

Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
 Data File : BF083103.D  
 Acq On : 21 Nov 2015 4:31  
 Operator : UM/IZ  
 Sample : G4486-01  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-13

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-BF111915.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         | 5.230       | 259           | 275         | 276          | rVV      | 279927         | 1798155       | 1.71%           | 0.654%        |
| 2         | 5.287       | 278           | 280         | 283          | rVV      | 1562820        | 2303110       | 2.18%           | 0.838%        |
| 3         | 5.687       | 299           | 315         | 317          | rBV      | 325066         | 1857150       | 1.76%           | 0.676%        |
| 4         | 6.350       | 371           | 373         | 376          | rBV2     | 1403796        | 1797478       | 1.70%           | 0.654%        |
| 5         | 6.476       | 381           | 384         | 385          | rVV      | 2769114        | 4039994       | 3.83%           | 1.470%        |
| 6         | 6.533       | 385           | 389         | 391          | rVB      | 532577         | 1258003       | 1.19%           | 0.458%        |
| 7         | 6.853       | 415           | 417         | 419          | rBV      | 2632919        | 2837727       | 2.69%           | 1.033%        |
| 8         | 7.139       | 439           | 442         | 444          | rVB      | 1926337        | 2530963       | 2.40%           | 0.921%        |
| 9         | 7.264       | 451           | 453         | 456          | rVB      | 2268454        | 2311481       | 2.19%           | 0.841%        |
| 10        | 7.813       | 490           | 501         | 503          | rBV      | 605482         | 2681223       | 2.54%           | 0.976%        |
| 11        | 7.984       | 503           | 516         | 518          | rVV      | 2694332        | 11606735      | 11.01%          | 4.225%        |
| 12        | 8.065       | 518           | 523         | 524          | rVV      | 4470435        | 14582758      | 13.83%          | 5.308%        |
| 13        | 8.282       | 524           | 542         | 543          | rVV2     | 14883330       | 105425445     | 100.00%         | 38.372%       |
| 14        | 8.316       | 543           | 545         | 547          | rVV      | 15963636       | 34936481      | 33.14%          | 12.716%       |
| 15        | 8.362       | 548           | 549         | 551          | rVV      | 10227820       | 7728056       | 7.33%           | 2.813%        |
| 16        | 8.522       | 559           | 563         | 564          | rVV      | 1965288        | 2466642       | 2.34%           | 0.898%        |
| 17        | 9.070       | 601           | 611         | 614          | rBV2     | 6597893        | 18159146      | 17.22%          | 6.610%        |
| 18        | 9.745       | 667           | 670         | 672          | rBV      | 1508268        | 1343868       | 1.27%           | 0.489%        |
| 19        | 10.065      | 687           | 698         | 700          | rBV      | 5559898        | 16891489      | 16.02%          | 6.148%        |
| 20        | 10.533      | 735           | 739         | 741          | rBV2     | 2667734        | 2584652       | 2.45%           | 0.941%        |
| 21        | 10.762      | 757           | 759         | 761          | rVV2     | 900708         | 1202503       | 1.14%           | 0.438%        |
| 22        | 10.956      | 772           | 776         | 777          | rVV2     | 3666997        | 6777975       | 6.43%           | 2.467%        |
| 23        | 11.219      | 797           | 799         | 800          | rVV      | 1197197        | 1112466       | 1.06%           | 0.405%        |
| 24        | 11.253      | 800           | 802         | 808          | rVB      | 1917128        | 2386539       | 2.26%           | 0.869%        |
| 25        | 11.825      | 845           | 852         | 854          | rBV      | 7186857        | 15044954      | 14.27%          | 5.476%        |
| 26        | 12.511      | 907           | 912         | 916          | rBV      | 893579         | 2638457       | 2.50%           | 0.960%        |
| 27        | 12.579      | 916           | 918         | 919          | rBV      | 851858         | 1079032       | 1.02%           | 0.393%        |
| 28        | 12.808      | 936           | 938         | 941          | rVB      | 3107716        | 2885388       | 2.74%           | 1.050%        |
| 29        | 12.922      | 946           | 948         | 950          | rBV2     | 794936         | 1236551       | 1.17%           | 0.450%        |
| 30        | 13.848      | 1027          | 1029        | 1033         | rBV      | 998151         | 1237802       | 1.17%           | 0.451%        |

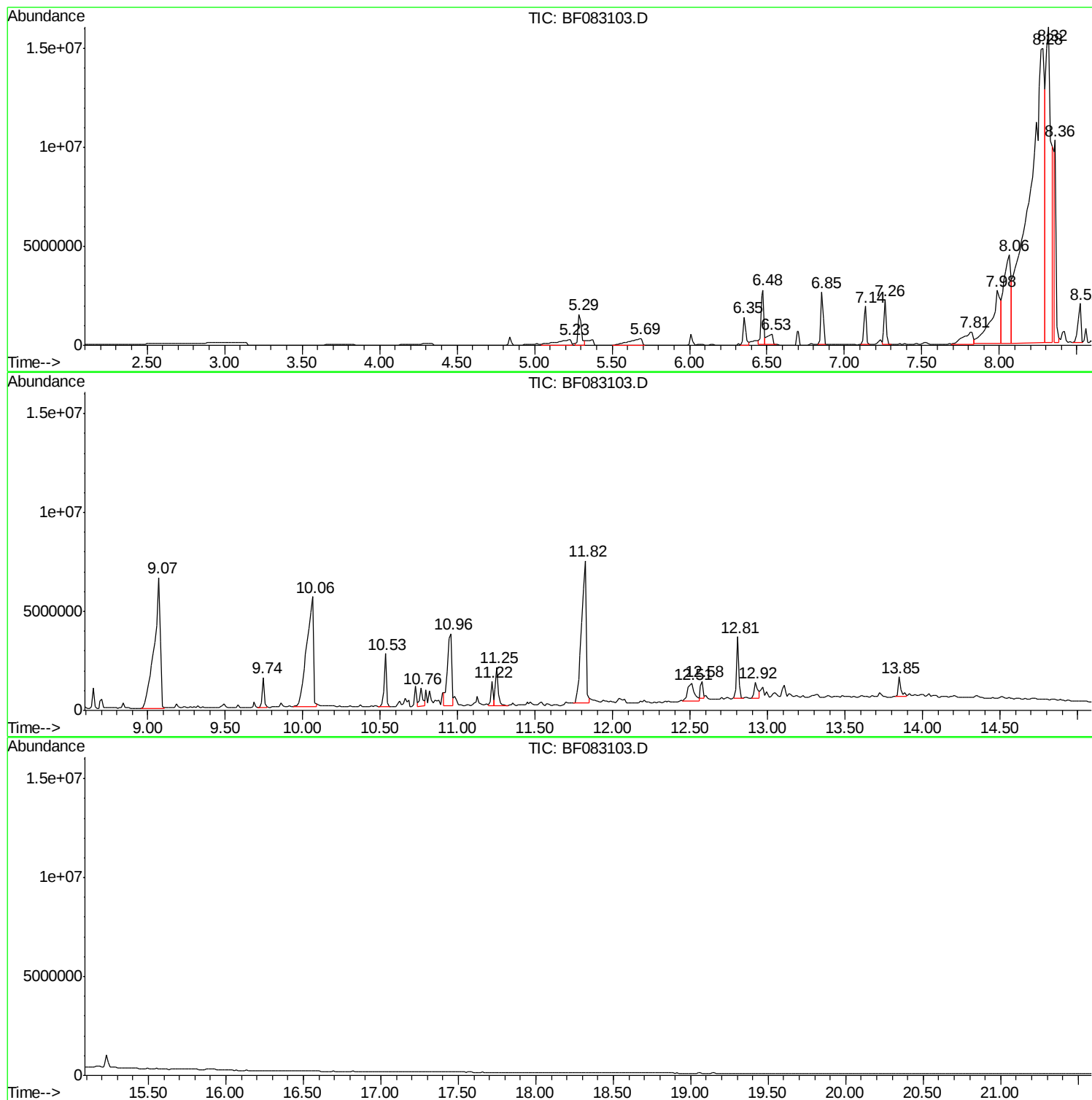
Sum of corrected areas: 274742223

Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

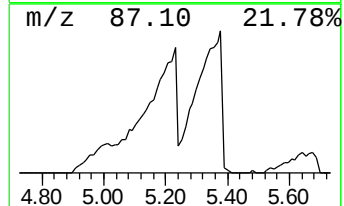
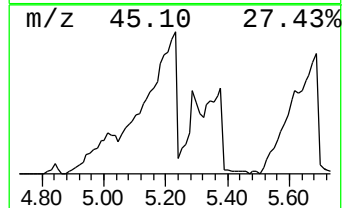
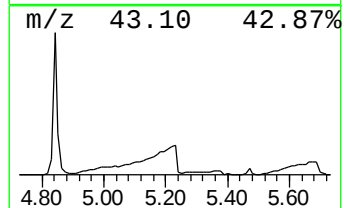
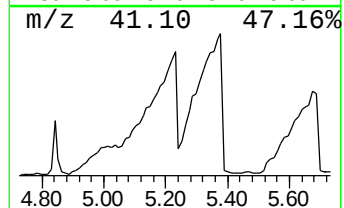
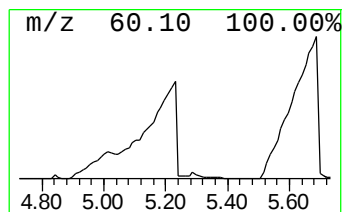
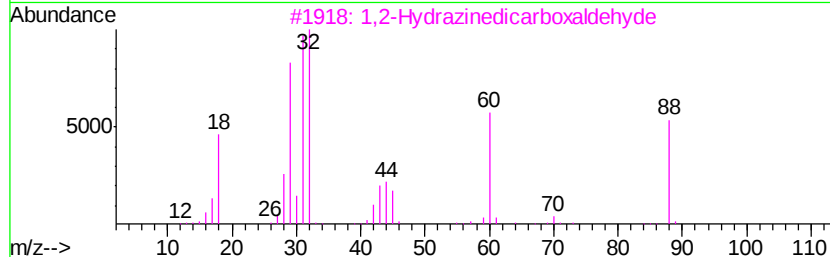
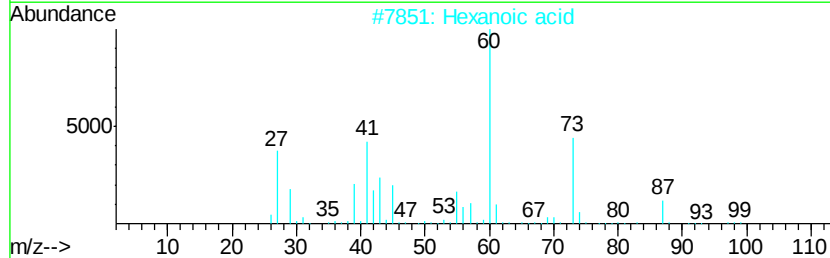
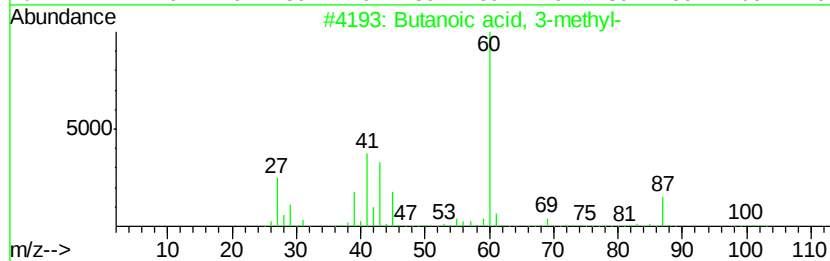
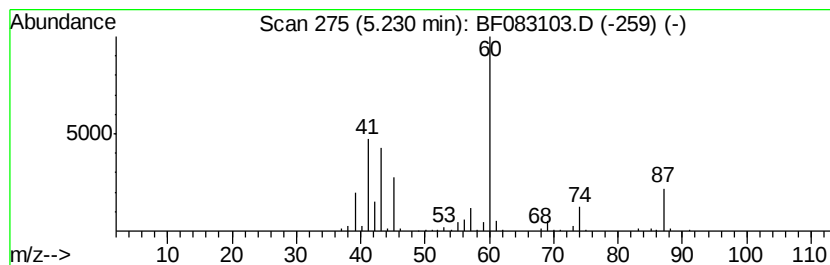
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 15

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.23 | 12.67 ng | 1798160 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | Tentative ID                  | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------|-----|----------|-------------|------|
| 1    | 5  | Butanoic acid, 3-methyl-      | 102 | C5H10O2  | 000503-74-2 | 59   |
| 2    |    | Hexanoic acid                 | 116 | C6H12O2  | 000142-62-1 | 53   |
| 3    |    | 1,2-Hydrazinedicarboxaldehyde | 88  | C2H4N2O2 | 000628-36-4 | 36   |
| 4    |    | Pentanoic acid                | 102 | C5H10O2  | 000109-52-4 | 33   |
| 5    |    | 1-Pentanamine                 | 87  | C5H13N   | 000110-58-7 | 9    |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

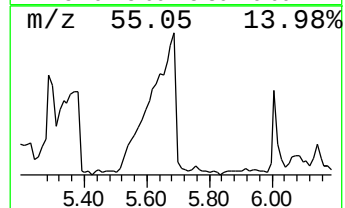
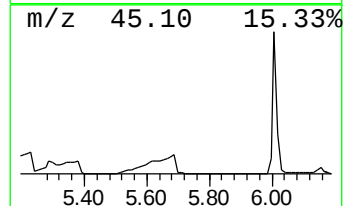
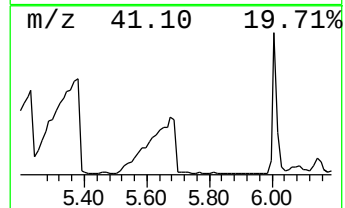
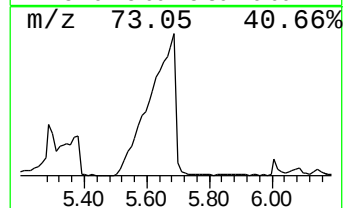
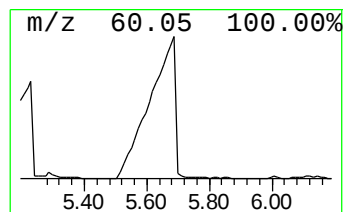
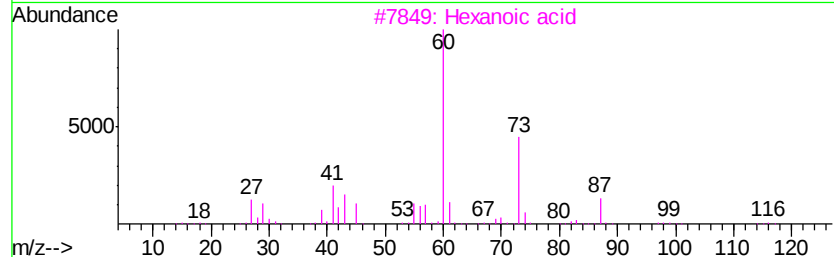
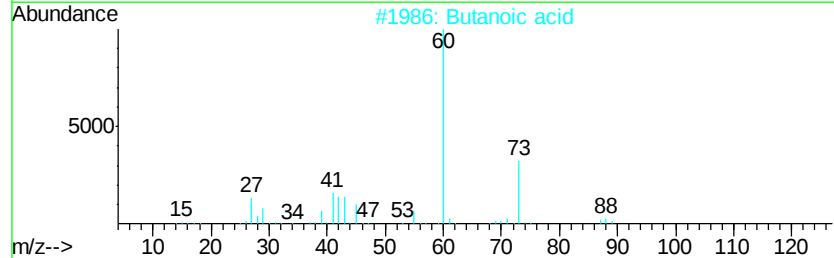
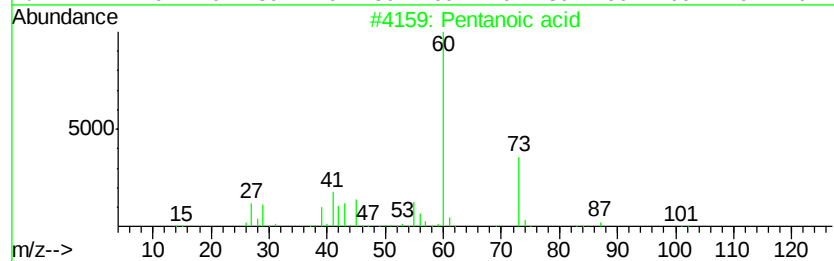
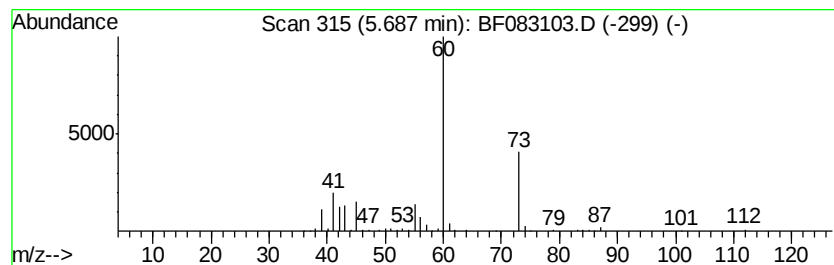
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Pentanoic acid Concentration Rank 14

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.69 | 13.09 ng | 1857150 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentanoic acid             | 102 | C5H10O2 | 000109-52-4 | 90   |
| 2    |    | Butanoic acid              | 88  | C4H8O2  | 000107-92-6 | 78   |
| 3    |    | Hexanoic acid              | 116 | C6H12O2 | 000142-62-1 | 64   |
| 4    |    | Propanedioic acid, propyl- | 146 | C6H10O4 | 000616-62-6 | 53   |
| 5    |    | Butanoic acid, 3-methyl-   | 102 | C5H10O2 | 000503-74-2 | 50   |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

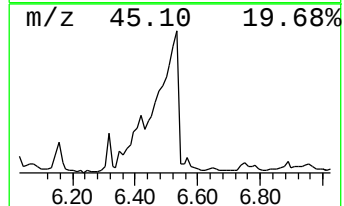
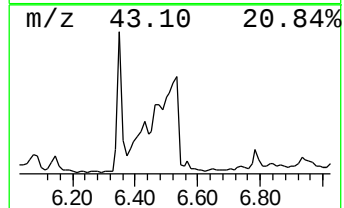
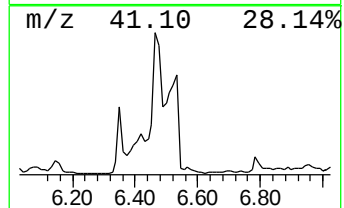
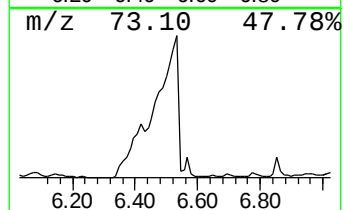
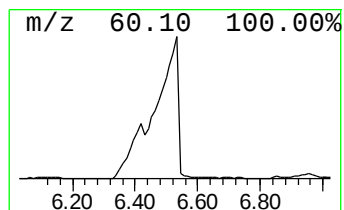
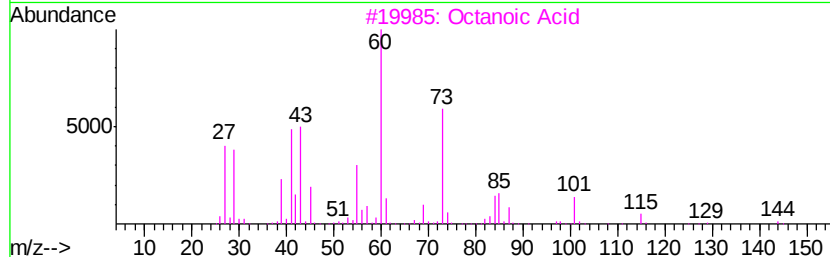
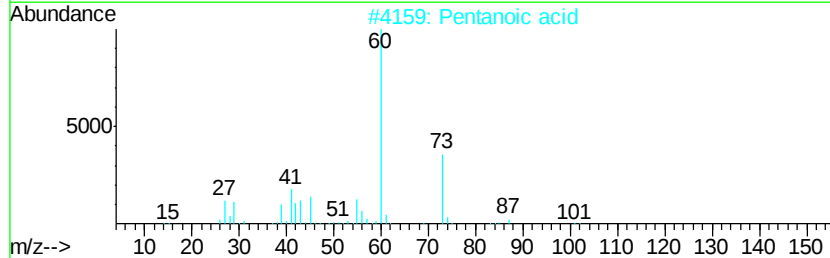
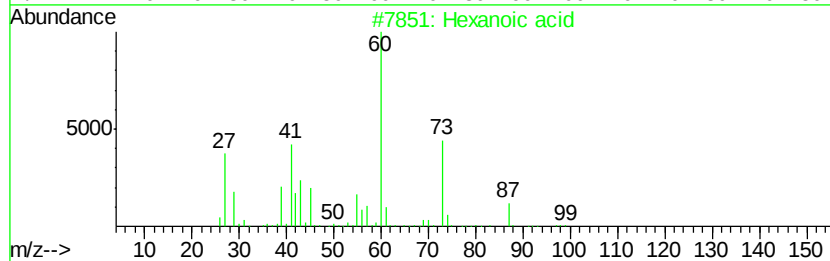
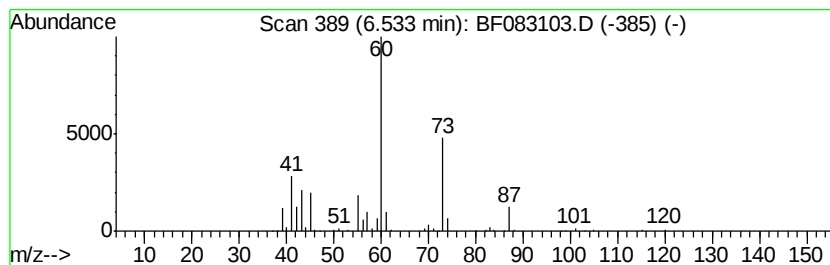
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Hexanoic acid Concentration Rank 16

| R.T. | EstConc | Area    | Relative to ISTD       | R.T. |
|------|---------|---------|------------------------|------|
| 6.53 | 8.87 ng | 1258000 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexanoic acid            | 116 | C6H12O2 | 000142-62-1 | 78   |
| 2    |    |   | Pentanoic acid           | 102 | C5H10O2 | 000109-52-4 | 72   |
| 3    |    |   | Octanoic Acid            | 144 | C8H16O2 | 000124-07-2 | 64   |
| 4    |    |   | Heptanoic acid           | 130 | C7H14O2 | 000111-14-8 | 64   |
| 5    |    |   | Butanoic acid, 3-methyl- | 102 | C5H10O2 | 000503-74-2 | 45   |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

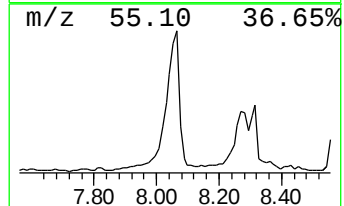
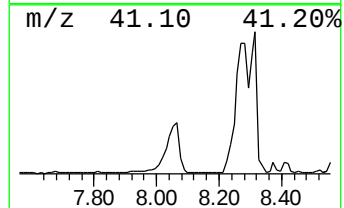
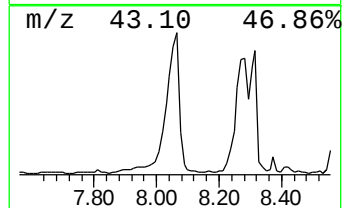
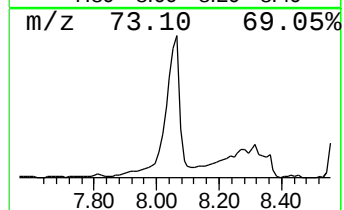
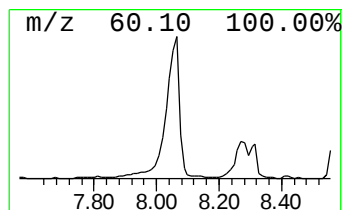
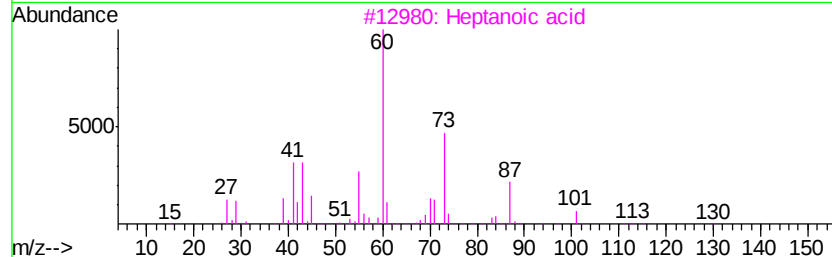
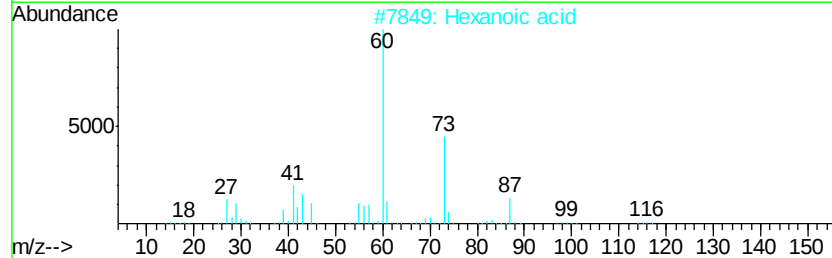
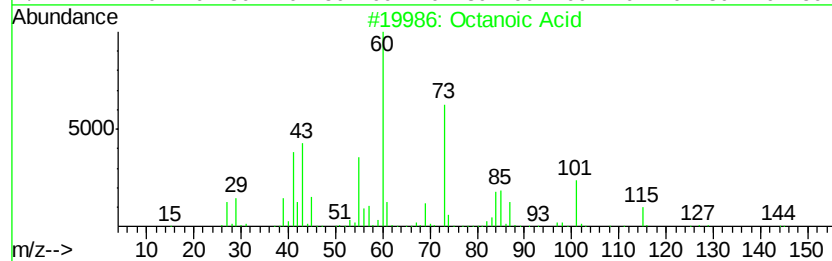
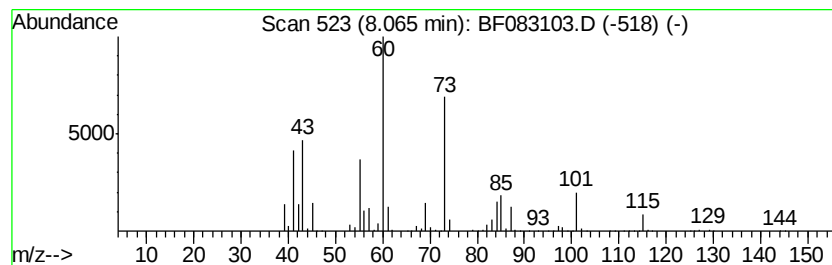
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Octanoic Acid Concentration Rank 8

| R.T. | EstConc  | Area     | Relative to ISTD | R.T. |
|------|----------|----------|------------------|------|
| 8.06 | 25.13 ng | 14582800 | Naphthalene-d8   | 7.98 |

| Hit# of | 5                          | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|----------------------------|--------------|-----|---------|-------------|------|
| 1       | Octanoic Acid              |              | 144 | C8H16O2 | 000124-07-2 | 94   |
| 2       | Hexanoic acid              |              | 116 | C6H12O2 | 000142-62-1 | 64   |
| 3       | Heptanoic acid             |              | 130 | C7H14O2 | 000111-14-8 | 40   |
| 4       | Butanoic acid, 3-methyl-   |              | 102 | C5H10O2 | 000503-74-2 | 40   |
| 5       | Propanedioic acid, propyl- |              | 146 | C6H10O4 | 000616-62-6 | 40   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

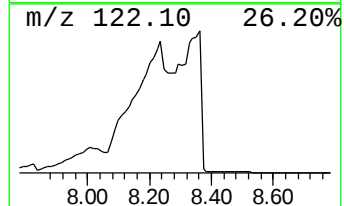
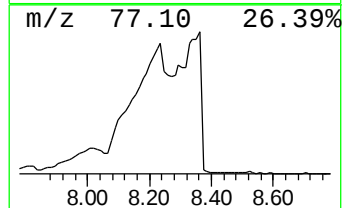
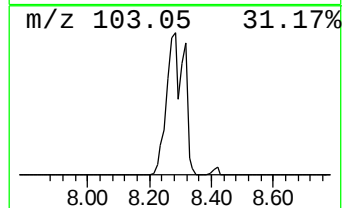
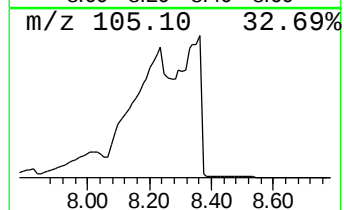
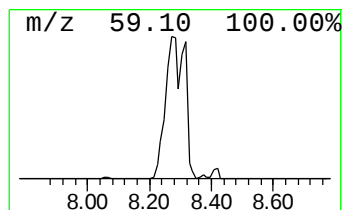
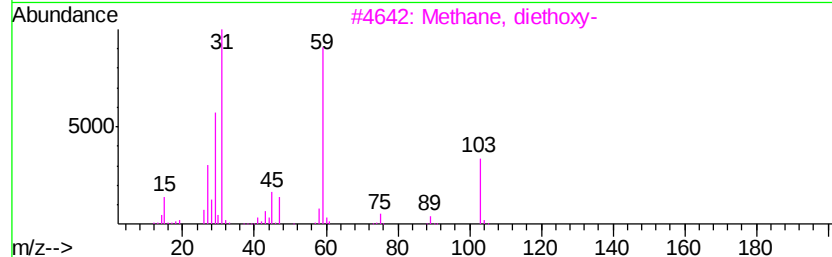
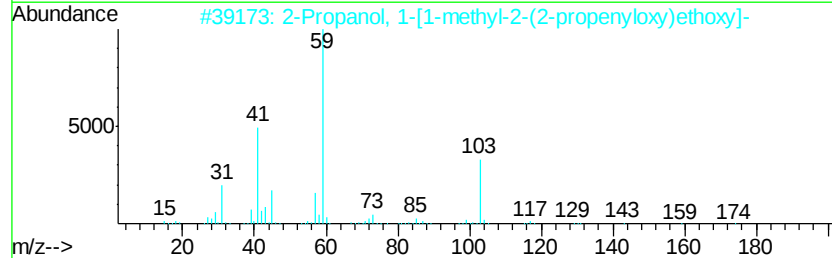
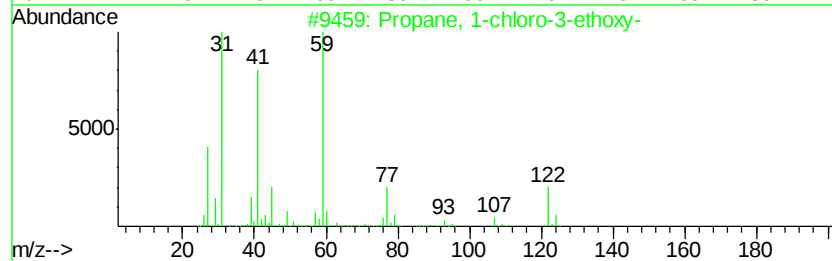
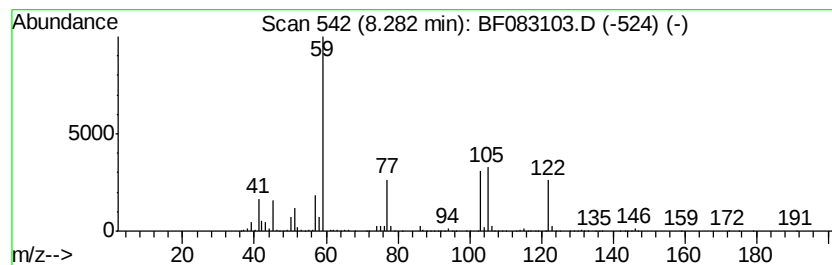
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Propane, 1-chloro-3-ethoxy- Concentration Rank 3

| R.T. | EstConc   | Area      | Relative to ISTD | R.T. |
|------|-----------|-----------|------------------|------|
| 8.28 | 181.66 ng | 105425000 | Naphthalene-d8   | 7.98 |

| Hit# of | 5                                   | Tentative ID | MW       | MolForm     | CAS# | Qual |
|---------|-------------------------------------|--------------|----------|-------------|------|------|
| 1       | Propane, 1-chloro-3-ethoxy-         | 122          | C5H11ClO | 036865-38-0 | 52   |      |
| 2       | 2-Propanol, 1-[1-methyl-2-(2-pro... | 174          | C9H18O3  | 055956-25-7 | 47   |      |
| 3       | Methane, diethoxy-                  | 104          | C5H12O2  | 000462-95-3 | 43   |      |
| 4       | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134          | C6H14O3  | 000106-62-7 | 43   |      |
| 5       | Ethanol, 1-methoxy-, benzoate       | 180          | C10H12O3 | 051835-44-0 | 36   |      |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

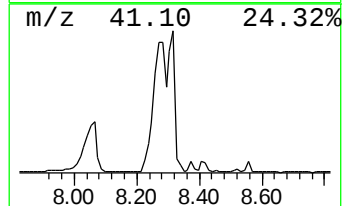
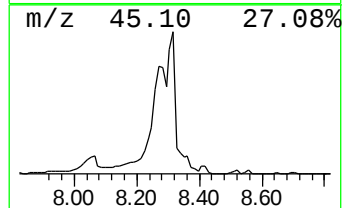
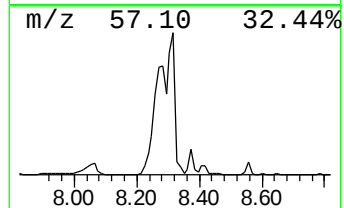
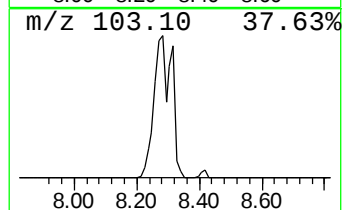
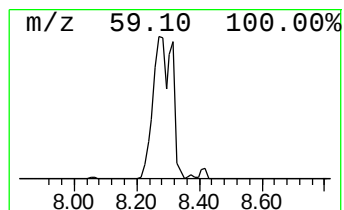
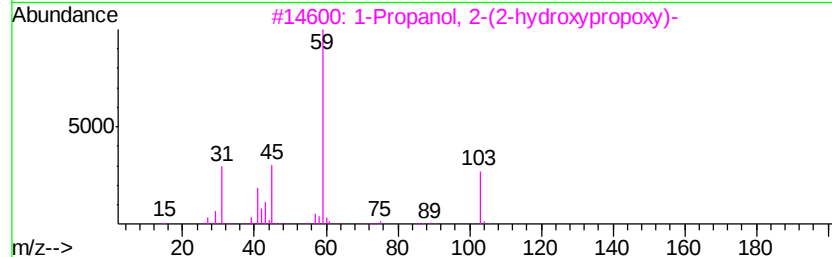
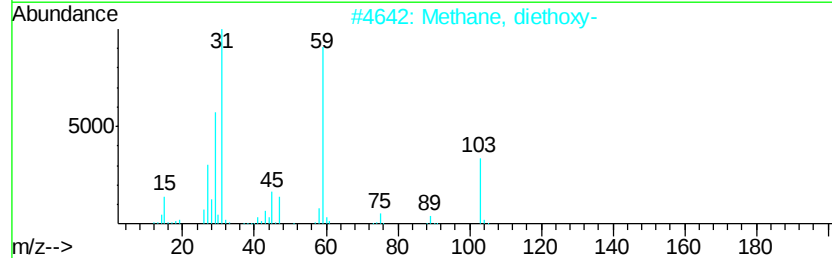
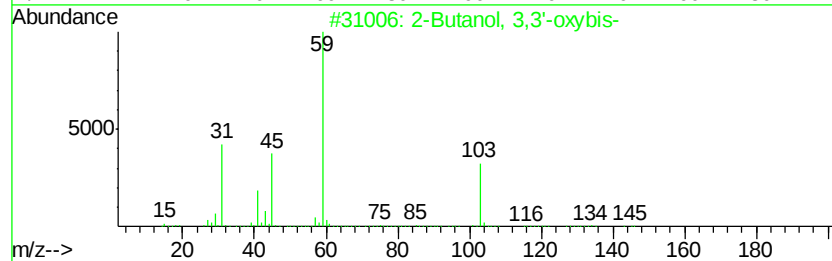
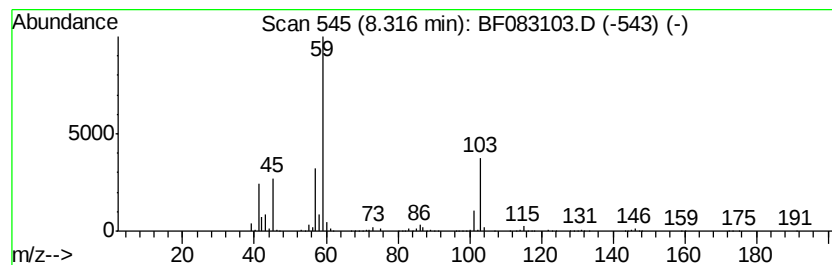
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 2-Butanol, 3,3'-oxybis- Concentration Rank 5

| R.T. | EstConc  | Area     | Relative to ISTD | R.T. |
|------|----------|----------|------------------|------|
| 8.32 | 60.20 ng | 34936500 | Naphthalene-d8   | 7.98 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Butanol, 3,3'-oxybis-           | 162 | C8H18O3 | 054305-61-2 | 72   |
| 2    |    | Methane, diethoxy-                | 104 | C5H12O2 | 000462-95-3 | 72   |
| 3    |    | 1-Propanol, 2-(2-hydroxypropoxy)- | 134 | C6H14O3 | 000106-62-7 | 64   |
| 4    |    | 1-Propanol, 2,2'-oxybis-          | 134 | C6H14O3 | 000108-61-2 | 64   |
| 5    |    | Dipropylene glycol                | 134 | C6H14O3 | 025265-71-8 | 56   |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

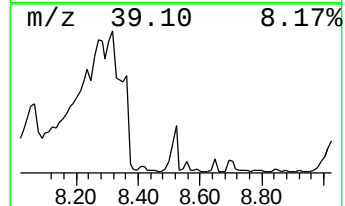
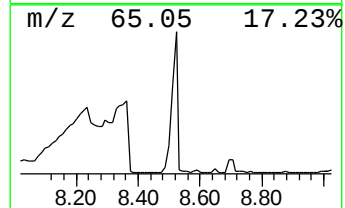
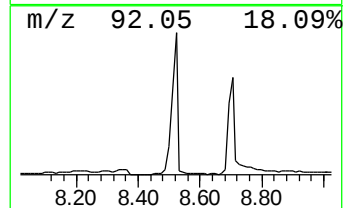
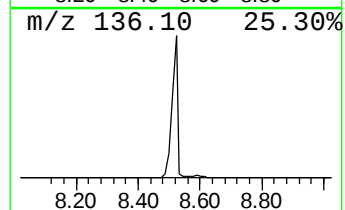
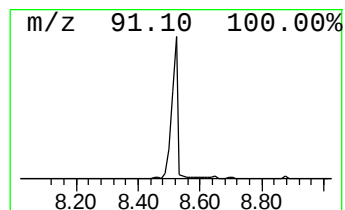
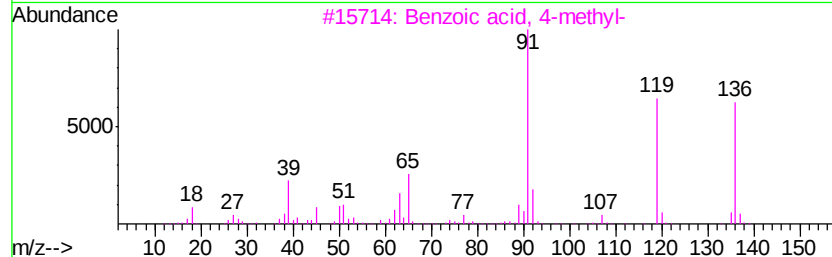
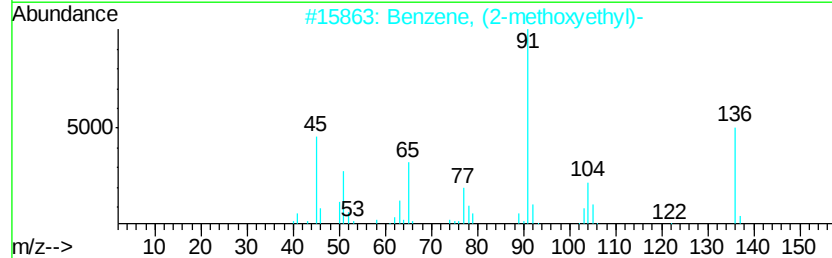
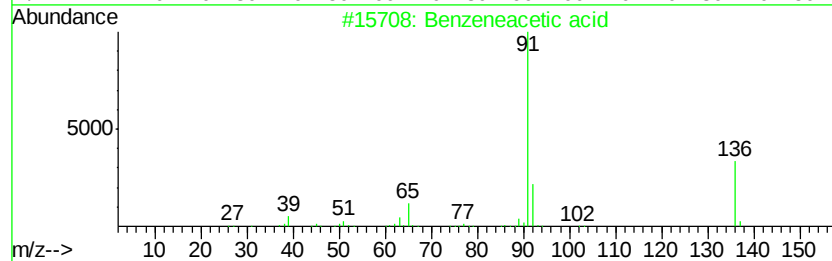
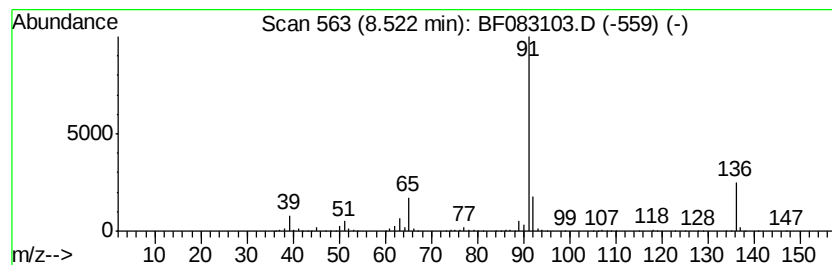
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzeneacetic acid Concentration Rank 17

| R.T. | EstConc | Area    | Relative to ISTD | R.T. |
|------|---------|---------|------------------|------|
| 8.52 | 4.25 ng | 2466640 | Naphthalene-d8   | 7.98 |

| Hit# | of | Tentative ID                 | MW  | MolForm    | CAS#        | Qual |
|------|----|------------------------------|-----|------------|-------------|------|
| 1    | 5  | Benzeneacetic acid           | 136 | C8H8O2     | 000103-82-2 | 87   |
| 2    |    | Benzene, (2-methoxyethyl)-   | 136 | C9H12O     | 003558-60-9 | 72   |
| 3    |    | Benzoic acid, 4-methyl-      | 136 | C8H8O2     | 000099-94-5 | 59   |
| 4    |    | Propanedioic acid, phenyl-   | 180 | C9H8O4     | 002613-89-0 | 59   |
| 5    |    | Benzyl 2-chloroethyl sulfone | 218 | C9H11ClO2S | 066998-67-2 | 50   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

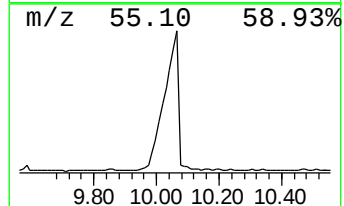
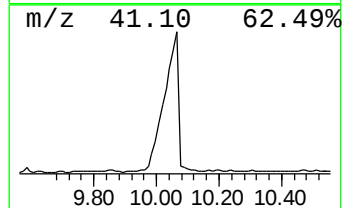
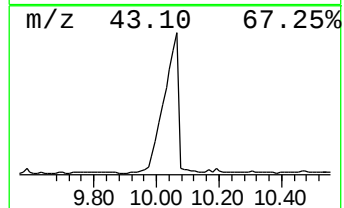
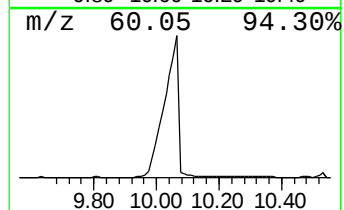
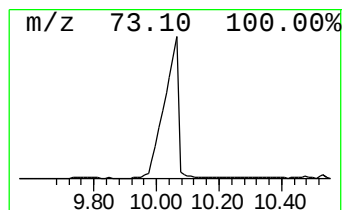
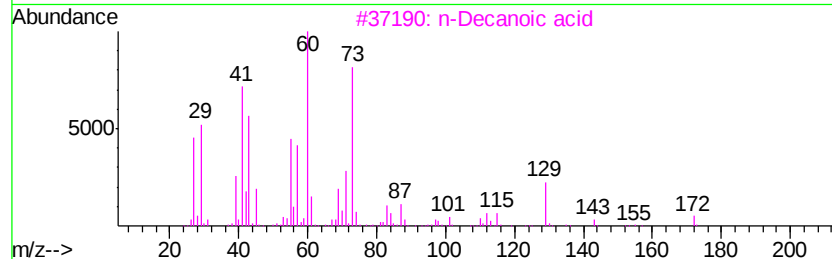
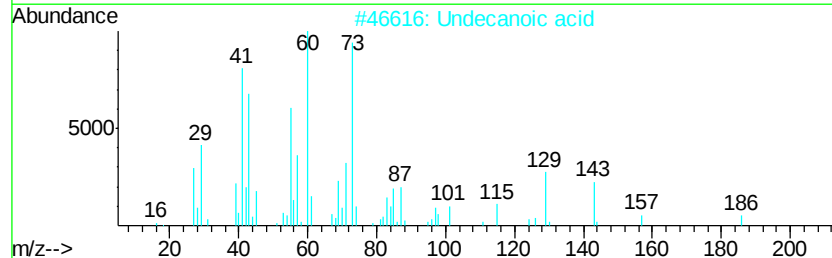
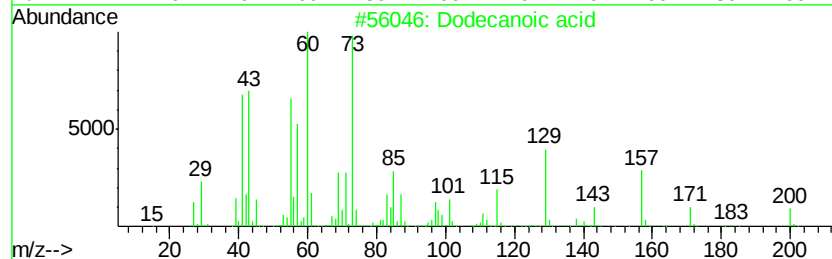
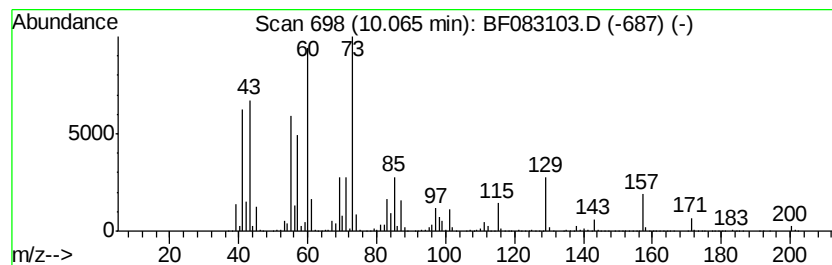
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Dodecanoic acid Concentration Rank 2

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T. |
|-------|-----------|----------|------------------|------|
| 10.06 | 251.39 ng | 16891500 | Acenaphthene-d10 | 9.74 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Dodecanoic acid                     | 200 | C12H24O2 | 000143-07-7 | 98   |
| 2    |    |   | Undecanoic acid                     | 186 | C11H22O2 | 000112-37-8 | 86   |
| 3    |    |   | n-Decanoic acid                     | 172 | C10H20O2 | 000334-48-5 | 81   |
| 4    |    |   | Nonanoic acid                       | 158 | C9H18O2  | 000112-05-0 | 58   |
| 5    |    |   | Ethanone, 1-(4,5-dihydro-2-thiaz... | 129 | C5H7NOS  | 029926-41-8 | 43   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

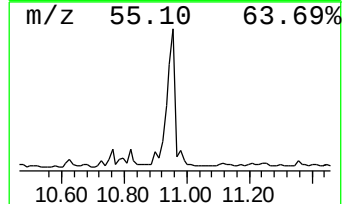
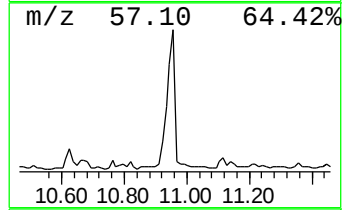
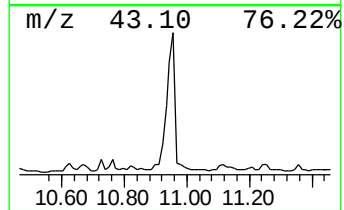
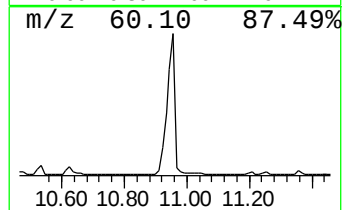
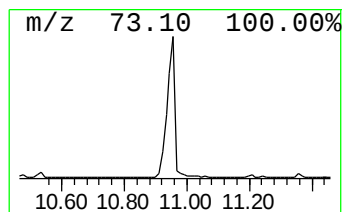
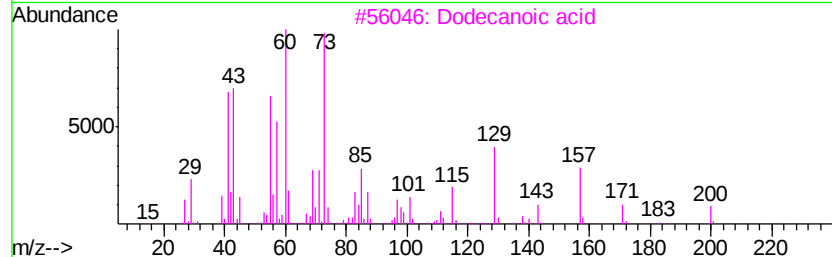
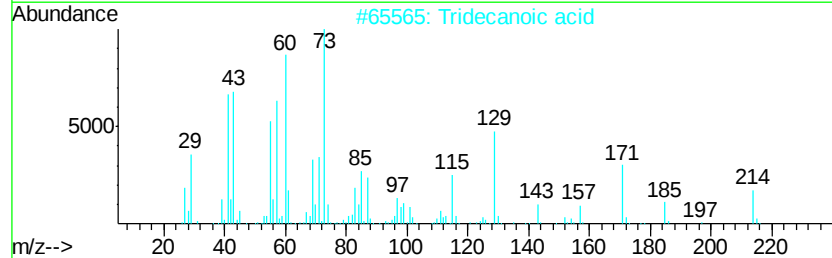
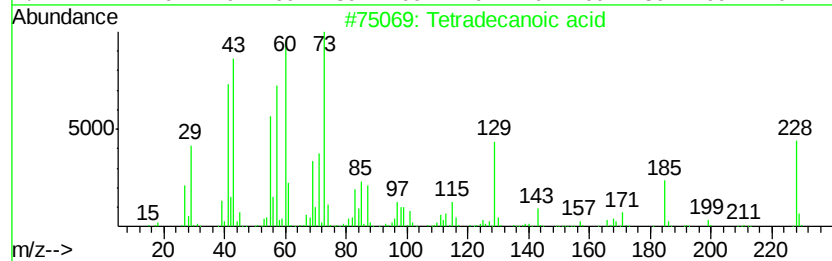
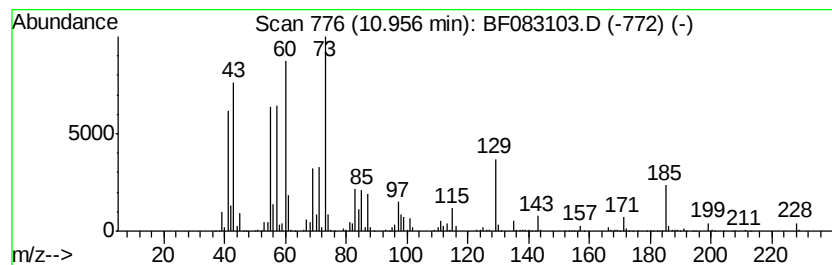
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Tetradecanoic acid Concentration Rank 4

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 10.96 | 121.86 ng | 6777980 | Phenanthrene-d10 | 11.22 |

| Hit# of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|---------|---|--------------------|-----|----------|-------------|------|
| 1       |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 95   |
| 2       |   | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 80   |
| 3       |   | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 64   |
| 4       |   | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 53   |
| 5       |   | Butanoic acid      | 88  | C4H8O2   | 000107-92-6 | 27   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

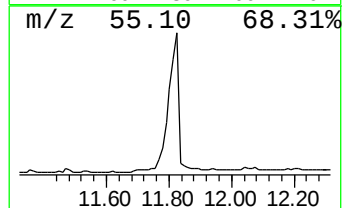
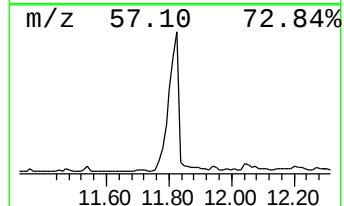
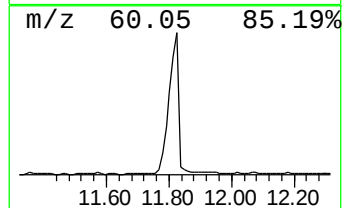
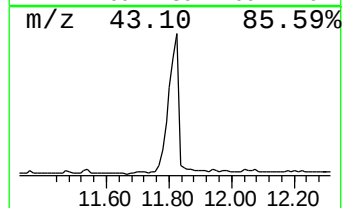
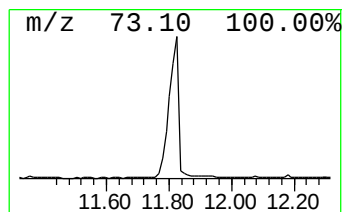
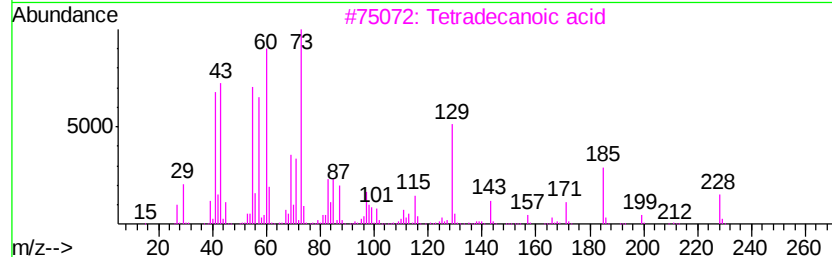
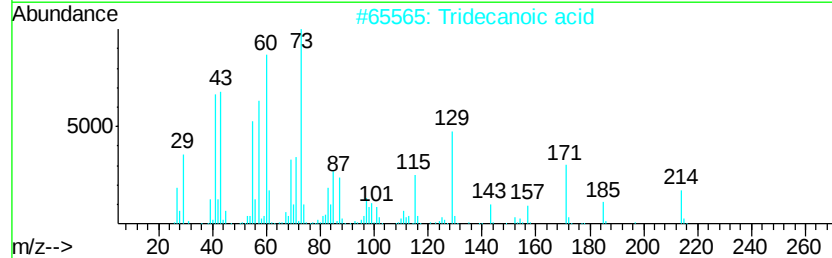
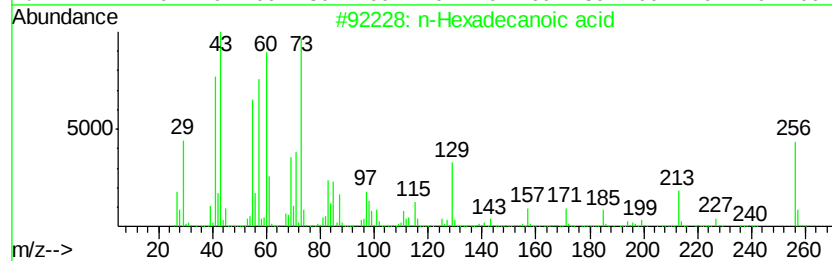
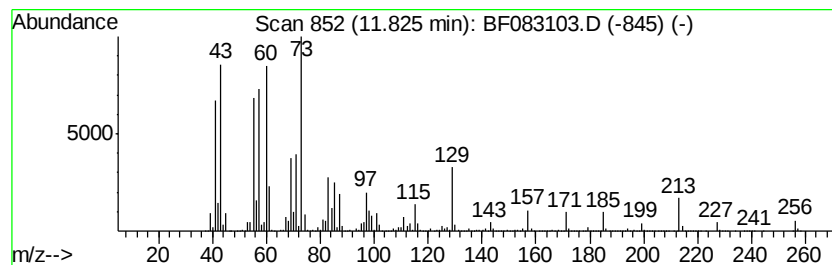
Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 n-Hexadecanoic acid Concentration Rank 1

| R.T.    | EstConc   | Area                                | Relative to ISTD | R.T.      |              |      |
|---------|-----------|-------------------------------------|------------------|-----------|--------------|------|
| 11.82   | 270.48 ng | 15045000                            | Phenanthrene-d10 | 11.22     |              |      |
| Hit# of | 5         | Tentative ID                        | MW               | MolForm   | CAS#         | Qual |
| 1       |           | n-Hexadecanoic acid                 | 256              | C16H32O2  | 000057-10-3  | 98   |
| 2       |           | Tridecanoic acid                    | 214              | C13H26O2  | 000638-53-9  | 95   |
| 3       |           | Tetradecanoic acid                  | 228              | C14H28O2  | 000544-63-8  | 87   |
| 4       |           | Estra-1,3,5(10)-trien-17.beta.-ol   | 256              | C18H24O   | 002529-64-8  | 25   |
| 5       |           | 6-Methyl(pentamethylene)silyloxy... | 326              | C20H42OSi | 1000278-79-4 | 18   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

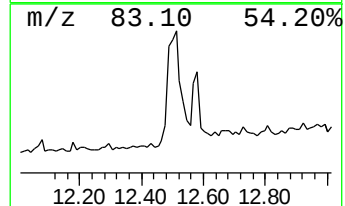
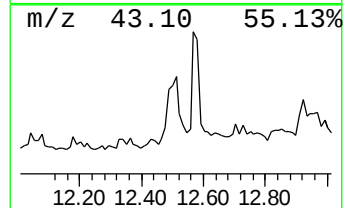
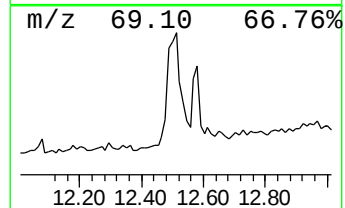
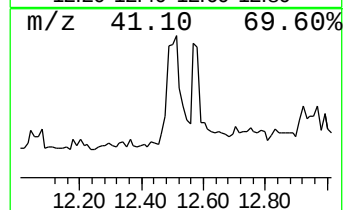
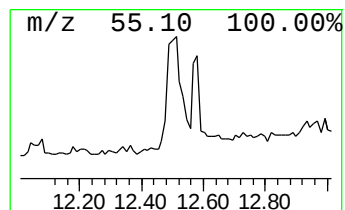
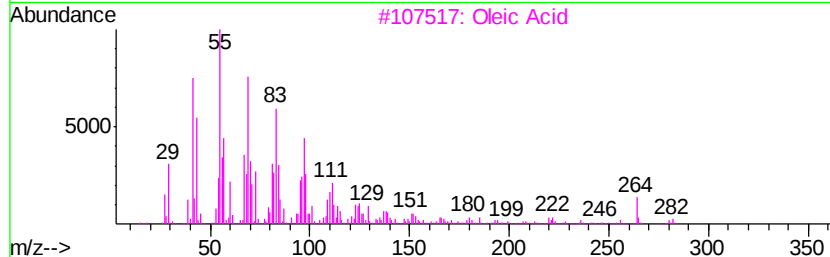
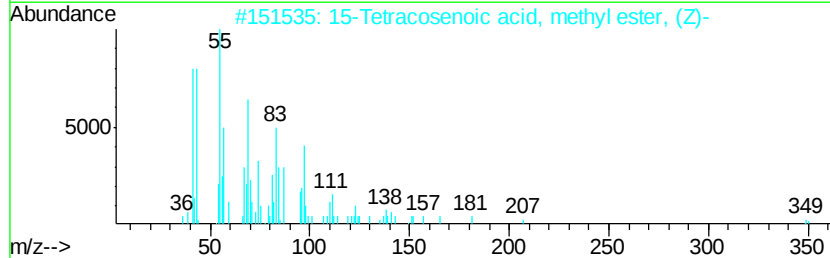
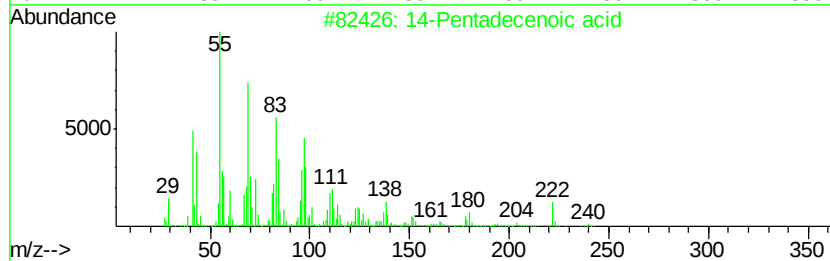
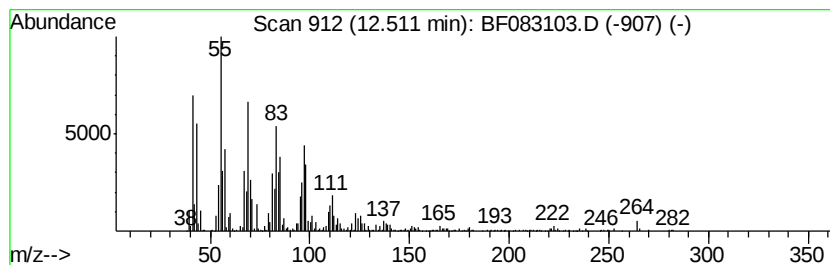
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 14-Pentadecenoic acid Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.51 | 47.43 ng | 2638460 | Phenanthrene-d10 | 11.22 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 14-Pentadecenoic acid               | 240 | C15H28O2 | 017351-34-7 | 83   |
| 2    |    | 15-Tetracosenoic acid, methyl es... | 380 | C25H48O2 | 002733-88-2 | 81   |
| 3    |    | Oleic Acid                          | 282 | C18H34O2 | 000112-80-1 | 80   |
| 4    |    | 6-Octadecenoic acid, (Z)-           | 282 | C18H34O2 | 000593-39-5 | 80   |
| 5    |    | Cyclopropane, 1-(1,2-dimethylpro... | 252 | C18H36   | 041977-42-8 | 64   |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

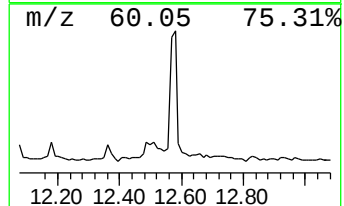
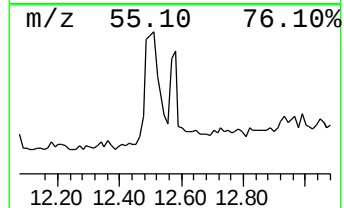
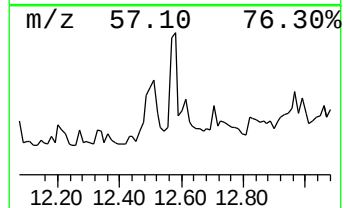
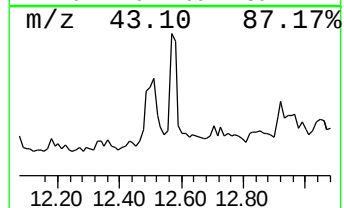
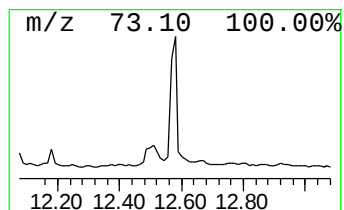
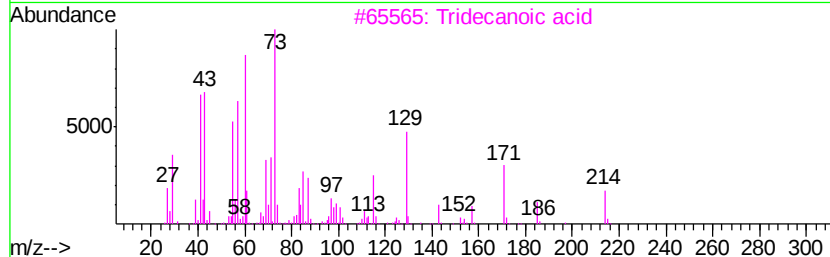
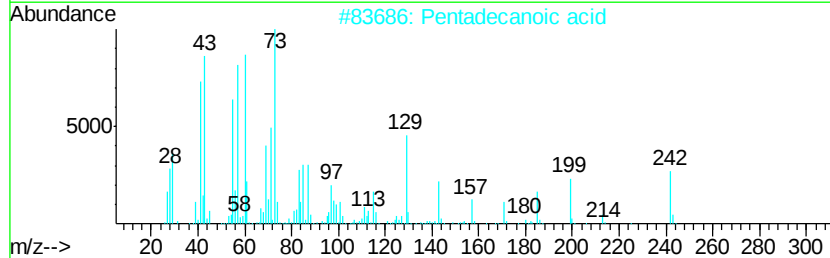
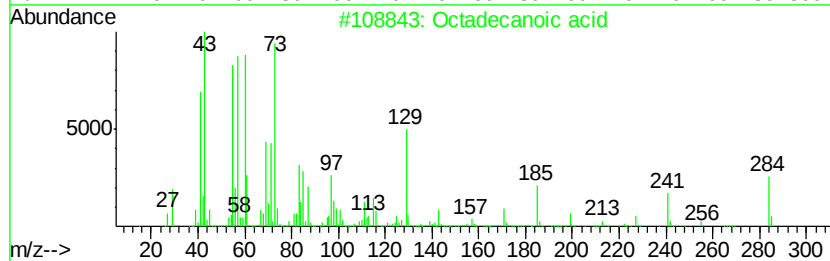
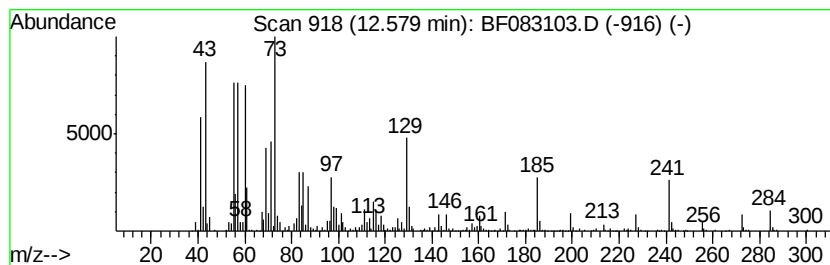
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Octadecanoic acid Concentration Rank 12

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.58 | 17.43 ng | 1079030 | Chrysene-d12     | 13.85 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid                  | 284 | C18H36O2 | 000057-11-4 | 94   |
| 2    |    | Pentadecanoic acid                 | 242 | C15H30O2 | 001002-84-2 | 90   |
| 3    |    | Tridecanoic acid                   | 214 | C13H26O2 | 000638-53-9 | 72   |
| 4    |    | Tetradecanoic acid                 | 228 | C14H28O2 | 000544-63-8 | 64   |
| 5    |    | Octadecanoic acid, 2-(2-hydroxy... | 372 | C22H44O4 | 000106-11-6 | 64   |



Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\  
Data File : BF083103.D  
Acq On : 21 Nov 2015 4:31  
Operator : UM/IZ  
Sample : G4486-01  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

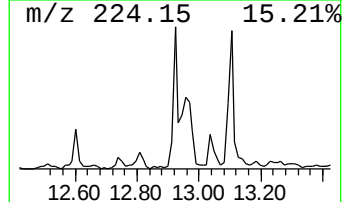
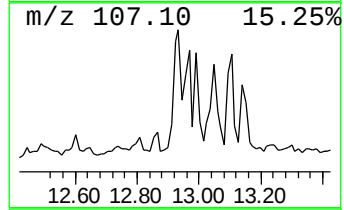
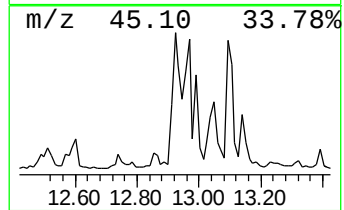
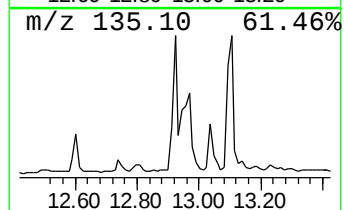
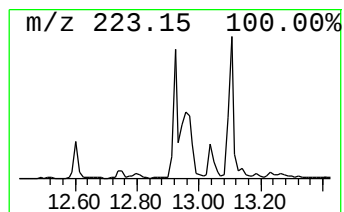
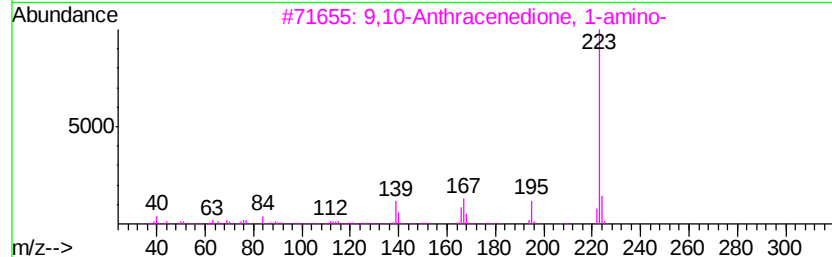
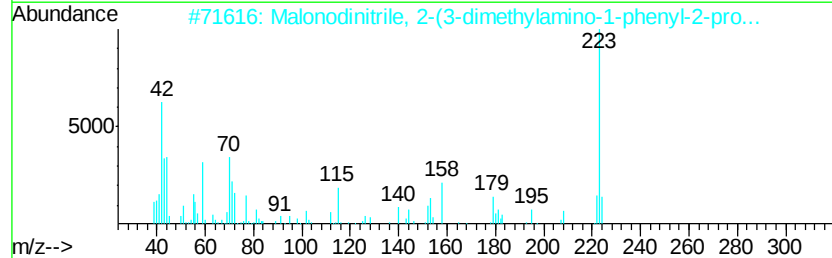
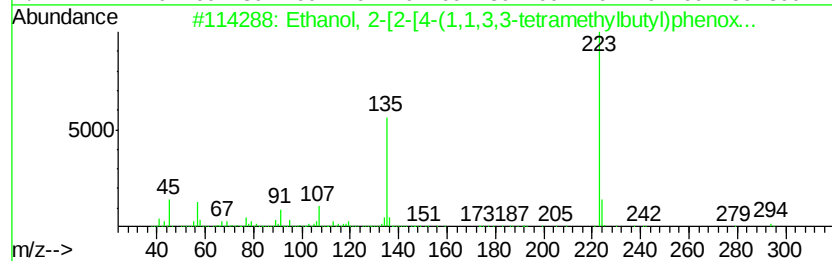
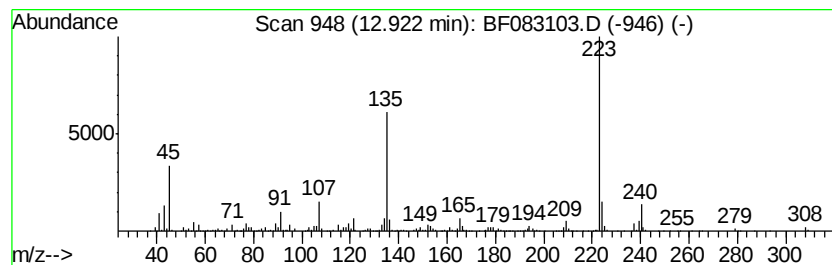
Instrument :  
BNA\_F  
ClientSampleId :  
MW-13

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 11

| R.T.    | EstConc  | Area                                | Relative to ISTD | R.T.       |              |      |
|---------|----------|-------------------------------------|------------------|------------|--------------|------|
| 12.92   | 19.98 ng | 1236550                             | Chrysene-d12     | 13.85      |              |      |
| Hit# of | 5        | Tentative ID                        | MW               | MolForm    | CAS#         | Qual |
| 1       |          | Ethanol, 2-[2-[4-(1,1,3,3-tetram... | 294              | C18H30O3   | 002315-61-9  | 72   |
| 2       |          | Malonodinitrile, 2-(3-dimethylam... | 223              | C14H13N3   | 065996-13-6  | 46   |
| 3       |          | 9,10-Anthracenedione, 1-amino-      | 223              | C14H9NO2   | 000082-45-1  | 35   |
| 4       |          | 2-[p-Aminobenzyl]benzimidazole      | 223              | C14H13N3   | 099206-51-6  | 35   |
| 5       |          | (4-Oxo-2-phenyl-1,2,3,4-tetrahyd... | 282              | C16H14N2O3 | 1000190-14-1 | 25   |



Data Path : V:\HPCHEM1\BNA\_F\DATA\BF112015\  
 Data File : BF083103.D  
 Acq On : 21 Nov 2015 4:31  
 Operator : UM/IZ  
 Sample : G4486-01  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-13

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF111915.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 |
| Butanoic acid, 3-... | 5.23  | 12.7    | ng    | 1798160   | 1                     | 6.70  | 2837730  | 20.0 |
| Pentanoic acid       | 5.69  | 13.1    | ng    | 1857150   | 1                     | 6.70  | 2837730  | 20.0 |
| Hexanoic acid        | 6.53  | 8.9     | ng    | 1258000   | 1                     | 6.70  | 2837730  | 20.0 |
| Octanoic Acid        | 8.06  | 25.1    | ng    | 14582800  | 2                     | 7.98  | 11606700 | 20.0 |
| Propane, 1-chloro... | 8.28  | 181.7   | ng    | 105425000 | 2                     | 7.98  | 11606700 | 20.0 |
| 2-Butanol, 3,3'-o... | 8.32  | 60.2    | ng    | 34936500  | 2                     | 7.98  | 11606700 | 20.0 |
| Benzeneacetic acid   | 8.52  | 4.3     | ng    | 2466640   | 2                     | 7.98  | 11606700 | 20.0 |
| Dodecanoic acid      | 10.06 | 251.4   | ng    | 16891500  | 3                     | 9.74  | 1343870  | 20.0 |
| Tetradecanoic acid   | 10.96 | 121.9   | ng    | 6777980   | 4                     | 11.22 | 1112470  | 20.0 |
| n-Hexadecanoic acid  | 11.82 | 270.5   | ng    | 15045000  | 4                     | 11.22 | 1112470  | 20.0 |
| 14-Pentadecenoic ... | 12.51 | 47.4    | ng    | 2638460   | 4                     | 11.22 | 1112470  | 20.0 |
| Octadecanoic acid    | 12.58 | 17.4    | ng    | 1079030   | 5                     | 13.85 | 1237800  | 20.0 |
| Ethanol, 2-[2-[4-... | 12.92 | 20.0    | ng    | 1236550   | 5                     | 13.85 | 1237800  | 20.0 |