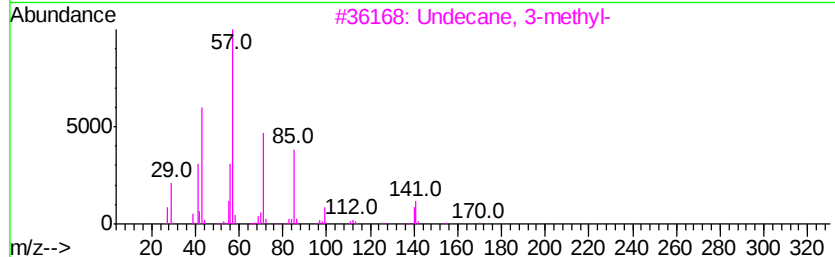
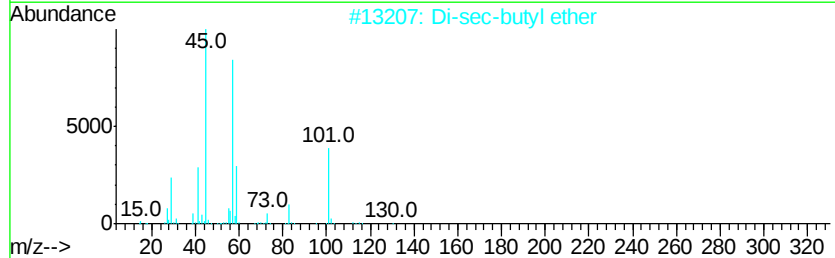
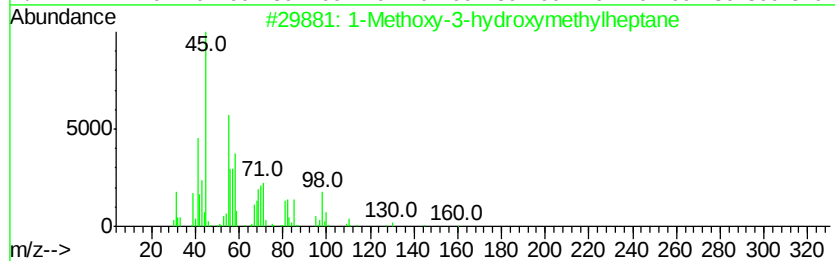
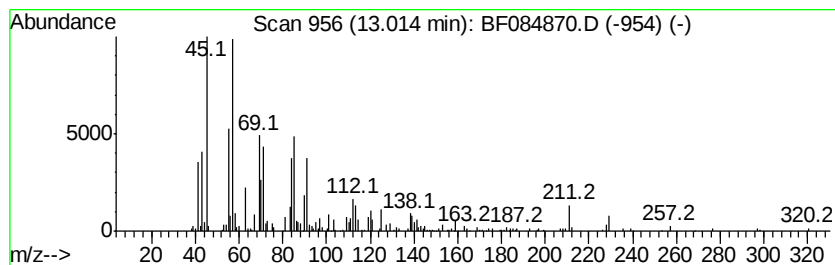
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown13.01 Concentration Rank 19

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.01 | 30.92 ng | 3314010 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#         | Qual |
|------|----|----------------------------------|-----|---------|--------------|------|
| 1    | 5  | 1-Methoxy-3-hydroxymethylheptane | 160 | C9H20O2 | 071052-95-4  | 35   |
| 2    |    | Di-sec-butyl ether               | 130 | C8H18O  | 006863-58-7  | 22   |
| 3    |    | Undecane, 3-methyl-              | 170 | C12H26  | 001002-43-3  | 22   |
| 4    |    | 3,3-Dimethylbutane-2-ol          | 102 | C6H14O  | 000464-07-3  | 22   |
| 5    |    | 5-Methoxymethoxyhexa-2,3-diene   | 142 | C8H14O2 | 1000190-45-1 | 22   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

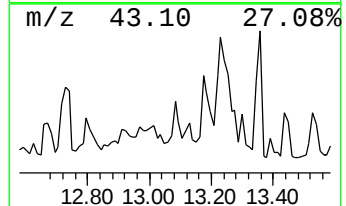
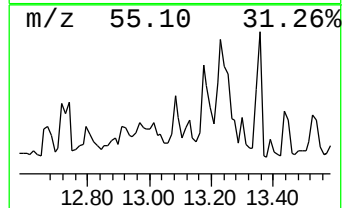
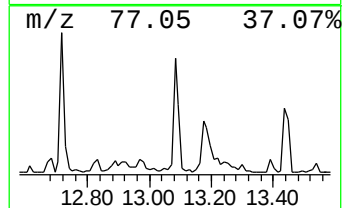
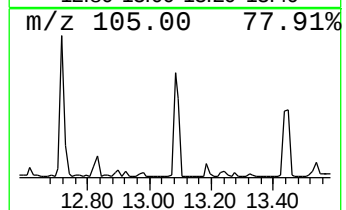
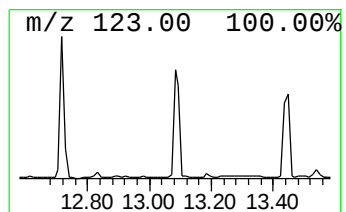
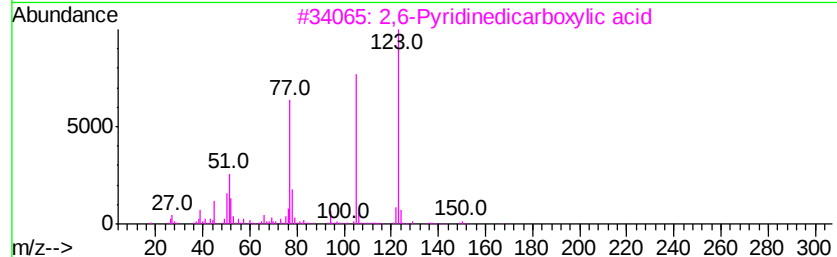
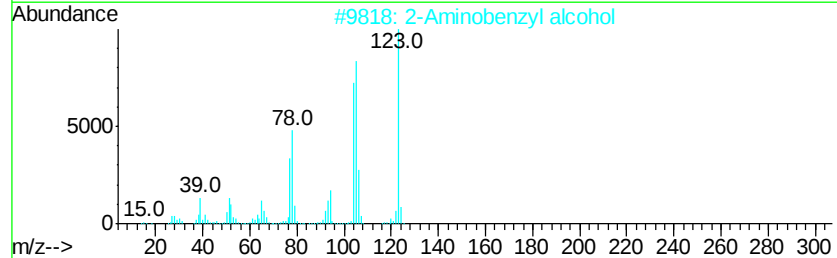
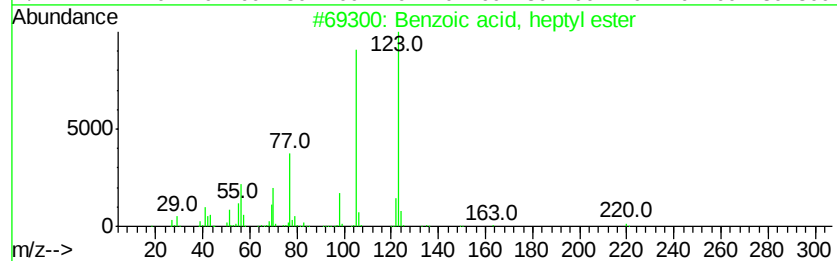
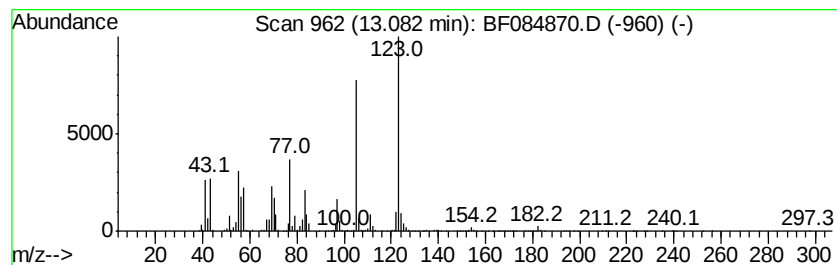
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benzoic acid, heptyl ester Concentration Rank 18

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.08 | 36.47 ng | 3909610 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Benzoic acid, heptyl ester          | 220 | C14H20O2  | 007155-12-6  | 72   |
| 2    |    | 2-Aminobenzyl alcohol               | 123 | C7H9NO    | 005344-90-1  | 64   |
| 3    |    | 2,6-Pyridinedicarboxylic acid       | 167 | C7H5N04   | 000499-83-2  | 59   |
| 4    |    | 2-Fluorobenzoic acid, 4-pentadec... | 350 | C22H35FO2 | 1000280-61-7 | 53   |
| 5    |    | 2-Fluorobenzoic acid, 4-tetradec... | 336 | C21H33FO2 | 1000280-61-4 | 50   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

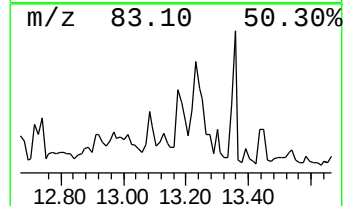
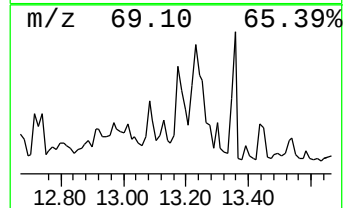
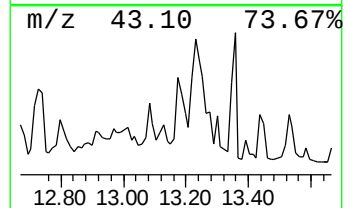
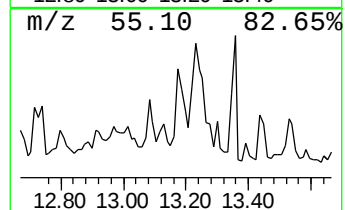
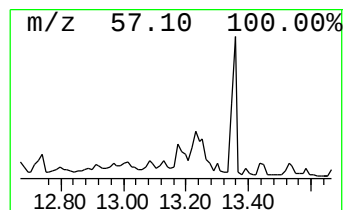
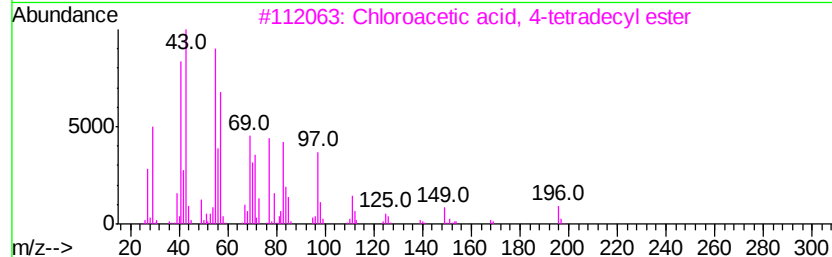
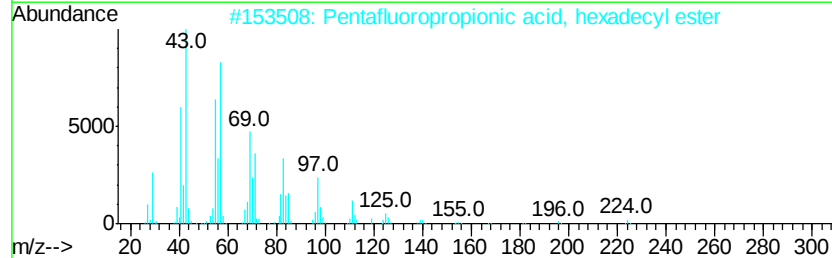
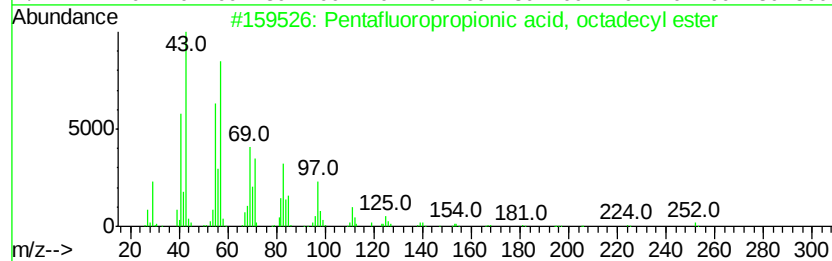
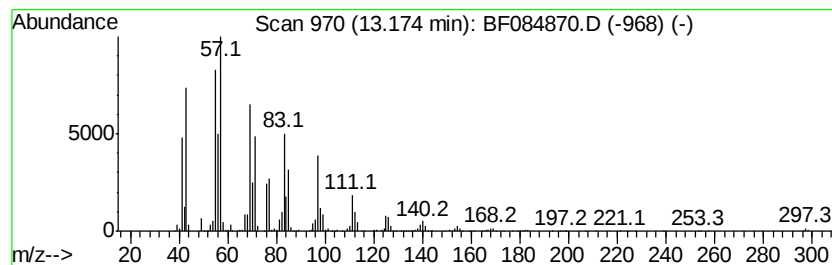
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Pentafluoropropionic acid, ... Concentration Rank 13

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.17 | 48.71 ng | 5221210 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                               | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Pentafluoropropionic acid, octadecyl ester | 416 | C21H37F5O2 | 1000280-07-7 | 83   |
| 2    |    | Pentafluoropropionic acid, hexadecyl ester | 388 | C19H33F5O2 | 006222-07-7  | 83   |
| 3    |    | Chloroacetic acid, 4-tetradecyl ester      | 290 | C16H31ClO2 | 1000280-08-6 | 80   |
| 4    |    | 1-Hexadecanol, 2-methyl-                   | 256 | C17H36O    | 002490-48-4  | 74   |
| 5    |    | 1-Hexacosanol                              | 382 | C26H54O    | 000506-52-5  | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

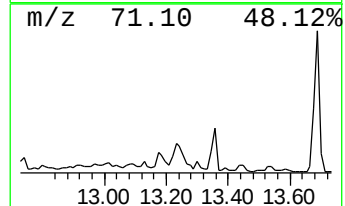
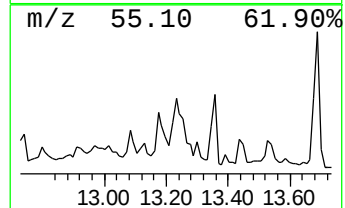
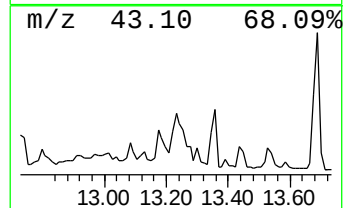
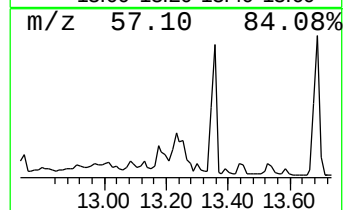
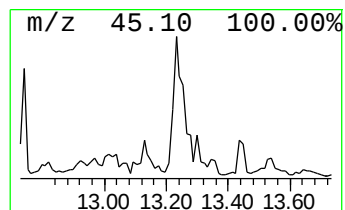
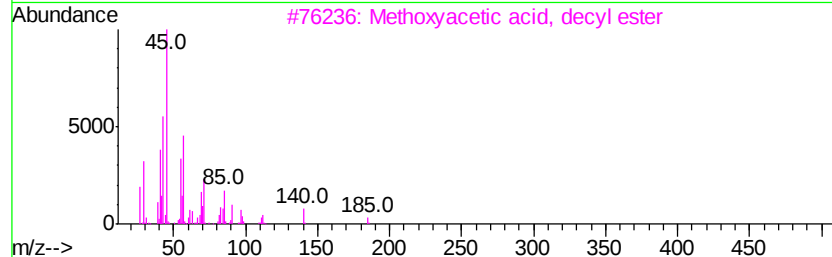
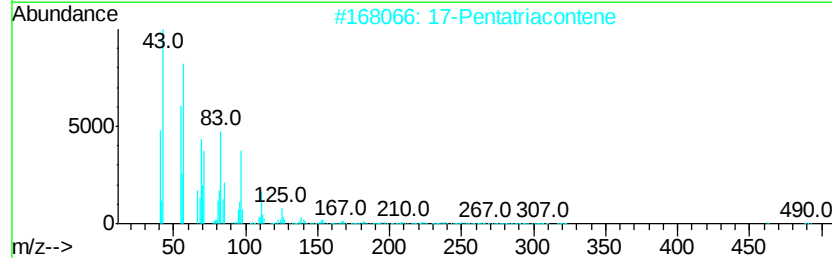
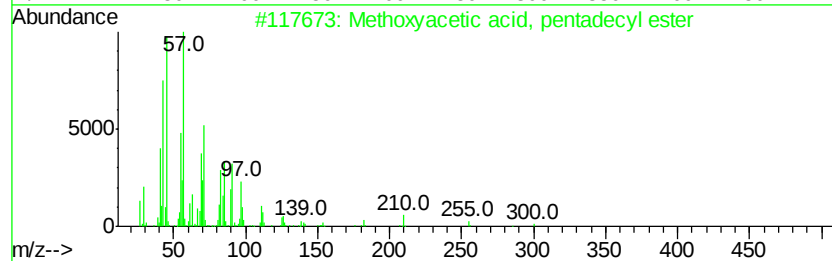
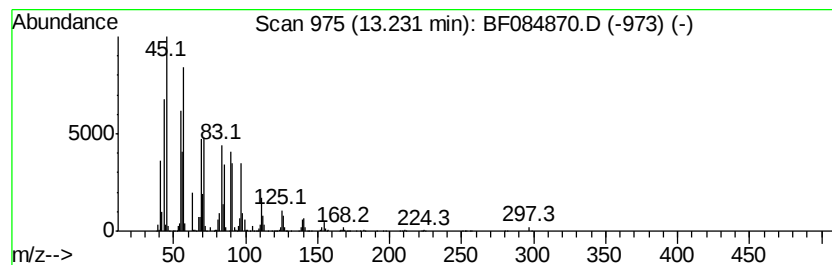
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Methoxyacetic acid, pentade... Concentration Rank 6

| R.T.  | EstConc  | Area     | Relative to ISTD | R.T.  |
|-------|----------|----------|------------------|-------|
| 13.23 | 93.74 ng | 10048000 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Methoxvacetic acid, pentadecyl e... | 300 | C18H36O3   | 1000281-82-2 | 53   |
| 2    |    | 17-Pentatriacontene                 | 491 | C35H70     | 006971-40-0  | 49   |
| 3    |    | Methoxvacetic acid, decyl ester     | 230 | C13H26O3   | 259141-02-1  | 46   |
| 4    |    | Pentafluoropropionic acid, hexad... | 388 | C19H33F5O2 | 006222-07-7  | 46   |
| 5    |    | 2-Pentadecanol                      | 228 | C15H32O    | 001653-34-5  | 43   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

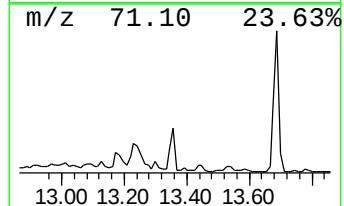
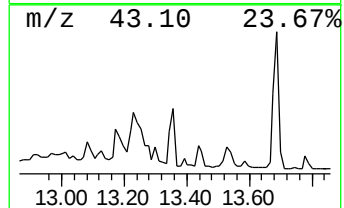
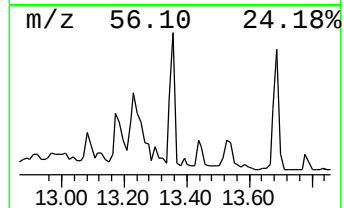
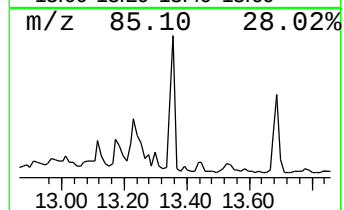
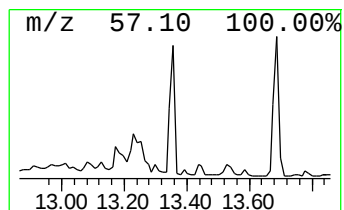
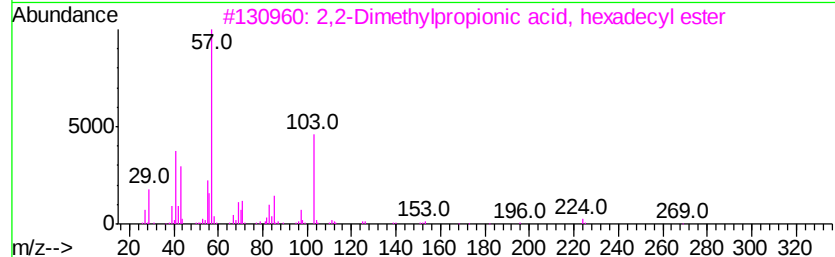
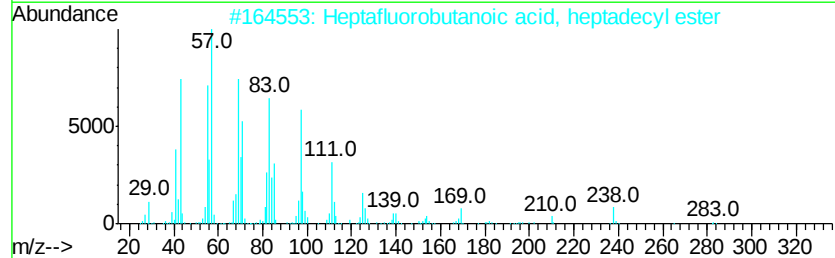
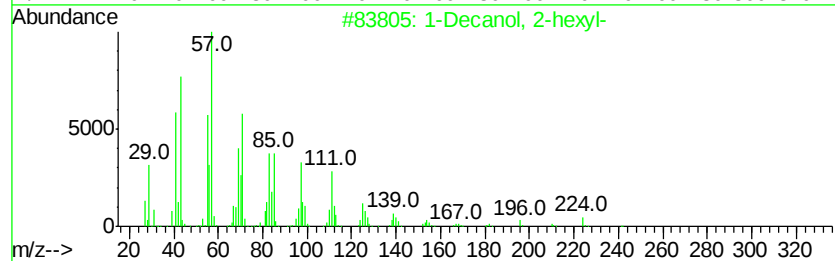
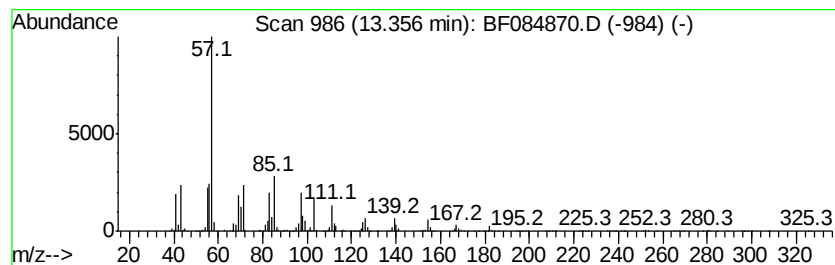
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 1-Decanol, 2-hexyl- Concentration Rank 11

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.36 | 55.41 ng | 5939880 | Chrysene-d12     | 13.80 |

| Hit# | of | 5 | Tentative ID                                 | MW  | MolForm    | CAS#         | Qual |
|------|----|---|----------------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Decanol, 2-hexyl-                          | 242 | C16H34O    | 002425-77-6  | 60   |
| 2    |    |   | Heptafluorobutanoic acid, heptadecyl ester   | 452 | C21H35F7O2 | 1000282-97-3 | 50   |
| 3    |    |   | 2,2-Dimethylpropionic acid, hexadecyl ester  | 326 | C21H42O2   | 032767-89-8  | 47   |
| 4    |    |   | Ethanol, 2-(tetradecyloxy)-                  | 258 | C16H34O2   | 002136-70-1  | 47   |
| 5    |    |   | 2,2-Dimethylpropionic acid, tetradecyl ester | 298 | C19H38O2   | 144610-93-5  | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

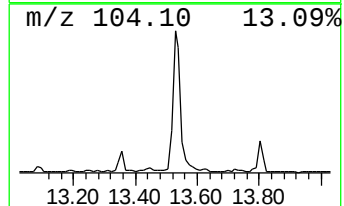
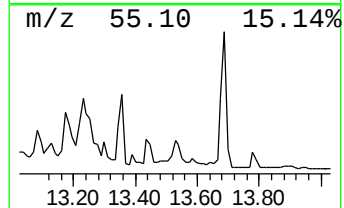
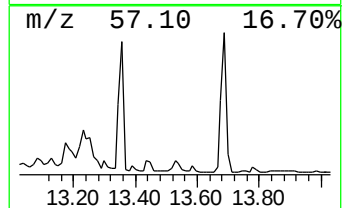
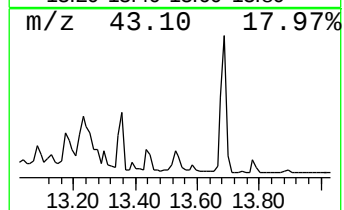
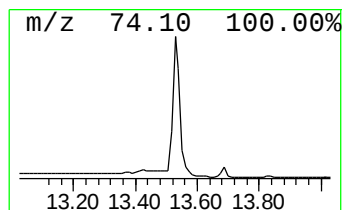
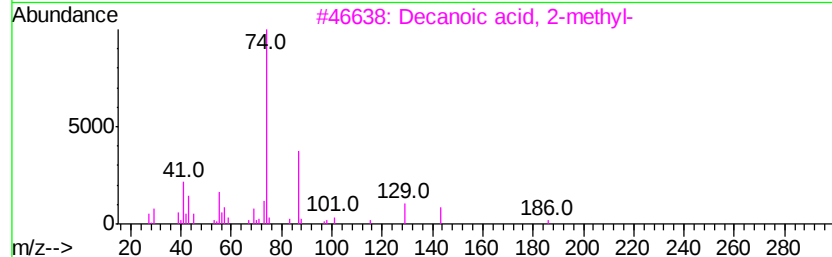
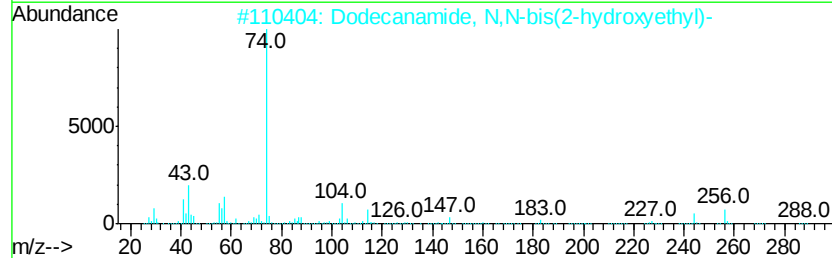
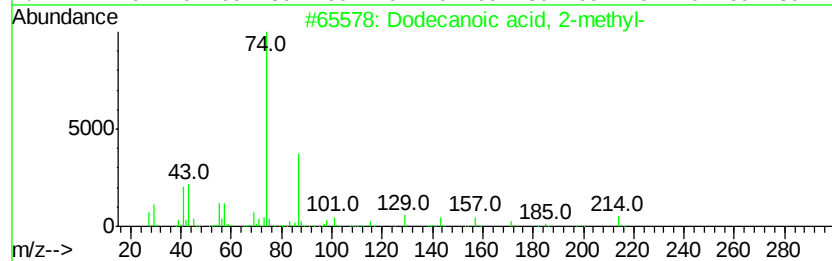
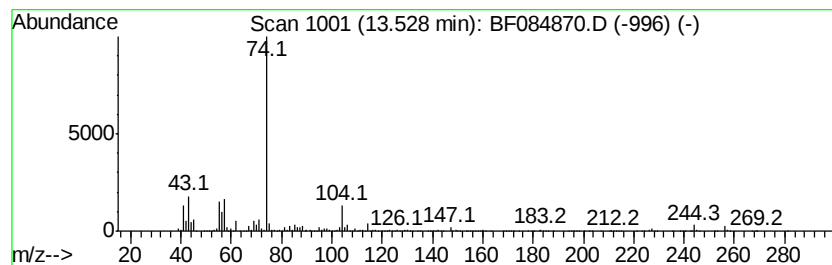
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown13.53 Concentration Rank 16

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.53 | 42.08 ng | 4510130 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Dodecanoic acid, 2-methyl-           | 214 | C13H26O2  | 002874-74-0 | 42   |
| 2    |    | Dodecanamide, N,N-bis(2-hydroxyve... | 287 | C16H33NO3 | 000120-40-1 | 38   |
| 3    |    | Decanoic acid, 2-methyl-             | 186 | C11H22O2  | 024323-23-7 | 38   |
| 4    |    | Pentanoic acid, methyl ester         | 116 | C6H12O2   | 000624-24-8 | 36   |
| 5    |    | Hexanoic acid, 2-methyl-             | 130 | C7H14O2   | 004536-23-6 | 36   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

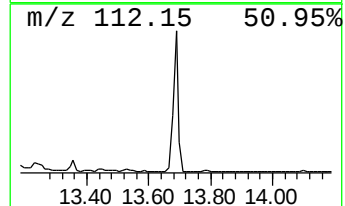
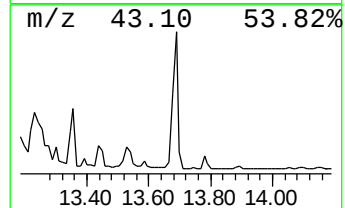
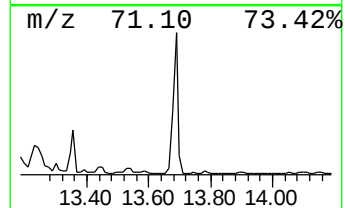
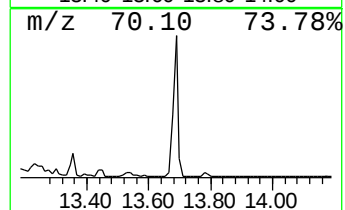
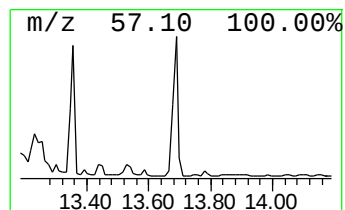
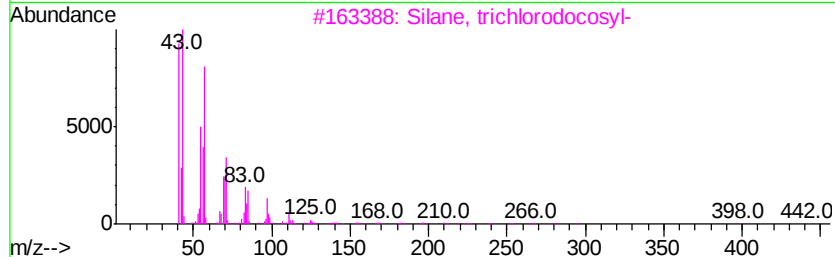
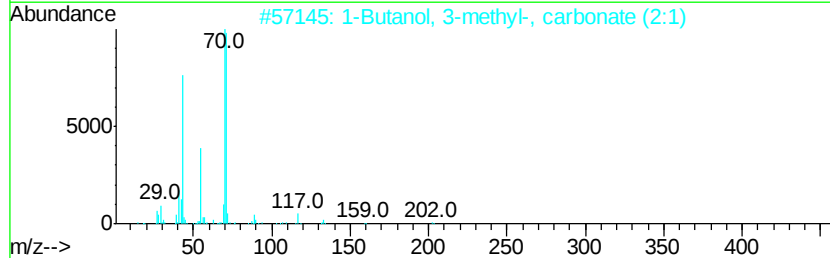
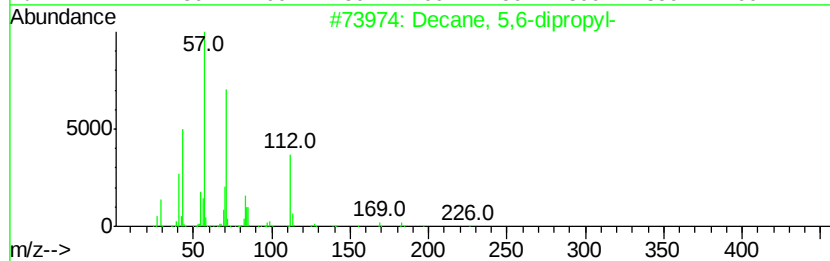
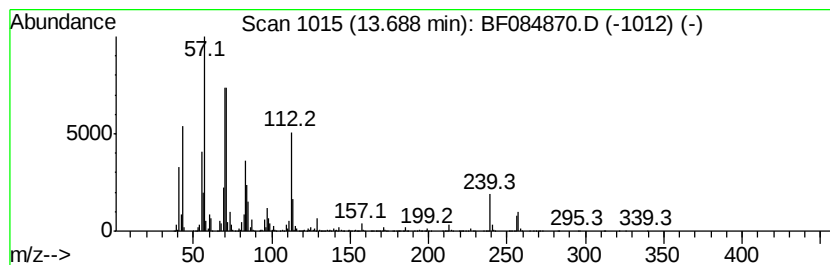
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown13.69 Concentration Rank 5

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T.  |
|-------|-----------|----------|------------------|-------|
| 13.69 | 103.95 ng | 11142800 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Decane, 5,6-dipropyl-               | 226 | C16H34      | 119209-20-0  | 47   |
| 2    |    | 1-Butanol, 3-methyl-, carbonate ... | 202 | C11H22O3    | 002050-95-5  | 43   |
| 3    |    | Silane, trichlorodocosyl-           | 442 | C22H45Cl3Si | 007325-84-0  | 38   |
| 4    |    | Cyclohexane, (1,1-dimethylpropyl)-  | 154 | C11H22      | 031797-64-5  | 35   |
| 5    |    | Heptane, 3-ethyl-5-methylene-       | 140 | C10H20      | 1000160-41-7 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
Data File : BF084870.D  
Acq On : 5 Feb 2016 20:06  
Operator : SJ/IZ  
Sample : H1357-02  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

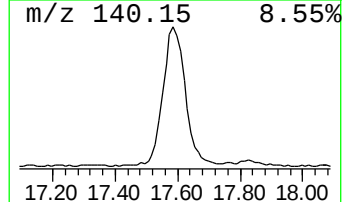
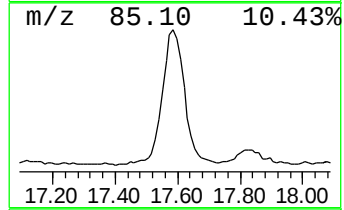
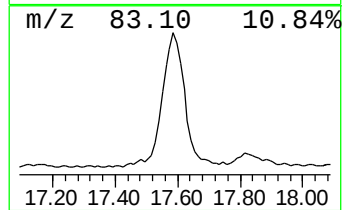
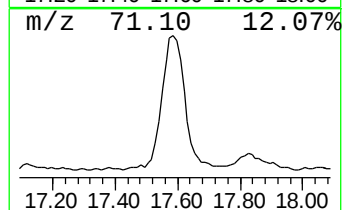
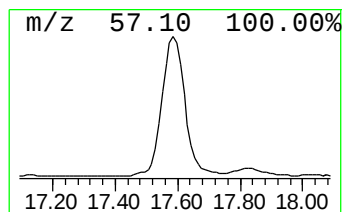
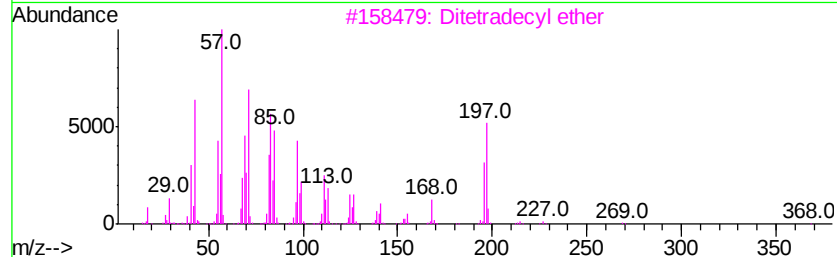
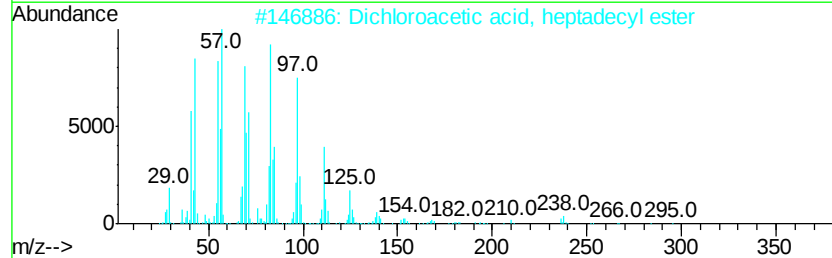
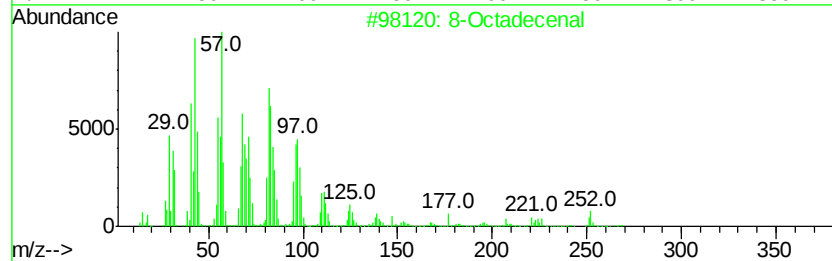
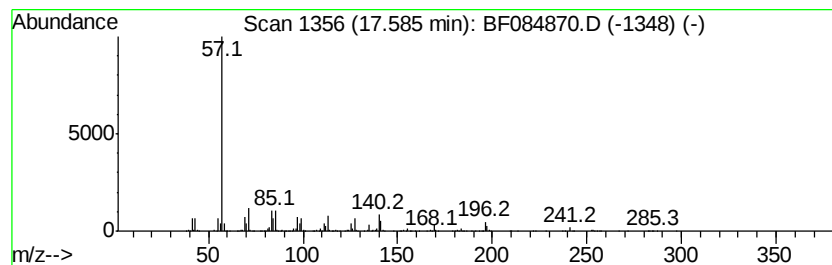
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 8-Octadecenal Concentration Rank 4

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 17.59 | 118.41 ng | 6273830 | Perylene-d12     | 15.17 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 8-Octadecenal                       | 266 | C18H34O     | 056554-94-0  | 56   |
| 2    |    | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 55   |
| 3    |    | Ditetradecyl ether                  | 410 | C28H58O     | 005412-98-6  | 50   |
| 4    |    | Acetic acid, chloro-, octadecyl ... | 346 | C20H39ClO2  | 005348-82-3  | 49   |
| 5    |    | Octane, 2,5,6-trimethyl-            | 156 | C11H24      | 062016-14-2  | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020516\  
 Data File : BF084870.D  
 Acq On : 5 Feb 2016 20:06  
 Operator : SJ/IZ  
 Sample : H1357-02  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EFF-WW

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.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 |
| 2-Pentanone, 4-hy... | 4.81  | 29.1 ng  |       | 1624260  | 1                     | 6.66  | 1116870 | 20.0 |
| (S)-Isopropyl lac... | 4.97  | 46.1 ng  |       | 2573500  | 1                     | 6.66  | 1116870 | 20.0 |
| unknown6.43          | 6.43  | 69.6 ng  |       | 3886510  | 1                     | 6.66  | 1116870 | 20.0 |
| Hexanoic acid        | 6.53  | 56.3 ng  |       | 3143900  | 1                     | 6.66  | 1116870 | 20.0 |
| 1-Hexanol, 2-ethyl-  | 6.78  | 441.6 ng |       | 24662800 | 1                     | 6.66  | 1116870 | 20.0 |
| Cyclopropane, pen... | 7.07  | 52.6 ng  |       | 2937620  | 1                     | 6.66  | 1116870 | 20.0 |
| Hexanoic acid, 2-... | 7.82  | 274.6 ng |       | 85718000 | 2                     | 7.95  | 6242100 | 20.0 |
| Octanoic Acid        | 8.04  | 61.7 ng  |       | 19251600 | 2                     | 7.95  | 6242100 | 20.0 |
| Dodecanoic acid      | 10.02 | 137.5 ng |       | 15513500 | 3                     | 9.70  | 2256560 | 20.0 |
| 17-Pentatriacontene  | 12.02 | 36.7 ng  |       | 15430700 | 4                     | 11.17 | 8408180 | 20.0 |
| Isopropyl stearate   | 12.50 | 43.5 ng  |       | 4658060  | 5                     | 13.80 | 2143830 | 20.0 |
| unknown13.01         | 13.01 | 30.9 ng  |       | 3314010  | 5                     | 13.80 | 2143830 | 20.0 |
| Benzoic acid, hep... | 13.08 | 36.5 ng  |       | 3909610  | 5                     | 13.80 | 2143830 | 20.0 |
| Pentafluoropropio... | 13.17 | 48.7 ng  |       | 5221210  | 5                     | 13.80 | 2143830 | 20.0 |
| Methoxyacetic aci... | 13.23 | 93.7 ng  |       | 10048000 | 5                     | 13.80 | 2143830 | 20.0 |
| 1-Decanol, 2-hexyl-  | 13.36 | 55.4 ng  |       | 5939880  | 5                     | 13.80 | 2143830 | 20.0 |
| unknown13.53         | 13.53 | 42.1 ng  |       | 4510130  | 5                     | 13.80 | 2143830 | 20.0 |
| unknown13.69         | 13.69 | 104.0 ng |       | 11142800 | 5                     | 13.80 | 2143830 | 20.0 |
| 8-Octadecenal        | 17.59 | 118.4 ng |       | 6273830  | 6                     | 15.17 | 1059650 | 20.0 |