

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF043015\  
 Data File : BF078879.D  
 Acq On : 1 May 2015 6:21  
 Operator : TP/IZ  
 Sample : G2032-04  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B2(5-7)

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-BF043015.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         | 1.541       | 28            | 31          | 34           | rVB      | 6696514        | 8625752       | 100.00%         | 21.322%       |
| 2         | 1.678       | 40            | 43          | 50           | rVB      | 225409         | 539304        | 6.25%           | 1.333%        |
| 3         | 1.781       | 50            | 52          | 57           | rVV      | 404932         | 815944        | 9.46%           | 2.017%        |
| 4         | 1.873       | 57            | 60          | 62           | rVV      | 235628         | 405514        | 4.70%           | 1.002%        |
| 5         | 1.907       | 62            | 63          | 91           | rVB      | 3085565        | 5728556       | 66.41%          | 14.161%       |
| 6         | 5.187       | 348           | 350         | 359          | rVB      | 223305         | 326794        | 3.79%           | 0.808%        |
| 7         | 5.667       | 390           | 392         | 395          | rBV      | 1463015        | 2075490       | 24.06%          | 5.130%        |
| 8         | 6.925       | 499           | 502         | 504          | rBV      | 2331770        | 2301501       | 26.68%          | 5.689%        |
| 9         | 7.085       | 513           | 516         | 518          | rBV      | 2325049        | 2308792       | 26.77%          | 5.707%        |
| 10        | 7.370       | 539           | 541         | 543          | rBV      | 590825         | 541540        | 6.28%           | 1.339%        |
| 11        | 7.565       | 555           | 558         | 560          | rVB      | 1265431        | 1709605       | 19.82%          | 4.226%        |
| 12        | 8.056       | 599           | 601         | 604          | rBV      | 1091938        | 1304158       | 15.12%          | 3.224%        |
| 13        | 8.948       | 677           | 679         | 682          | rBV      | 880558         | 753903        | 8.74%           | 1.864%        |
| 14        | 10.296      | 794           | 797         | 799          | rBV      | 2909480        | 2570857       | 29.80%          | 6.355%        |
| 15        | 10.799      | 839           | 841         | 843          | rBV      | 102955         | 94907         | 1.10%           | 0.235%        |
| 16        | 11.131      | 867           | 870         | 872          | rBV      | 715207         | 822145        | 9.53%           | 2.032%        |
| 17        | 11.474      | 898           | 900         | 903          | rBV      | 97737          | 105619        | 1.22%           | 0.261%        |
| 18        | 12.102      | 953           | 955         | 958          | rBV2     | 1494356        | 1703527       | 19.75%          | 4.211%        |
| 19        | 12.971      | 1029          | 1031        | 1033         | rBV      | 836028         | 876946        | 10.17%          | 2.168%        |
| 20        | 13.679      | 1091          | 1093        | 1096         | rBV      | 63835          | 87931         | 1.02%           | 0.217%        |
| 21        | 14.480      | 1161          | 1163        | 1166         | rBV      | 129673         | 138809        | 1.61%           | 0.343%        |
| 22        | 14.571      | 1168          | 1171        | 1173         | rVB      | 74029          | 91001         | 1.05%           | 0.225%        |
| 23        | 14.765      | 1185          | 1188        | 1189         | rBV      | 125046         | 143868        | 1.67%           | 0.356%        |
| 24        | 14.834      | 1192          | 1194        | 1197         | rBV      | 313717         | 402643        | 4.67%           | 0.995%        |
| 25        | 14.971      | 1203          | 1206        | 1209         | rBV      | 2895341        | 2787096       | 32.31%          | 6.890%        |
| 26        | 16.137      | 1306          | 1308        | 1313         | rBV      | 251660         | 397646        | 4.61%           | 0.983%        |
| 27        | 16.274      | 1317          | 1320        | 1322         | rVV      | 788296         | 1051713       | 12.19%          | 2.600%        |
| 28        | 16.308      | 1322          | 1323        | 1326         | rVB2     | 76356          | 89800         | 1.04%           | 0.222%        |
| 29        | 16.983      | 1380          | 1382        | 1387         | rVB      | 109042         | 149205        | 1.73%           | 0.369%        |
| 30        | 17.611      | 1435          | 1437        | 1439         | rBV      | 57561          | 109297        | 1.27%           | 0.270%        |
| 31        | 17.794      | 1451          | 1453        | 1459         | rBV2     | 67404          | 205573        | 2.38%           | 0.508%        |
| 32        | 18.080      | 1475          | 1478        | 1483         | rVB      | 682198         | 1025958       | 11.89%          | 2.536%        |
| 33        | 20.252      | 1665          | 1668        | 1674         | rBV5     | 59462          | 162700        | 1.89%           | 0.402%        |

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BNA\_F  
ClientSampleId :  
B2(5-7)

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-BF043015.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

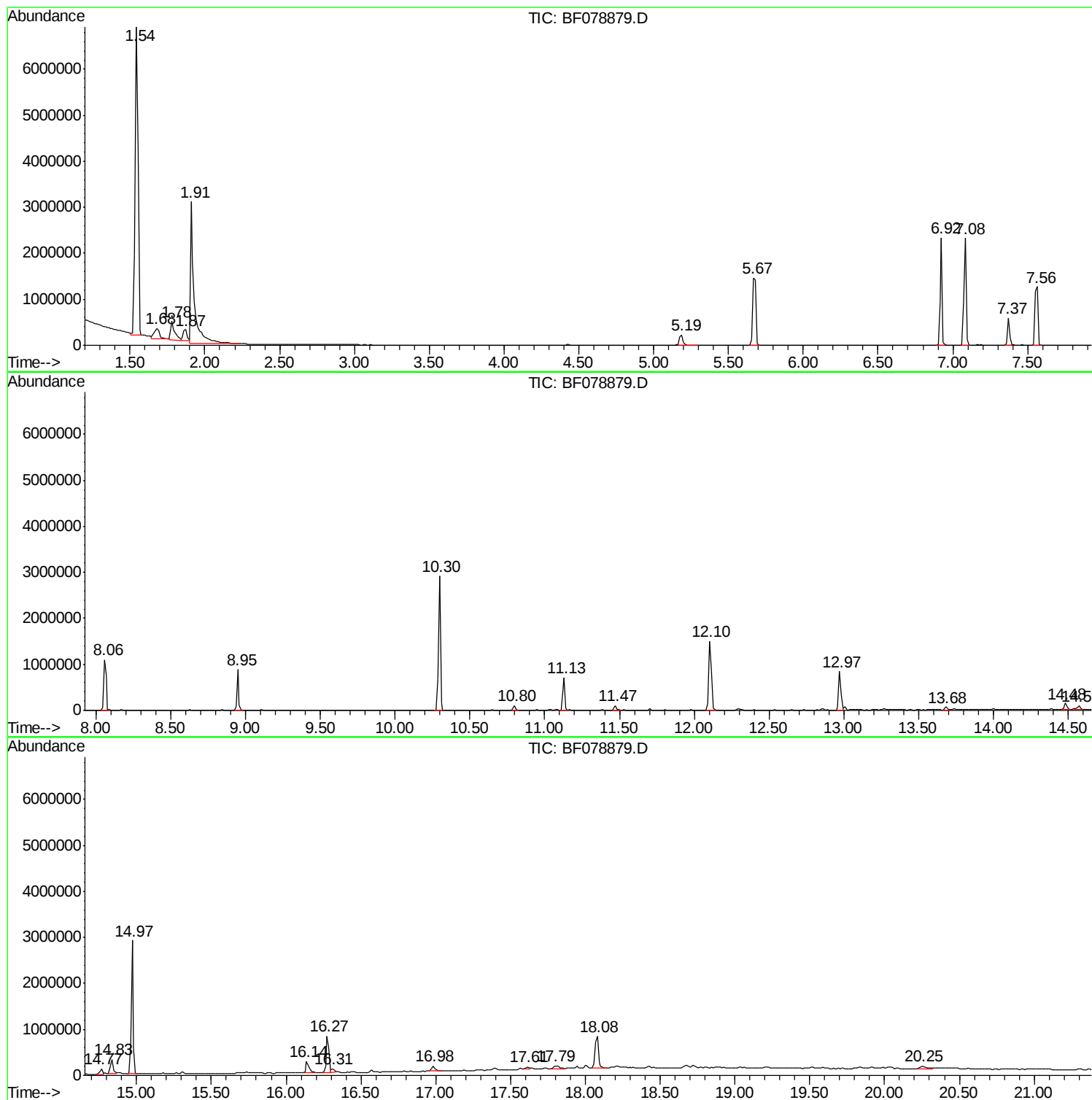
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :  
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B2(5-7)

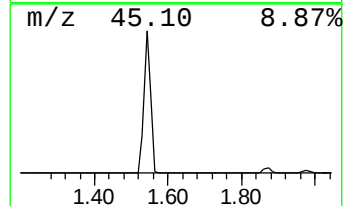
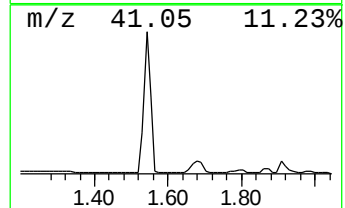
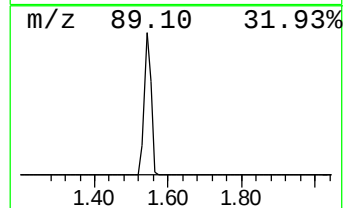
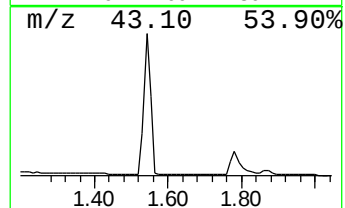
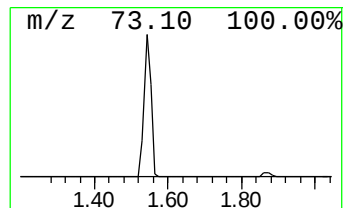
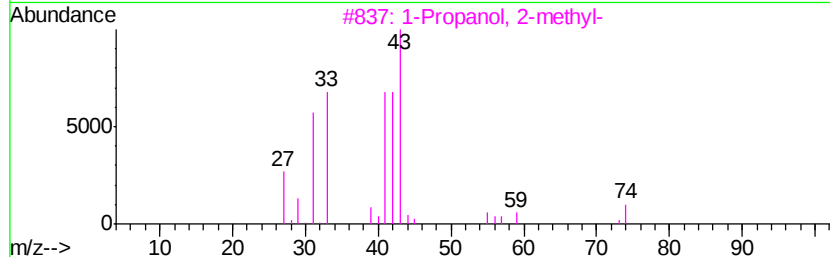
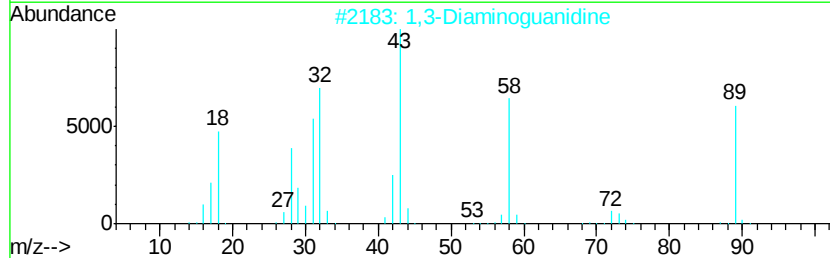
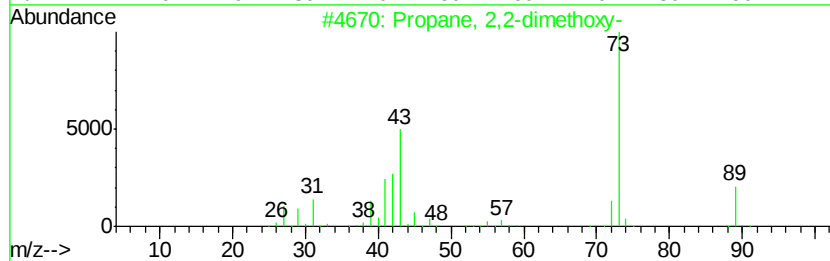
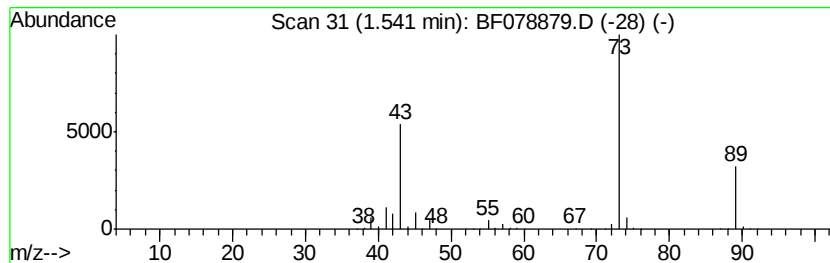
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 unknown1.54 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 1.54 | 318.56 ng | 8625750 | 1,4-Dichlorobenzene-d4 | 7.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Propane, 2,2-dimethoxy-             | 104 | C5H12O2 | 000077-76-9 | 38   |
| 2    |    |   | 1,3-Diaminoguanidine                | 89  | CH7N5   | 004364-78-7 | 9    |
| 3    |    |   | 1-Propanol, 2-methyl-               | 74  | C4H10O  | 000078-83-1 | 9    |
| 4    |    |   | Propane, 2-methoxy-2-methyl-        | 88  | C5H12O  | 001634-04-4 | 9    |
| 5    |    |   | Propanoic acid, 2-methyl-, propy... | 130 | C7H14O2 | 000644-49-5 | 9    |



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ClientSampleId :  
B2(5-7)

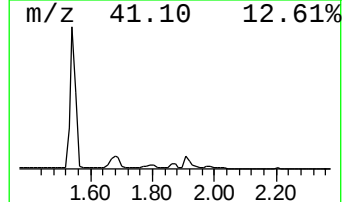
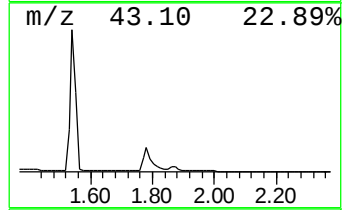
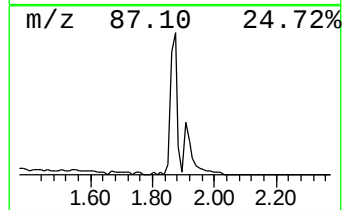
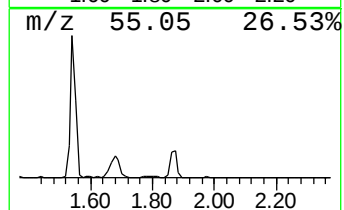
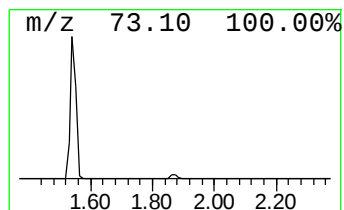
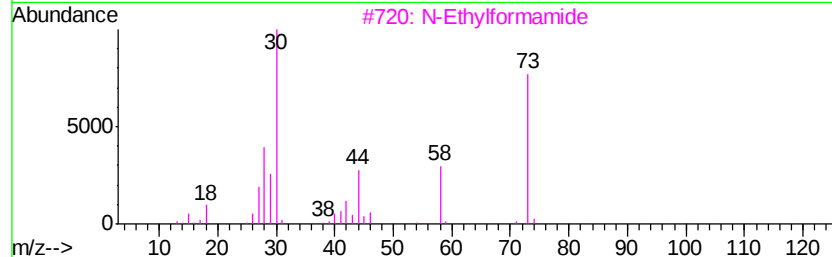
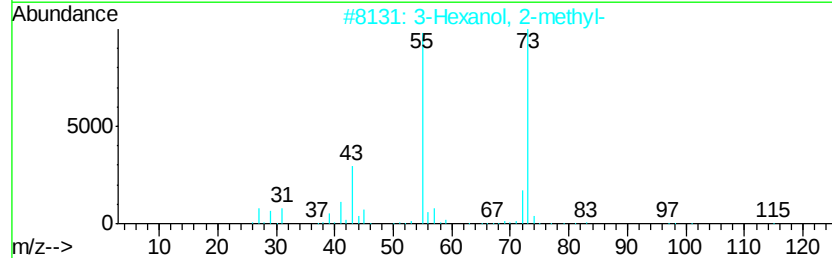
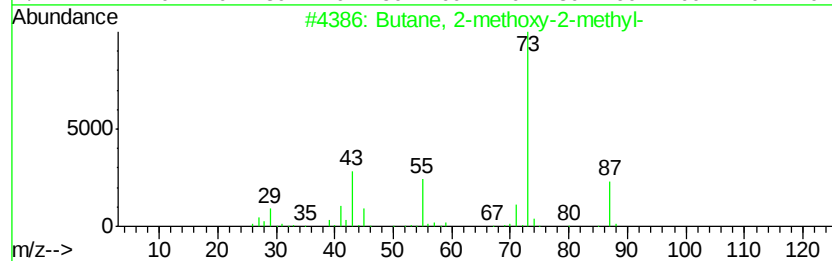
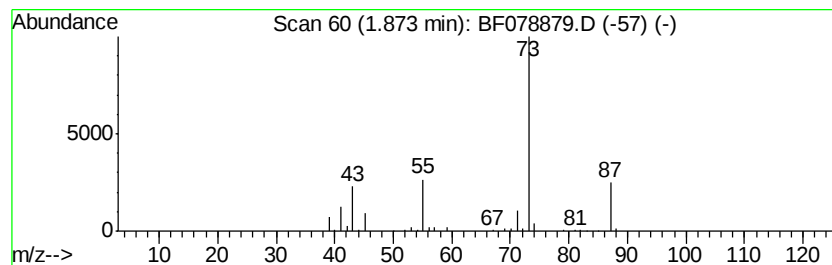
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
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TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 7

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 1.87 | 14.98 ng | 405514 | 1,4-Dichlorobenzene-d4 | 7.37 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 90   |
| 2    |    | 3-Hexanol, 2-methyl-        | 116 | C7H16O  | 000617-29-8 | 23   |
| 3    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 4    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



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Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

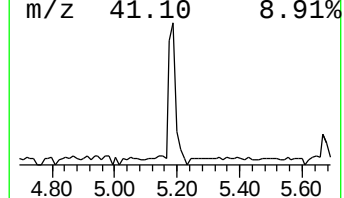
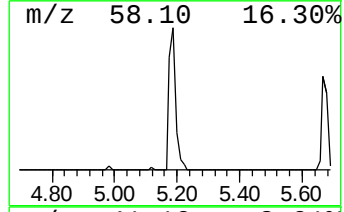
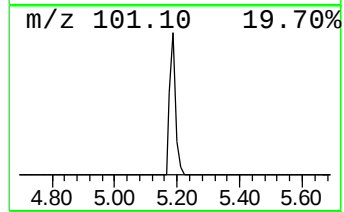
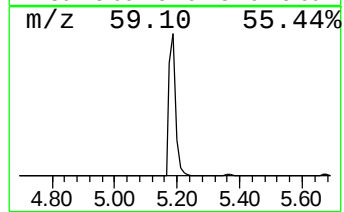
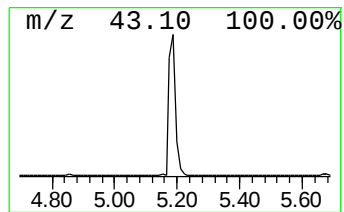
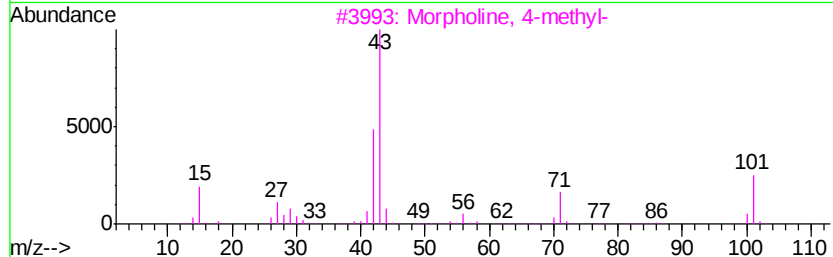
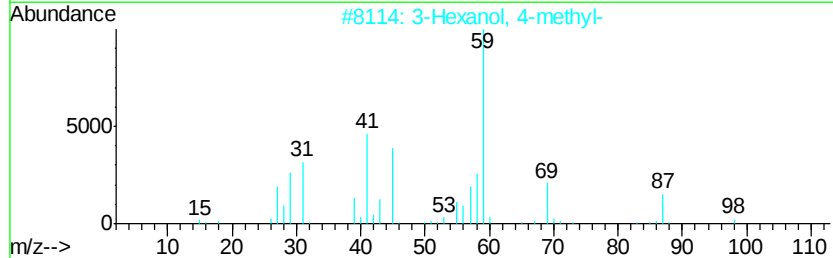
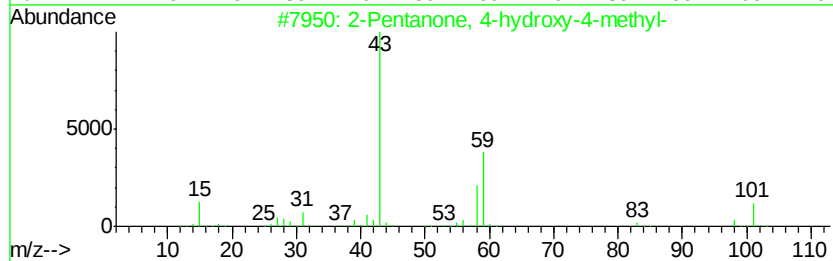
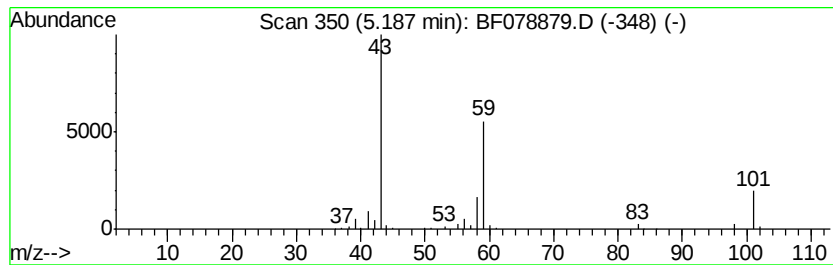
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.19 | 12.07 ng | 326794 | 1,4-Dichlorobenzene-d4 | 7.37 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2 | 000123-42-2 | 72   |
| 2    |    | 3-Hexanol, 4-methyl-                 | 116 | C7H16O  | 000615-29-2 | 28   |
| 3    |    | Morpholine, 4-methyl-                | 101 | C5H11NO | 000109-02-4 | 9    |
| 4    |    | 4-Penten-2-one, 4-methyl-            | 98  | C6H10O  | 003744-02-3 | 9    |
| 5    |    | Propane, 2-methyl-2-(1-methylethyl)- | 116 | C7H16O  | 017348-59-3 | 9    |



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ALS Vial : 29 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

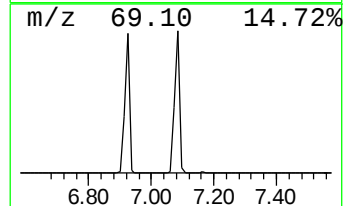
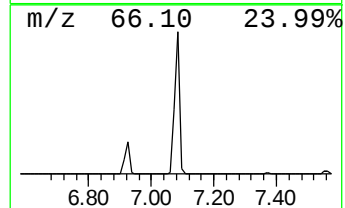
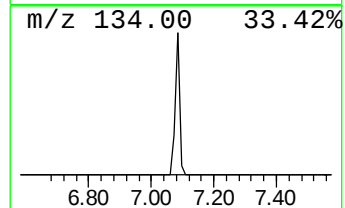
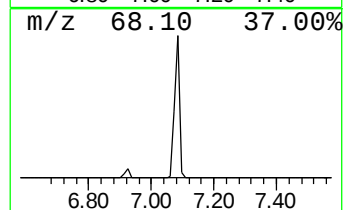
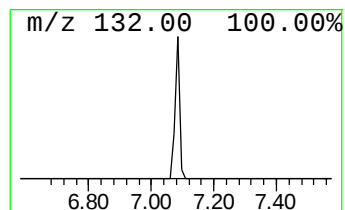
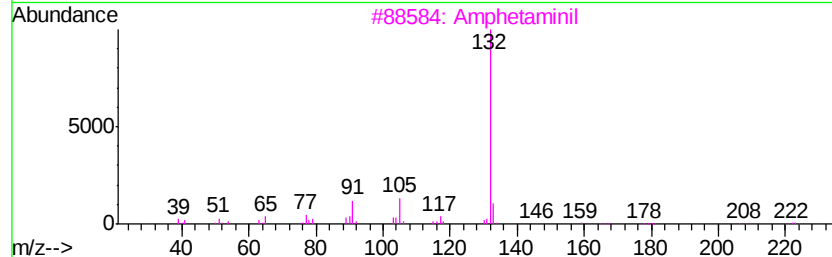
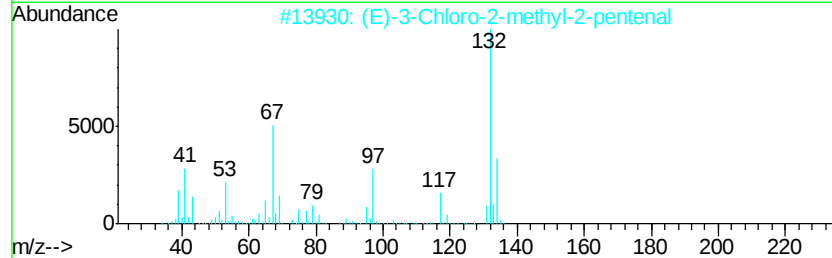
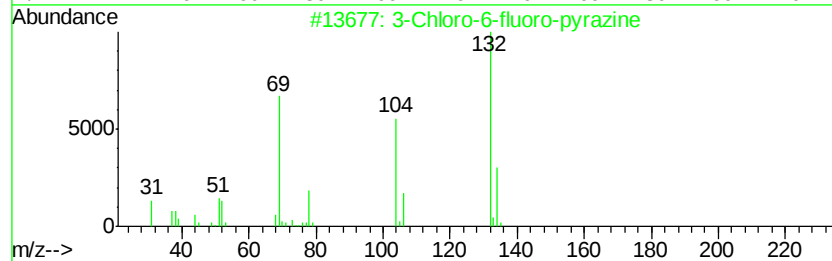
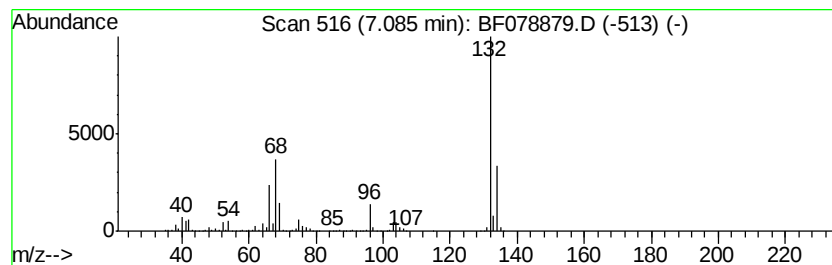
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown7.08 Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.08 | 85.27 ng | 2308790 | 1,4-Dichlorobenzene-d4 | 7.37 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 12   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 5    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |



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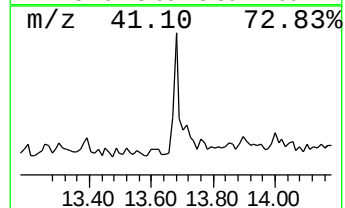
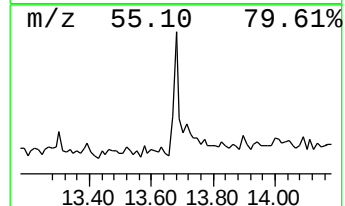
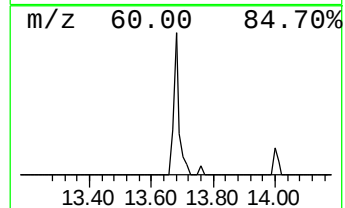
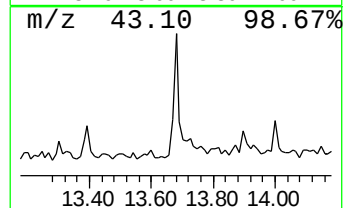
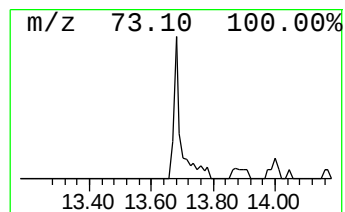
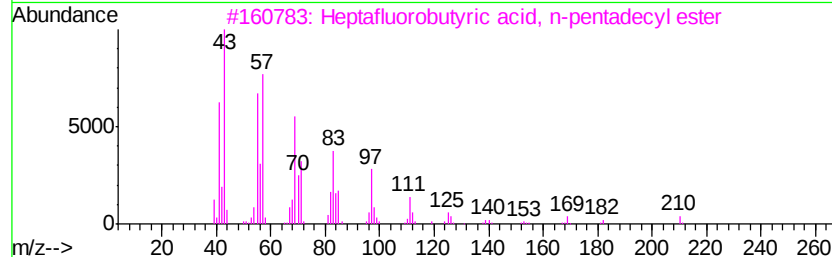
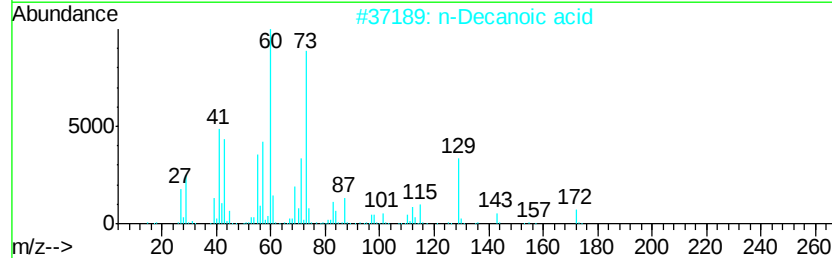
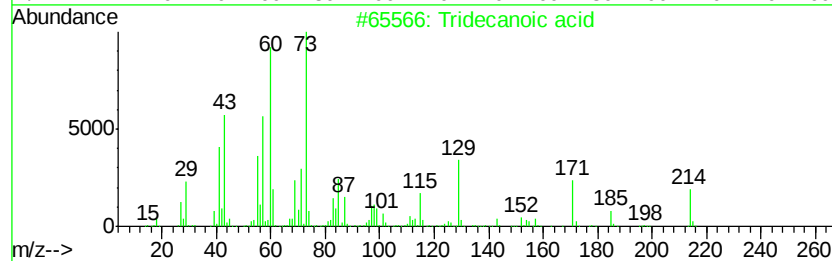
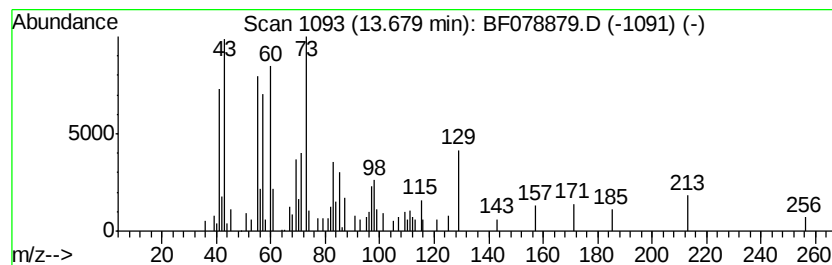
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Tridecanoic acid Concentration Rank 16

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.68 | 2.01 ng | 87931 | Phenanthrene-d10 | 12.97 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Tridecanoic acid                    | 214 | C13H26O2   | 000638-53-9  | 86   |
| 2    |    |   | n-Decanoic acid                     | 172 | C10H20O2   | 000334-48-5  | 58   |
| 3    |    |   | Heptafluorobutyric acid, n-penta... | 424 | C19H31F7O2 | 1000216-79-3 | 15   |
| 4    |    |   | 17-Pentatriacontene                 | 491 | C35H70     | 006971-40-0  | 15   |
| 5    |    |   | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2 | 079392-43-1  | 11   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

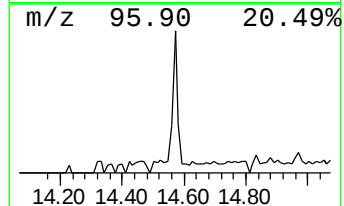
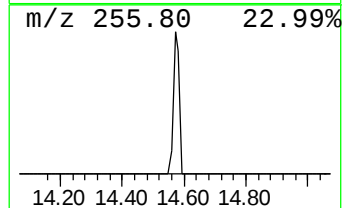
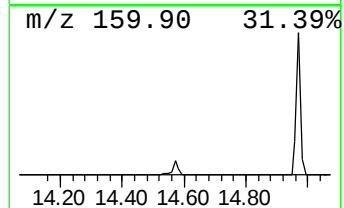
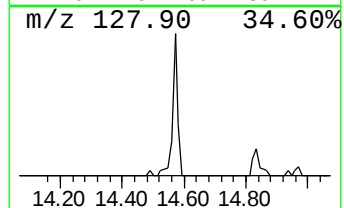
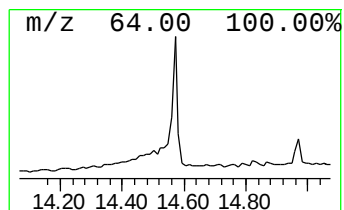
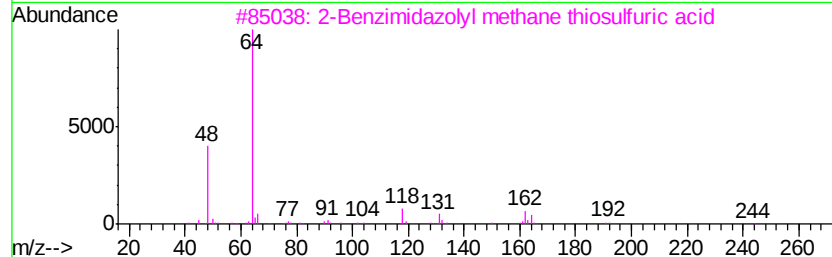
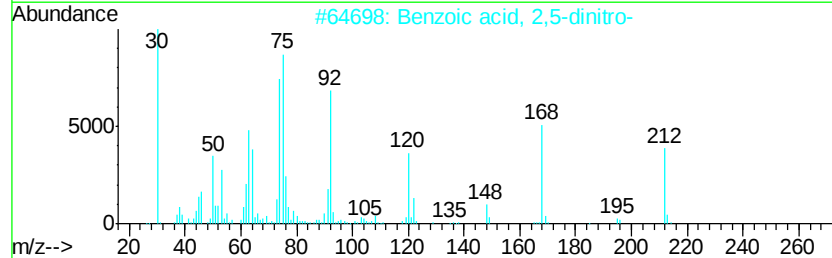
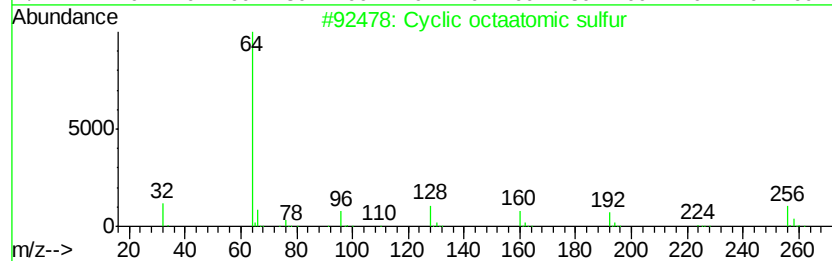
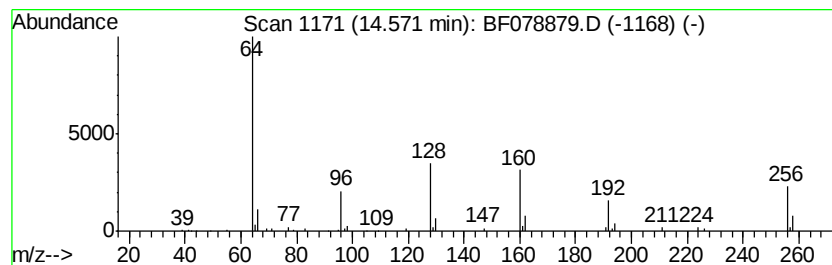
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Cyclic octaatomic sulfur Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.57 | 2.08 ng | 91001 | Phenanthrene-d10 | 12.97 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Cyclic octaatomic sulfur            | 256 | S8         | 010544-50-0  | 90   |
| 2    |    |   | Benzoic acid, 2,5-dinitro-          | 212 | C7H4N2O6   | 000610-28-6  | 53   |
| 3    |    |   | 2-Benzimidazolyl methane thiosul... | 244 | C8H8N2O3S2 | 1000256-39-2 | 38   |
| 4    |    |   | 1,2,5,6-Tetrathiocane               | 184 | C4H8S4     | 001940-01-8  | 9    |
| 5    |    |   | 2-Norbornanone, 1-(epithioethyl)... | 228 | C11H16O3S  | 001782-93-0  | 9    |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

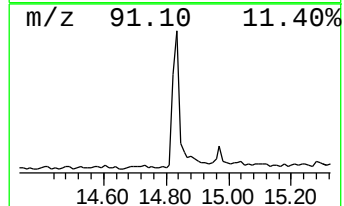
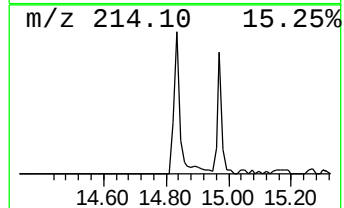
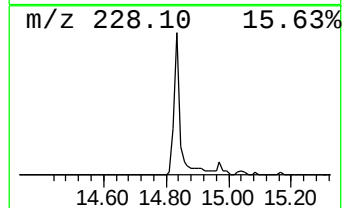
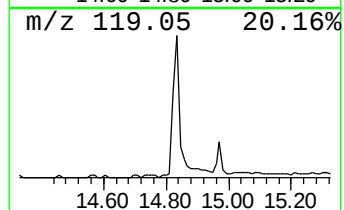
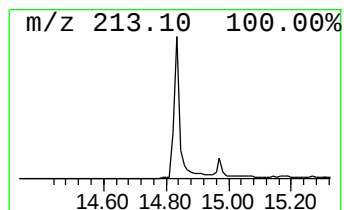
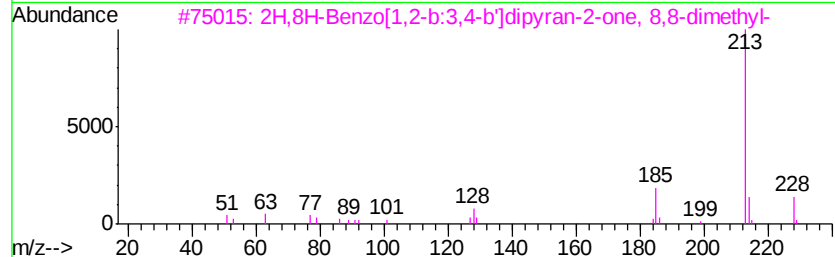
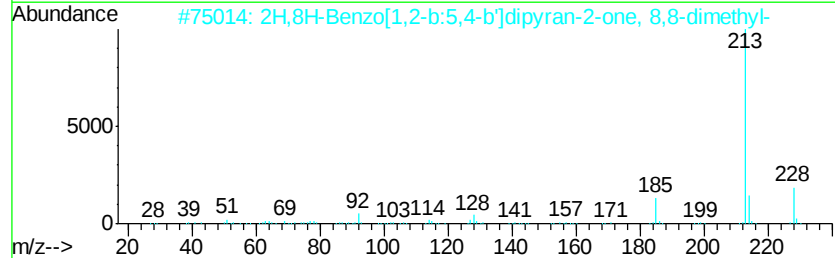
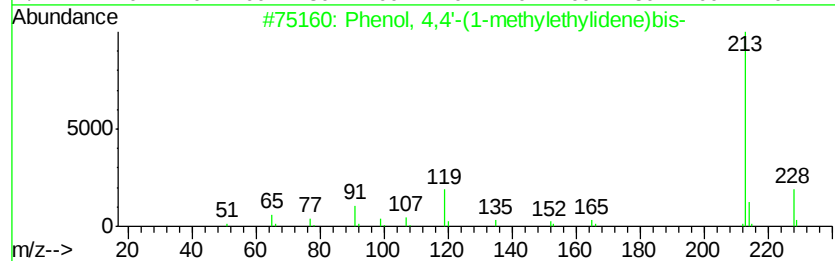
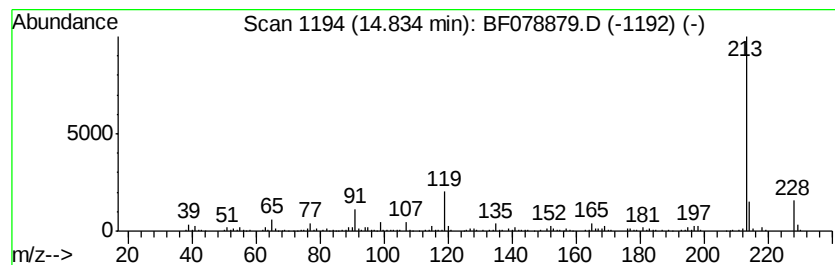
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Phenol, 4,4'-(1-methylethyl... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.83 | 7.66 ng | 402643 | Chrysene-d12     | 16.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2   | 000080-05-7  | 98   |
| 2    |    | 2H,8H-Benzo[1,2-b:5,4-b']dipyran... | 228 | C14H12O3   | 000553-19-5  | 52   |
| 3    |    | 2H,8H-Benzo[1,2-b:3,4-b']dipyran... | 228 | C14H12O3   | 000523-59-1  | 50   |
| 4    |    | Trimethyl[5-(trimethylsilyl)-2-t... | 228 | C10H20SSi2 | 017906-71-7  | 42   |
| 5    |    | 1,4-Puran, 5,6-dihydro-3-[3-indo... | 213 | C14H15NO   | 1000128-72-9 | 40   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

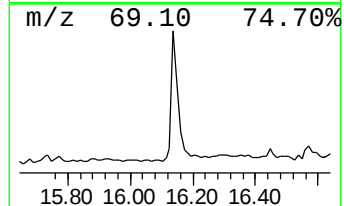
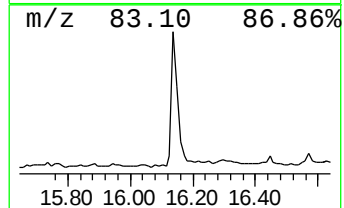
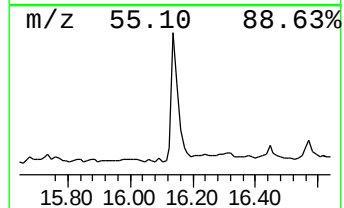
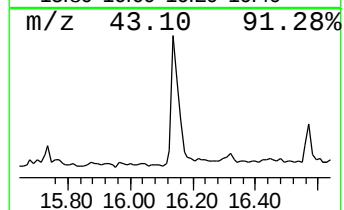
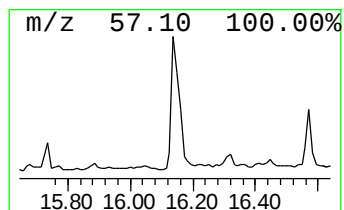
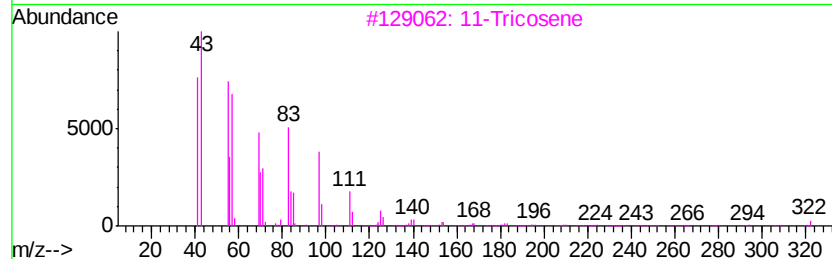
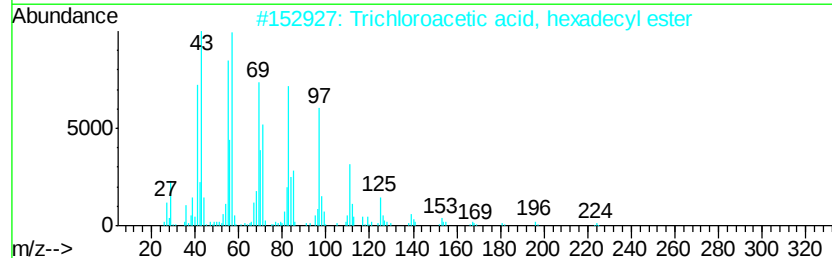
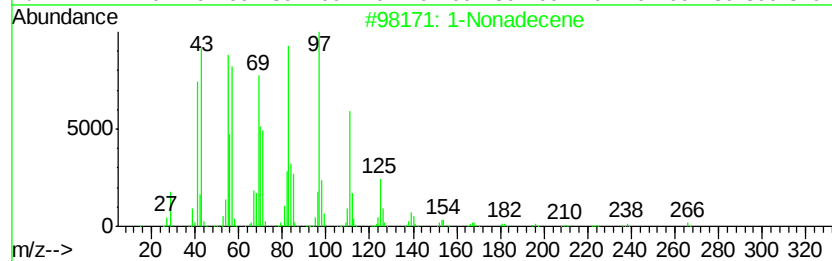
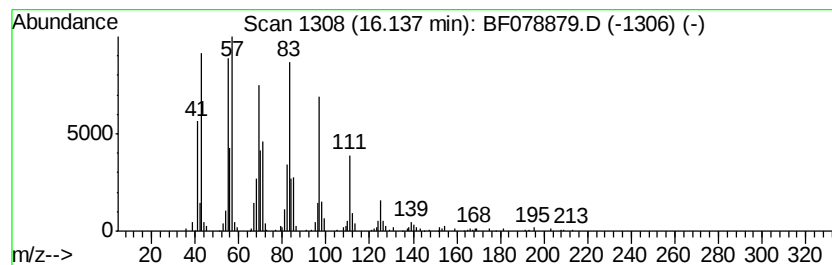
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 1-Nonadecene Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.14 | 7.56 ng | 397646 | Chrysene-d12     | 16.27 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 1-Nonadecene                        | 266 | C19H38      | 018435-45-5 | 91   |
| 2    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 90   |
| 3    |    |   | 11-Tricosene                        | 322 | C23H46      | 052078-56-5 | 90   |
| 4    |    |   | 1-Heptadecanol                      | 256 | C17H36O     | 001454-85-9 | 87   |
| 5    |    |   | 3-Eicosene, (E)-                    | 280 | C20H40      | 074685-33-9 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

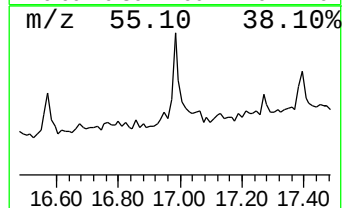
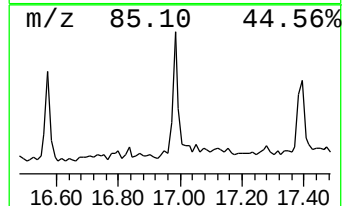
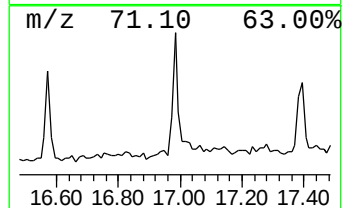
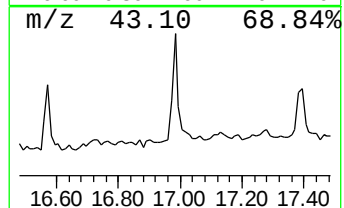
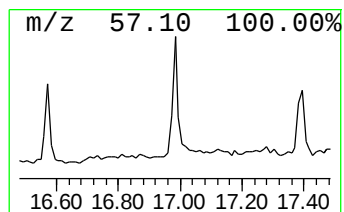
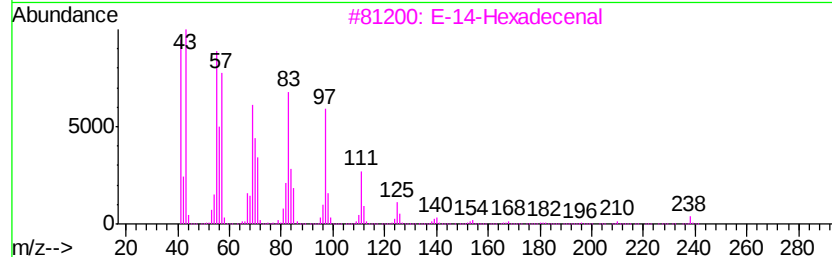
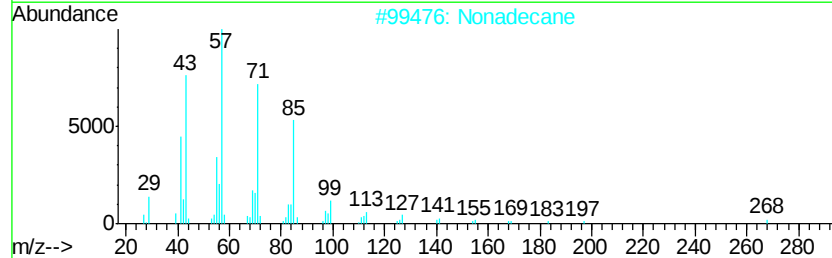
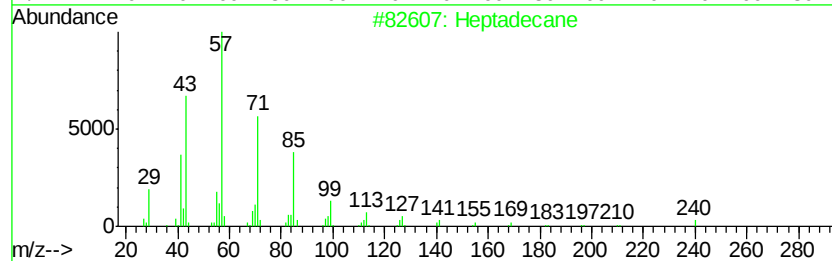
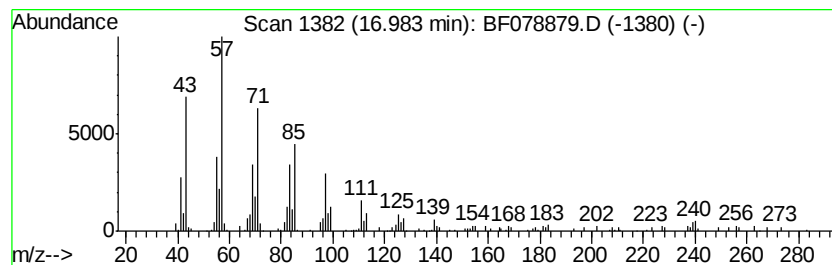
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Heptadecane Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.98 | 2.84 ng | 149205 | Chrysene-d12     | 16.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Heptadecane                         | 240 | C17H36     | 000629-78-7 | 95   |
| 2    |    | Nonadecane                          | 268 | C19H40     | 000629-92-5 | 93   |
| 3    |    | E-14-Hexadecenal                    | 238 | C16H30O    | 330207-53-9 | 89   |
| 4    |    | Nonahexacontanoic acid              | 999 | C69H138O2  | 040710-32-5 | 87   |
| 5    |    | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2 | 079392-43-1 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

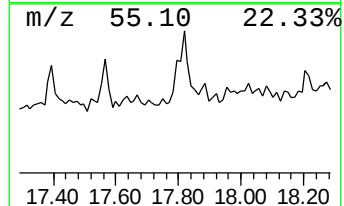
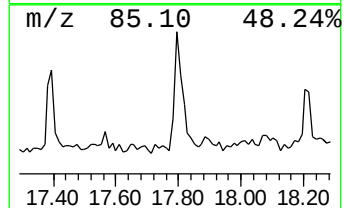
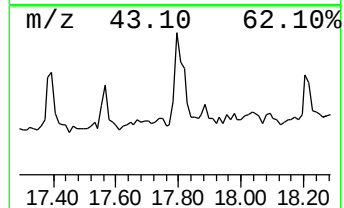
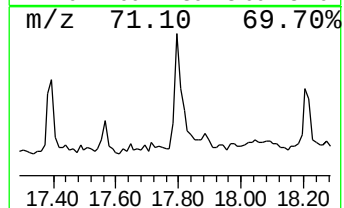
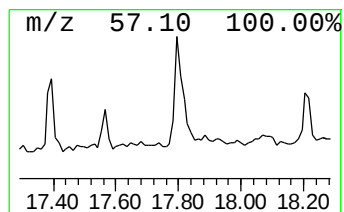
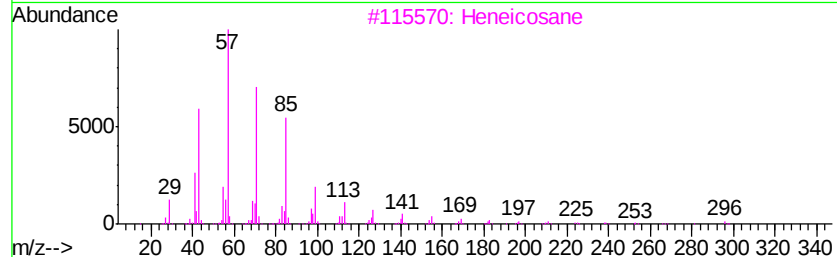
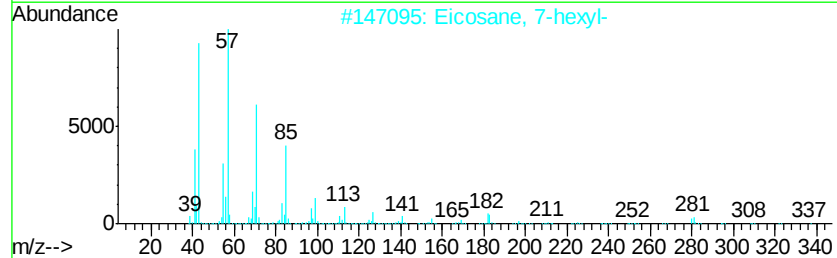
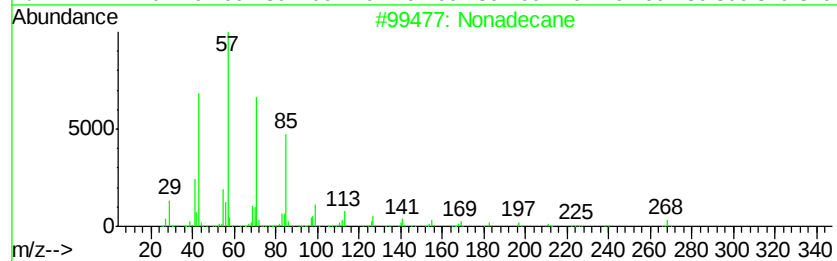
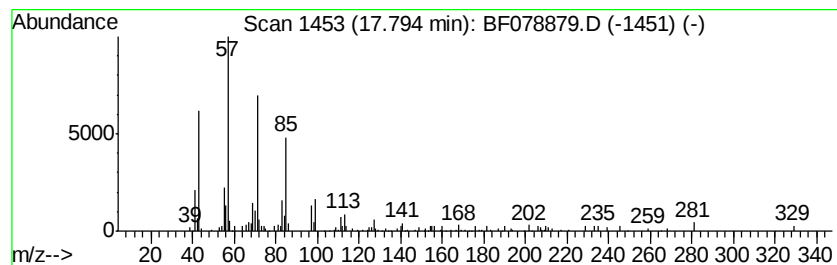
Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.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 15 Nonadecane Concentration Rank 11

| R.T.      | EstConc            | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------------|--------|------------------|-------------|-------|
| 17.79     | 4.01 ng            | 205573 | Perylene-d12     |             | 18.08 |
| Hit# of 5 | Tentative ID       | MW     | MolForm          | CAS#        | Qual  |
| 1         | Nonadecane         | 268    | C19H40           | 000629-92-5 | 91    |
| 2         | Eicosane, 7-hexyl- | 366    | C26H54           | 055333-99-8 | 86    |
| 3         | Heneicosane        | 296    | C21H44           | 000629-94-7 | 83    |
| 4         | Tetracosane        | 338    | C24H50           | 000646-31-1 | 83    |
| 5         | Hexatriacontane    | 507    | C36H74           | 000630-06-8 | 83    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF043015\  
Data File : BF078879.D  
Acq On : 1 May 2015 6:21  
Operator : TP/IZ  
Sample : G2032-04  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B2(5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

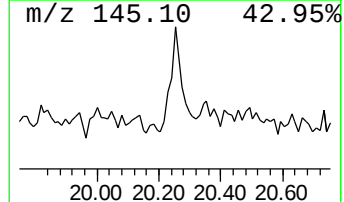
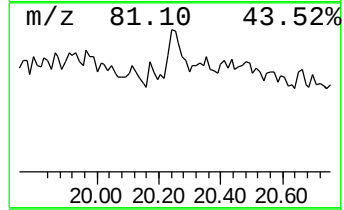
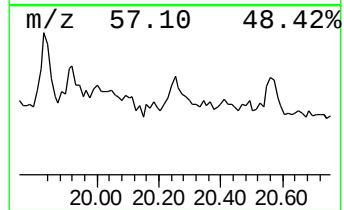
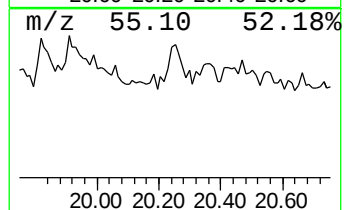
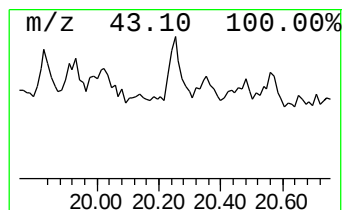
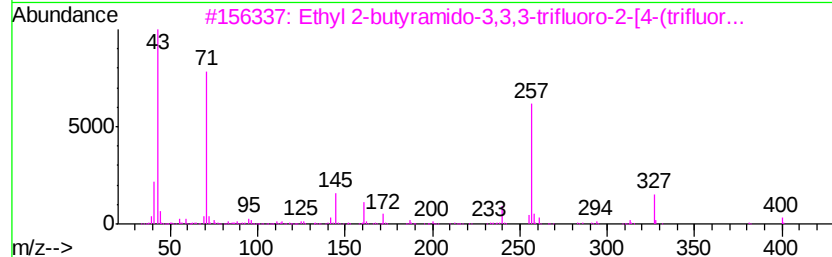
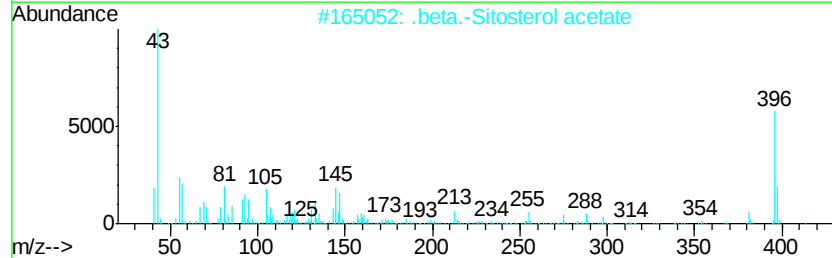
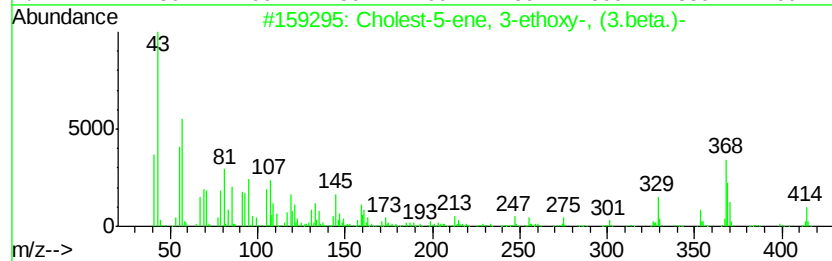
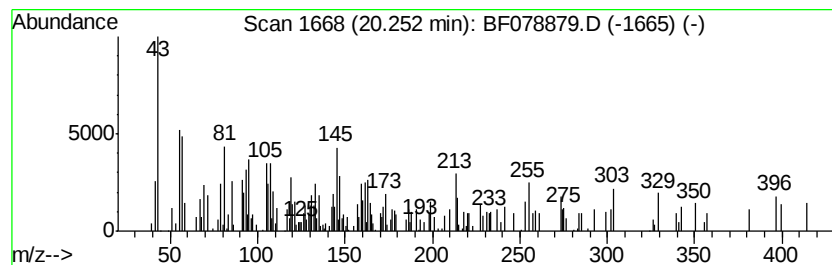
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown20.25 Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.25 | 3.17 ng | 162700 | Perylene-d12     | 18.08 |

| Hit# | of | Tentative ID                         | MW  | MolForm      | CAS#         | Qual |
|------|----|--------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Cholest-5-ene, 3-ethoxy-, (3.beta... | 414 | C29H50O      | 000986-19-6  | 38   |
| 2    |    | .beta.-Sitosterol acetate            | 456 | C31H52O2     | 000915-05-9  | 10   |
| 3    |    | Ethyl 2-butyramido-3,3,3-trifluo...  | 400 | C16H18F6N2O3 | 1000224-15-9 | 9    |
| 4    |    | 3.beta.-Acetoxy-16-isothiocyana...   | 415 | C24H33N3O3S  | 006084-08-8  | 7    |
| 5    |    | Stigmastan-3,5-diene                 | 396 | C29H48       | 1000214-16-4 | 2    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF043015\  
 Data File : BF078879.D  
 Acq On : 1 May 2015 6:21  
 Operator : TP/IZ  
 Sample : G2032-04  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B2(5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF043015.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 |
| unknown1.54          | 1.54  | 318.6   | ng    | 8625750  | 1                     | 7.37  | 541540  | 20.0 |
| Butane, 2-methoxy... | 1.87  | 15.0    | ng    | 405514   | 1                     | 7.37  | 541540  | 20.0 |
| 2-Pentanone, 4-hy... | 5.19  | 12.1    | ng    | 326794   | 1                     | 7.37  | 541540  | 20.0 |
| unknown7.08          | 7.08  | 85.3    | ng    | 2308790  | 1                     | 7.37  | 541540  | 20.0 |
| Tridecanoic acid     | 13.68 | 2.0     | ng    | 87931    | 4                     | 12.97 | 876946  | 20.0 |
| Cyclic octaatomic... | 14.57 | 2.1     | ng    | 91001    | 4                     | 12.97 | 876946  | 20.0 |
| Phenol, 4,4'-(1-m... | 14.83 | 7.7     | ng    | 402643   | 5                     | 16.27 | 1051710 | 20.0 |
| 1-Nonadecene         | 16.14 | 7.6     | ng    | 397646   | 5                     | 16.27 | 1051710 | 20.0 |
| Heptadecane          | 16.98 | 2.8     | ng    | 149205   | 5                     | 16.27 | 1051710 | 20.0 |
| Nonadecane           | 17.79 | 4.0     | ng    | 205573   | 6                     | 18.08 | 1025960 | 20.0 |
| unknown20.25         | 20.25 | 3.2     | ng    | 162700   | 6                     | 18.08 | 1025960 | 20.0 |