

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
 Data File : BF080111.D  
 Acq On : 23 Jun 2015 1:31  
 Operator : TP/IZ  
 Sample : G2715-05  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 76-INFIELD(0-2)-3

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-BF061715.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         | 4.481       | 119           | 122         | 124          | rVB      | 31963          | 42611         | 2.53%           | 0.266%        |
| 2         | 5.533       | 212           | 214         | 217          | rBV      | 324576         | 310896        | 18.49%          | 1.940%        |
| 3         | 5.933       | 246           | 249         | 251          | rBV      | 1333550        | 1298495       | 77.24%          | 8.102%        |
| 4         | 6.333       | 281           | 284         | 286          | rBV      | 23845          | 21964         | 1.31%           | 0.137%        |
| 5         | 6.550       | 301           | 303         | 308          | rBV      | 34502          | 41065         | 2.44%           | 0.256%        |
| 6         | 6.927       | 333           | 336         | 338          | rBV      | 1096023        | 1304451       | 77.59%          | 8.139%        |
| 7         | 7.087       | 348           | 350         | 352          | rBV      | 1715542        | 1418333       | 84.37%          | 8.849%        |
| 8         | 7.133       | 352           | 354         | 360          | rVB      | 15605          | 17561         | 1.04%           | 0.110%        |
| 9         | 7.316       | 368           | 370         | 373          | rBV      | 255586         | 340895        | 20.28%          | 2.127%        |
| 10        | 7.476       | 382           | 384         | 387          | rBV      | 938053         | 996376        | 59.27%          | 6.217%        |
| 11        | 7.876       | 417           | 419         | 421          | rBV      | 1040599        | 826202        | 49.15%          | 5.155%        |
| 12        | 8.607       | 481           | 483         | 486          | rBV      | 434319         | 463529        | 27.57%          | 2.892%        |
| 13        | 9.156       | 529           | 531         | 533          | rBV      | 39075          | 34736         | 2.07%           | 0.217%        |
| 14        | 9.556       | 564           | 566         | 569          | rBV      | 21566          | 24696         | 1.47%           | 0.154%        |
| 15        | 9.693       | 575           | 578         | 580          | rVB      | 1207721        | 1483599       | 88.25%          | 9.257%        |
| 16        | 10.070      | 610           | 611         | 614          | rBV3     | 104893         | 115161        | 6.85%           | 0.719%        |
| 17        | 10.379      | 636           | 638         | 641          | rVB      | 459636         | 558093        | 33.20%          | 3.482%        |
| 18        | 11.088      | 698           | 700         | 703          | rBV      | 124869         | 117626        | 7.00%           | 0.734%        |
| 19        | 11.179      | 705           | 708         | 710          | rBV      | 942116         | 1010916       | 60.13%          | 6.307%        |
| 20        | 11.888      | 768           | 770         | 772          | rBV      | 789530         | 662678        | 39.42%          | 4.135%        |
| 21        | 13.156      | 880           | 881         | 883          | rBV      | 17104          | 17890         | 1.06%           | 0.112%        |
| 22        | 13.419      | 899           | 904         | 907          | rBV3     | 30137          | 62047         | 3.69%           | 0.387%        |
| 23        | 13.557      | 914           | 916         | 919          | rBV      | 1462128        | 1681147       | 100.00%         | 10.489%       |
| 24        | 13.785      | 933           | 936         | 937          | rBV3     | 16710          | 22800         | 1.36%           | 0.142%        |
| 25        | 14.094      | 958           | 963         | 965          | rBV2     | 55869          | 93936         | 5.59%           | 0.586%        |
| 26        | 14.608      | 1005          | 1008        | 1015         | rBV2     | 159333         | 328703        | 19.55%          | 2.051%        |
| 27        | 14.814      | 1024          | 1026        | 1029         | rBV      | 601568         | 790644        | 47.03%          | 4.933%        |
| 28        | 15.020      | 1041          | 1044        | 1049         | rBV3     | 25317          | 58094         | 3.46%           | 0.362%        |
| 29        | 15.420      | 1077          | 1079        | 1083         | rBV      | 154843         | 251170        | 14.94%          | 1.567%        |
| 30        | 16.243      | 1149          | 1151        | 1153         | rBV      | 142313         | 231360        | 13.76%          | 1.444%        |
| 31        | 16.677      | 1186          | 1189        | 1192         | rBV      | 459232         | 657567        | 39.11%          | 4.103%        |
| 32        | 16.985      | 1214          | 1216        | 1219         | rBV      | 67217          | 118683        | 7.06%           | 0.740%        |
| 33        | 17.294      | 1240          | 1243        | 1246         | rBV      | 102757         | 204389        | 12.16%          | 1.275%        |
| 34        | 17.351      | 1246          | 1248        | 1252         | rVV4     | 41430          | 91463         | 5.44%           | 0.571%        |

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Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |       |        |        |        |
|----|--------|------|------|------|------|-------|--------|--------|--------|
| 35 | 17.591 | 1267 | 1269 | 1280 | rVB4 | 65061 | 218943 | 13.02% | 1.366% |
| 36 | 18.311 | 1329 | 1332 | 1337 | rVB  | 46877 | 108907 | 6.48%  | 0.679% |

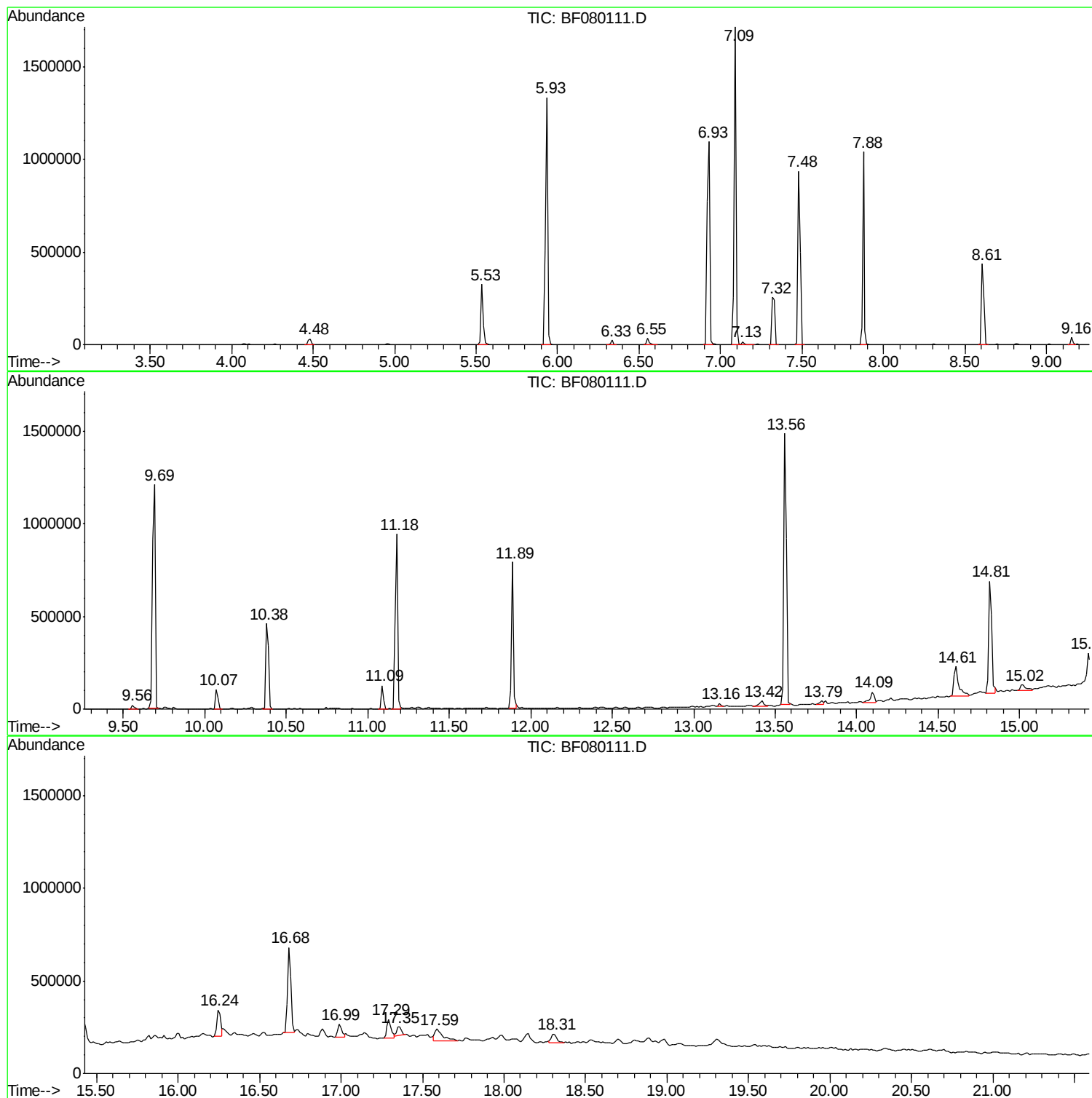
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

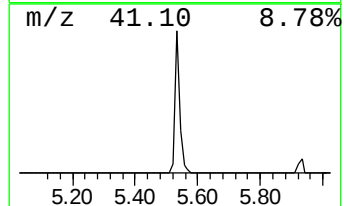
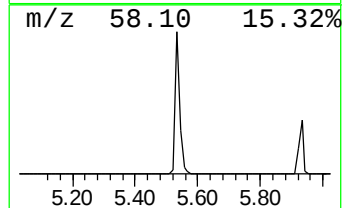
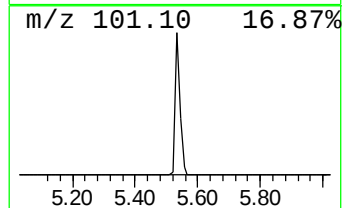
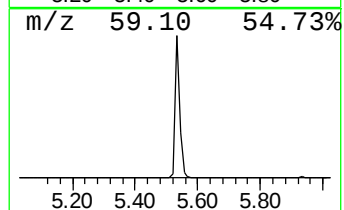
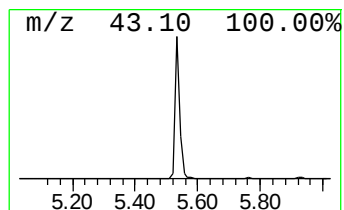
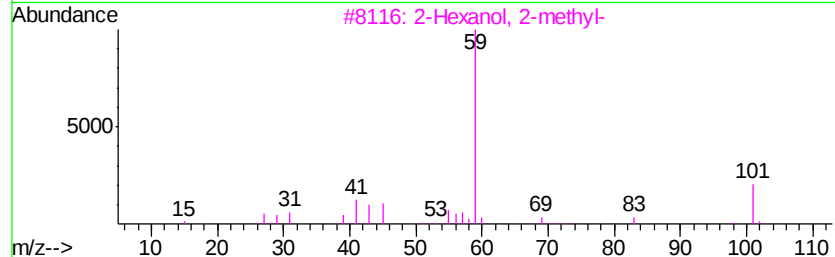
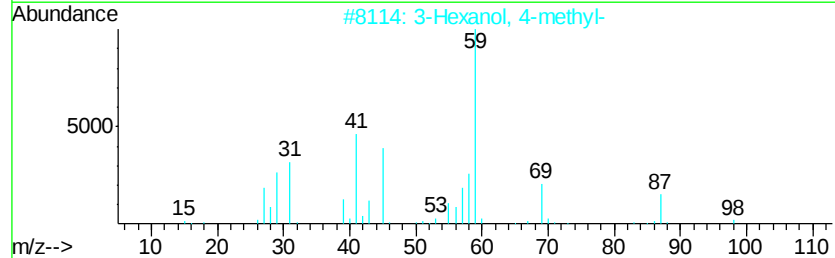
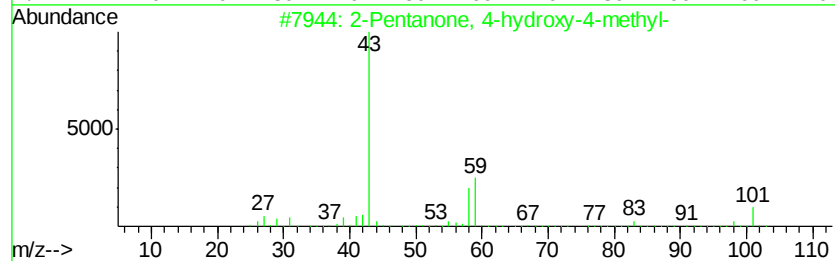
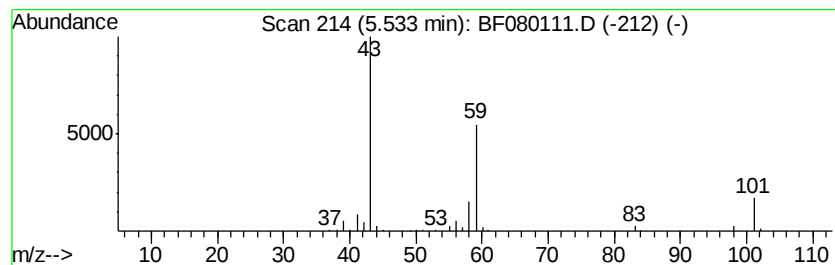
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.53 | 18.24 ng | 310896 | 1,4-Dichlorobenzene-d4 | 7.33 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 3    |    |   | 2-Hexanol, 2-methyl-                | 116 | C7H16O   | 000625-23-0 | 28   |
| 4    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 25   |
| 5    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 25   |



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ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

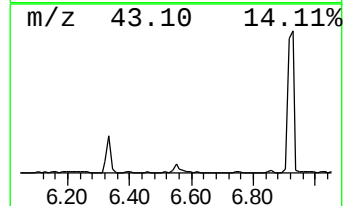
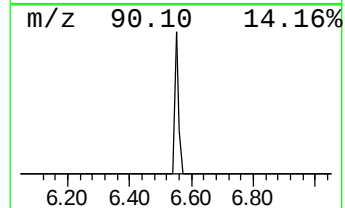
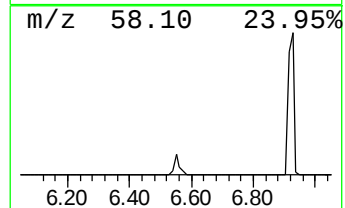
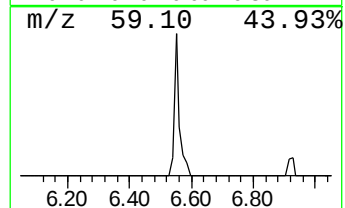
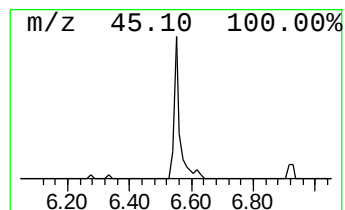
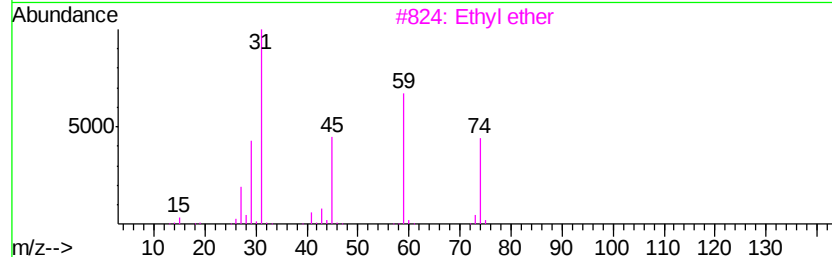
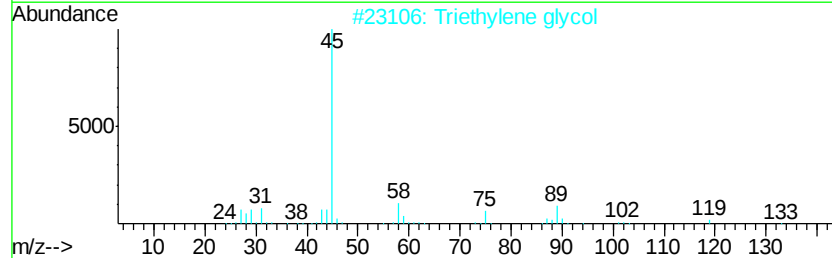
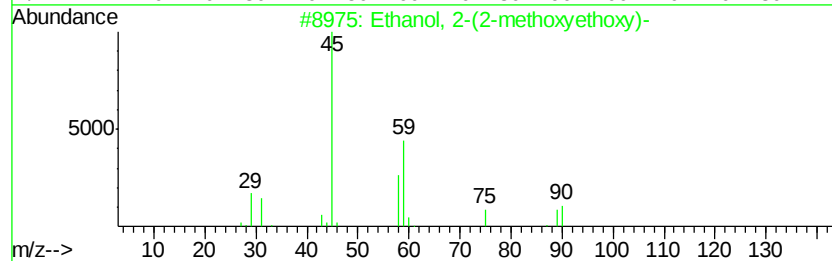
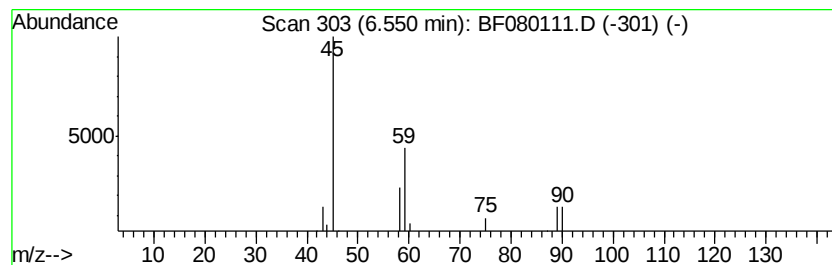
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Ethanol, 2-(2-methoxyethoxy)- Concentration Rank 14

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 6.55 | 2.41 ng | 41065 | 1,4-Dichlorobenzene-d4 | 7.33 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Ethanol, 2-(2-methoxyethoxy)- | 120 | C5H12O3 | 000111-77-3 | 83   |
| 2    |    |   | Triethylene glycol            | 150 | C6H14O4 | 000112-27-6 | 28   |
| 3    |    |   | Ethyl ether                   | 74  | C4H10O  | 000060-29-7 | 9    |
| 4    |    |   | Carbonic acid, dimethyl ester | 90  | C3H6O3  | 000616-38-6 | 9    |
| 5    |    |   | Butanoic acid, 4-methoxy-     | 118 | C5H10O3 | 029006-02-8 | 9    |



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Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

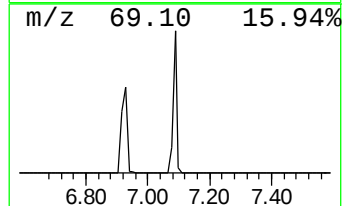
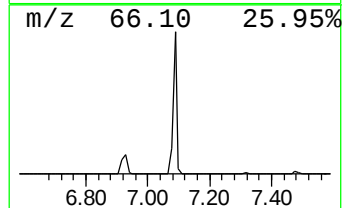
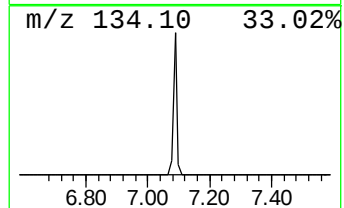
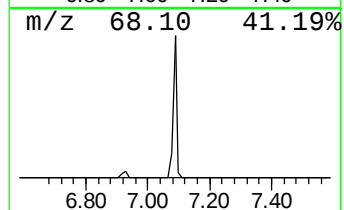
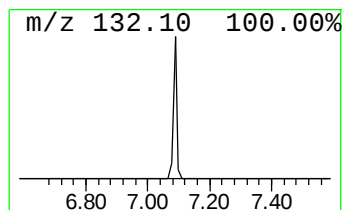
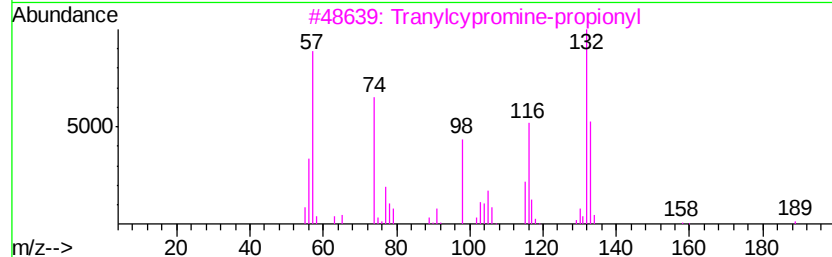
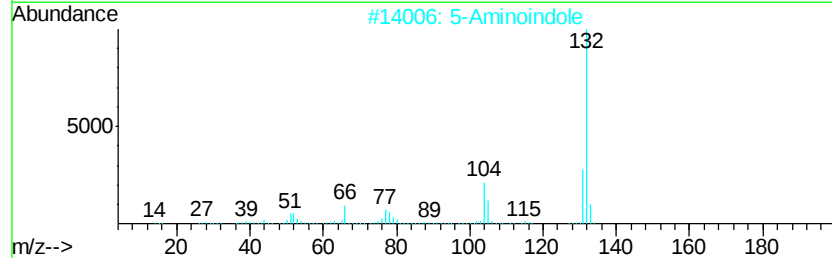
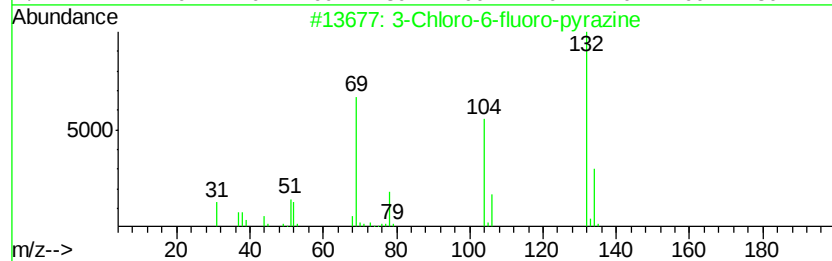
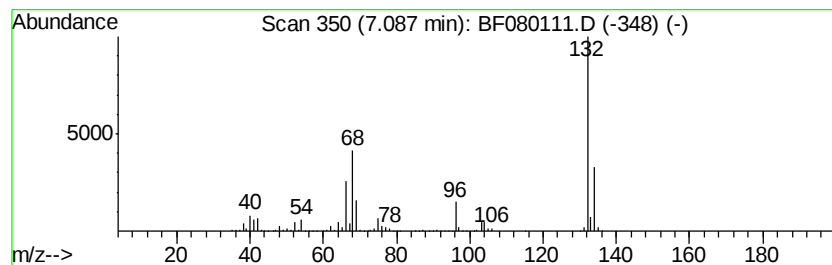
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 unknown7.09 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.09 | 83.21 ng | 1418330 | 1,4-Dichlorobenzene-d4 | 7.33 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    |   | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    |   | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    |   | Benzene, 1,3,5-trifluoro-           | 132 | C6H3F3    | 000372-38-3  | 9    |
| 5    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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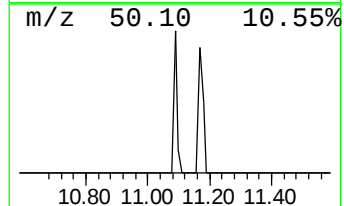
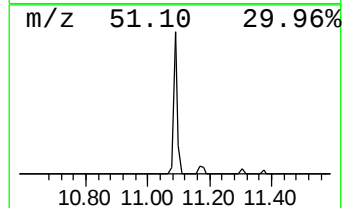
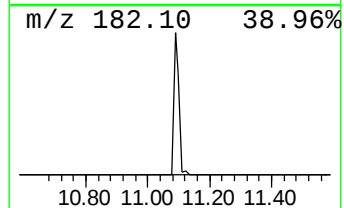
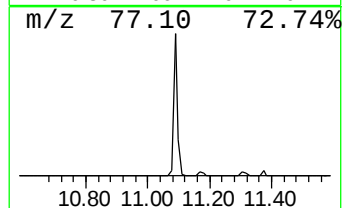
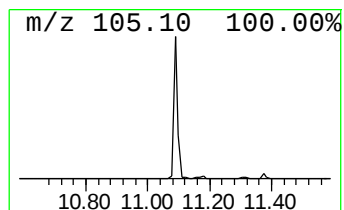
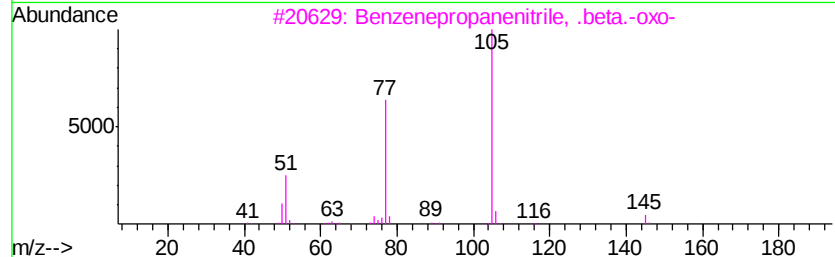
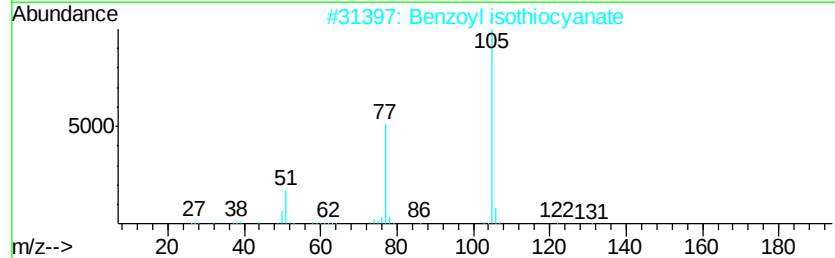
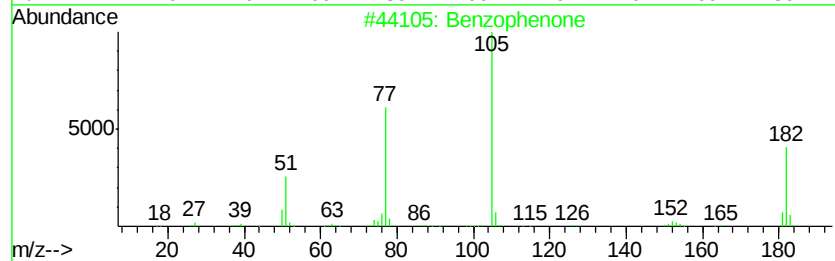
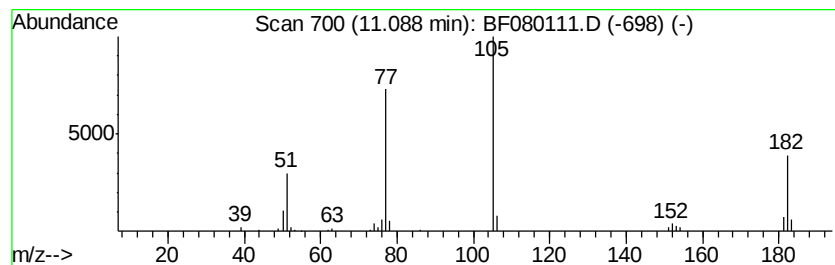
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzophenone Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.09 | 4.22 ng | 117626 | Acenaphthene-d10 | 10.38 |

| Hit# of | 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|---------|---|------------------------------------|-----|----------|-------------|------|
| 1       |   | Benzophenone                       | 182 | C13H10O  | 000119-61-9 | 96   |
| 2       |   | Benzoyl isothiocyanate             | 163 | C8H5NOS  | 000532-55-8 | 50   |
| 3       |   | Benzenepropanenitrile, .beta.-oxo- | 145 | C9H7NO   | 000614-16-4 | 50   |
| 4       |   | Benzoic acid, phenyl ester         | 198 | C13H10O2 | 000093-99-2 | 50   |
| 5       |   | Benzenebutanoic acid, .gamma.-oxo- | 178 | C10H10O3 | 002051-95-8 | 50   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

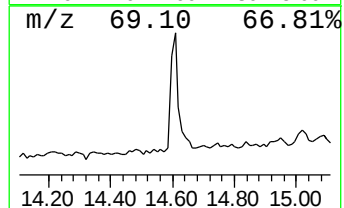
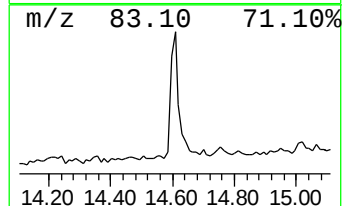
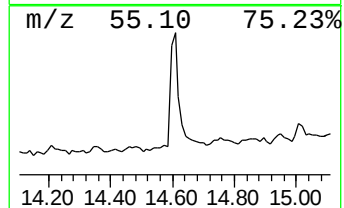
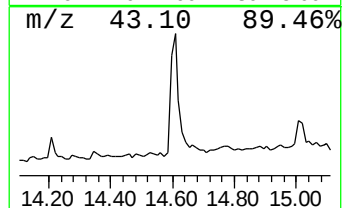
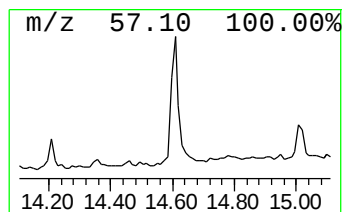
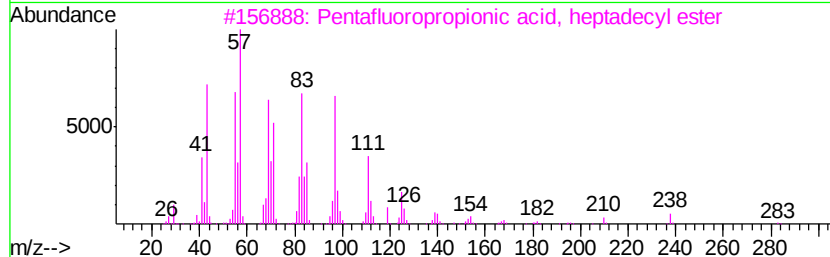
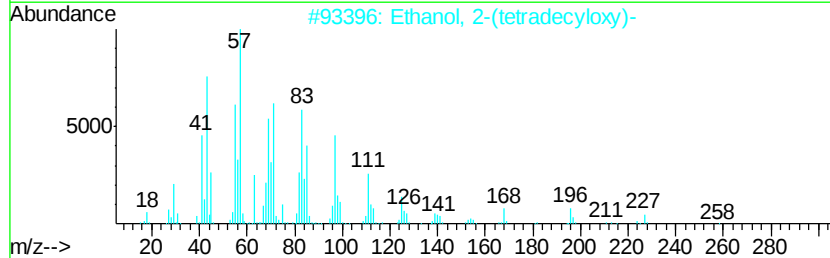
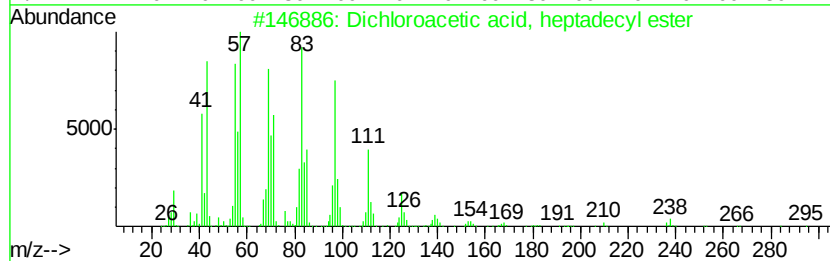
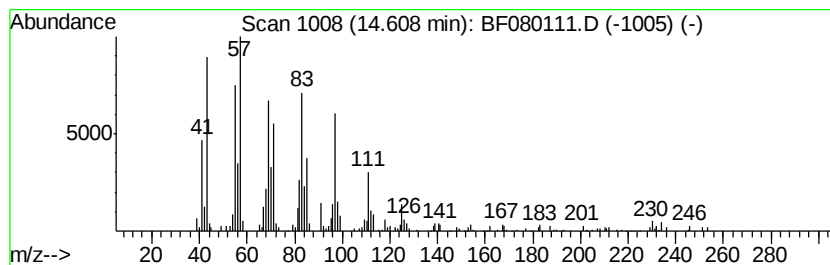
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Dichloroacetic acid, heptad... Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.61 | 8.31 ng | 328703 | Chrysene-d12     | 14.81 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 93   |
| 2    |    | Ethanol, 2-(tetradecyloxy)-         | 258 | C16H34O2    | 002136-70-1  | 91   |
| 3    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2  | 1000283-04-2 | 91   |
| 4    |    | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2  | 000400-57-7  | 91   |
| 5    |    | Heptafluorobutanoic acid, heptad... | 452 | C21H35F7O2  | 1000282-97-3 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

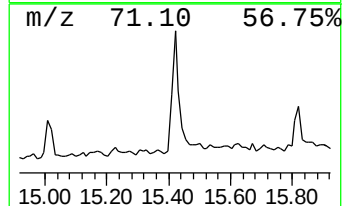
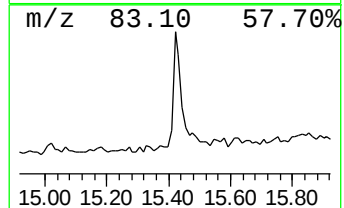
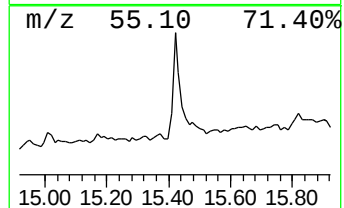
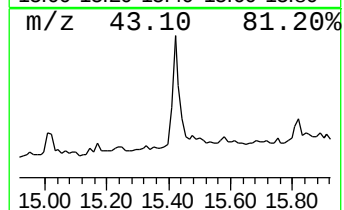
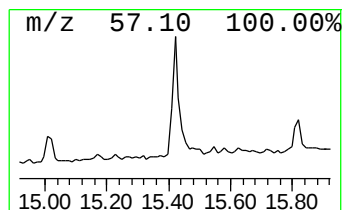
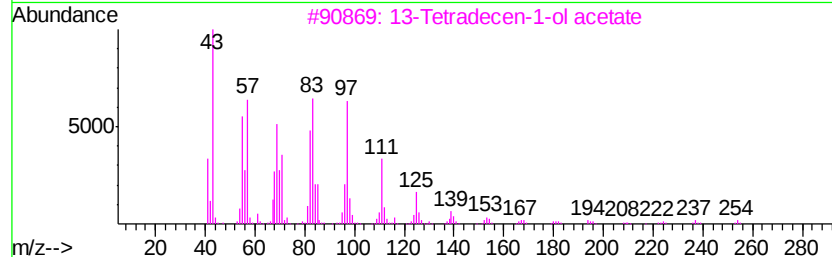
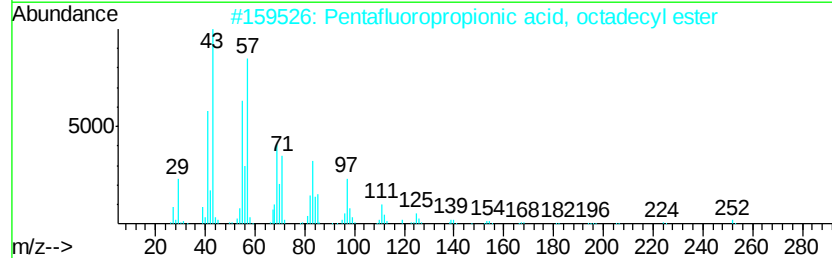
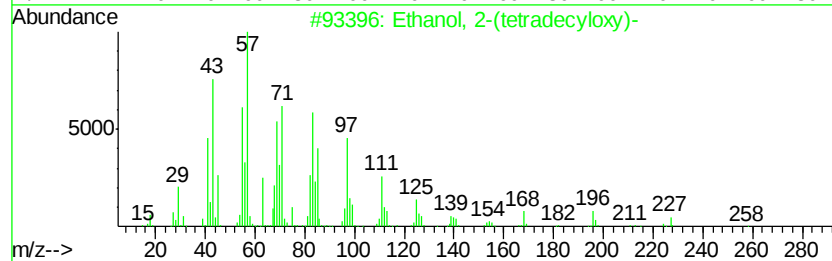
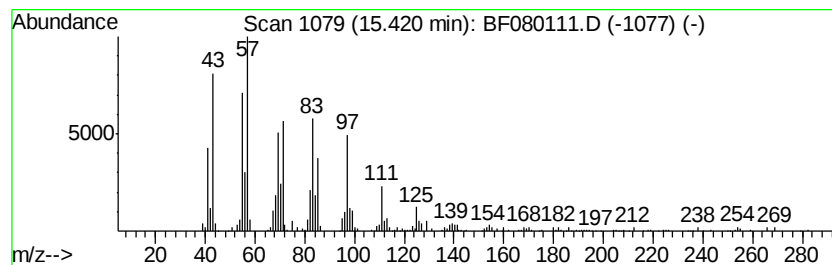
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ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.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 Ethanol, 2-(tetradecyloxy)- Concentration Rank 7

| R.T.      | EstConc                                                                                                                                                                                                                                                                                                                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--------|------------------|--------------|------|
| 15.42     | 6.35 ng                                                                                                                                                                                                                                                                                                                             | 251170 | Chrysene-d12     | 14.81        |      |
| Hit# of 5 | Tentative ID                                                                                                                                                                                                                                                                                                                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Ethanol, 2-(tetradecyloxy)-                                                                                                                                                                                                                                                                                                         | 258    | C16H34O2         | 002136-70-1  | 91   |
| 2         | Pentafluoropropionic acid, octadecanoic acid, heptadecanoic acid, hexadecanoic acid, pentadecanoic acid, tetradecanoic acid, tridecanoic acid, dodecanoic acid, undecanoic acid, decanoic acid, nonanoic acid, octanoic acid, heptanoic acid, hexanoic acid, pentanoic acid, butyric acid, propionic acid, acetic acid, formic acid | 416    | C21H37F5O2       | 1000280-07-7 | 91   |
| 3         | 13-Tetradecen-1-ol acetate                                                                                                                                                                                                                                                                                                          | 254    | C16H30O2         | 056221-91-1  | 89   |
| 4         | Heptafluorobutanoic acid, heptadecanoic acid, hexadecanoic acid, pentadecanoic acid, tetradecanoic acid, tridecanoic acid, dodecanoic acid, undecanoic acid, decanoic acid, nonanoic acid, octanoic acid, heptanoic acid, hexanoic acid, pentanoic acid, butyric acid, propionic acid, acetic acid, formic acid                     | 452    | C21H35F7O2       | 1000282-97-3 | 87   |
| 5         | 1-Nonadecanol                                                                                                                                                                                                                                                                                                                       | 284    | C19H40O          | 001454-84-8  | 74   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

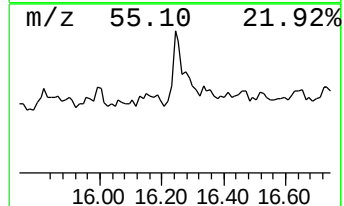
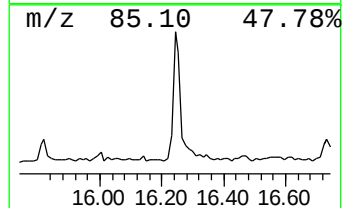
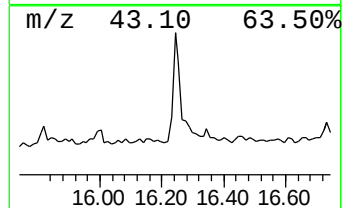
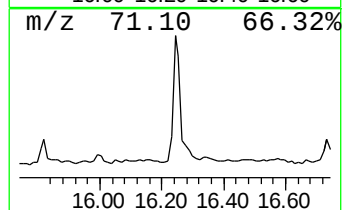
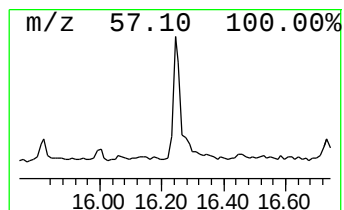
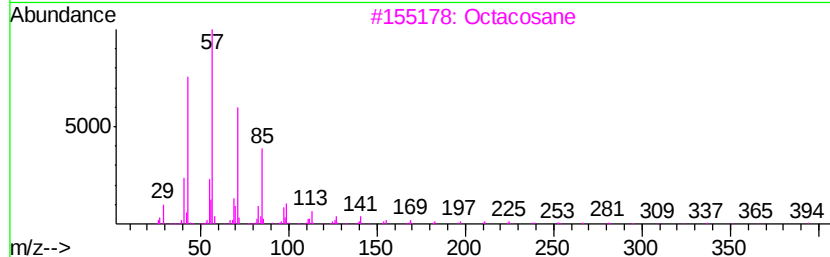
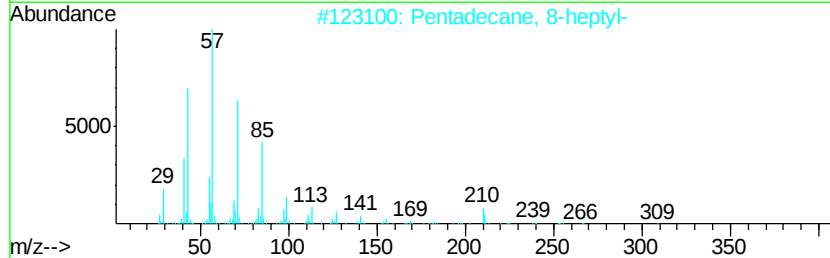
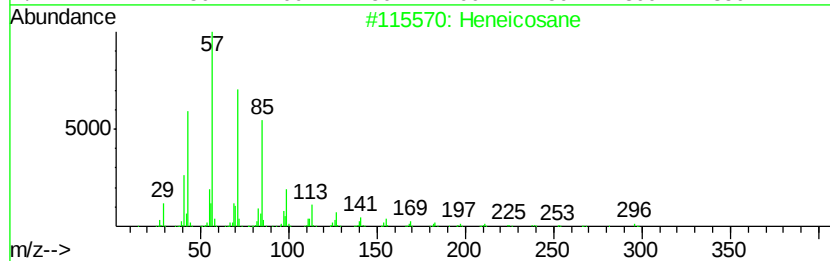
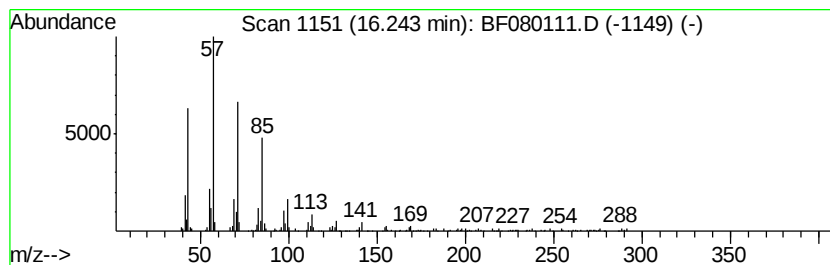
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Heneicosane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.24 | 7.04 ng | 231360 | Perylene-d12     | 16.68 |

| Hit# of | 5                      | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane            |              | 296 | C21H44  | 000629-94-7 | 90   |
| 2       | Pentadecane, 8-heptyl- |              | 310 | C22H46  | 071005-15-7 | 87   |
| 3       | Octacosane             |              | 394 | C28H58  | 000630-02-4 | 87   |
| 4       | Pentacosane            |              | 352 | C25H52  | 000629-99-2 | 86   |
| 5       | Tetracosane            |              | 338 | C24H50  | 000646-31-1 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

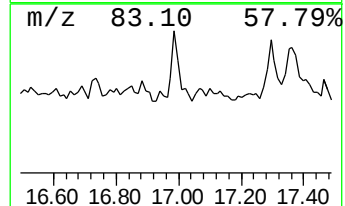
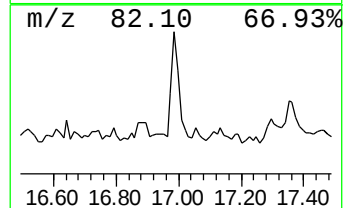
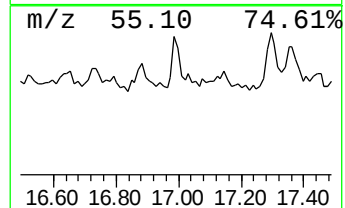
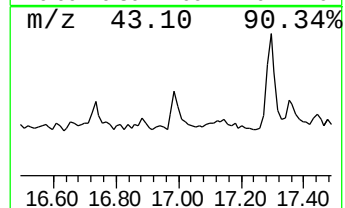
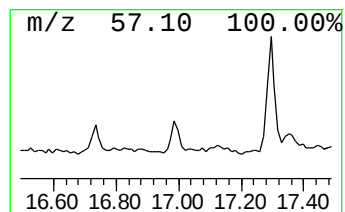
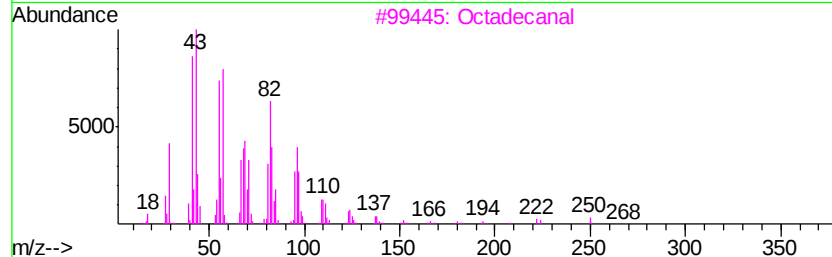
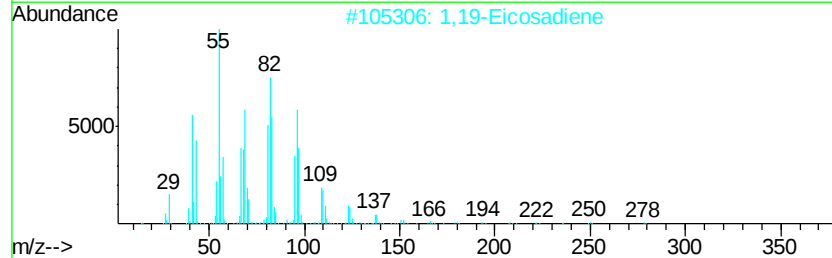
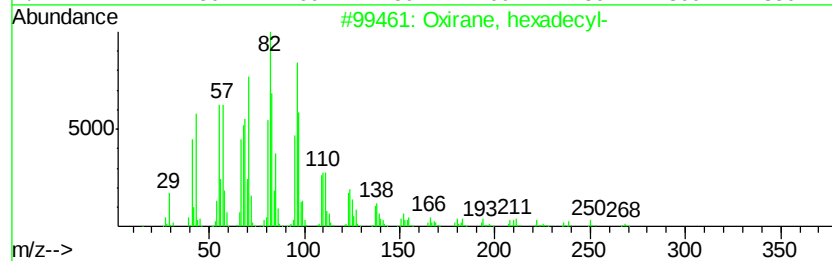
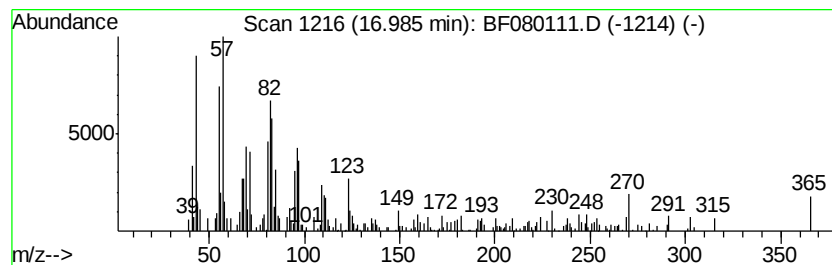
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Oxirane, hexadecyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.99 | 3.61 ng | 118683 | Perylene-d12     | 16.68 |

| Hit# | of | Tentative ID              | MW  | MolForm  | CAS#         | Qual |
|------|----|---------------------------|-----|----------|--------------|------|
| 1    | 5  | Oxirane, hexadecyl-       | 268 | C18H36O  | 007390-81-0  | 55   |
| 2    |    | 1,19-Eicosadiene          | 278 | C20H38   | 014811-95-1  | 49   |
| 3    |    | Octadecanal               | 268 | C18H36O  | 000638-66-4  | 47   |
| 4    |    | (Z)-14-Tricosenyl formate | 366 | C24H46O2 | 077899-10-6  | 47   |
| 5    |    | 18-Nonadecen-1-ol         | 282 | C19H38O  | 1000142-89-2 | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

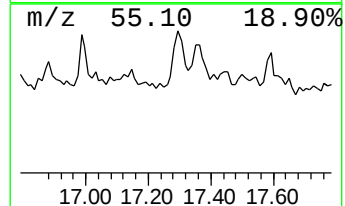
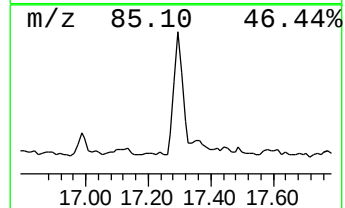
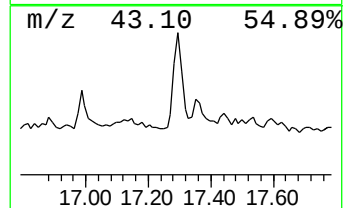
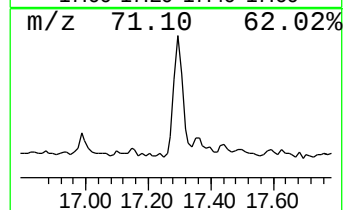
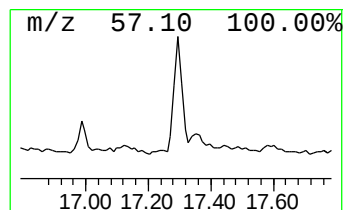
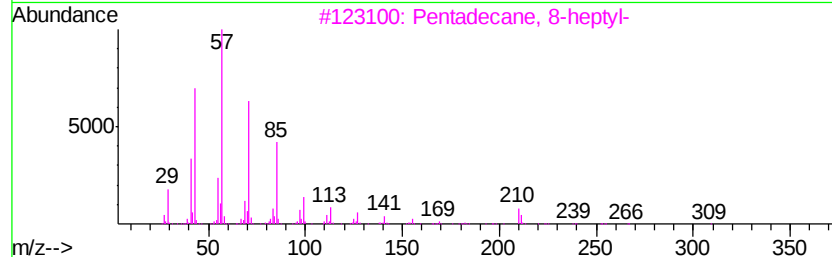
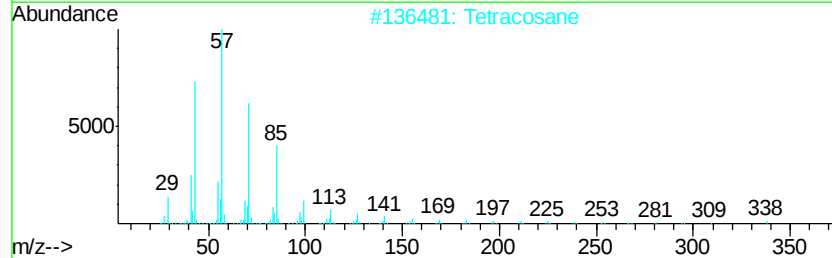
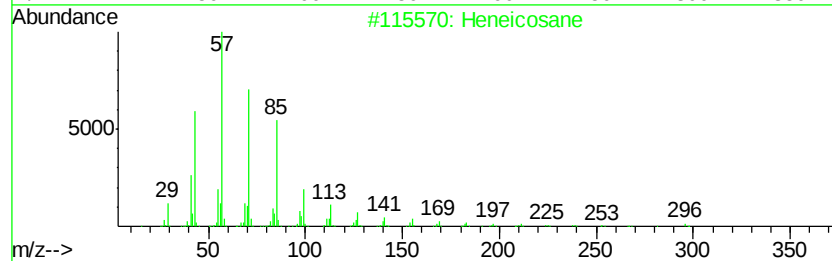
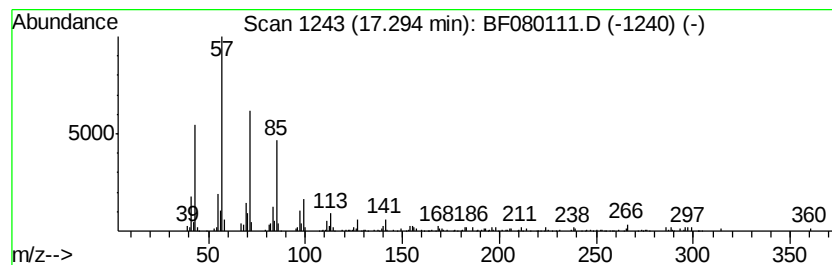
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Tetracosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.29 | 6.22 ng | 204389 | Perylene-d12     | 16.68 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7 | 99   |
| 2    |    | Tetracosane            | 338 | C24H50  | 000646-31-1 | 87   |
| 3    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 87   |
| 4    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 87   |
| 5    |    | Docosane               | 310 | C22H46  | 000629-97-0 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

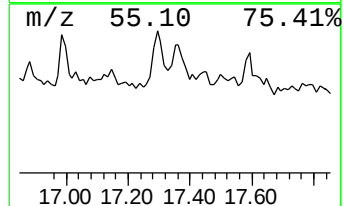
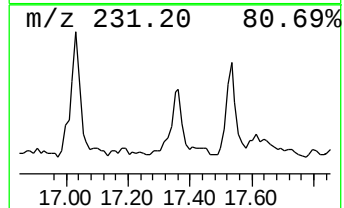
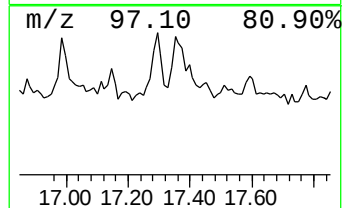
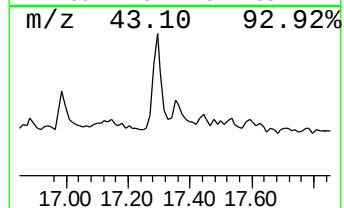
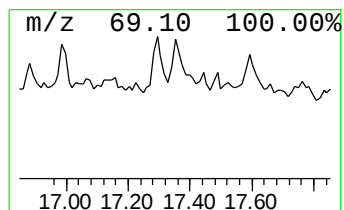
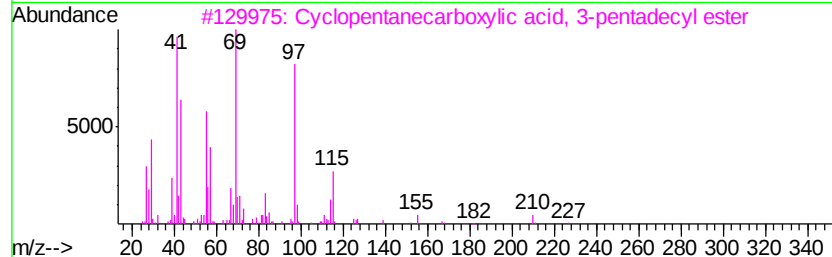
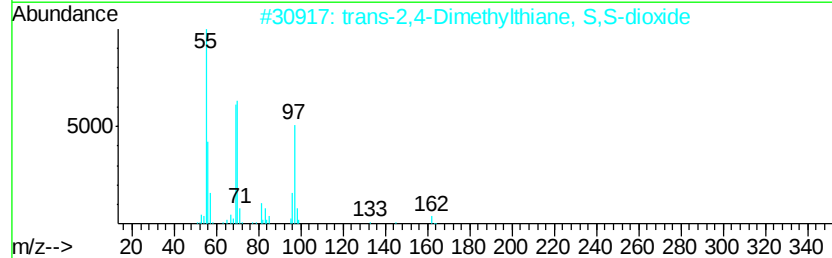
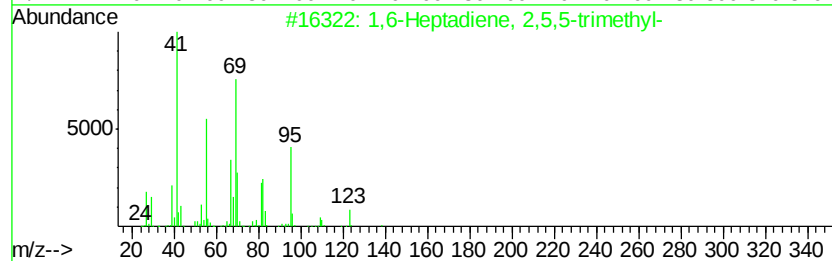
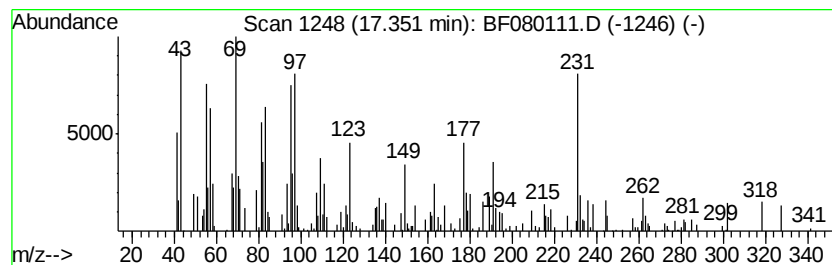
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown17.35 Concentration Rank 12

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 17.35 | 2.78 ng | 91463 | Perylene-d12     | 16.68 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1,6-Heptadiene, 2,5,5-trimethyl-    | 138 | C10H18    | 062238-28-2  | 22   |
| 2    |    |   | trans-2,4-Dimethylthiane, S,S-di... | 162 | C7H14O2S  | 1000215-67-6 | 22   |
| 3    |    |   | Cyclopentanecarboxylic acid, 3-p... | 324 | C21H40O2  | 1000280-59-1 | 14   |
| 4    |    |   | 1,2-Cyclohexanediol, cyclic sulf... | 162 | C6H10O3S  | 019456-19-0  | 14   |
| 5    |    |   | 12-Bromo-1-dodecanol                | 264 | C12H25BrO | 003344-77-2  | 14   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

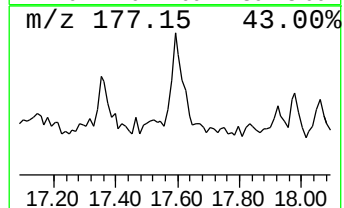
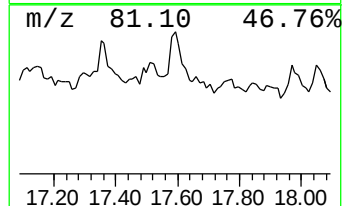
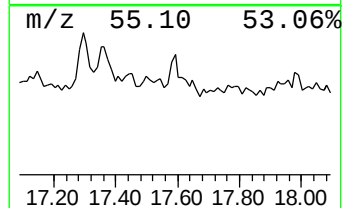
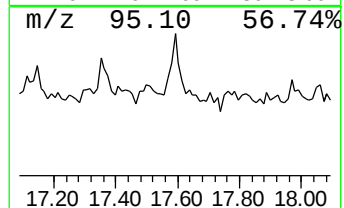
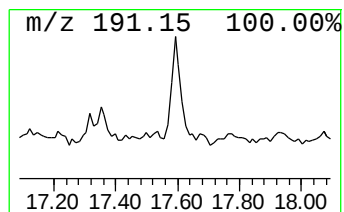
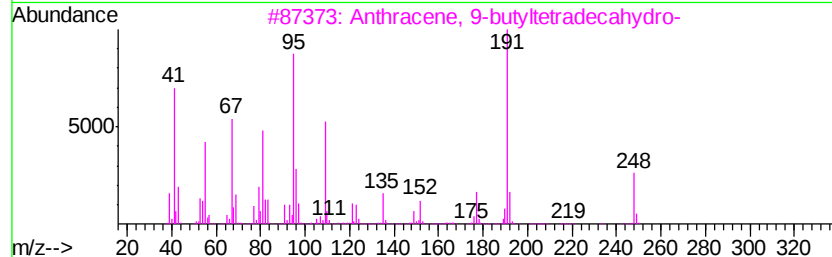
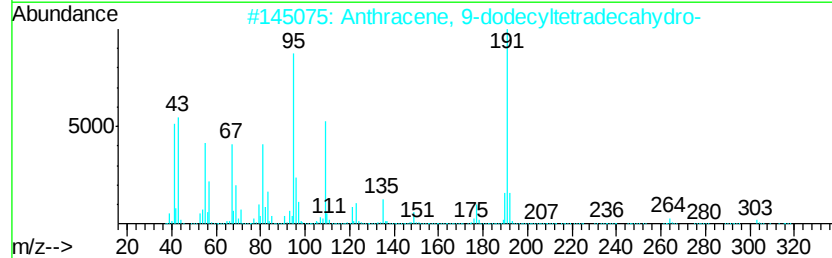
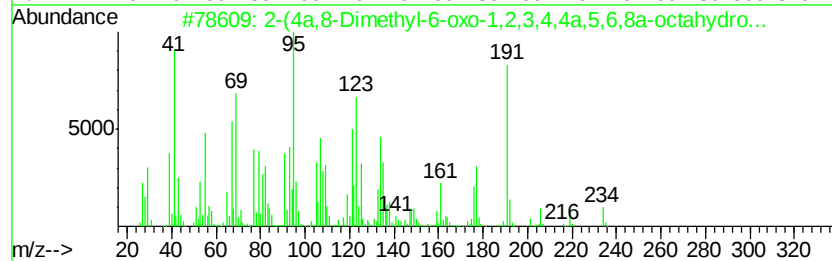
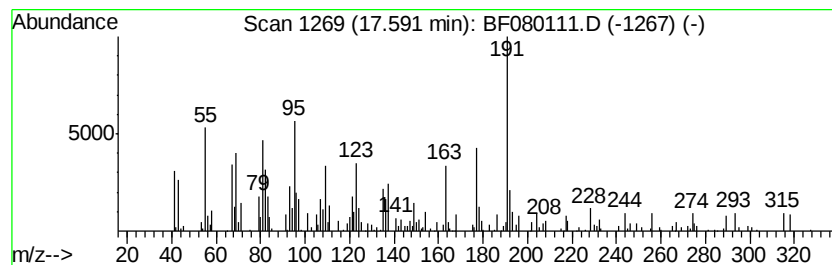
Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.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 unknown17.59 Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 17.59     | 6.66 ng                             | 218943 | Perylene-d12     | 16.68        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2-(4a,8-Dimethyl-6-oxo-1,2,3,4,4... | 234    | C15H22O2         | 1000190-21-8 | 32   |
| 2         | Anthracene, 9-dodecyltetradecahy... | 360    | C26H48           | 055401-75-7  | 32   |
| 3         | Anthracene, 9-butyltetradecahydro-  | 248    | C18H32           | 055133-89-6  | 32   |
| 4         | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206    | C14H22O          | 000127-43-5  | 27   |
| 5         | 2'-Fluoro-6'-(trifluoromethyl)-a... | 206    | C9H6F4O          | 174013-29-7  | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

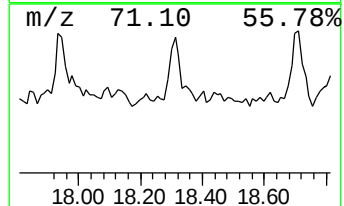
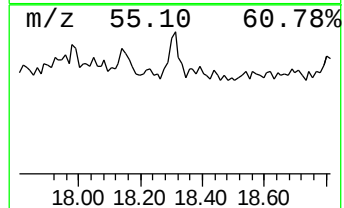
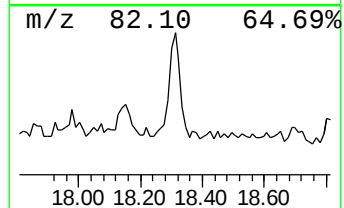
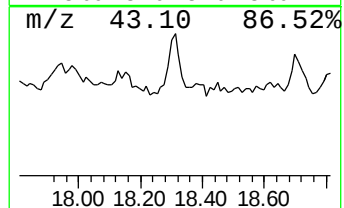
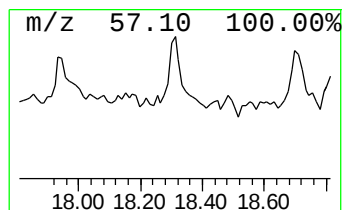
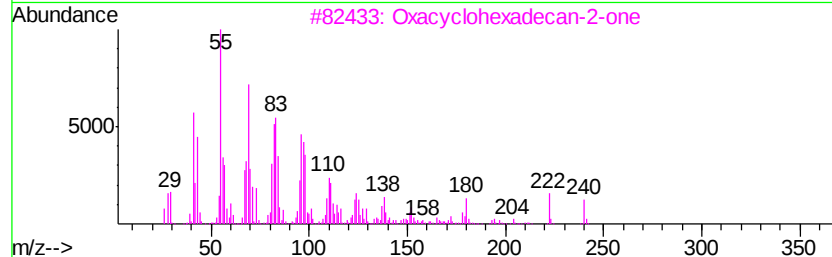
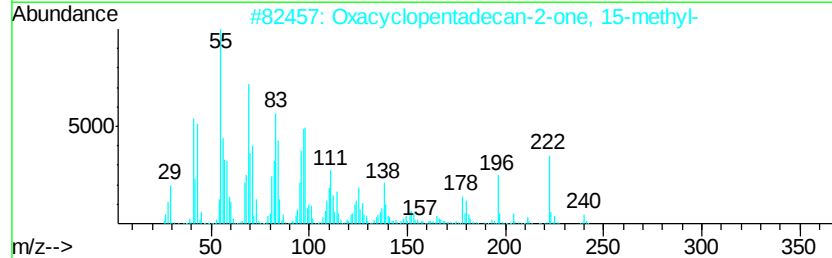
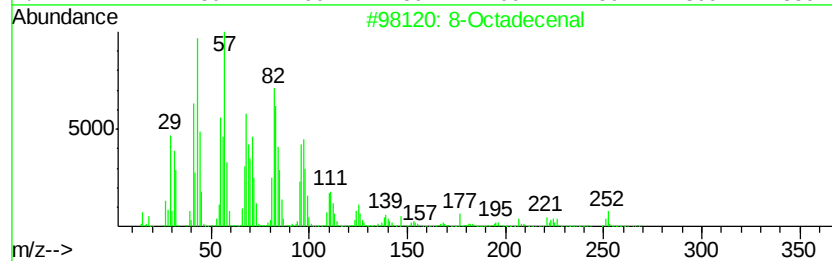
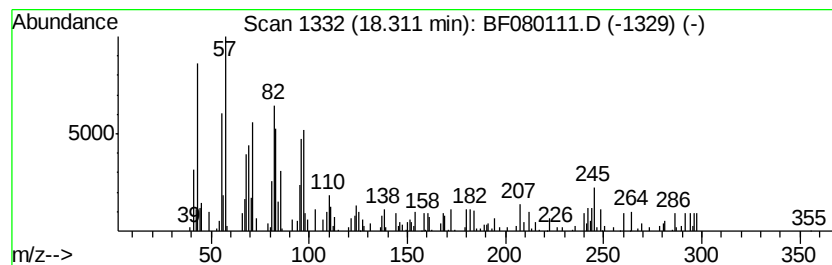
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown18.31 Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.31 | 3.31 ng | 108907 | Perylene-d12     | 16.68 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 8-Octadecenal                       | 266 | C18H34O  | 056554-94-0 | 38   |
| 2    |    | Oxacyclopentadecan-2-one, 15-met... | 240 | C15H28O2 | 032539-85-8 | 35   |
| 3    |    | Oxacyclohexadecan-2-one             | 240 | C15H28O2 | 000106-02-5 | 30   |
| 4    |    | Oxirane, heptadecyl-                | 282 | C19H38O  | 067860-04-2 | 25   |
| 5    |    | 1,19-Eicosadiene                    | 278 | C20H38   | 014811-95-1 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062215\  
Data File : BF080111.D  
Acq On : 23 Jun 2015 1:31  
Operator : TP/IZ  
Sample : G2715-05  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
76-INFIELD(0-2)-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF061715.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... | 5.53  | 18.2    | ng    | 310896   | 1                     | 7.33  | 340895 | 20.0 |
| Ethanol, 2-(2-met... | 6.55  | 2.4     | ng    | 41065    | 1                     | 7.33  | 340895 | 20.0 |
| unknown7.09          | 7.09  | 83.2    | ng    | 1418330  | 1                     | 7.33  | 340895 | 20.0 |
| Benzophenone         | 11.09 | 4.2     | ng    | 117626   | 3                     | 10.38 | 558093 | 20.0 |
| Dichloroacetic ac... | 14.61 | 8.3     | ng    | 328703   | 5                     | 14.81 | 790644 | 20.0 |
| Ethanol, 2-(tetra... | 15.42 | 6.3     | ng    | 251170   | 5                     | 14.81 | 790644 | 20.0 |
| Heneicosane          | 16.24 | 7.0     | ng    | 231360   | 6                     | 16.68 | 657567 | 20.0 |
| Oxirane, hexadecyl-  | 16.99 | 3.6     | ng    | 118683   | 6                     | 16.68 | 657567 | 20.0 |
| Tetracosane          | 17.29 | 6.2     | ng    | 204389   | 6                     | 16.68 | 657567 | 20.0 |
| unknown17.35         | 17.35 | 2.8     | ng    | 91463    | 6                     | 16.68 | 657567 | 20.0 |
| unknown17.59         | 17.59 | 6.7     | ng    | 218943   | 6                     | 16.68 | 657567 | 20.0 |
| unknown18.31         | 18.31 | 3.3     | ng    | 108907   | 6                     | 16.68 | 657567 | 20.0 |