

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011615\  
 Data File : BF076917.D  
 Acq On : 16 Jan 2015 20:13  
 Operator : IZ / TP  
 Sample : G1080-03  
 Misc : W  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 FIELDBLANK-1

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-BF011415.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         | 1.384       | 23            | 26          | 31           | rBV      | 9961           | 19368         | 1.25%           | 0.195%        |
| 2         | 1.452       | 31            | 32          | 46           | rVB      | 60680          | 138621        | 8.92%           | 1.396%        |
| 3         | 4.104       | 262           | 264         | 273          | rBV      | 45205          | 102790        | 6.62%           | 1.035%        |
| 4         | 5.030       | 342           | 345         | 358          | rVB      | 189212         | 348852        | 22.46%          | 3.514%        |
| 5         | 7.362       | 547           | 549         | 555          | rBV      | 115193         | 183961        | 11.84%          | 1.853%        |
| 6         | 7.465       | 555           | 558         | 563          | rVB      | 526433         | 824331        | 53.07%          | 8.303%        |
| 7         | 7.967       | 599           | 602         | 607          | rVB      | 117124         | 200696        | 12.92%          | 2.021%        |
| 8         | 8.299       | 627           | 631         | 634          | rBV      | 507643         | 827720        | 53.29%          | 8.337%        |
| 9         | 9.270       | 712           | 716         | 719          | rBV      | 453226         | 675486        | 43.49%          | 6.804%        |
| 10        | 10.882      | 853           | 857         | 860          | rBV      | 186964         | 285675        | 18.39%          | 2.877%        |
| 11        | 11.396      | 898           | 902         | 904          | rVB      | 21024          | 32994         | 2.12%           | 0.332%        |
| 12        | 12.139      | 965           | 967         | 971          | rBV      | 33396          | 50964         | 3.28%           | 0.513%        |
| 13        | 13.534      | 1085          | 1089        | 1092         | rBV      | 946167         | 1393728       | 89.72%          | 14.038%       |
| 14        | 13.957      | 1123          | 1126        | 1130         | rBV      | 193143         | 303379        | 19.53%          | 3.056%        |
| 15        | 14.757      | 1193          | 1196        | 1200         | rVB      | 53624          | 78339         | 5.04%           | 0.789%        |
| 16        | 15.008      | 1215          | 1218        | 1222         | rVB      | 213660         | 324189        | 20.87%          | 3.265%        |
| 17        | 15.260      | 1236          | 1240        | 1243         | rVB      | 71642          | 110626        | 7.12%           | 1.114%        |
| 18        | 15.362      | 1246          | 1249        | 1252         | rBV      | 45213          | 64755         | 4.17%           | 0.652%        |
| 19        | 16.311      | 1329          | 1332        | 1334         | rVB      | 38480          | 52435         | 3.38%           | 0.528%        |
| 20        | 16.471      | 1342          | 1346        | 1349         | rVB      | 308833         | 476844        | 30.70%          | 4.803%        |
| 21        | 16.677      | 1361          | 1364        | 1367         | rBV      | 648012         | 853750        | 54.96%          | 8.599%        |
| 22        | 17.774      | 1457          | 1460        | 1463         | rVB      | 309816         | 335217        | 21.58%          | 3.376%        |
| 23        | 17.957      | 1474          | 1476        | 1479         | rBV      | 29854          | 33462         | 2.15%           | 0.337%        |
| 24        | 18.460      | 1518          | 1520        | 1522         | rBV      | 21042          | 27045         | 1.74%           | 0.272%        |
| 25        | 20.003      | 1652          | 1655        | 1658         | rBV      | 1438945        | 1553356       | 100.00%         | 15.646%       |
| 26        | 21.272      | 1763          | 1766        | 1768         | rBV      | 245285         | 337094        | 21.70%          | 3.395%        |
| 27        | 22.906      | 1906          | 1909        | 1912         | rBV      | 250791         | 292657        | 18.84%          | 2.948%        |

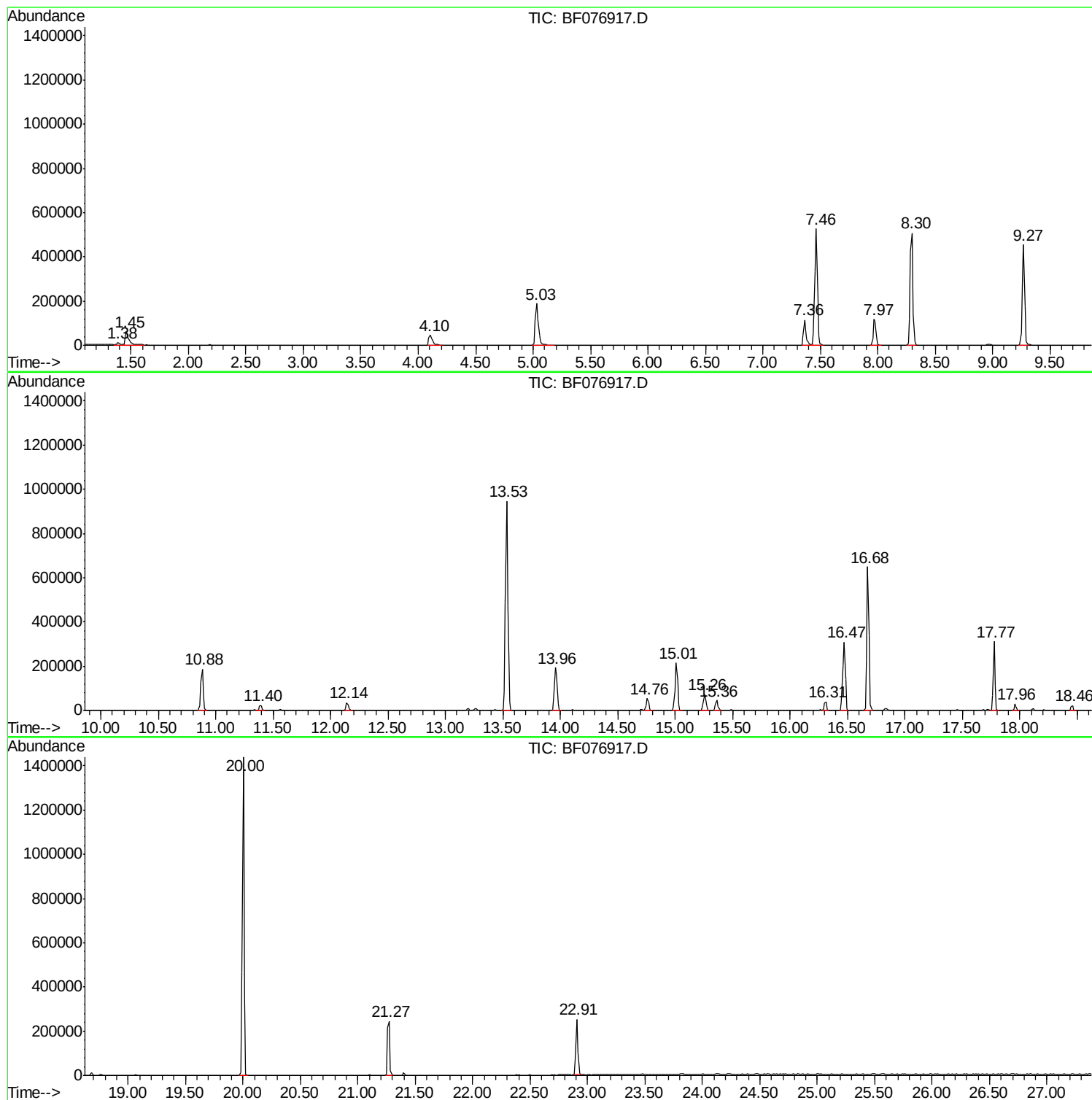
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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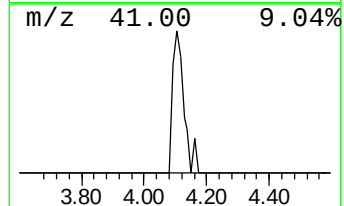
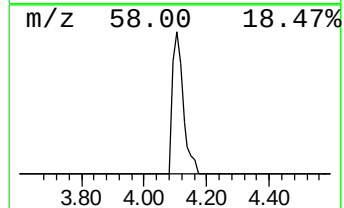
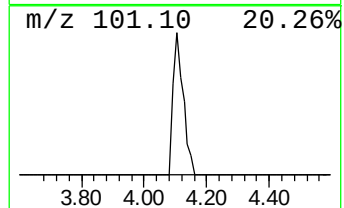
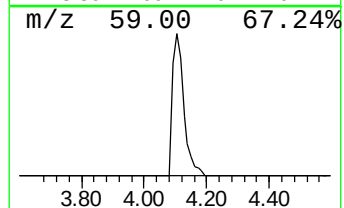
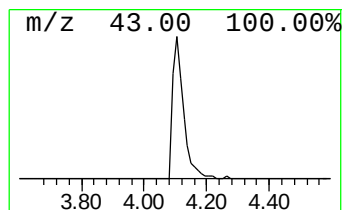
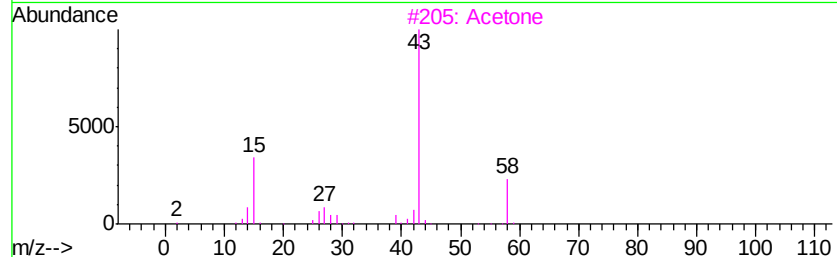
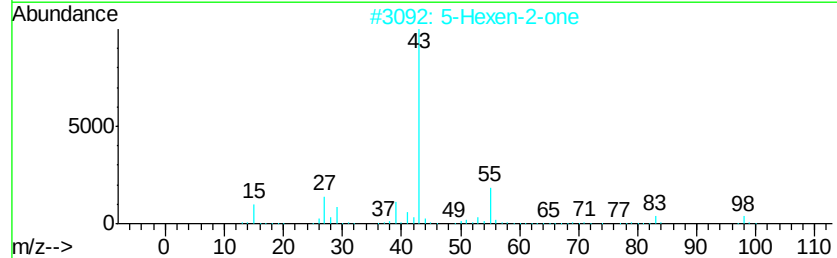
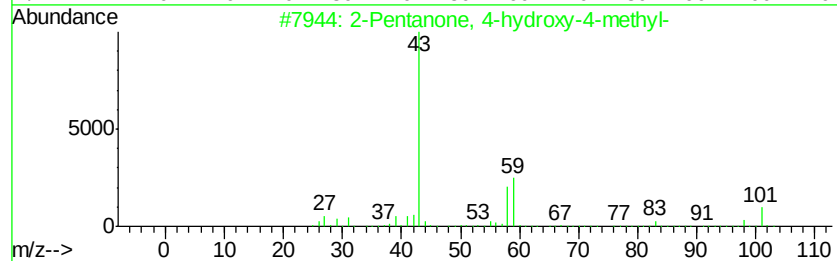
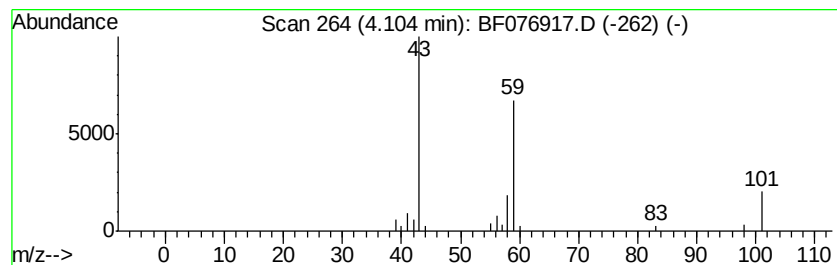
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.10 | 10.24 ng | 102790 | 1,4-Dichlorobenzene-d4 | 7.98 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 64   |
| 2    |    | 5-Hexen-2-one                    | 98  | C6H10O  | 000109-49-9 | 9    |
| 3    |    | Acetone                          | 58  | C3H6O   | 000067-64-1 | 9    |
| 4    |    | Guanidine                        | 59  | CH5N3   | 000113-00-8 | 9    |
| 5    |    | 4-Hydroxy-3-hexanone             | 116 | C6H12O2 | 004984-85-4 | 9    |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

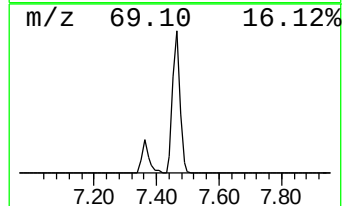
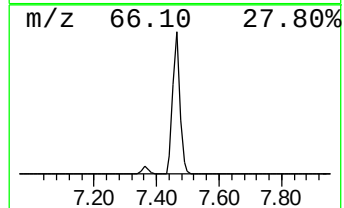
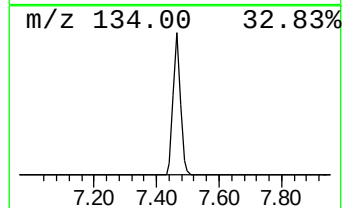
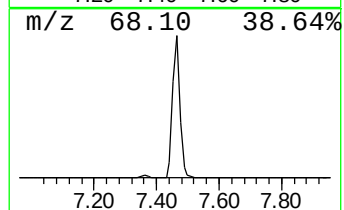
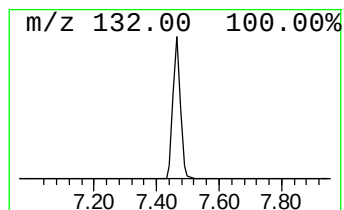
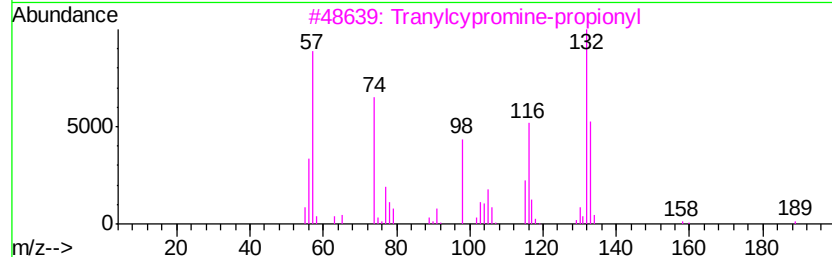
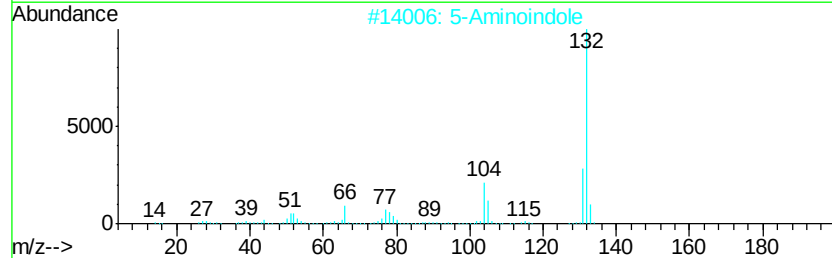
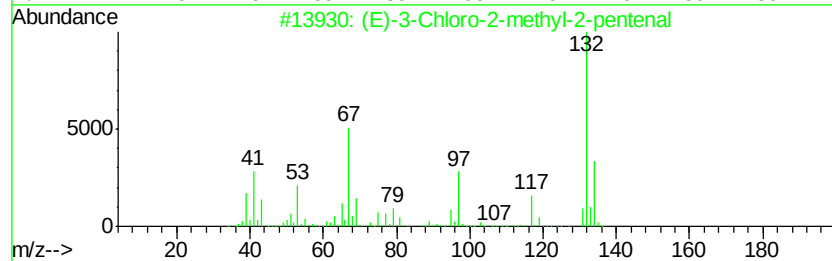
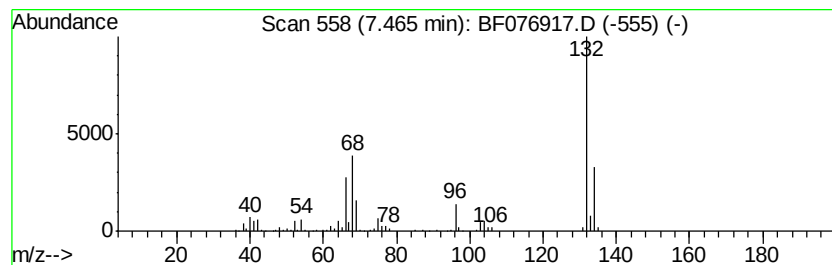
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown 7.46 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.46 | 82.15 ng | 824331 | 1,4-Dichlorobenzene-d4 | 7.98 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 2    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 5    |    | 4-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-60-2  | 9    |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
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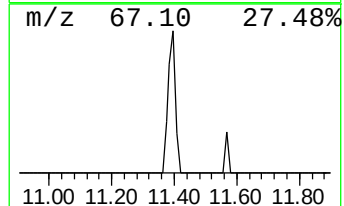
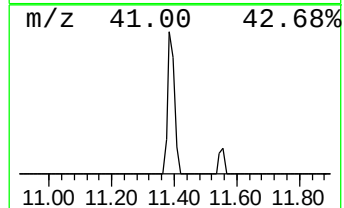
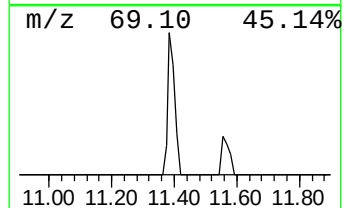
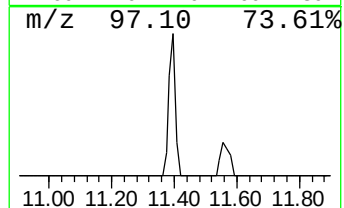
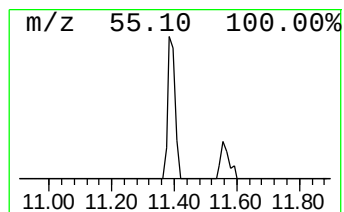
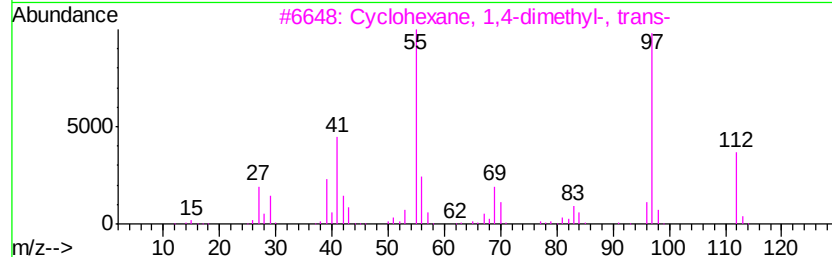
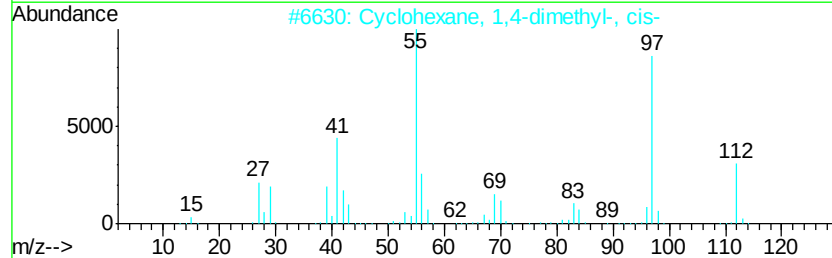
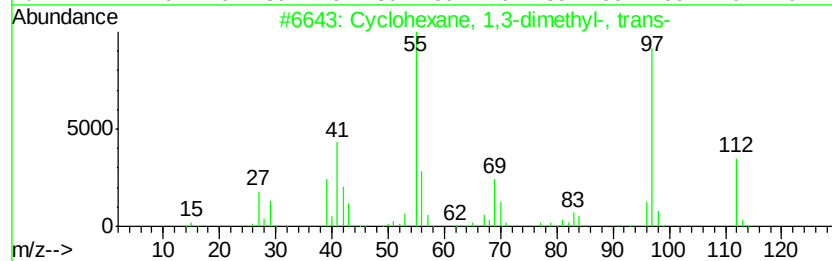
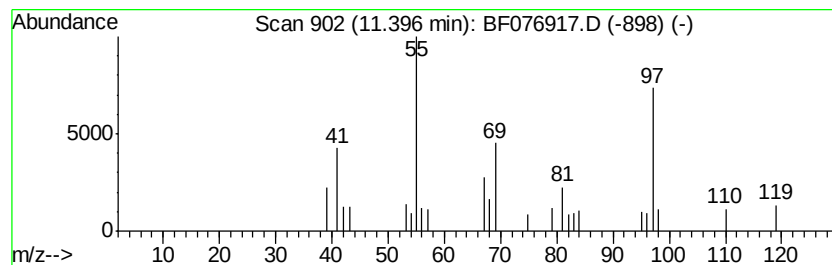
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 12

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.40 | 2.31 ng | 32994 | Napthalene-d8    | 10.88 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 59   |
| 2    |    | Cyclohexane, 1,4-dimethyl-, cis-   | 112 | C8H16   | 000624-29-3 | 53   |
| 3    |    | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 53   |
| 4    |    | Cyclohexane, 1,2-dimethyl-, trans- | 112 | C8H16   | 006876-23-9 | 50   |
| 5    |    | 2-Pentene, 1-bromo-3,4-dimethyl-   | 176 | C7H13Br | 000922-99-6 | 50   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

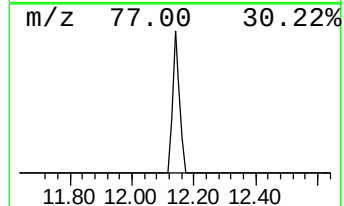
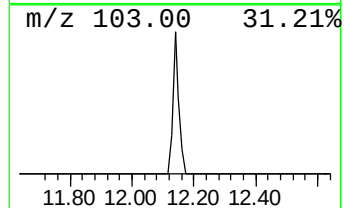
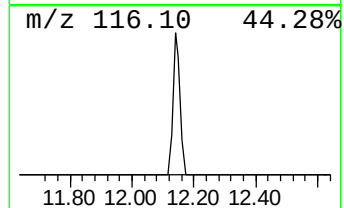
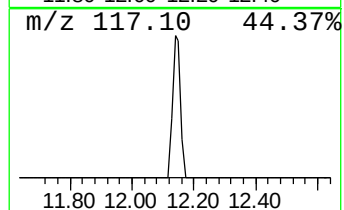
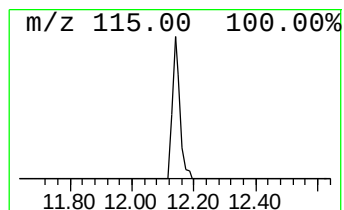
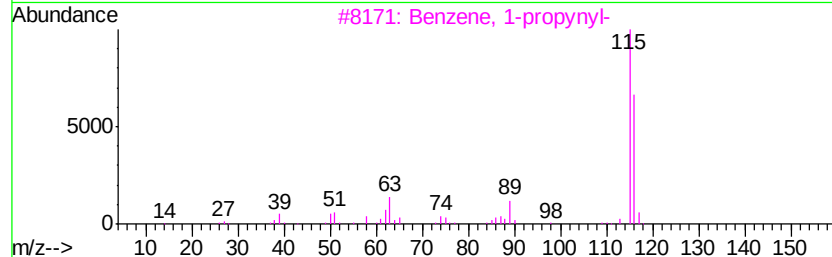
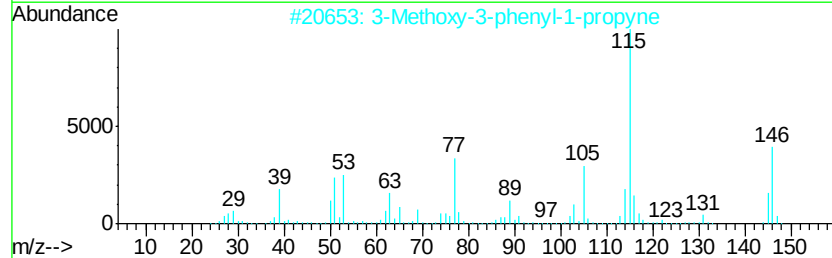
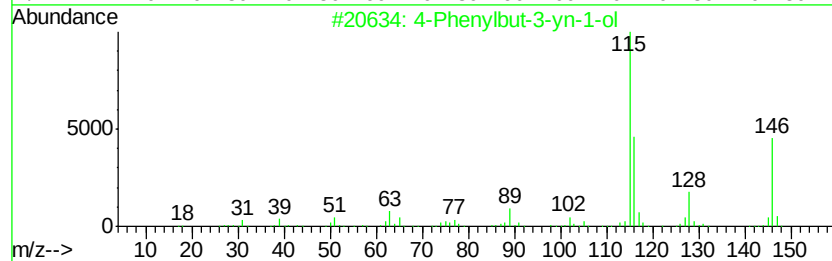
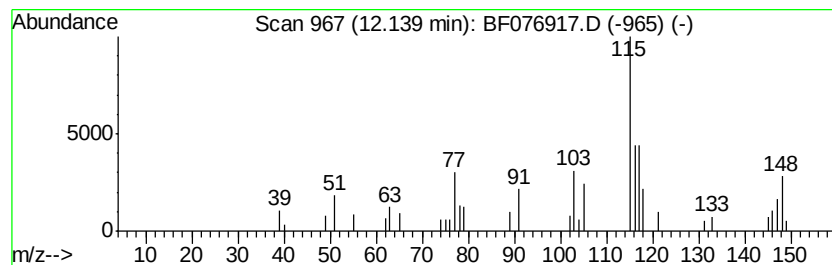
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown12.14 Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.14 | 3.57 ng | 50964 | Naphthalene-d8   | 10.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 4-Phenylbut-3-yn-1-ol               | 146 | C10H10O | 010229-11-5  | 38   |
| 2    |    | 3-Methoxy-3-phenyl-1-propyne        | 146 | C10H10O | 1000217-10-1 | 38   |
| 3    |    | Benzene, 1-propynyl-                | 116 | C9H8    | 000673-32-5  | 35   |
| 4    |    | Cyclohexane, 2,4-cyclopentadien-... | 146 | C11H14  | 003141-04-6  | 23   |
| 5    |    | 1,4-Epoxy naphthalene, 1,4-dihydro- | 144 | C10H8O  | 000573-57-9  | 23   |



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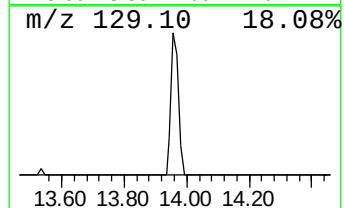
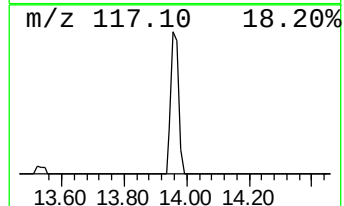
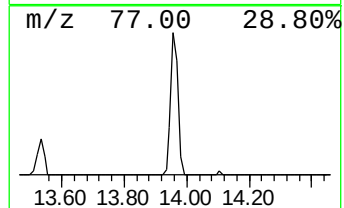
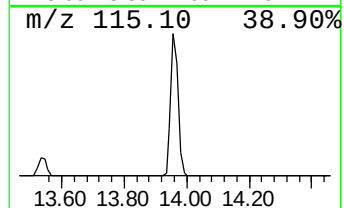
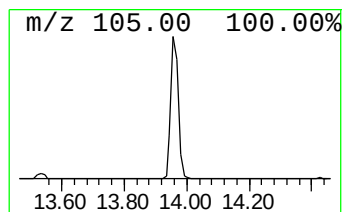
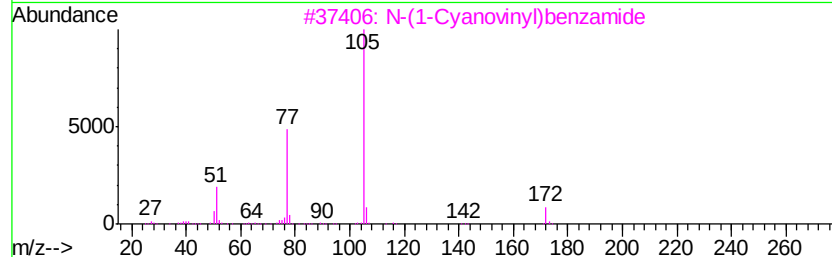
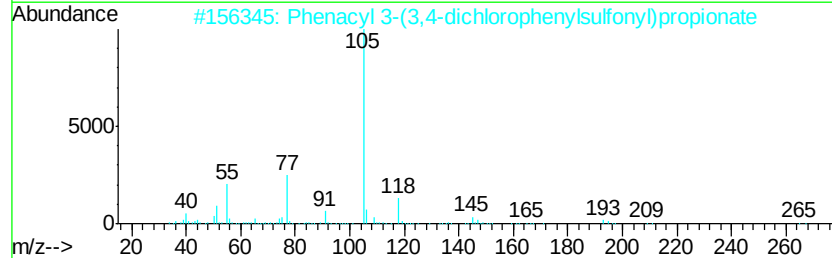
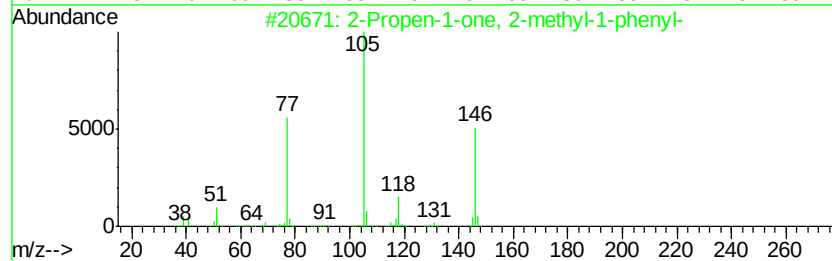
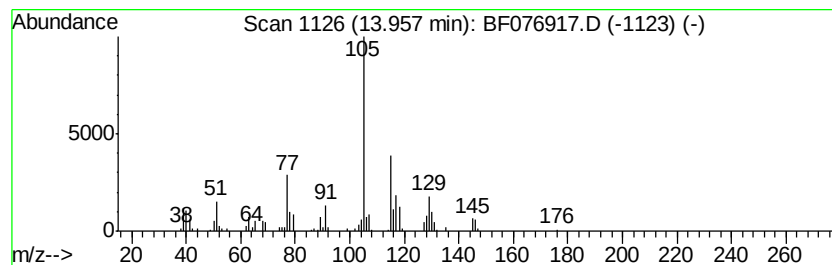
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown13.96 Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.96 | 18.72 ng | 303379 | Acenaphthene-d10 | 15.01 |

| Hit# | of | Tentative ID                                      | MW  | MolForm      | CAS#         | Qual |
|------|----|---------------------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 2-Propen-1-one, 2-methyl-1-phenyl-                | 146 | C10H10O      | 000769-60-8  | 27   |
| 2    |    | Phenacyl 3-(3,4-dichlorophenyl)sulfonylpropionate | 400 | C17H14Cl2O5S | 1000225-01-1 | 22   |
| 3    |    | N-(1-Cyanovinyl)benzamide                         | 172 | C10H8N2O     | 1000186-19-1 | 10   |
| 4    |    | Benzene, (1-bromoethyl)-                          | 184 | C8H9Br       | 000585-71-7  | 10   |
| 5    |    | 1,2(4H)-Oxazine-3-ol, 5,6-dihydro-                | 219 | C11H9N2O4    | 1000253-25-6 | 10   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011615\  
Data File : BF076917.D  
Acq On : 16 Jan 2015 20:13  
Operator : IZ / TP  
Sample : G1080-03  
Misc : W  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

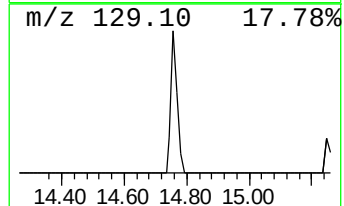
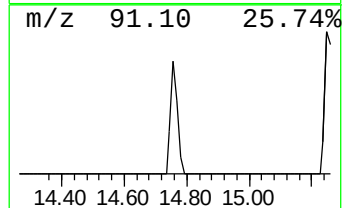
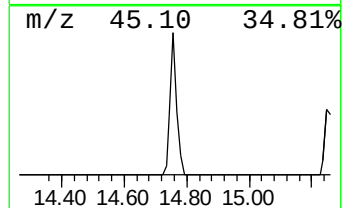
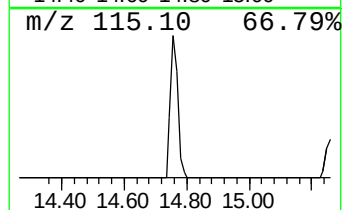
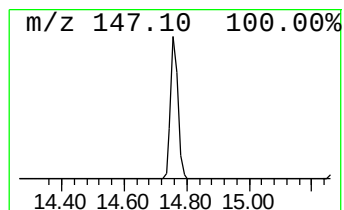
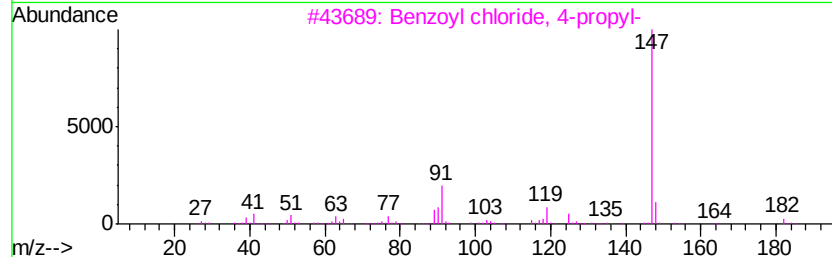
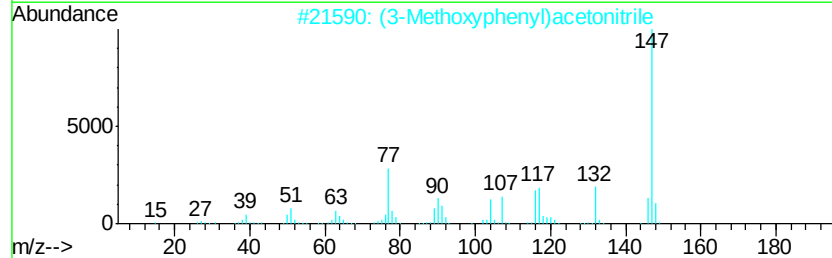
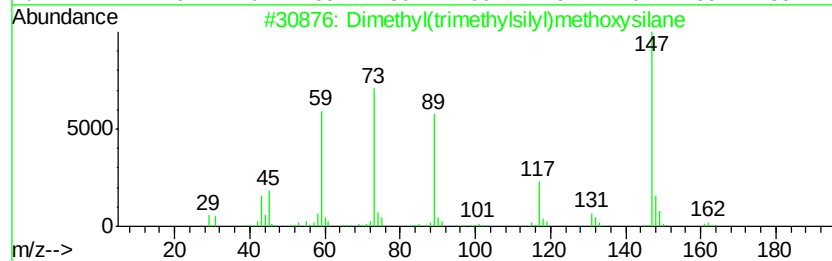
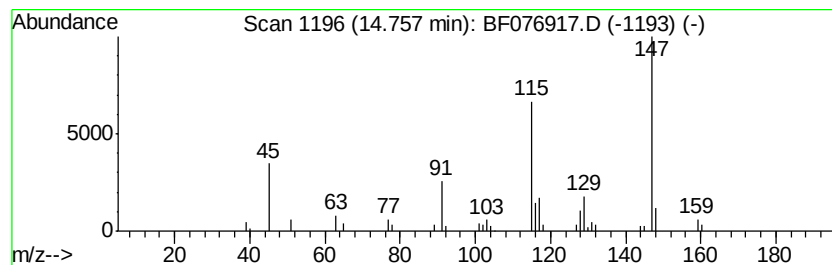
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown14.76 Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.76 | 4.83 ng | 78339 | Acenaphthene-d10 | 15.01 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|---------|---|-------------------------------------|-----|-------------|--------------|------|
| 1       |   | Dimethyl(trimethylsilyl)methoxys... | 162 | C6H18OSi2   | 018107-29-4  | 43   |
| 2       |   | (3-Methoxyphenyl)acetonitrile       | 147 | C9H9NO      | 019924-43-7  | 43   |
| 3       |   | Benzoyl chloride, 4-propyl-         | 182 | C10H11ClO   | 052710-27-7  | 35   |
| 4       |   | Benzoic acid, 4-propyl-, 4-cyano... | 265 | C17H15NO2   | 056131-49-8  | 25   |
| 5       |   | 2-(5-Methyl-[1,3,4]thiadiazol-2-... | 292 | C14H16N2OS2 | 1000296-87-8 | 25   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF011615\  
Data File : BF076917.D  
Acq On : 16 Jan 2015 20:13  
Operator : IZ / TP  
Sample : G1080-03  
Misc : W  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

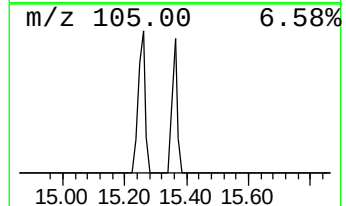
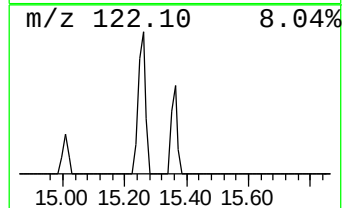
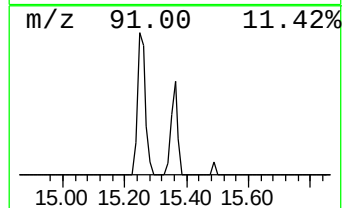
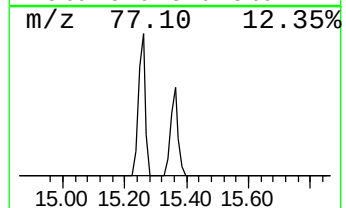
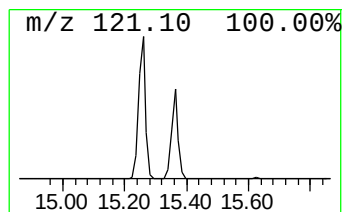
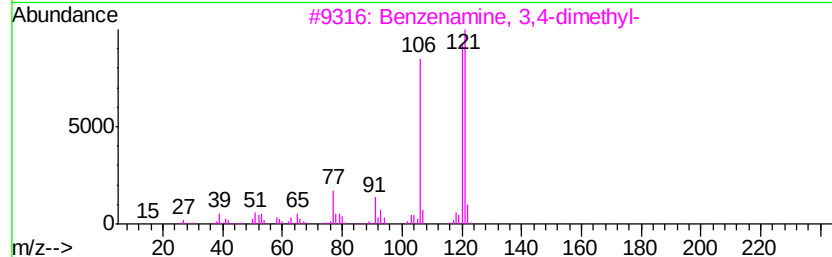
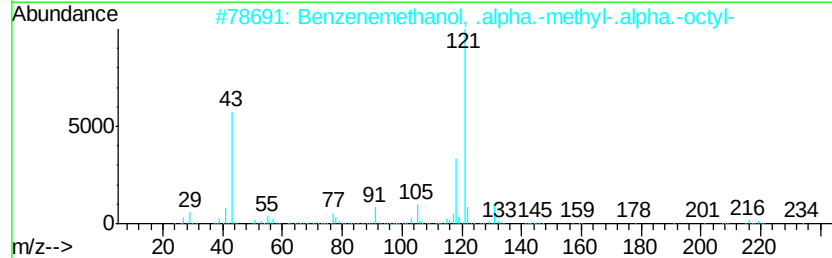
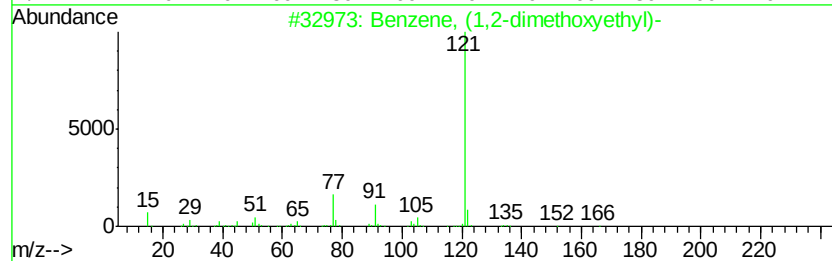
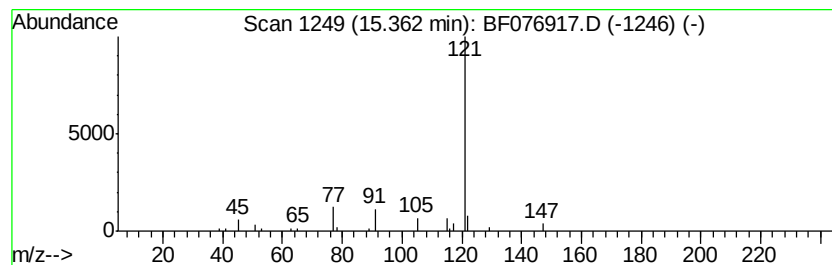
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzenemethanol, .alpha.-me... Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.36 | 3.99 ng | 64755 | Acenaphthene-d10 | 15.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzene, (1,2-dimethoxyethyl)-      | 166 | C10H14O2 | 004013-37-0 | 72   |
| 2    |    | Benzenemethanol, .alpha.-methyl-... | 234 | C16H26O  | 021078-96-6 | 56   |
| 3    |    | Benzenamine, 3,4-dimethyl-          | 121 | C8H11N   | 000095-64-7 | 50   |
| 4    |    | Benzaldehyde dimethyl acetal        | 152 | C9H12O2  | 001125-88-8 | 50   |
| 5    |    | Benzenamine, 3,5-dimethyl-          | 121 | C8H11N   | 000108-69-0 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011615\  
Data File : BF076917.D  
Acq On : 16 Jan 2015 20:13  
Operator : IZ / TP  
Sample : G1080-03  
Misc : W  
ALS Vial : 10 Sample Multiplier: 1

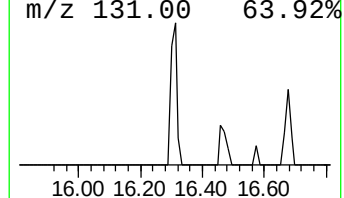
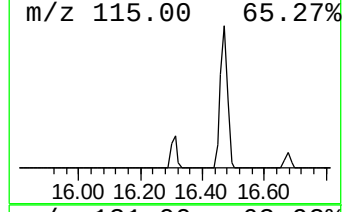
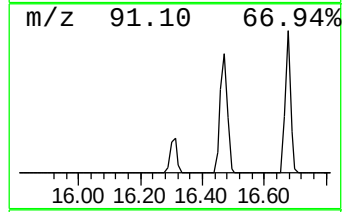
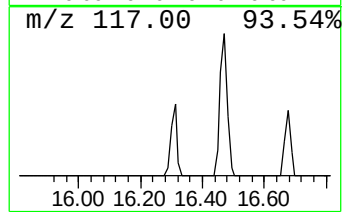
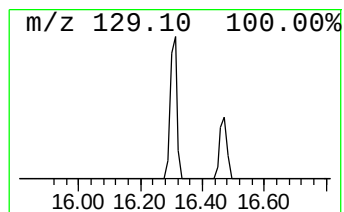
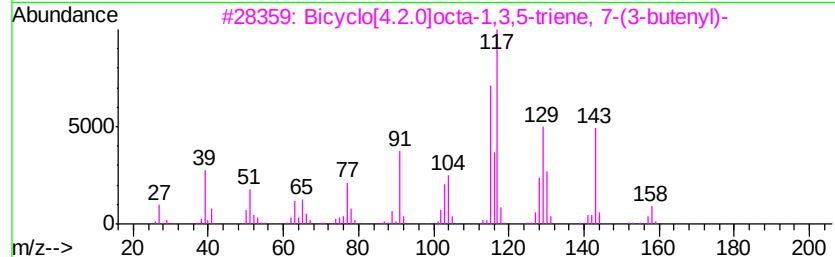
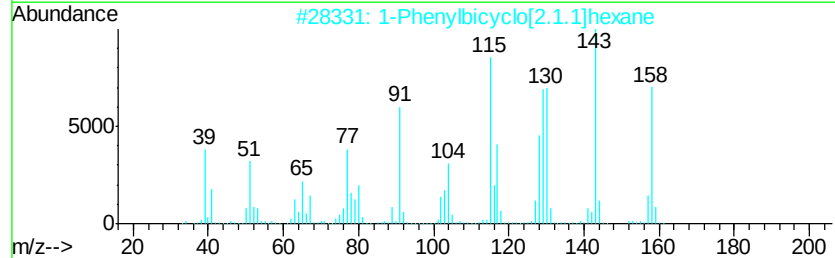
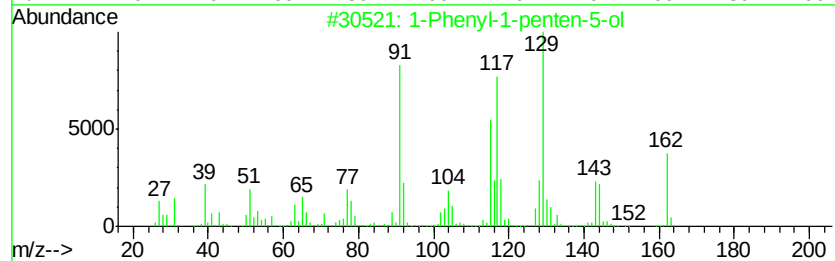
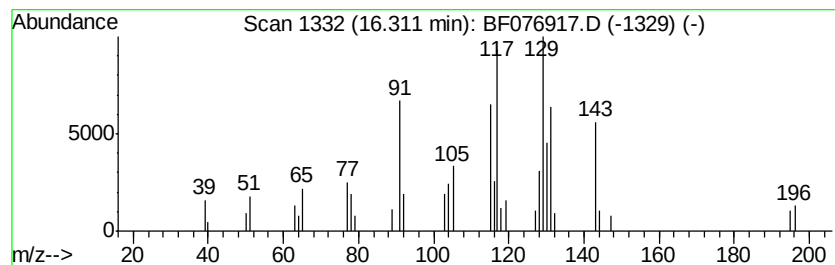
Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 1-Phenyl-1-penten-5-ol Concentration Rank 11

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 16.31   | 3.23 ng                             | 52435        | Acenaphthene-d10 | 15.01     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1-Phenyl-1-penten-5-ol              | 162 C11H14O  | 037464-86-1      | 59        |
| 2       | 1-Phenylbicyclo[2.1.1]hexane        | 158 C12H14   | 029508-78-9      | 43        |
| 3       | Bicyclo[4.2.0]octa-1,3,5-triene,... | 158 C12H14   | 122057-61-8      | 43        |
| 4       | Bicyclo[4.2.0]octa-1,3,5-triene,... | 144 C11H12   | 071721-26-1      | 30        |
| 5       | 5H-Benzocycloheptene, 6,7-dihydro-  | 144 C11H12   | 007125-62-4      | 27        |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011615\  
Data File : BF076917.D  
Acq On : 16 Jan 2015 20:13  
Operator : IZ / TP  
Sample : G1080-03  
Misc : W  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
FIELDBLANK-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

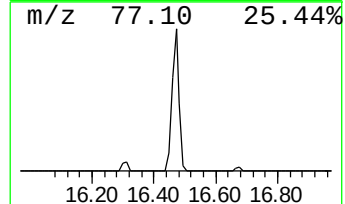
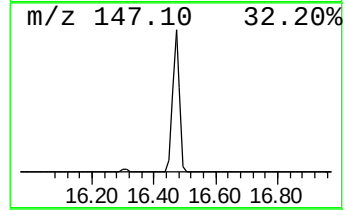
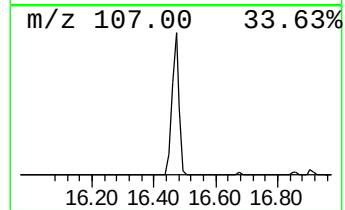
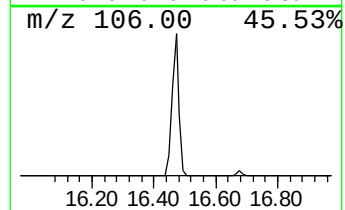
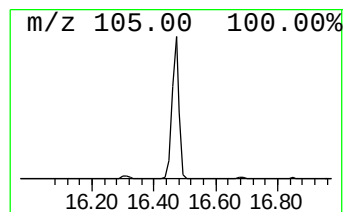
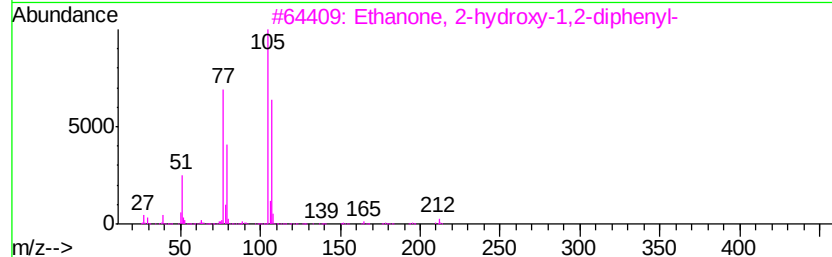
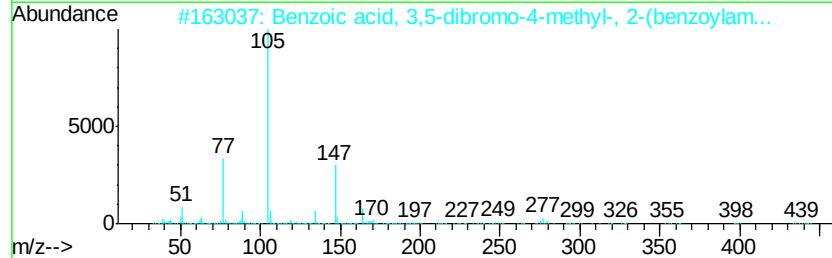
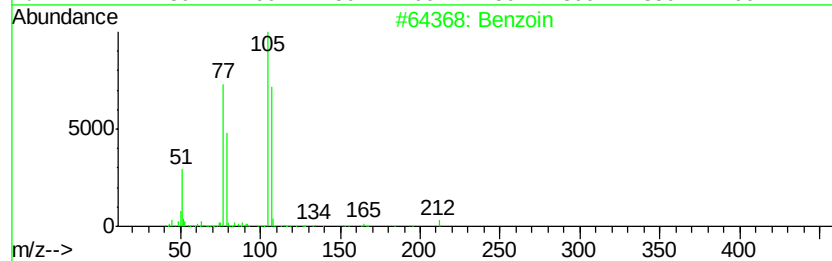
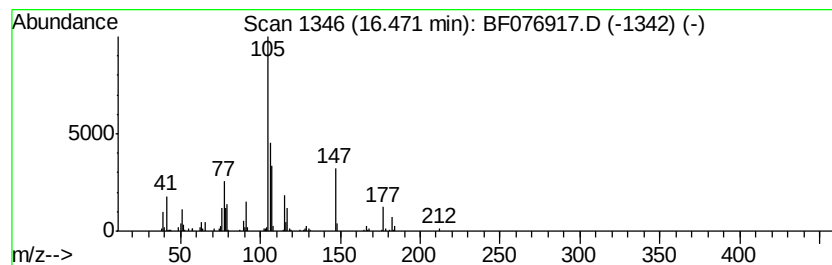
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown16.47 Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.47 | 28.45 ng | 476844 | Phenanthrene-d10 | 17.77 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Benzoin                             | 212 | C14H12O2     | 000579-44-2  | 27   |
| 2    |    | Benzoic acid, 3,5-dibromo-4-meth... | 439 | C17H15Br2NO3 | 1000258-73-5 | 27   |
| 3    |    | Ethanone, 2-hydroxy-1,2-diphenyl-   | 212 | C14H12O2     | 000119-53-9  | 27   |
| 4    |    | 2-Cyclopropyl-2-nitro-1-phenyl-e... | 207 | C11H13NO3    | 1000194-91-5 | 25   |
| 5    |    | Benzamide, N-(1,1-dimethylethyl)-   | 177 | C11H15NO     | 005894-65-5  | 22   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF011615\  
Data File : BF076917.D  
Acq On : 16 Jan 2015 20:13  
Operator : IZ / TP  
Sample : G1080-03  
Misc : W  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.10  | 10.2    | ng    | 102790   | 1                     | 7.98  | 200696 | 20.0 |
| unknown7.46          | 7.46  | 82.2    | ng    | 824331   | 1                     | 7.98  | 200696 | 20.0 |
| Cyclohexane, 1,3-... | 11.40 | 2.3     | ng    | 32994    | 2                     | 10.88 | 285675 | 20.0 |
| unknown12.14         | 12.14 | 3.6     | ng    | 50964    | 2                     | 10.88 | 285675 | 20.0 |
| unknown13.96         | 13.96 | 18.7    | ng    | 303379   | 3                     | 15.01 | 324189 | 20.0 |
| unknown14.76         | 14.76 | 4.8     | ng    | 78339    | 3                     | 15.01 | 324189 | 20.0 |
| Benzenemethanol, ... | 15.36 | 4.0     | ng    | 64755    | 3                     | 15.01 | 324189 | 20.0 |
| 1-Phenyl-1-penten... | 16.31 | 3.2     | ng    | 52435    | 3                     | 15.01 | 324189 | 20.0 |
| unknown16.47         | 16.47 | 28.4    | ng    | 476844   | 4                     | 17.77 | 335217 | 20.0 |