

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031915\  
 Data File : BF077858.D  
 Acq On : 20 Mar 2015 6:13  
 Operator : TP/IZ  
 Sample : G1580-01  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DGA-DRH-01

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-BF031115.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.218       | 27            | 29          | 34           | rVB2     | 48847          | 74583         | 1.47%           | 0.386%        |
| 2         | 1.310       | 34            | 37          | 40           | rVB      | 4654088        | 5061503       | 100.00%         | 26.222%       |
| 3         | 1.424       | 45            | 47          | 60           | rVB      | 88515          | 224526        | 4.44%           | 1.163%        |
| 4         | 1.595       | 60            | 62          | 65           | rVB      | 342854         | 383777        | 7.58%           | 1.988%        |
| 5         | 1.664       | 66            | 68          | 77           | rVB      | 156620         | 260690        | 5.15%           | 1.351%        |
| 6         | 1.778       | 77            | 78          | 102          | rVB      | 865231         | 1560275       | 30.83%          | 8.083%        |
| 7         | 4.910       | 350           | 352         | 358          | rVB      | 130169         | 158831        | 3.14%           | 0.823%        |
| 8         | 5.447       | 396           | 399         | 401          | rBV      | 973410         | 1097515       | 21.68%          | 5.686%        |
| 9         | 6.727       | 509           | 511         | 514          | rBV      | 1065284        | 1047030       | 20.69%          | 5.424%        |
| 10        | 6.864       | 521           | 523         | 526          | rBV      | 1306505        | 1192956       | 23.57%          | 6.180%        |
| 11        | 7.150       | 546           | 548         | 551          | rBV      | 278821         | 245206        | 4.84%           | 1.270%        |
| 12        | 7.333       | 562           | 564         | 567          | rBV      | 914777         | 926168        | 18.30%          | 4.798%        |
| 13        | 7.847       | 606           | 609         | 611          | rBV      | 811262         | 726626        | 14.36%          | 3.764%        |
| 14        | 8.727       | 684           | 686         | 688          | rBV      | 367472         | 333329        | 6.59%           | 1.727%        |
| 15        | 10.076      | 802           | 804         | 806          | rBV      | 1663029        | 1476250       | 29.17%          | 7.648%        |
| 16        | 10.899      | 874           | 876         | 878          | rBV      | 406180         | 385970        | 7.63%           | 2.000%        |
| 17        | 11.882      | 959           | 962         | 964          | rBV      | 676250         | 875407        | 17.30%          | 4.535%        |
| 18        | 12.076      | 977           | 979         | 983          | rBV      | 65379          | 86777         | 1.71%           | 0.450%        |
| 19        | 12.739      | 1034          | 1037        | 1039         | rBV      | 386303         | 407629        | 8.05%           | 2.112%        |
| 20        | 14.728      | 1209          | 1211        | 1214         | rBV      | 1314165        | 1702909       | 33.64%          | 8.822%        |
| 21        | 15.905      | 1310          | 1314        | 1322         | rBV2     | 79983          | 153296        | 3.03%           | 0.794%        |
| 22        | 16.020      | 1322          | 1324        | 1327         | rBV      | 442485         | 457953        | 9.05%           | 2.372%        |
| 23        | 17.757      | 1473          | 1476        | 1479         | rBV      | 356284         | 463568        | 9.16%           | 2.402%        |

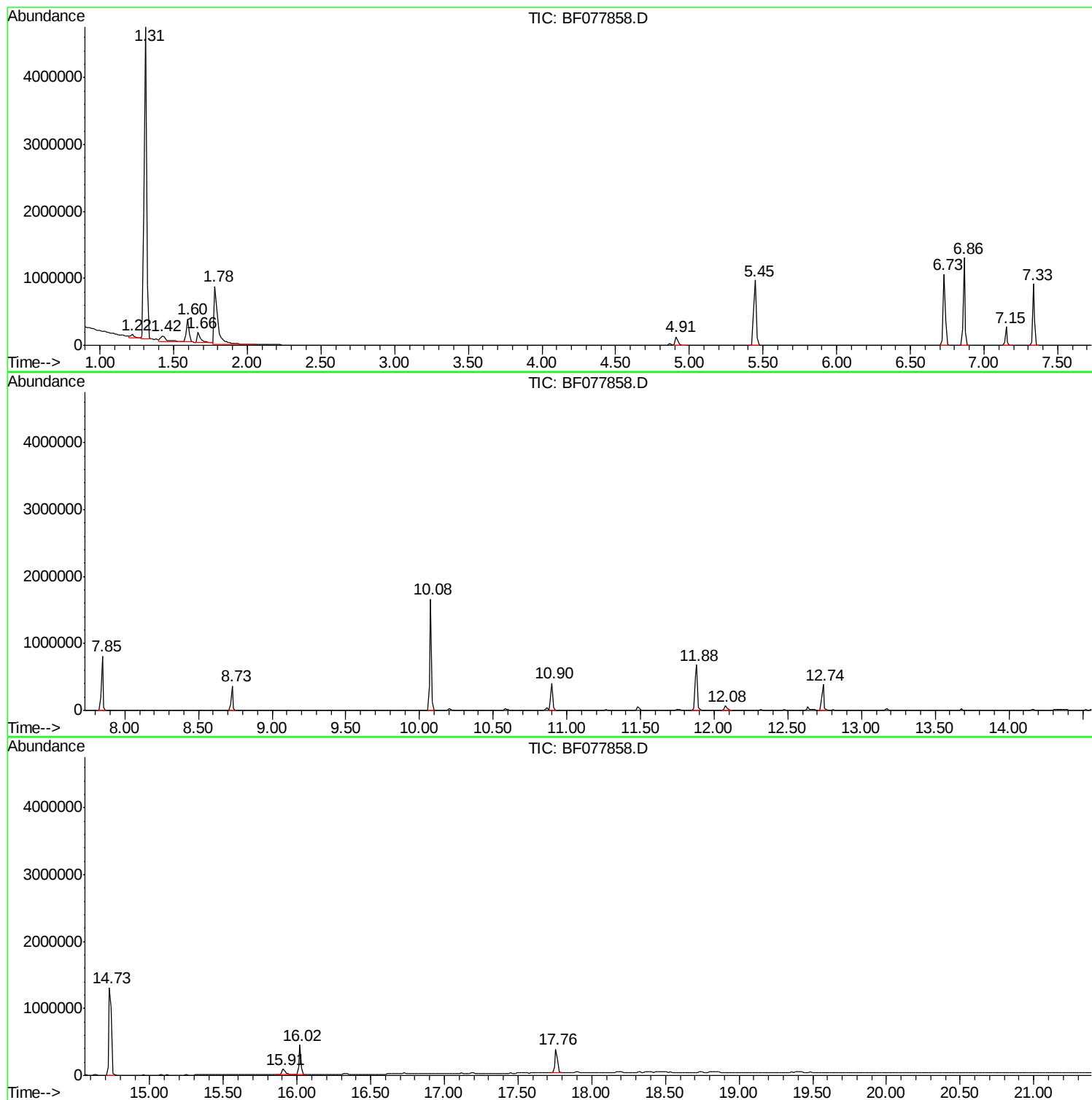
Sum of corrected areas: 19302774

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

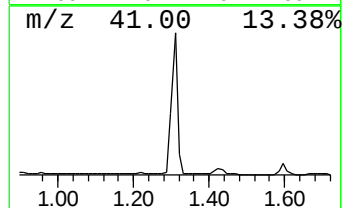
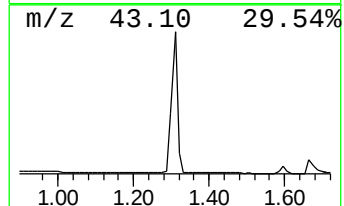
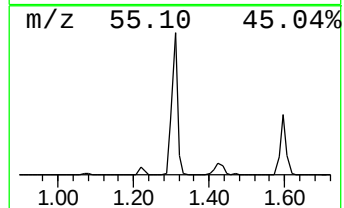
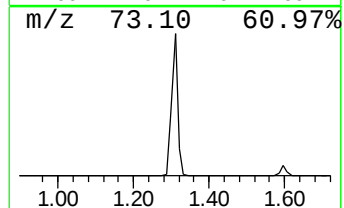
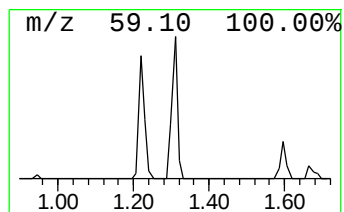
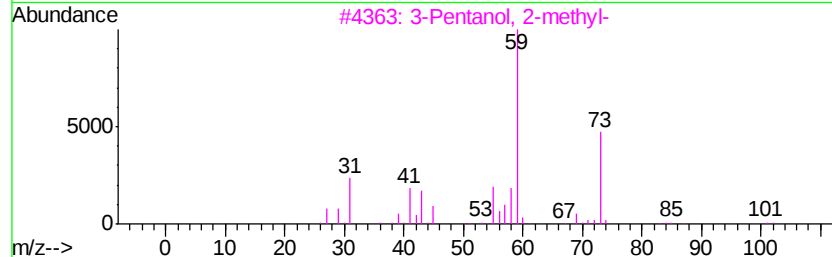
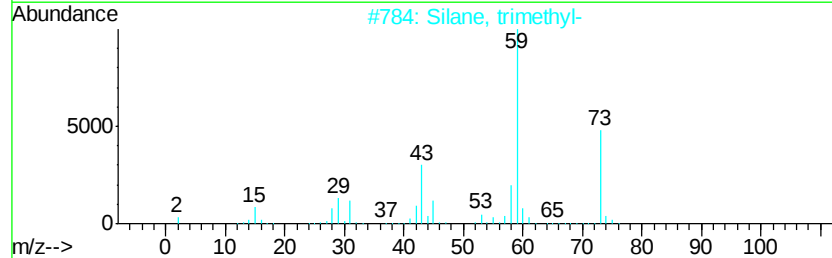
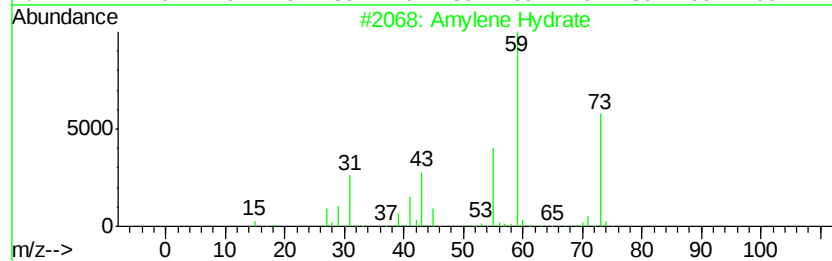
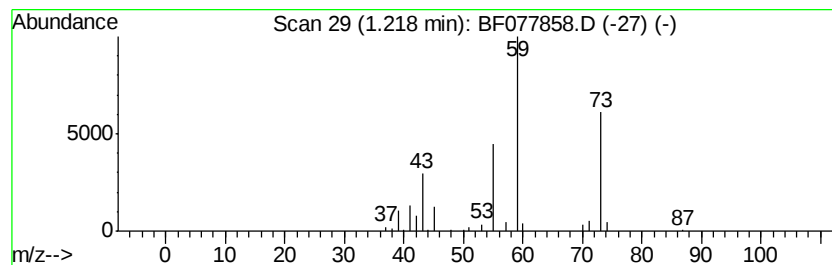
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Amylene Hydrate Concentration Rank 10

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 1.22 | 6.08 ng | 74583 | 1,4-Dichlorobenzene-d4 | 7.15 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Amylene Hydrate       | 88  | C5H12O  | 000075-85-4 | 83   |
| 2    |    |   | Silane, trimethyl-    | 74  | C3H10Si | 000993-07-7 | 64   |
| 3    |    |   | 3-Pentanol, 2-methyl- | 102 | C6H14O  | 000565-67-3 | 39   |
| 4    |    |   | 2-Methyl-5-hexen-3-ol | 114 | C7H14O  | 032815-70-6 | 10   |
| 5    |    |   | Hexanamide            | 115 | C6H13NO | 000628-02-4 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

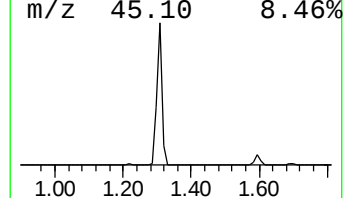
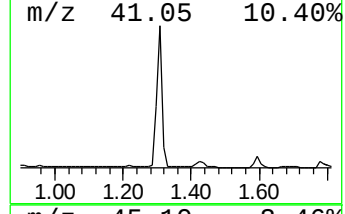
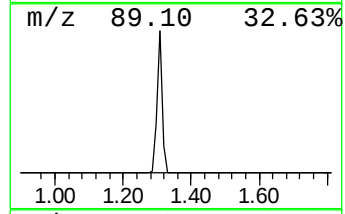
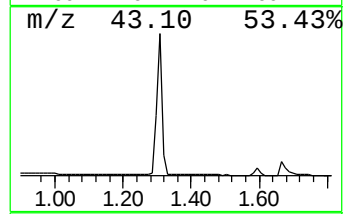
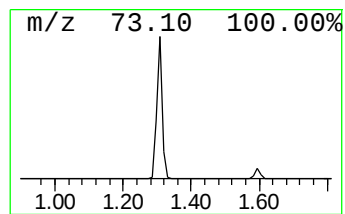
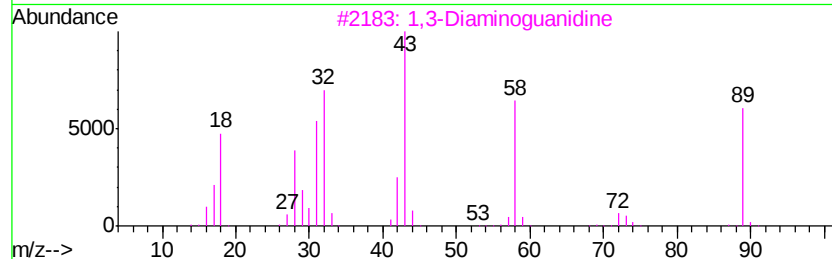
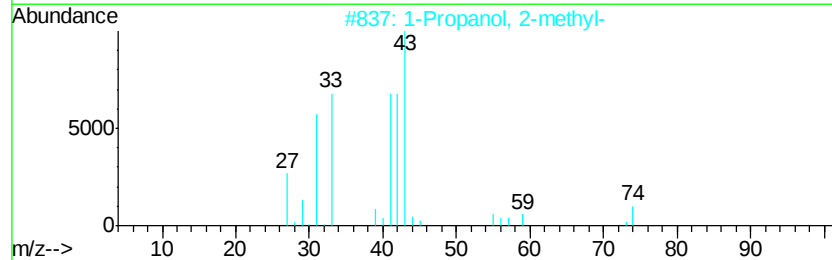
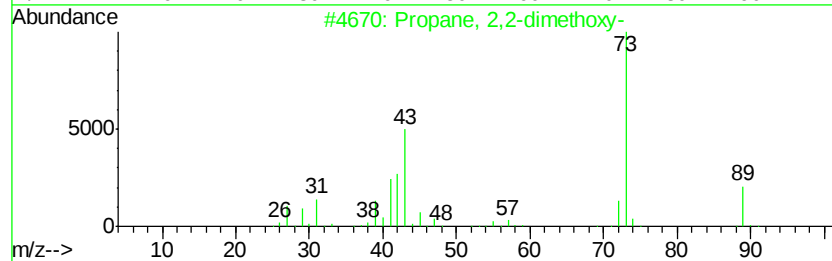
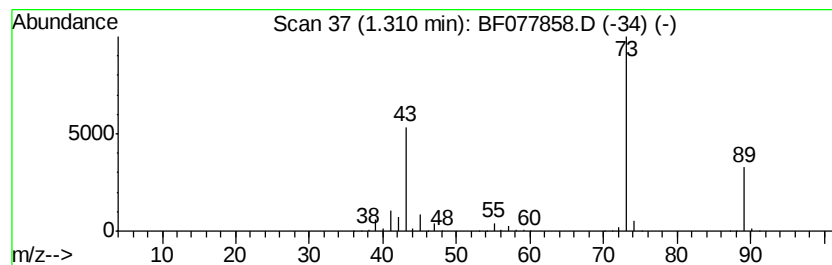
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 1.31 | 412.84 ng | 5061500 | 1,4-Dichlorobenzene-d4 | 7.15 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm       | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------------|--------------|------|
| 1    |    |   | Propane, 2,2-dimethoxy-             | 104 | C5H12O2       | 000077-76-9  | 56   |
| 2    |    |   | 1-Propanol, 2-methyl-               | 74  | C4H10O        | 000078-83-1  | 9    |
| 3    |    |   | 1,3-Diaminoguanidine                | 89  | CH7N5         | 004364-78-7  | 9    |
| 4    |    |   | 1,3-Dioxolane, 2-(3-bromo-5,5,5-... | 352 | C10H16BrCl3O2 | 1000115-31-4 | 9    |
| 5    |    |   | Propane, 2-methoxy-2-methyl-        | 88  | C5H12O        | 001634-04-4  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

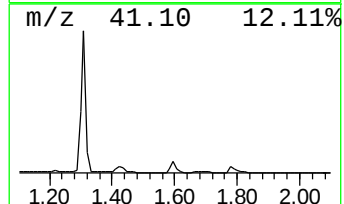
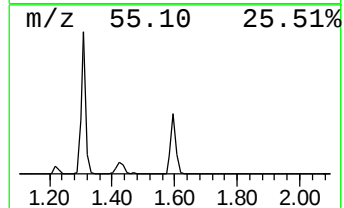
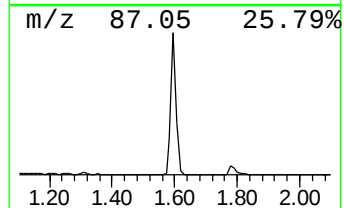
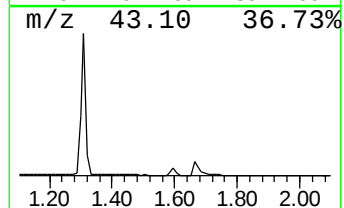
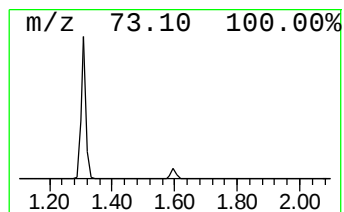
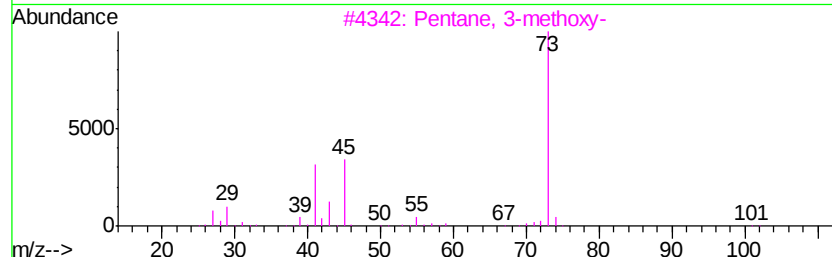
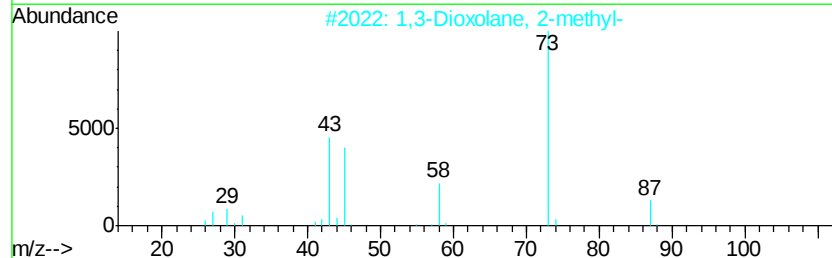
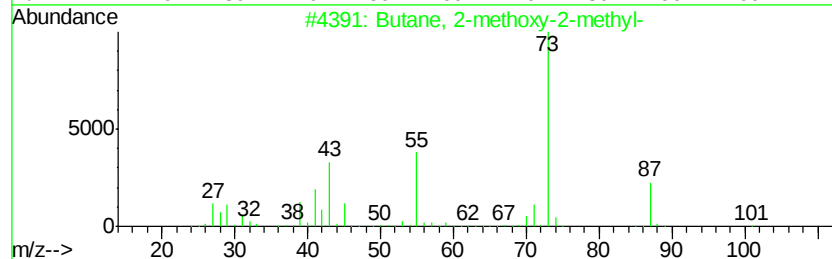
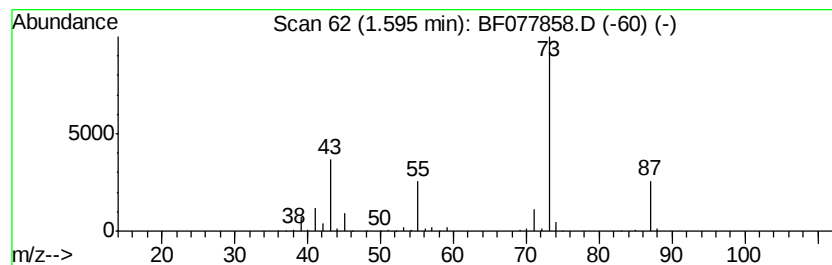
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 1.60 | 31.30 ng | 383777 | 1,4-Dichlorobenzene-d4 | 7.15 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2       |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 37   |
| 3       |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4       |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |
| 5       |   | Borane, dimethoxy-          | 74  | C2H7BO2 | 004542-61-4 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

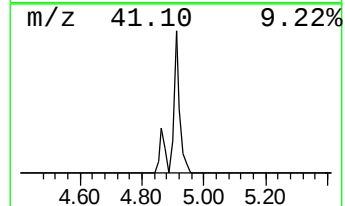
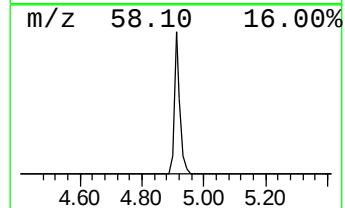
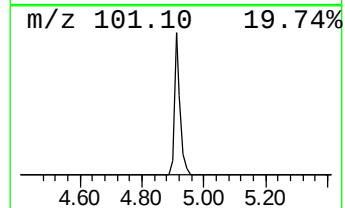
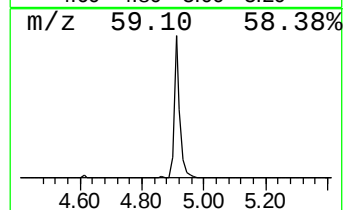
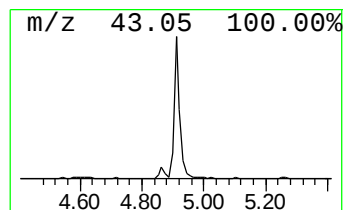
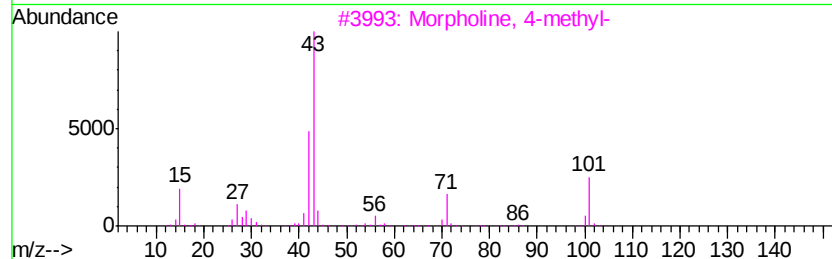
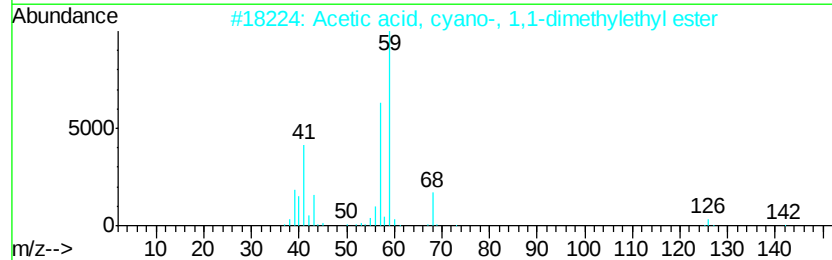
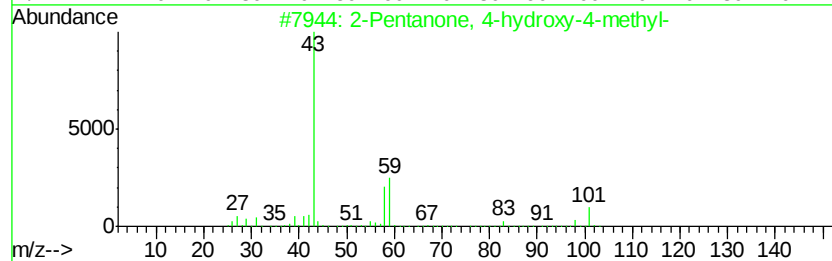
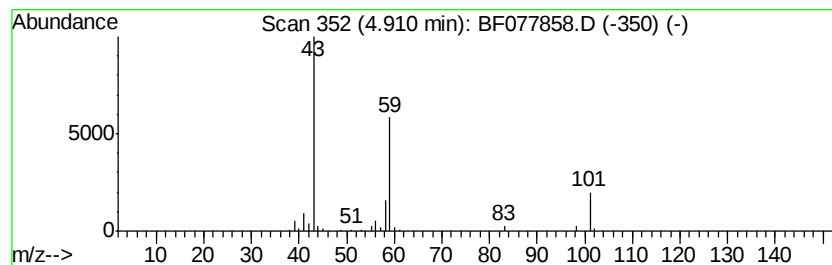
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.91 | 12.95 ng | 158831 | 1,4-Dichlorobenzene-d4 | 7.15 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 50   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 3    |    | Morpholine, 4-methyl-                | 101 | C5H11NO  | 000109-02-4 | 9    |
| 4    |    | Butane, 1-ethoxy-                    | 102 | C6H14O   | 000628-81-9 | 9    |
| 5    |    | 5-Hexen-2-one                        | 98  | C6H10O   | 000109-49-9 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

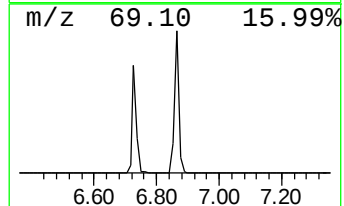
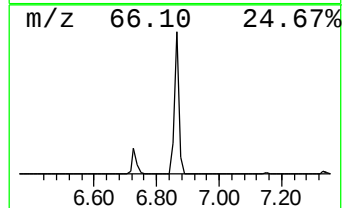
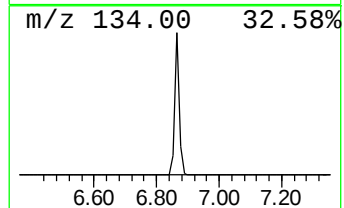
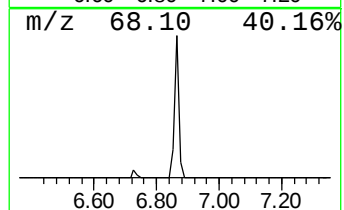
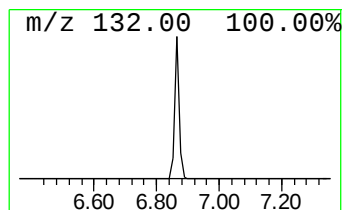
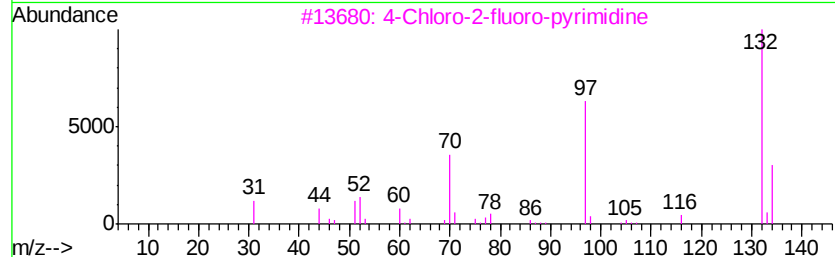
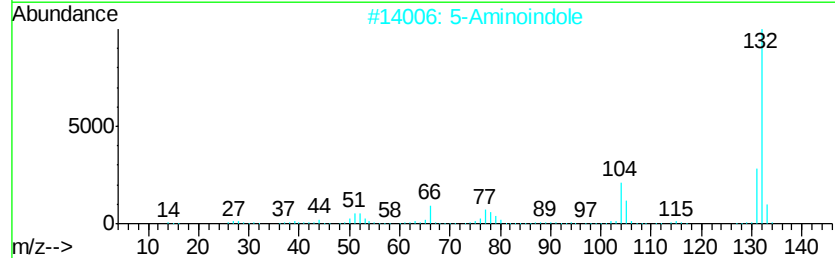
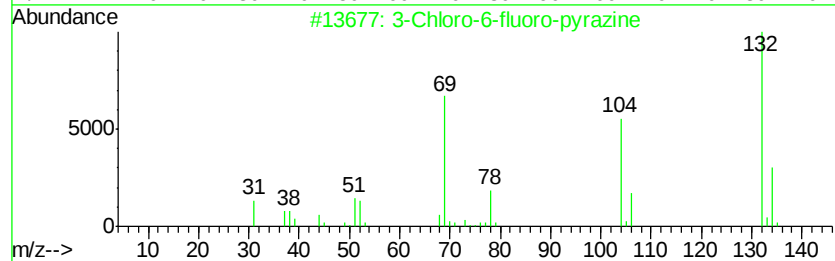
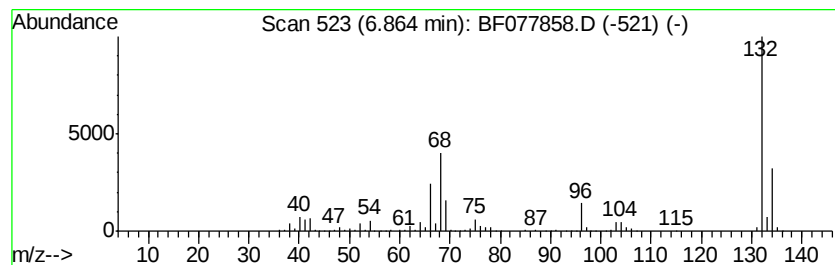
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 unknown6.86 Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.86 | 97.30 ng | 1192960 | 1,4-Dichlorobenzene-d4 | 7.15 |

| Hit# | of | 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    | 4  | Chloro-2-fluoro-pyrimidine   | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 4    | 4  | Chloro-6-fluoro-pyrimidine   | 132 | C4H2ClFN2 | 051422-01-6  | 9    |
| 5    | 1  | Methylpyrrolo[1,2-a]pyrazine | 132 | C8H8N2    | 064608-59-9  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

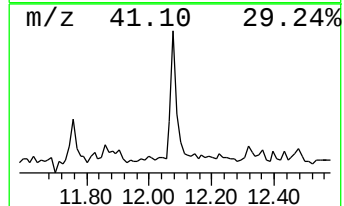
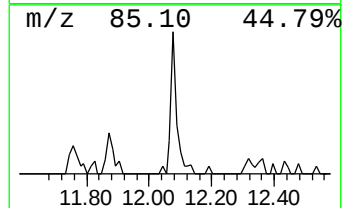
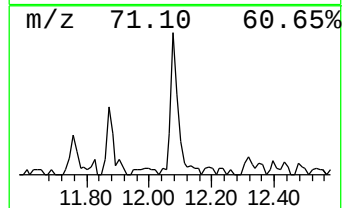
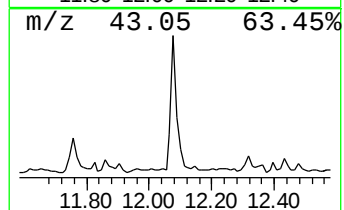
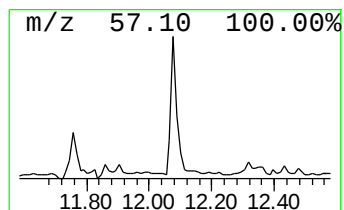
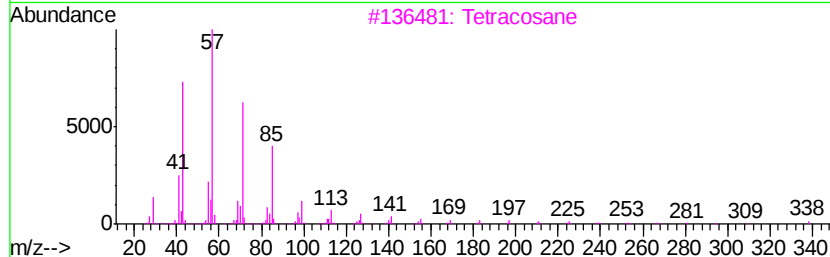
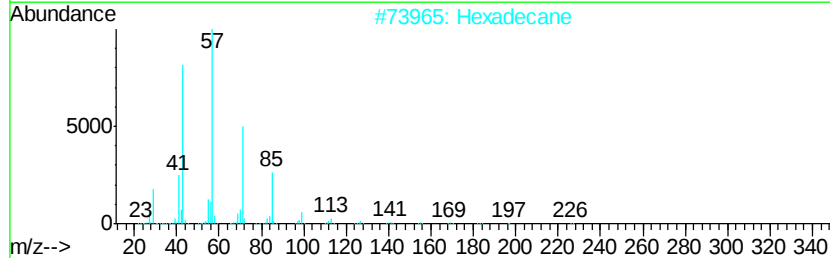
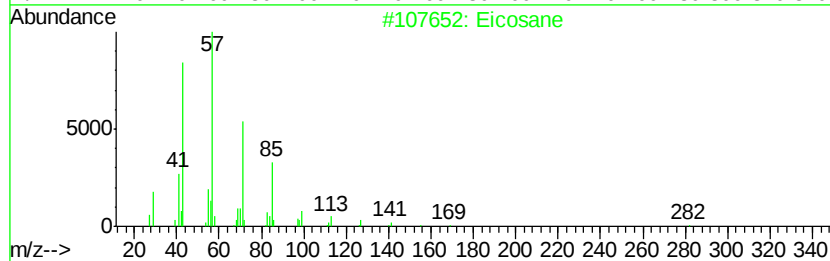
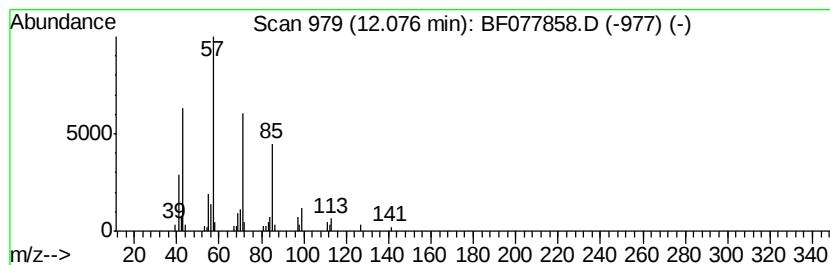
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Eicosane Concentration Rank 11

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.08 | 4.26 ng | 86777 | Phenanthrene-d10 | 12.74 |

| Hit# | of          | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------|---|--------------|-----|---------|-------------|------|
| 1    | Eicosane    |   |              | 282 | C20H42  | 000112-95-8 | 91   |
| 2    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 90   |
| 3    | Tetracosane |   |              | 338 | C24H50  | 000646-31-1 | 86   |
| 4    | Heptadecane |   |              | 240 | C17H36  | 000629-78-7 | 83   |
| 5    | Tetradecane |   |              | 198 | C14H30  | 000629-59-4 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

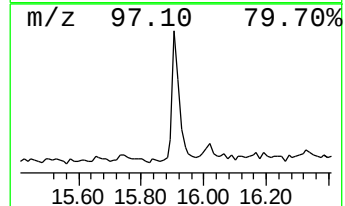
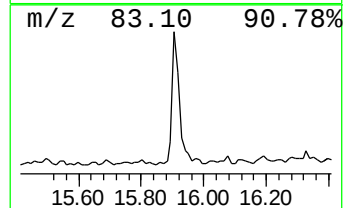
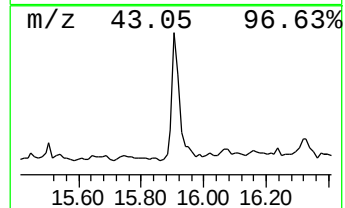
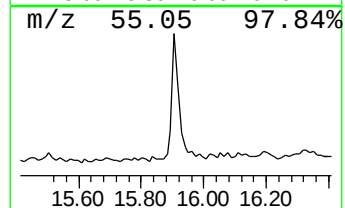
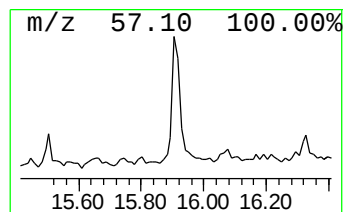
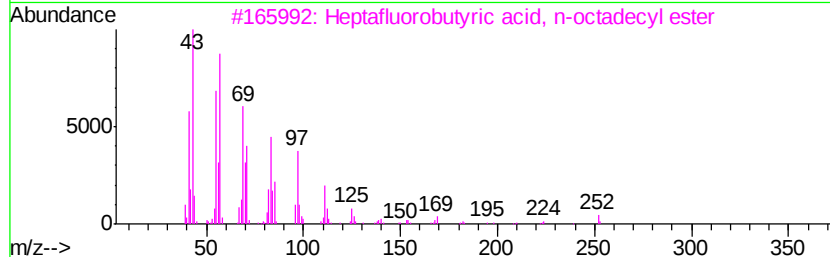
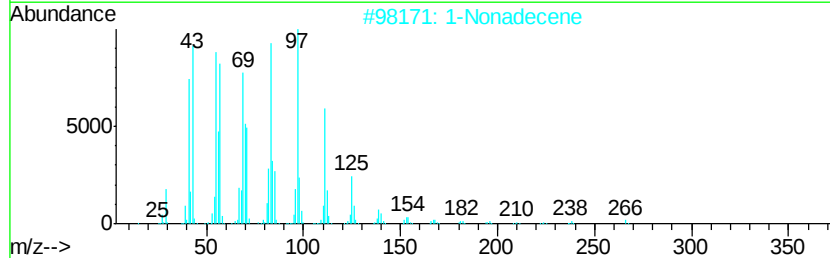
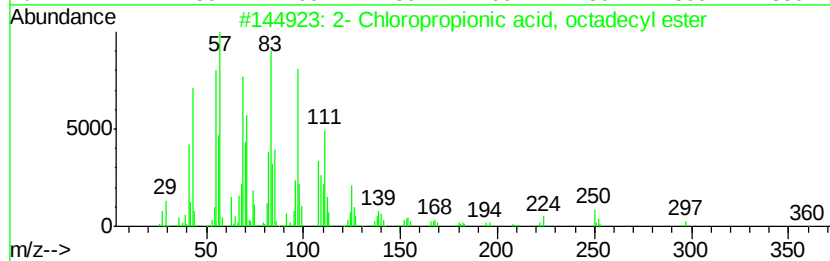
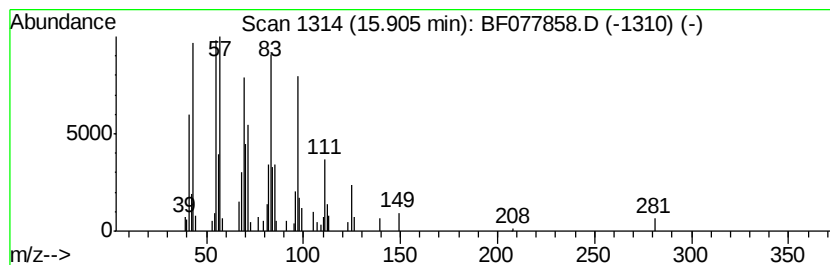
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 2- Chloropropionic acid, oc... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.91 | 6.69 ng | 153296 | Chrysene-d12     | 16.02 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|---------|---|-------------------------------------|-----|------------|--------------|------|
| 1       |   | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2 | 088104-31-8  | 94   |
| 2       |   | 1-Nonadecene                        | 266 | C19H38     | 018435-45-5  | 91   |
| 3       |   | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2 | 000400-57-7  | 91   |
| 4       |   | Heptafluorobutanoic acid, heptad... | 452 | C21H35F7O2 | 1000282-97-3 | 91   |
| 5       |   | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2 | 079392-43-1  | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031915\  
Data File : BF077858.D  
Acq On : 20 Mar 2015 6:13  
Operator : TP/IZ  
Sample : G1580-01  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DGA-DRH-01

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031115.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 |
| Amylene Hydrate      | 1.22  | 6.1     | ng    | 74583    | 1                     | 7.15  | 245206 | 20.0 |
| Propane, 2,2-dime... | 1.31  | 412.8   | ng    | 5061500  | 1                     | 7.15  | 245206 | 20.0 |
| Butane, 2-methoxy... | 1.60  | 31.3    | ng    | 383777   | 1                     | 7.15  | 245206 | 20.0 |
| 2-Pentanone, 4-hy... | 4.91  | 12.9    | ng    | 158831   | 1                     | 7.15  | 245206 | 20.0 |
| unknown6.86          | 6.86  | 97.3    | ng    | 1192960  | 1                     | 7.15  | 245206 | 20.0 |
| Eicosane             | 12.08 | 4.3     | ng    | 86777    | 4                     | 12.74 | 407629 | 20.0 |
| 2- Chloropropioni... | 15.91 | 6.7     | ng    | 153296   | 5                     | 16.02 | 457953 | 20.0 |