

Data Path : S:\HPCHEM1\BNA\_F\DATA\BF031016\  
 Data File : BF085365.D  
 Acq On : 11 Mar 2016 6:31  
 Operator : UM/SJ  
 Sample : H1731-12  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-28-COMP

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-BF030316.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         | 5.144       | 248           | 250         | 259          | rBV      | 201756         | 328003        | 6.74%           | 0.820%        |
| 2         | 5.544       | 282           | 285         | 288          | rBV      | 3636569        | 4052099       | 83.21%          | 10.125%       |
| 3         | 6.573       | 371           | 375         | 377          | rBV      | 3098654        | 4795741       | 98.48%          | 11.983%       |
| 4         | 6.710       | 384           | 387         | 389          | rBV      | 4176948        | 4779947       | 98.16%          | 11.944%       |
| 5         | 6.939       | 404           | 407         | 409          | rBV      | 1164895        | 963061        | 19.78%          | 2.406%        |
| 6         | 7.099       | 418           | 421         | 423          | rBV      | 3675692        | 3703148       | 76.05%          | 9.253%        |
| 7         | 7.499       | 453           | 456         | 459          | rBV      | 2330438        | 2905168       | 59.66%          | 7.259%        |
| 8         | 7.579       | 461           | 463         | 469          | rVB      | 57115          | 62965         | 1.29%           | 0.157%        |
| 9         | 8.219       | 517           | 519         | 522          | rBV      | 1123249        | 1112102       | 22.84%          | 2.779%        |
| 10        | 9.305       | 611           | 614         | 616          | rBV      | 4011972        | 4869640       | 100.00%         | 12.168%       |
| 11        | 9.693       | 645           | 648         | 649          | rBV      | 829459         | 777053        | 15.96%          | 1.942%        |
| 12        | 9.911       | 664           | 667         | 669          | rVB      | 115047         | 108199        | 2.22%           | 0.270%        |
| 13        | 9.979       | 671           | 673         | 675          | rVB      | 1358844        | 1196154       | 24.56%          | 2.989%        |
| 14        | 10.402      | 709           | 710         | 712          | rVB      | 119444         | 139599        | 2.87%           | 0.349%        |
| 15        | 10.631      | 727           | 730         | 733          | rBV      | 49202          | 94405         | 1.94%           | 0.236%        |
| 16        | 10.768      | 739           | 742         | 744          | rBV      | 2561267        | 2559144       | 52.55%          | 6.395%        |
| 17        | 10.882      | 750           | 752         | 758          | rBV      | 153106         | 195670        | 4.02%           | 0.489%        |
| 18        | 11.328      | 789           | 791         | 792          | rBV      | 96712          | 81900         | 1.68%           | 0.205%        |
| 19        | 11.465      | 800           | 803         | 805          | rBV      | 933428         | 1007956       | 20.70%          | 2.519%        |
| 20        | 11.991      | 847           | 849         | 858          | rBV      | 126000         | 210497        | 4.32%           | 0.526%        |
| 21        | 12.768      | 915           | 917         | 924          | rBV      | 68211          | 96910         | 1.99%           | 0.242%        |
| 22        | 13.042      | 939           | 941         | 944          | rBV      | 2364565        | 3354382       | 68.88%          | 8.382%        |
| 23        | 13.934      | 1017          | 1019        | 1027         | rBV      | 459348         | 640310        | 13.15%          | 1.600%        |
| 24        | 14.094      | 1031          | 1033        | 1036         | rVB      | 800692         | 757247        | 15.55%          | 1.892%        |
| 25        | 14.574      | 1072          | 1075        | 1081         | rBV2     | 41036          | 86515         | 1.78%           | 0.216%        |
| 26        | 14.951      | 1106          | 1108        | 1111         | rVB      | 148578         | 155953        | 3.20%           | 0.390%        |
| 27        | 15.557      | 1158          | 1161        | 1164         | rBV      | 585161         | 750025        | 15.40%          | 1.874%        |
| 28        | 16.083      | 1204          | 1207        | 1214         | rBV2     | 40314          | 123105        | 2.53%           | 0.308%        |
| 29        | 16.265      | 1220          | 1223        | 1230         | rVB5     | 16405          | 55199         | 1.13%           | 0.138%        |
| 30        | 17.614      | 1338          | 1341        | 1347         | rVB6     | 21584          | 58631         | 1.20%           | 0.147%        |

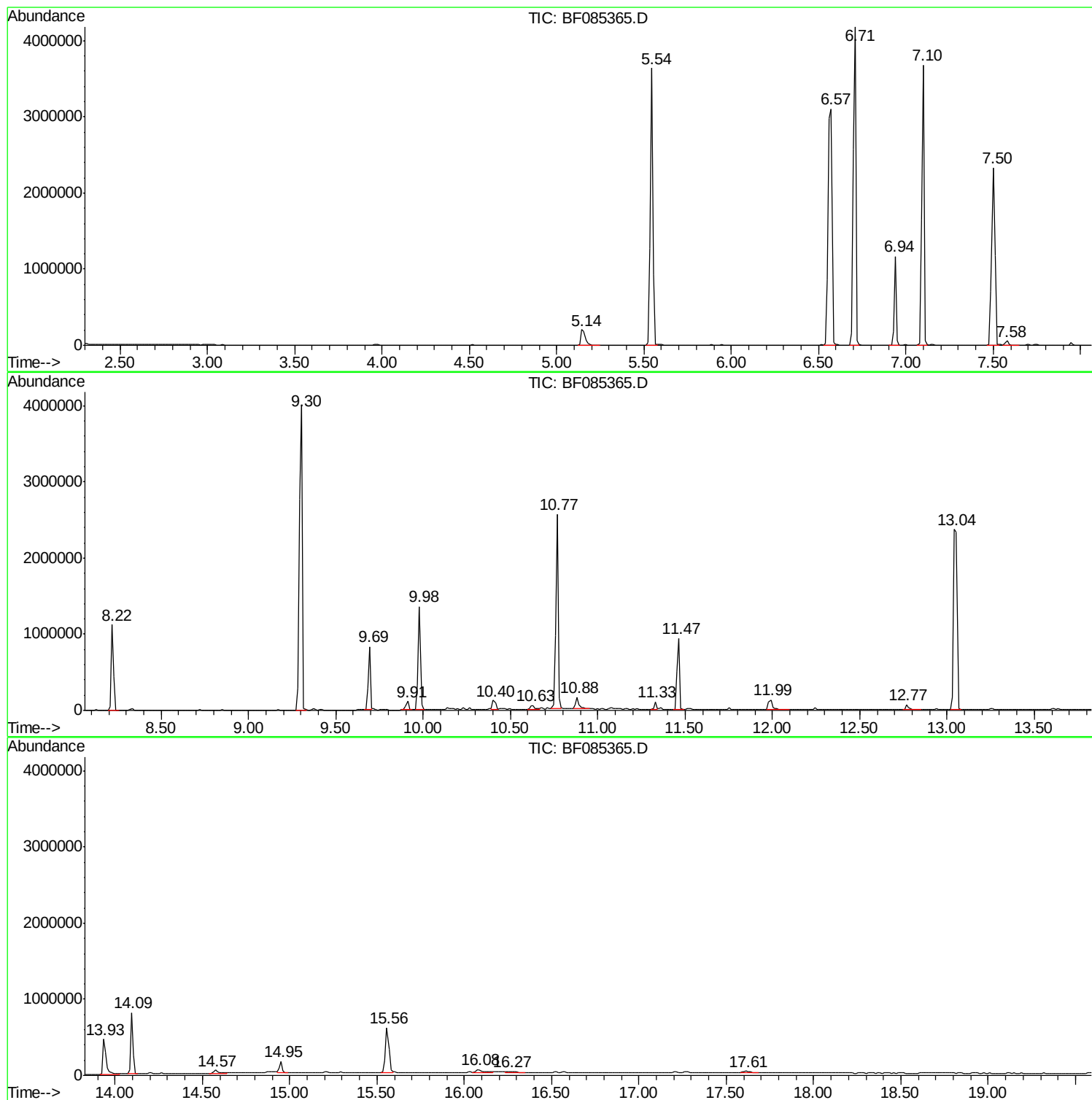
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SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.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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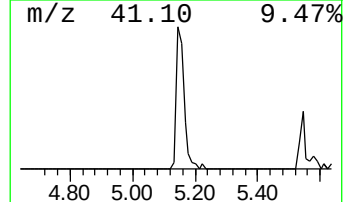
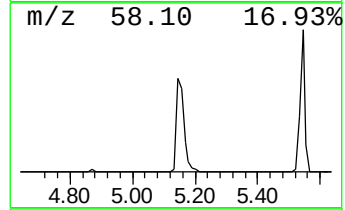
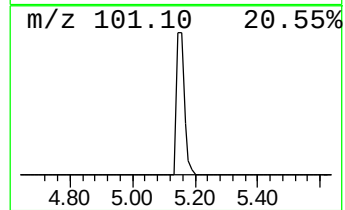
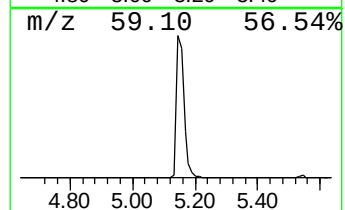
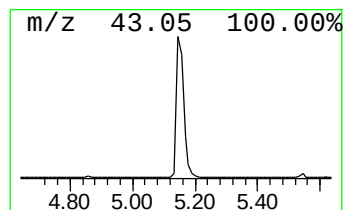
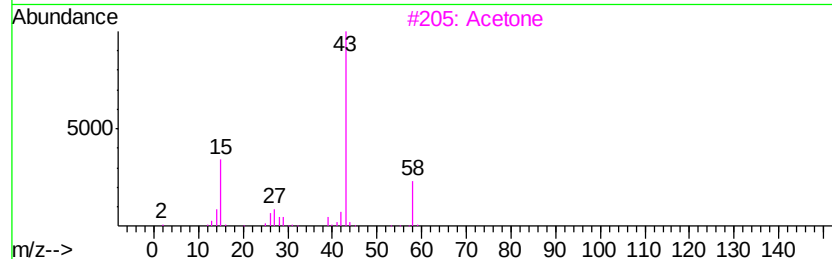
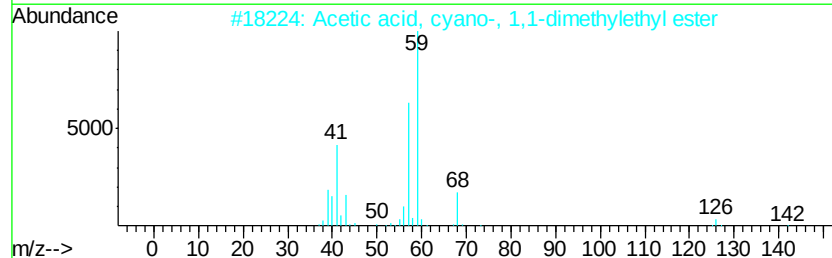
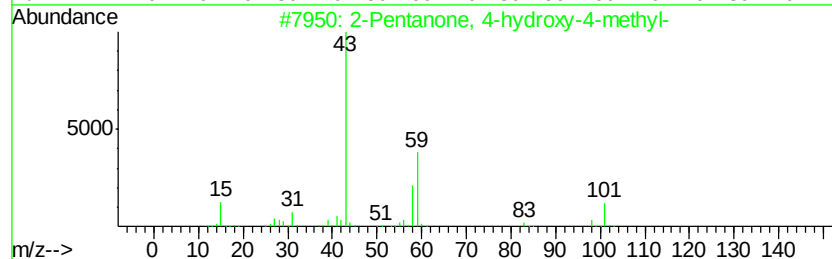
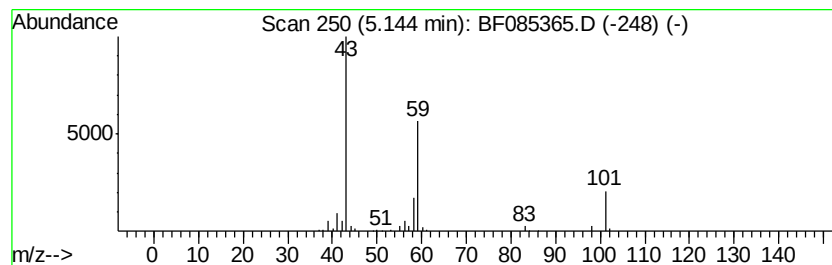
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.14 | 6.81 ng | 328003 | 1,4-Dichlorobenzene-d4 | 6.94 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 53   |
| 2    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 3    |    |   | Acetone                             | 58  | C3H6O    | 000067-64-1 | 10   |
| 4    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    |   | 5-Hexen-2-one                       | 98  | C6H10O   | 000109-49-9 | 9    |



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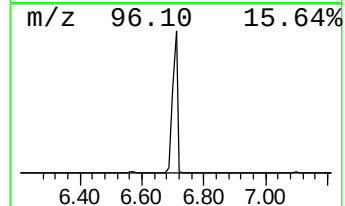
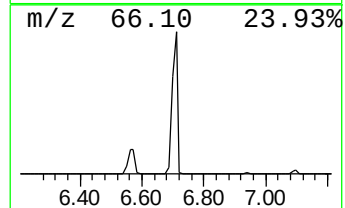
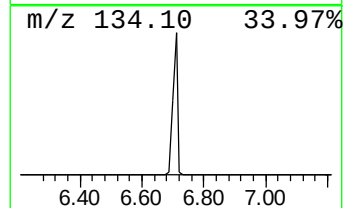
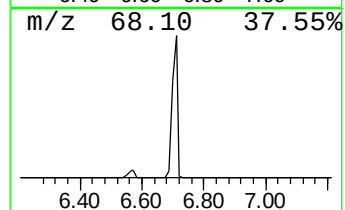
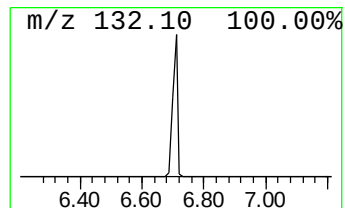
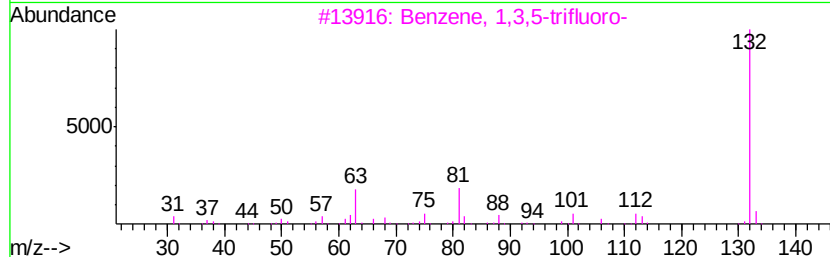
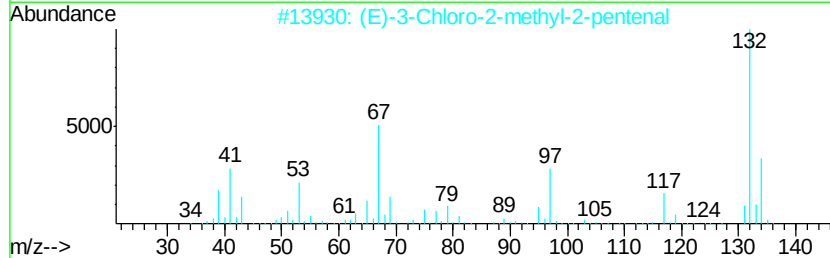
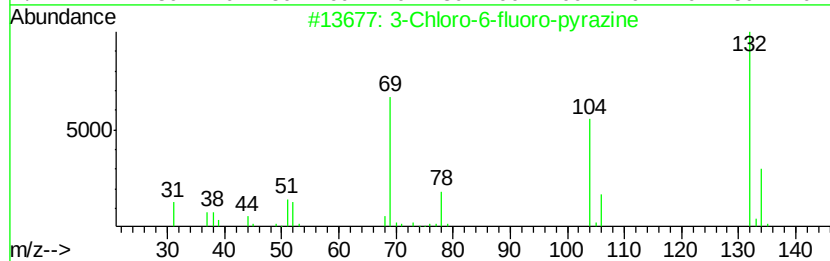
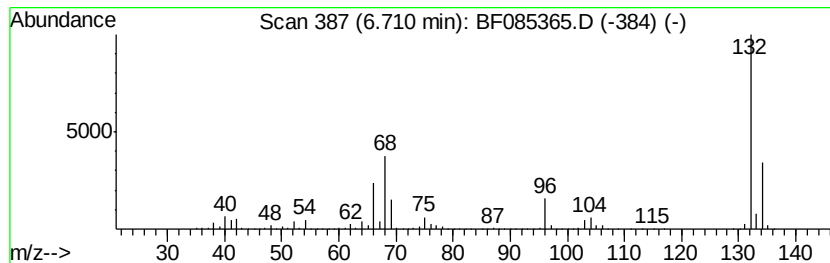
TIC Library : C:\DATABASE\NIST02.L  
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Peak Number 2 unknown6.71 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.71 | 99.27 ng | 4779950 | 1,4-Dichlorobenzene-d4 | 6.94 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 35   |
| 3    |    | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 9    |
| 4    |    | 1-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-59-9  | 9    |
| 5    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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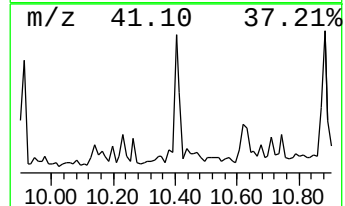
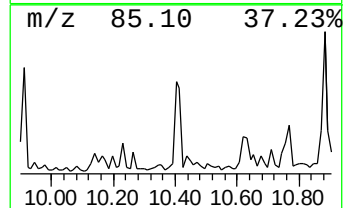
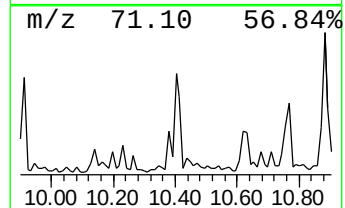
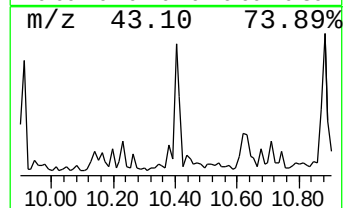
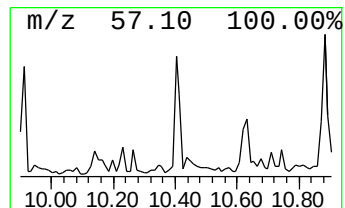
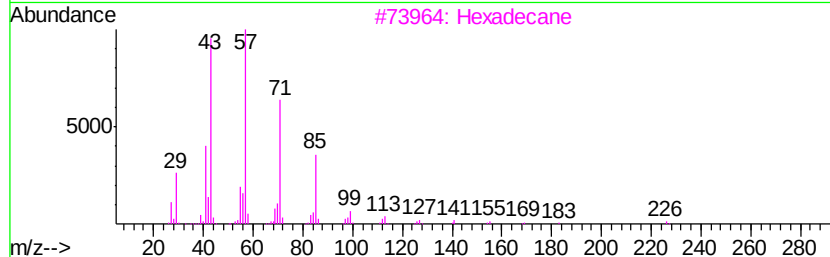
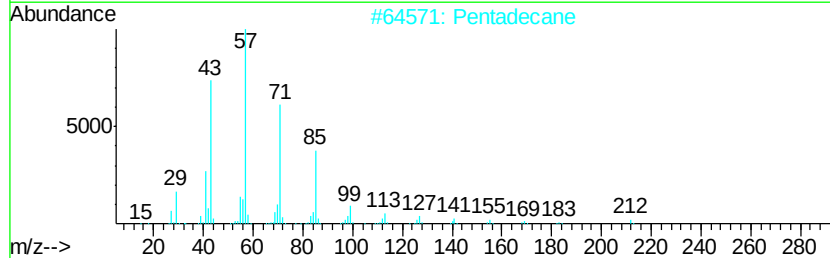
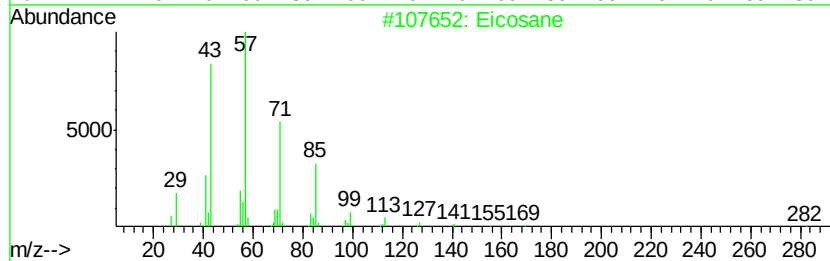
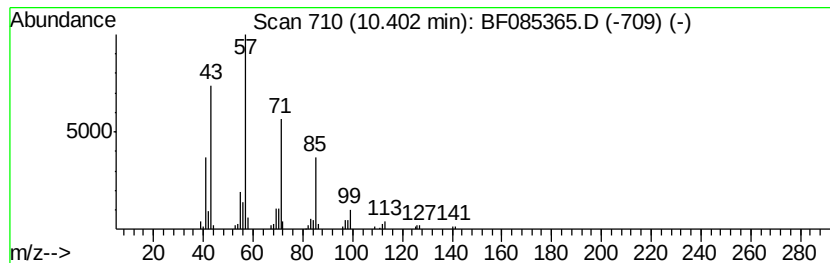
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Peak Number 4 Eicosane Concentration Rank 9

| R.T.    | EstConc     | Area         | Relative to ISTD | R.T.      |
|---------|-------------|--------------|------------------|-----------|
| 10.40   | 2.33 ng     | 139599       | Acenaphthene-d10 | 9.98      |
| Hit# of | 5           | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Eicosane    | 282 C20H42   | 000112-95-8      | 91        |
| 2       | Pentadecane | 212 C15H32   | 000629-62-9      | 90        |
| 3       | Hexadecane  | 226 C16H34   | 000544-76-3      | 90        |
| 4       | Nonadecane  | 268 C19H40   | 000629-92-5      | 86        |
| 5       | Tetradecane | 198 C14H30   | 000629-59-4      | 80        |



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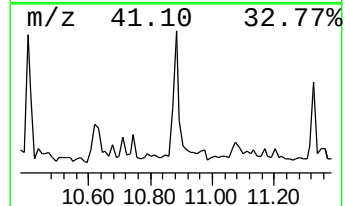
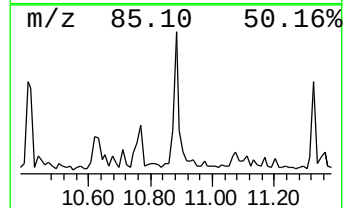
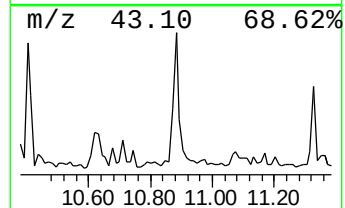
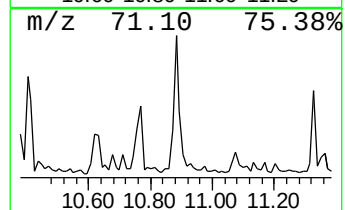
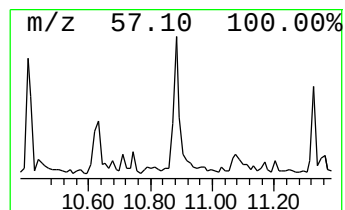
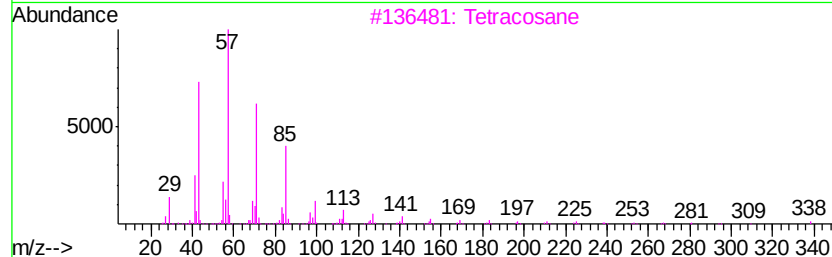
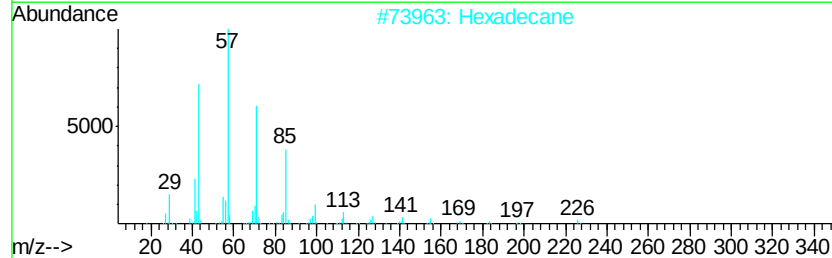
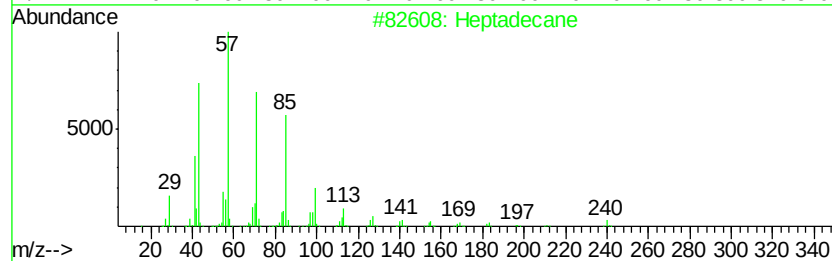
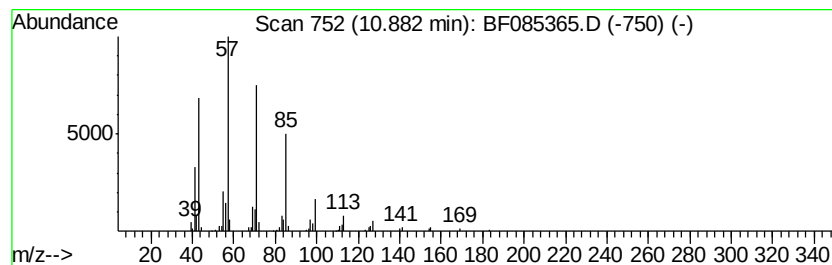
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Peak Number 5 Heptadecane Concentration Rank 7

| R.T.      | EstConc            | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------------|--------|------------------|-------------|-------|
| 10.88     | 3.88 ng            | 195670 | Phenanthrene-d10 |             | 11.47 |
| Hit# of 5 | Tentative ID       | MW     | MolForm          | CAS#        | Qual  |
| 1         | Heptadecane        | 240    | C17H36           | 000629-78-7 | 91    |
| 2         | Hexadecane         | 226    | C16H34           | 000544-76-3 | 90    |
| 3         | Tetracosane        | 338    | C24H50           | 000646-31-1 | 90    |
| 4         | Pentadecane        | 212    | C15H32           | 000629-62-9 | 90    |
| 5         | Tridecane, 1-iodo- | 310    | C13H27I          | 035599-77-0 | 86    |



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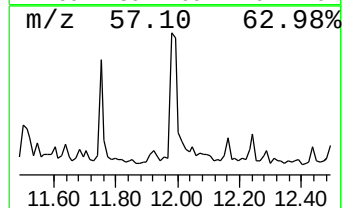
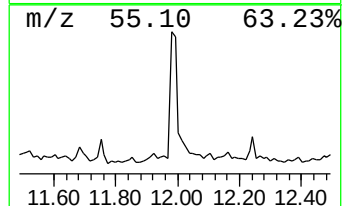
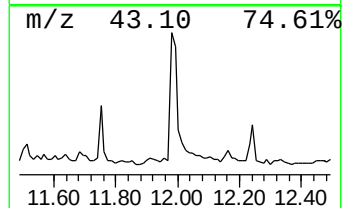
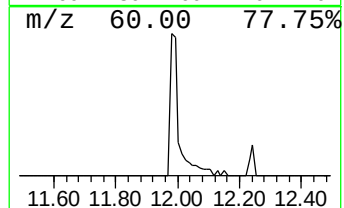
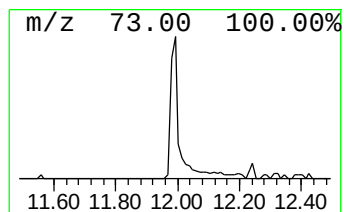
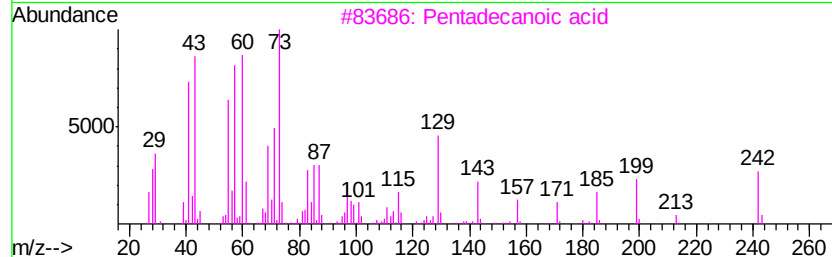
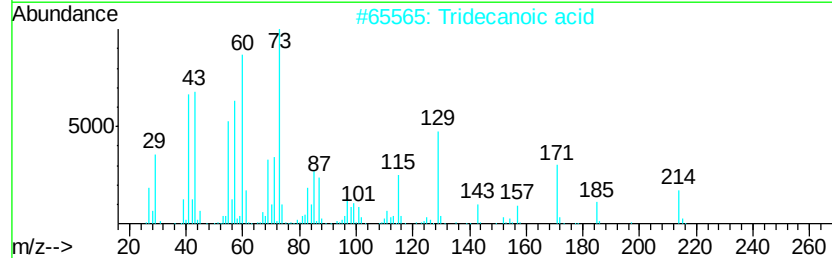
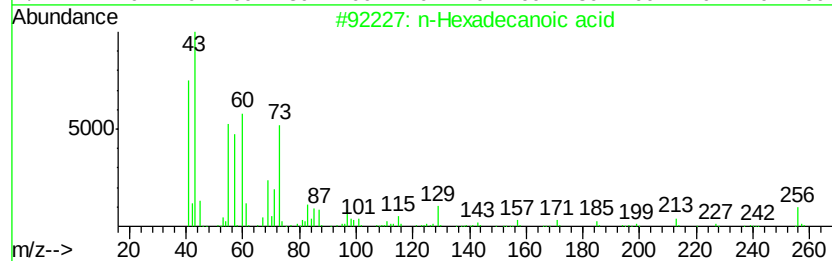
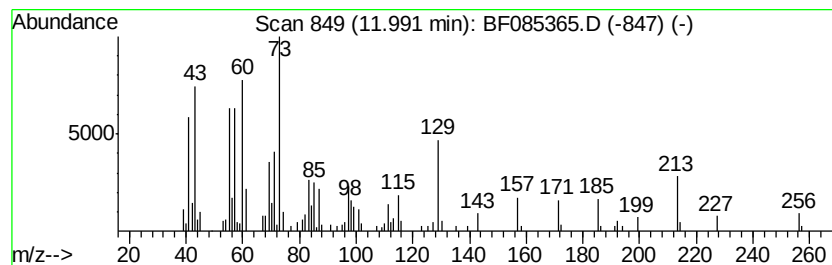
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\*\*\*\*\*  
Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 11.99     | 4.18 ng             | 210497 | Phenanthrene-d10 | 11.47       |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 96   |
| 2         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 91   |
| 3         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 83   |
| 4         | n-Decanoic acid     | 172    | C10H20O2         | 000334-48-5 | 76   |
| 5         | Undecanoic acid     | 186    | C11H22O2         | 000112-37-8 | 76   |



Data Path : S:\HPCHEM1\BNA\_F\DATA\BF031016\  
Data File : BF085365.D  
Acq On : 11 Mar 2016 6:31  
Operator : UM/SJ  
Sample : H1731-12  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

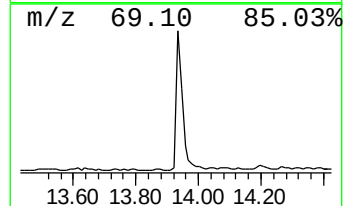
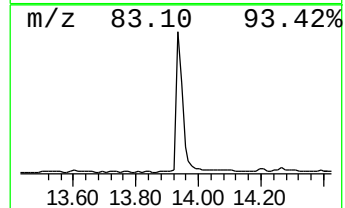
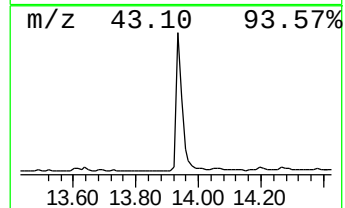
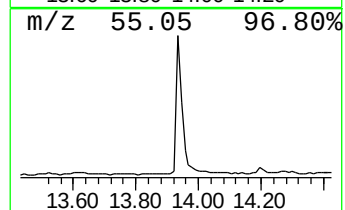
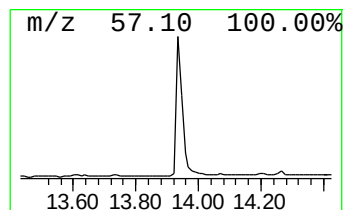
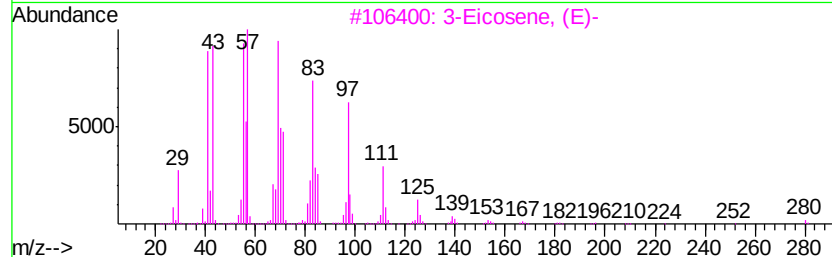
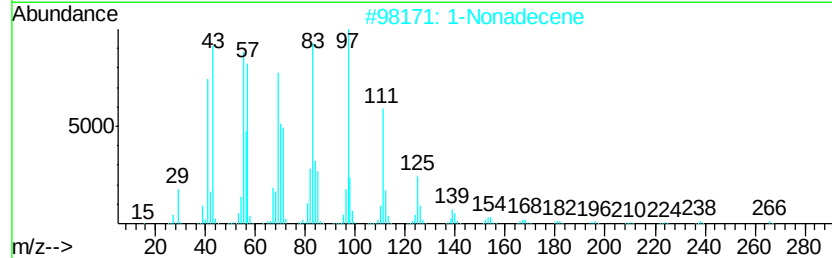
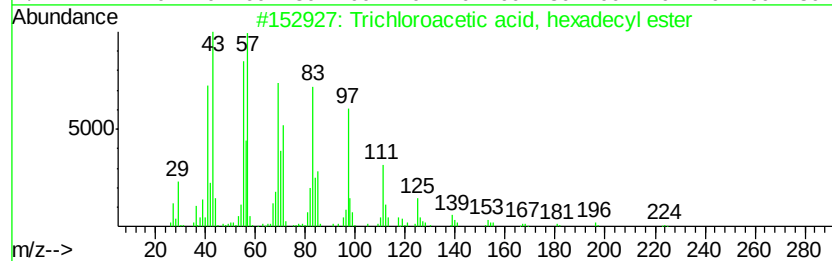
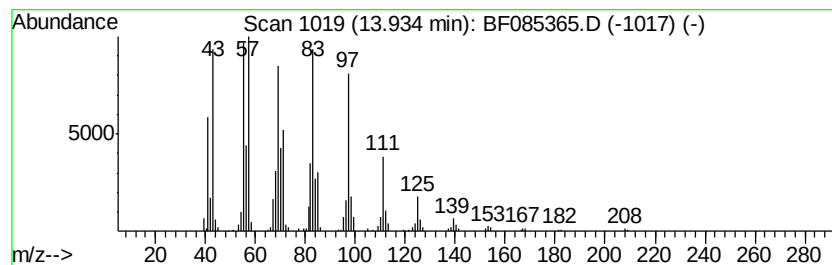
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Trichloroacetic acid, hexadecyl... Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.93 | 16.91 ng | 640310 | Chrysene-d12     | 14.09 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|-----------|-------------------------------------|-----|-------------|-------------|------|
| 1         | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 94   |
| 2         | 1-Nonadecene                        | 266 | C19H38      | 018435-45-5 | 91   |
| 3         | 3-Eicosene, (E)-                    | 280 | C20H40      | 074685-33-9 | 91   |
| 4         | 9-Eicosene, (E)-                    | 280 | C20H40      | 074685-29-3 | 91   |
| 5         | Trichloroacetic acid, tetradecyl... | 358 | C16H29Cl3O2 | 074339-52-9 | 90   |





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Data File : BF085365.D  
Acq On : 11 Mar 2016 6:31  
Operator : UM/SJ  
Sample : H1731-12  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

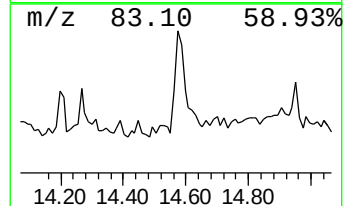
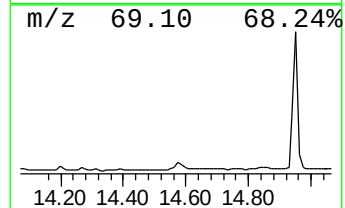
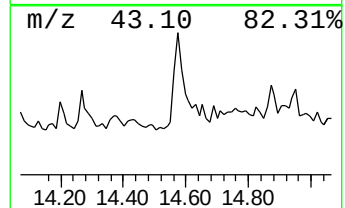
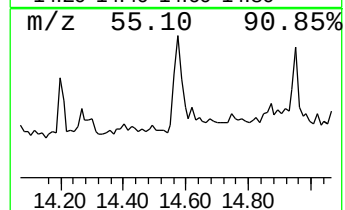
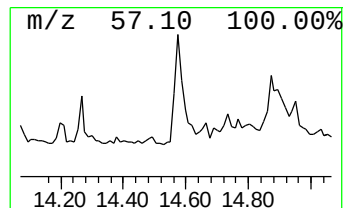
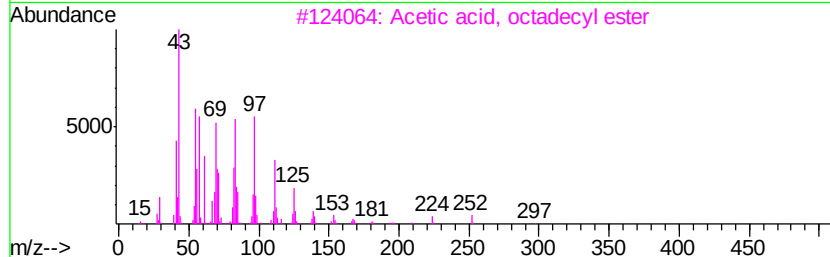
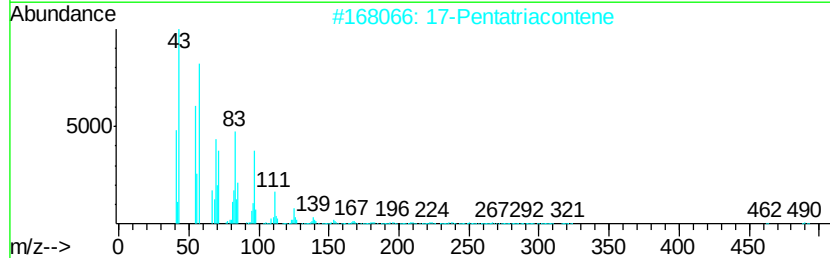
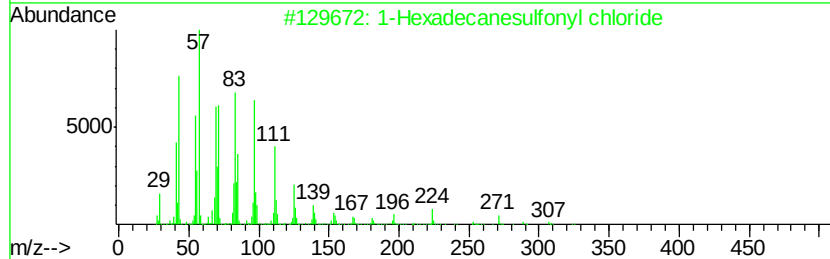
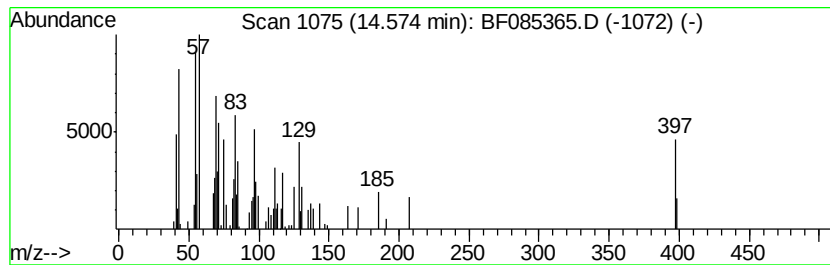
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown14.57 Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.57 | 2.28 ng | 86515 | Chrysene-d12     | 14.09 |

| Hit# | of | Tentative ID                  | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 1-Hexadecanesulfonyl chloride | 324 | C16H33ClO2S  | 038775-38-1  | 38   |
| 2    |    | 17-Pentatriacontene           | 491 | C35H70       | 006971-40-0  | 35   |
| 3    |    | Acetic acid, octadecyl ester  | 312 | C20H40O2     | 000822-23-1  | 30   |
| 4    |    | Bacteriochlorophyll-c-stearyl | 841 | C52H72MgN4O4 | 1000164-49-7 | 25   |
| 5    |    | 1-Hentetracontanol            | 593 | C41H84O      | 040710-42-7  | 25   |



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Operator : UM/SJ  
Sample : H1731-12  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

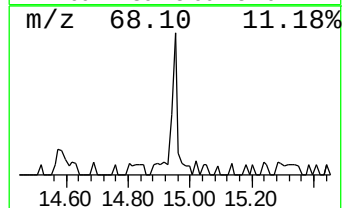
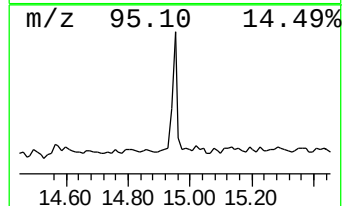
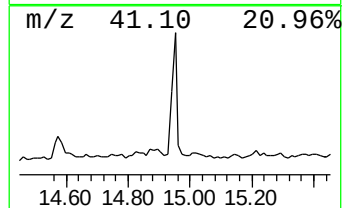
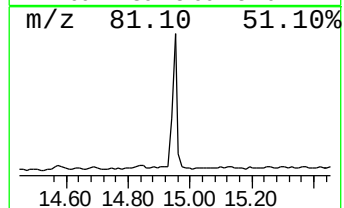
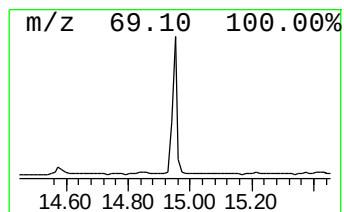
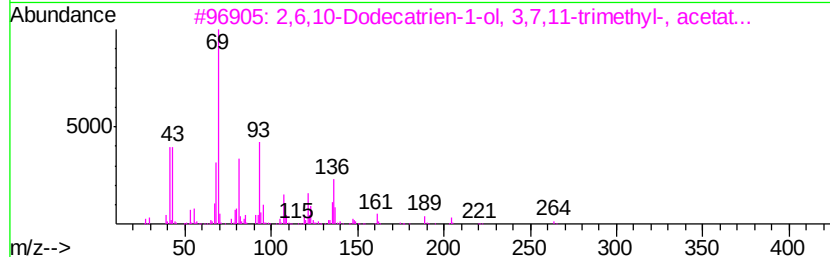
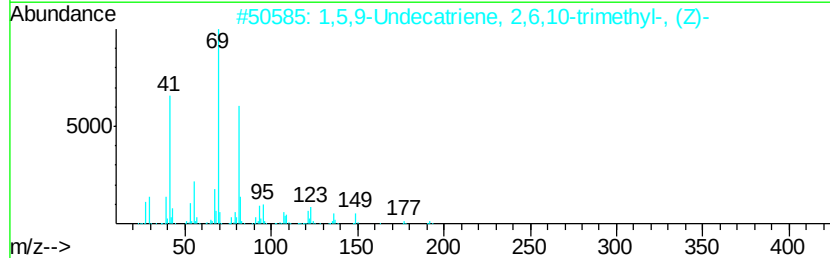
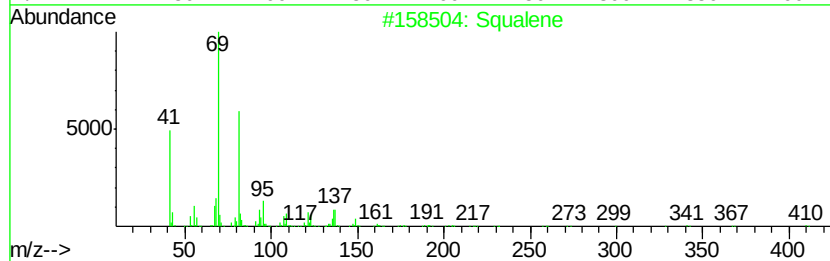
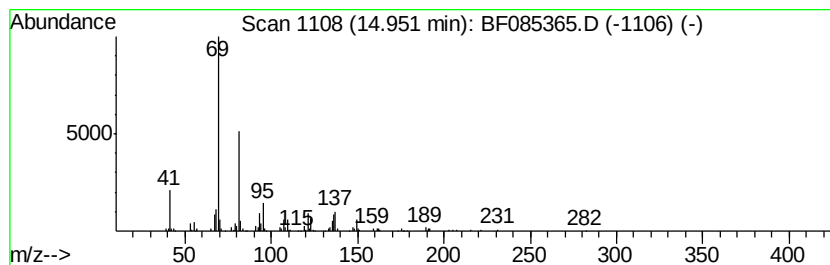
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Squalene Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.95 | 4.16 ng | 155953 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Squalene                            | 410 | C30H50   | 007683-64-9 | 91   |
| 2    |    | 1,5,9-Undecatriene, 2,6,10-trime... | 192 | C14H24   | 062951-96-6 | 80   |
| 3    |    | 2,6,10-Dodecatrien-1-ol, 3,7,11-... | 264 | C17H28O2 | 004128-17-0 | 64   |
| 4    |    | Furan, 3-(4-methyl-3-pentenyl)-     | 150 | C10H14O  | 000539-52-6 | 50   |
| 5    |    | 1,5-Heptadiene, 2,6-dimethyl-       | 124 | C9H16    | 006709-39-3 | 50   |



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Data File : BF085365.D  
Acq On : 11 Mar 2016 6:31  
Operator : UM/SJ  
Sample : H1731-12  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

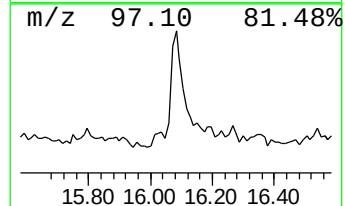
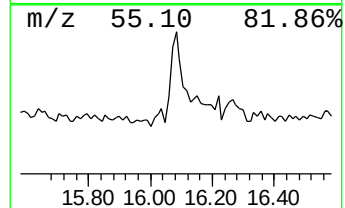
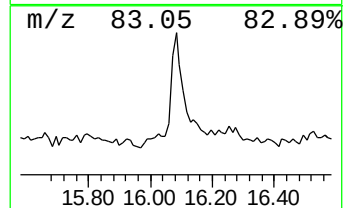
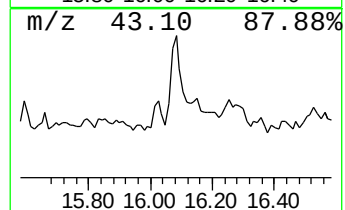
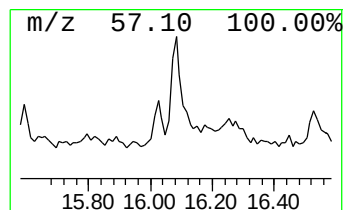
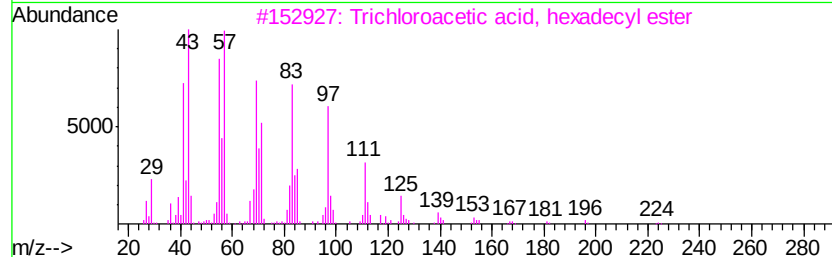
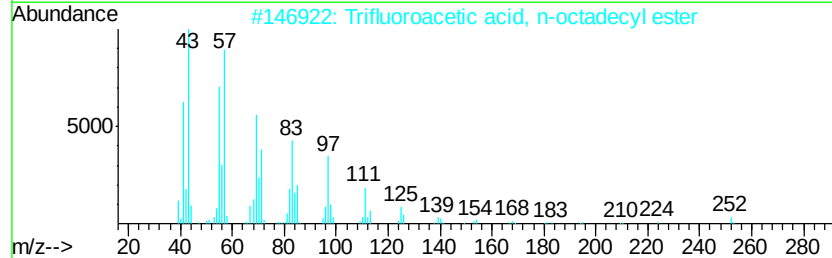
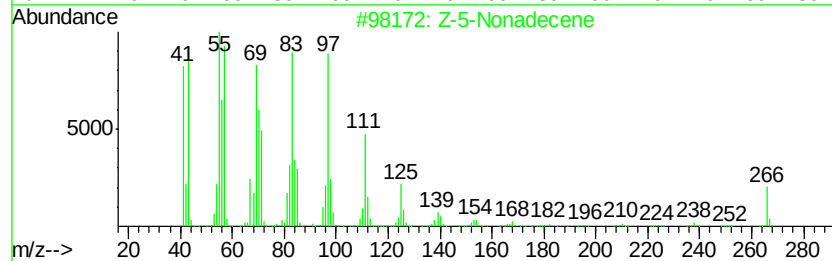
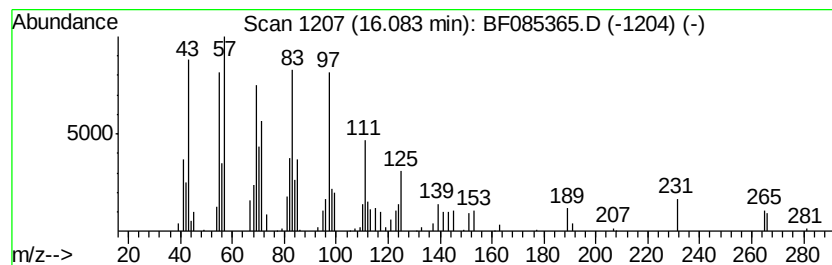
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Z-5-Nonadecene Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.08 | 3.28 ng | 123105 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Z-5-Nonadecene                      | 266 | C19H38      | 1000131-11-8 | 95   |
| 2    |    | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1  | 90   |
| 3    |    | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 87   |
| 4    |    | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2  | 000400-57-7  | 87   |
| 5    |    | 2- Chloropropionic acid, hexadec... | 332 | C19H37ClO2  | 086711-81-1  | 87   |



Data Path : S:\HPCHEM1\BNA\_F\DATA\BF031016\  
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Acq On : 11 Mar 2016 6:31  
Operator : UM/SJ  
Sample : H1731-12  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-28-COMP

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.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... | 5.14  | 6.8     | ng    | 328003   | 1                     | 6.94  | 963061  | 20.0 |
| unknown6.71          | 6.71  | 99.3    | ng    | 4779950  | 1                     | 6.94  | 963061  | 20.0 |
| Eicosane             | 10.40 | 2.3     | ng    | 139599   | 3                     | 9.98  | 1196150 | 20.0 |
| Heptadecane          | 10.88 | 3.9     | ng    | 195670   | 4                     | 11.47 | 1007960 | 20.0 |
| n-Hexadecanoic acid  | 11.99 | 4.2     | ng    | 210497   | 4                     | 11.47 | 1007960 | 20.0 |
| Trichloroacetic a... | 13.93 | 16.9    | ng    | 640310   | 5                     | 14.09 | 757247  | 20.0 |
| unknown14.57         | 14.57 | 2.3     | ng    | 86515    | 5                     | 14.09 | 757247  | 20.0 |
| Squalene             | 14.95 | 4.2     | ng    | 155953   | 6                     | 15.56 | 750025  | 20.0 |
| Z-5-Nonadecene       | 16.08 | 3.3     | ng    | 123105   | 6                     | 15.56 | 750025  | 20.0 |