

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF072017\  
 Data File : BF096916.D  
 Acq On : 20 Jul 2017 20:09  
 Operator : SJ/JU  
 Sample : I4312-05  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 72-11987

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-BF071317.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         | 2.798       | 136           | 138         | 146          | rBV      | 50755          | 99630         | 3.98%           | 0.508%        |
| 2         | 4.739       | 463           | 468         | 478          | rVB      | 174760         | 259538        | 10.36%          | 1.323%        |
| 3         | 5.181       | 538           | 543         | 546          | rBV      | 2136297        | 2178913       | 86.99%          | 11.109%       |
| 4         | 6.228       | 716           | 721         | 724          | rBV      | 2404502        | 2504891       | 100.00%         | 12.771%       |
| 5         | 6.345       | 736           | 741         | 744          | rBV      | 2378582        | 2407023       | 96.09%          | 12.272%       |
| 6         | 6.569       | 775           | 779         | 787          | rBV      | 739849         | 621143        | 24.80%          | 3.167%        |
| 7         | 6.728       | 801           | 806         | 809          | rBV      | 1951421        | 1802328       | 71.95%          | 9.189%        |
| 8         | 6.751       | 809           | 810         | 821          | rVB      | 37517          | 45000         | 1.80%           | 0.229%        |
| 9         | 7.133       | 870           | 875         | 878          | rBV      | 1623161        | 1592207       | 63.56%          | 8.118%        |
| 10        | 7.269       | 888           | 898         | 902          | rBV      | 62245          | 79445         | 3.17%           | 0.405%        |
| 11        | 7.769       | 979           | 983         | 988          | rBV      | 33952          | 35080         | 1.40%           | 0.179%        |
| 12        | 7.851       | 993           | 997         | 1000         | rBV      | 865401         | 702586        | 28.05%          | 3.582%        |
| 13        | 8.933       | 1175          | 1181        | 1184         | rBV      | 2478897        | 2441333       | 97.46%          | 12.447%       |
| 14        | 9.063       | 1199          | 1203        | 1206         | rVB      | 37016          | 35142         | 1.40%           | 0.179%        |
| 15        | 9.327       | 1244          | 1248        | 1251         | rBV      | 473204         | 370149        | 14.78%          | 1.887%        |
| 16        | 9.598       | 1290          | 1294        | 1298         | rBV      | 756784         | 627801        | 25.06%          | 3.201%        |
| 17        | 10.386      | 1423          | 1428        | 1431         | rBV      | 1061753        | 988229        | 39.45%          | 5.039%        |
| 18        | 10.521      | 1448          | 1451        | 1458         | rBV      | 43572          | 46442         | 1.85%           | 0.237%        |
| 19        | 11.074      | 1541          | 1545        | 1548         | rBV      | 548848         | 457725        | 18.27%          | 2.334%        |
| 20        | 11.321      | 1584          | 1587        | 1590         | rBV      | 59787          | 55338         | 2.21%           | 0.282%        |
| 21        | 12.657      | 1810          | 1814        | 1818         | rBV      | 1317119        | 1254034       | 50.06%          | 6.394%        |
| 22        | 13.568      | 1965          | 1969        | 1974         | rBV      | 73339          | 82637         | 3.30%           | 0.421%        |
| 23        | 13.698      | 1987          | 1991        | 1995         | rBV      | 482546         | 455984        | 18.20%          | 2.325%        |
| 24        | 15.068      | 2219          | 2224        | 2230         | rBV      | 430898         | 470680        | 18.79%          | 2.400%        |

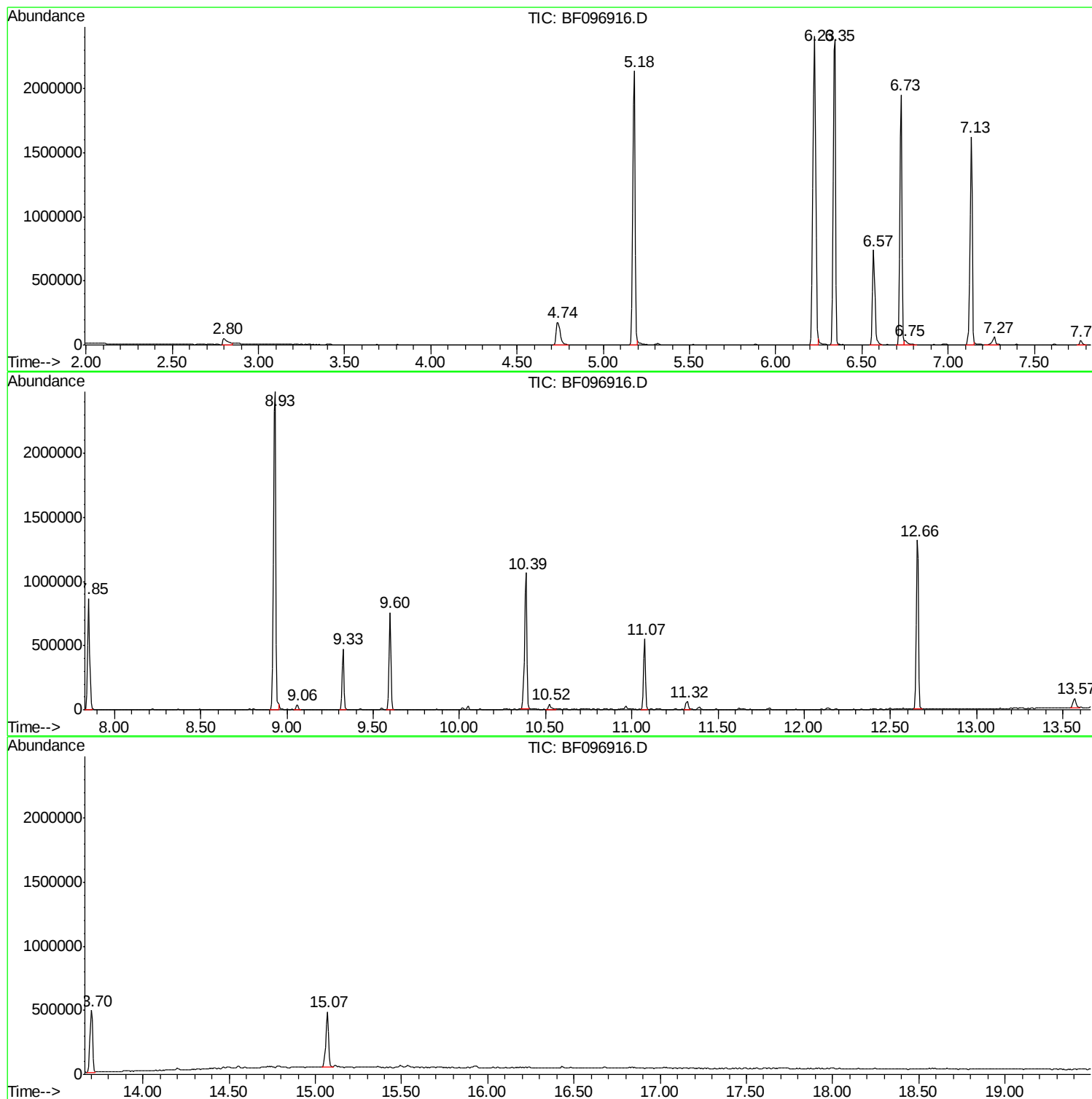
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ALS Vial : 15 Sample Multiplier: 1

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BNA\_F  
ClientSampleId :  
72-11987

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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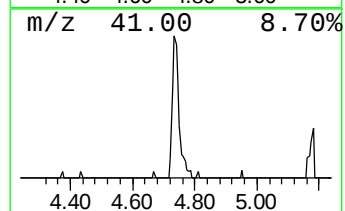
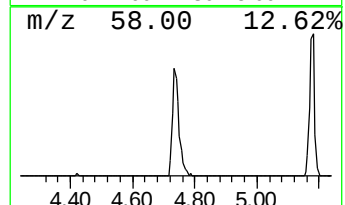
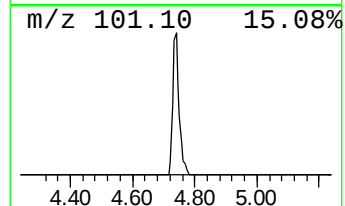
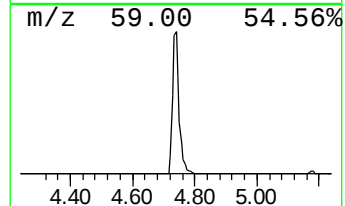
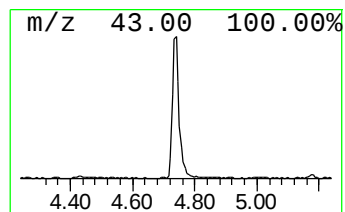
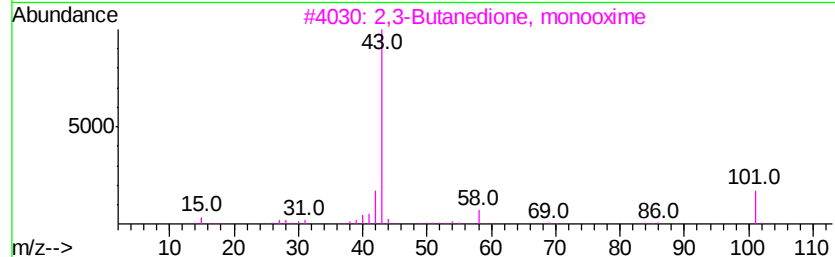
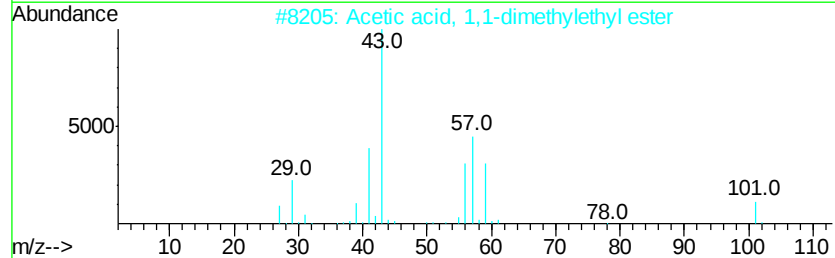
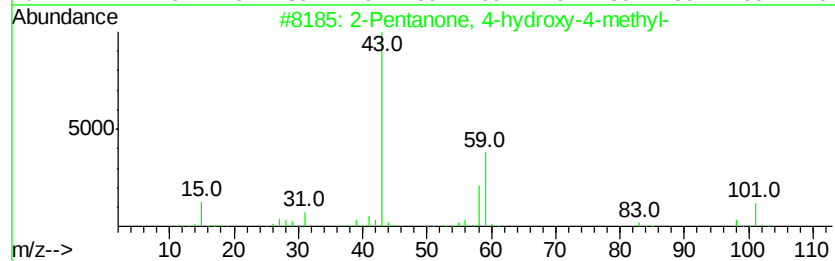
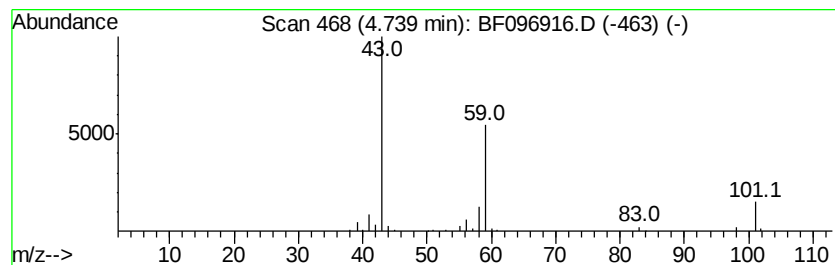
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.74 | 8.36 ng | 259538 | 1,4-Dichlorobenzene-d4 | 6.57 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 17   |
| 3    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 16   |
| 4    |    |   | 1,8-Nonanediol, 8-methyl-           | 174 | C10H22O2 | 054725-73-4 | 9    |
| 5    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 9    |



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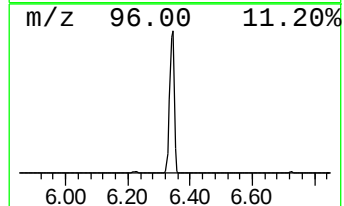
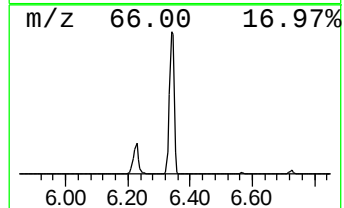
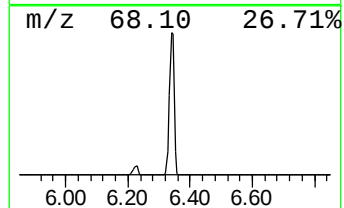
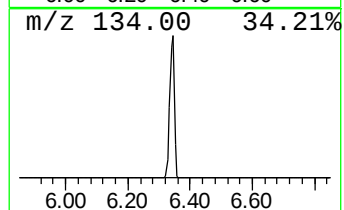
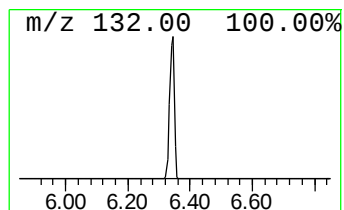
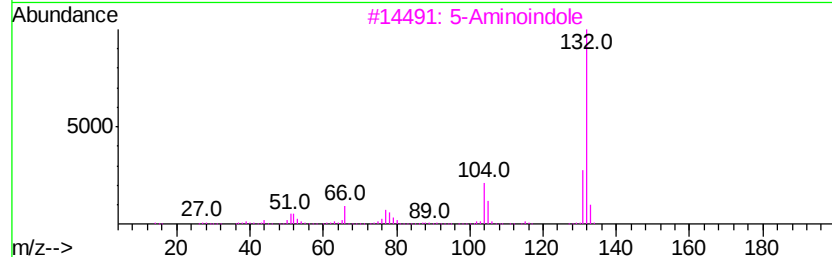
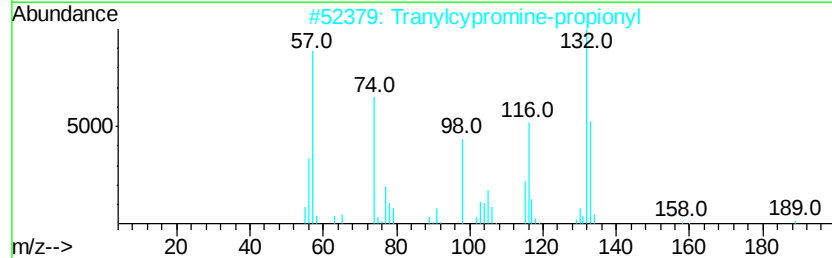
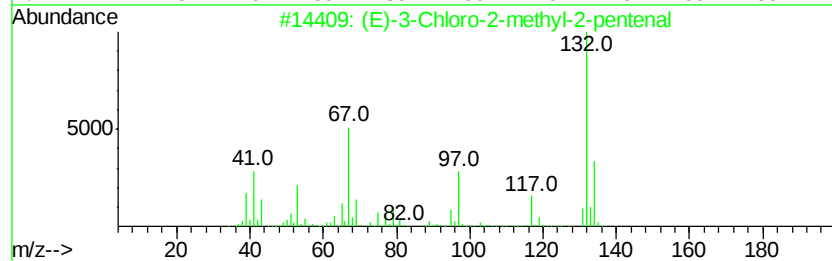
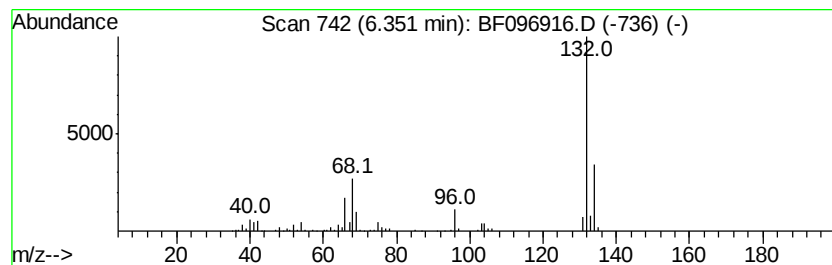
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.35 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.35 | 77.50 ng | 2407020 | 1,4-Dichlorobenzene-d4 | 6.57 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|----|----------------------------------|-----|----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO  | 031357-76-3  | 17   |
| 2    |    | Tranylcypromine-propionyl        | 189 | C12H15NO | 1000123-86-3 | 12   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2   | 005192-03-0  | 12   |
| 4    |    | Imidazole, 4,5-dicyano-1-methyl- | 132 | C6H4N4   | 019485-35-9  | 9    |
| 5    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3   | 000372-38-3  | 9    |



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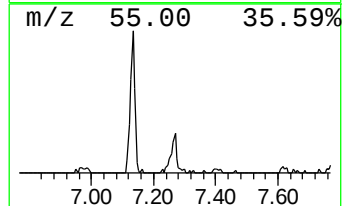
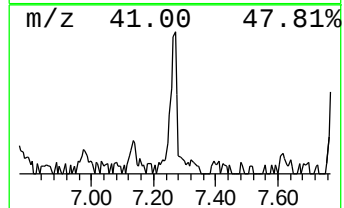
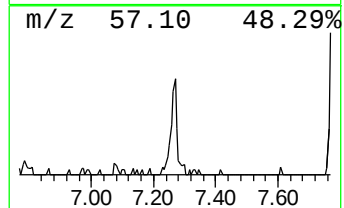
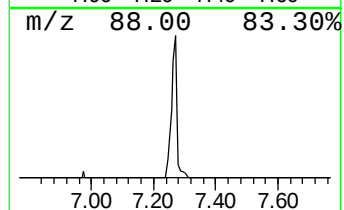
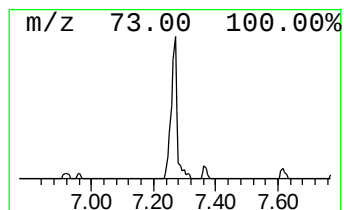
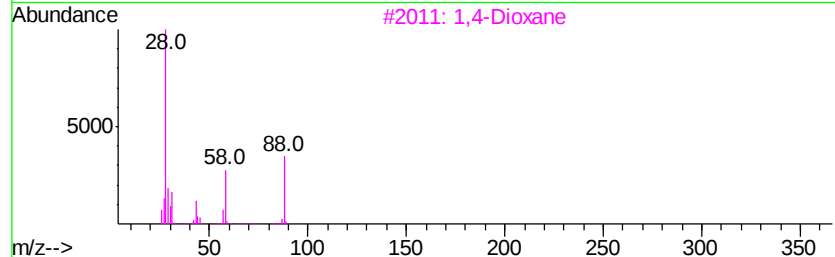
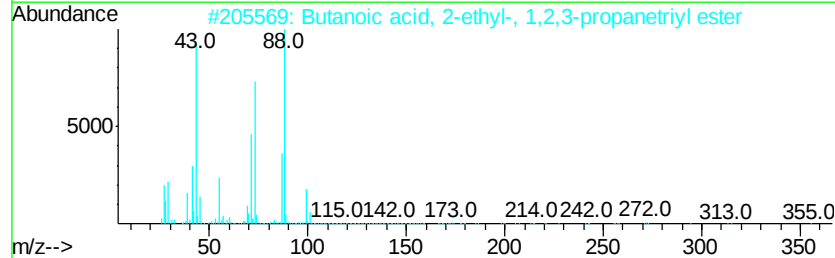
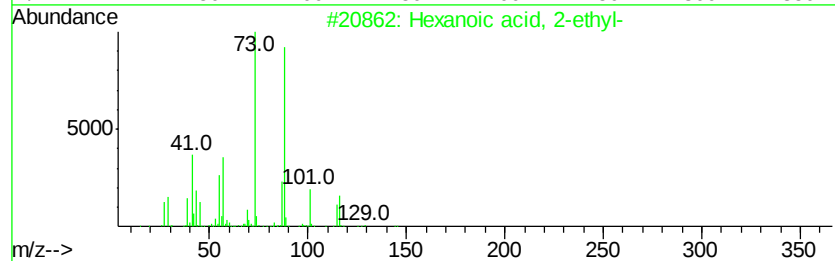
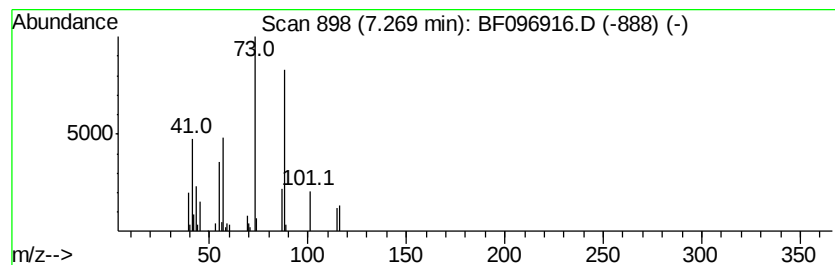
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 7

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 7.27 | 2.26 ng | 79445 | Naphthalene-d8   | 7.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2   | 000149-57-5 | 86   |
| 2    |    | Butanoic acid, 2-ethyl-, 1,2,3-p... | 386 | C21H38O6  | 056554-54-2 | 50   |
| 3    |    | 1,4-Dioxane                         | 88  | C4H8O2    | 000123-91-1 | 37   |
| 4    |    | Hexanamide, 6-(heptyloxy)-          | 229 | C13H27NO2 | 055191-08-7 | 33   |
| 5    |    | Methyl L-(+)-.beta.-hydroxyisobu... | 118 | C5H10O3   | 080657-57-4 | 25   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
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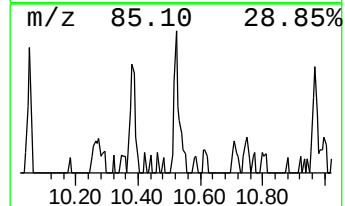
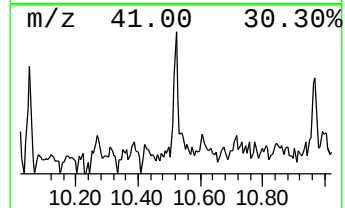
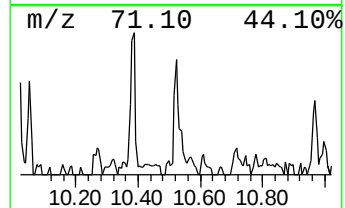
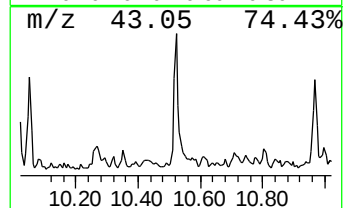
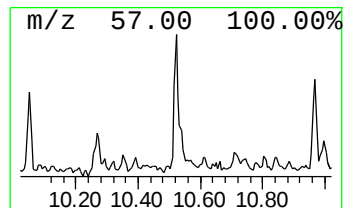
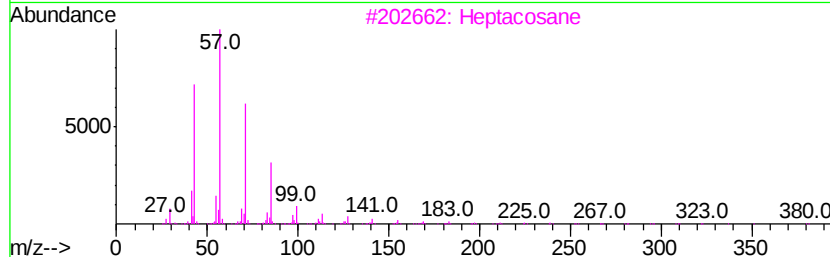
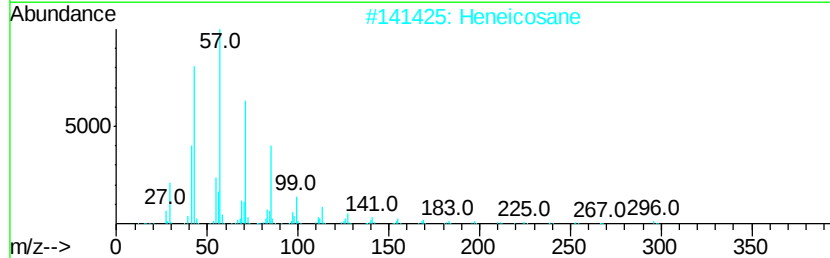
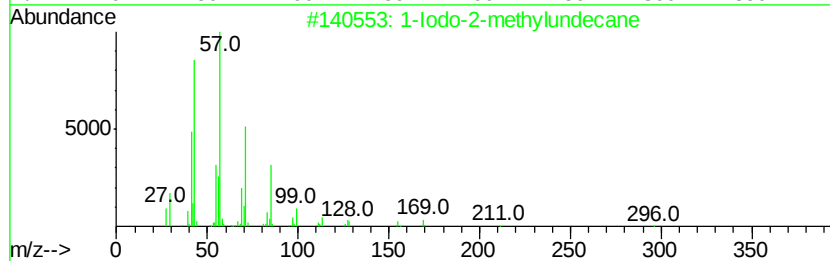
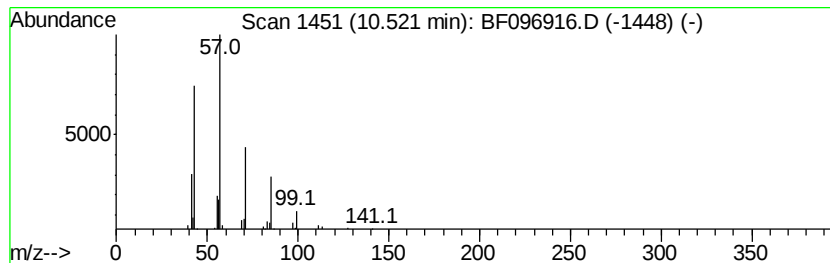
TIC Library : C:\DATABASE\NIST11.L  
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Peak Number 6 1-Iodo-2-methylundecane Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 10.52 | 2.03 ng | 46442 | Phenanthrene-d10 | 11.07 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 90   |
| 2    |    | Heneicosane             | 296 | C21H44  | 000629-94-7 | 83   |
| 3    |    | Heptacosane             | 380 | C27H56  | 000593-49-7 | 80   |
| 4    |    | Octadecane              | 254 | C18H38  | 000593-45-3 | 78   |
| 5    |    | Hexadecane              | 226 | C16H34  | 000544-76-3 | 78   |



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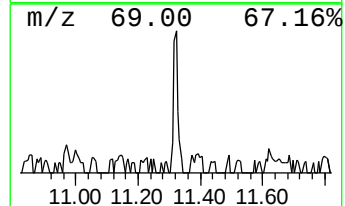
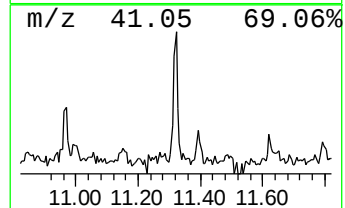
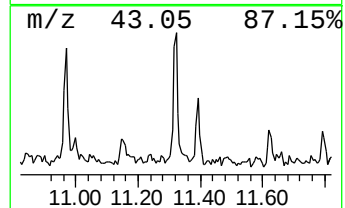
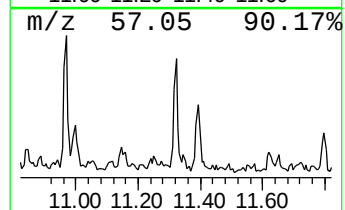
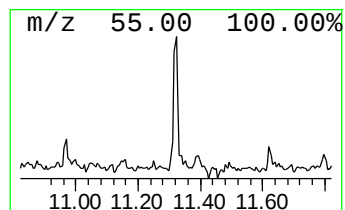
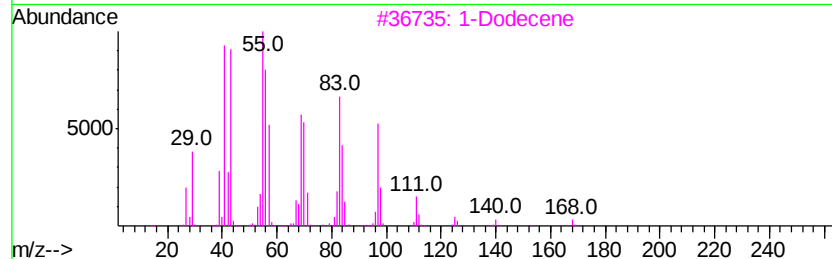
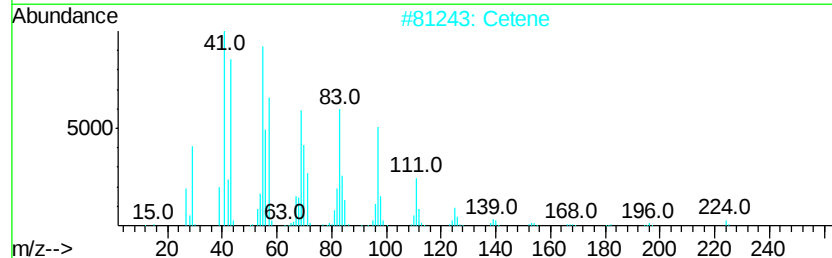
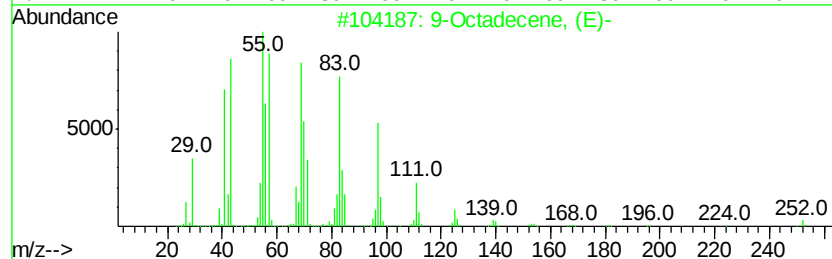
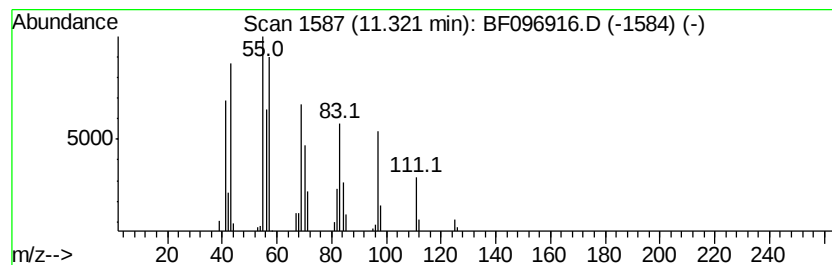
TIC Library : C:\DATABASE\NIST11.L  
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Peak Number 7 9-Octadecene, (E)- Concentration Rank 6

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.32 | 2.42 ng | 55338 | Phenanthrene-d10 | 11.07 |

| Hit# of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------------|-----|---------|-------------|------|
| 1       |   | 9-Octadecene, (E)- | 252 | C18H36  | 007206-25-9 | 91   |
| 2       |   | Cetene             | 224 | C16H32  | 000629-73-2 | 91   |
| 3       |   | 1-Dodecene         | 168 | C12H24  | 000112-41-4 | 86   |
| 4       |   | 5-Octadecene, (E)- | 252 | C18H36  | 007206-21-5 | 86   |
| 5       |   | 1-Tridecene        | 182 | C13H26  | 002437-56-1 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF072017\  
Data File : BF096916.D  
Acq On : 20 Jul 2017 20:09  
Operator : SJ/JU  
Sample : I4312-05  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
72-11987

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

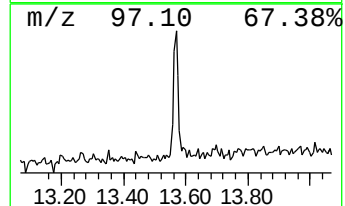
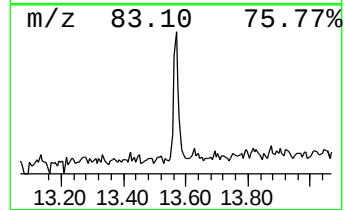
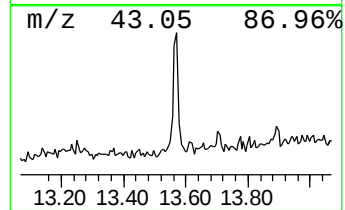
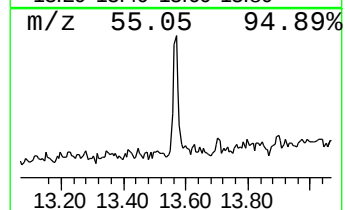
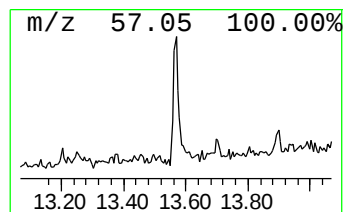
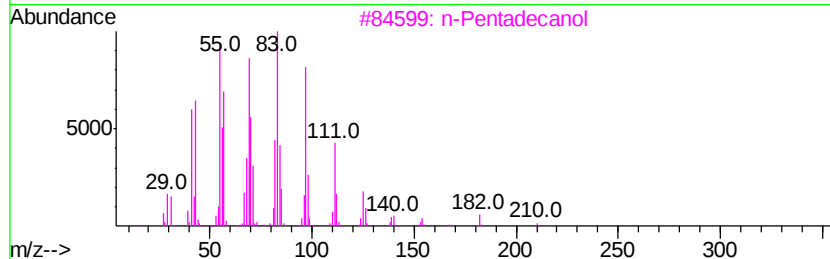
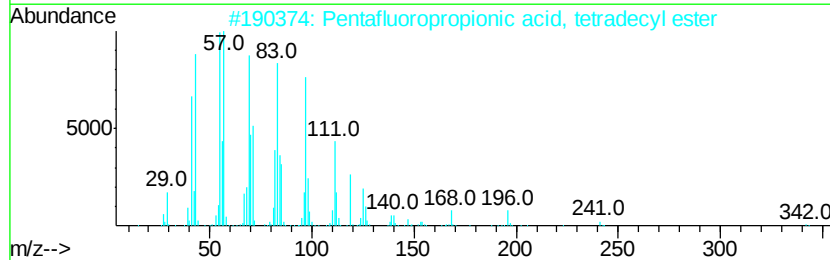
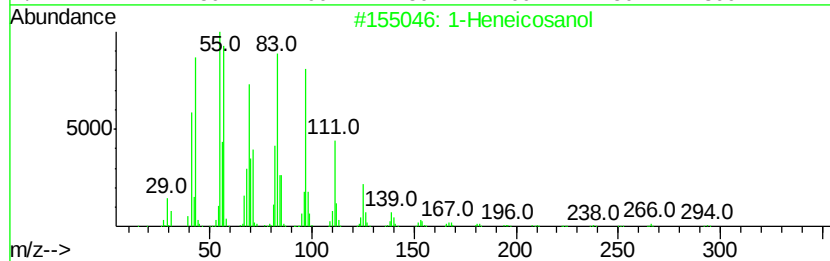
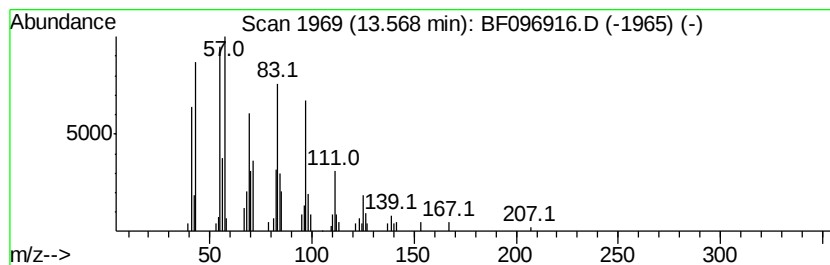
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 1-Heneicosanol Concentration Rank 4

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.57 | 3.62 ng | 82637 | Chrysene-d12     | 13.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8 | 93   |
| 2    |    | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6 | 91   |
| 3    |    | n-Pentadecanol                      | 228 | C15H32O    | 000629-76-5 | 91   |
| 4    |    | Octacosanol                         | 410 | C28H58O    | 000557-61-9 | 90   |
| 5    |    | n-Heptadecanol-1                    | 256 | C17H36O    | 001454-85-9 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF072017\  
Data File : BF096916.D  
Acq On : 20 Jul 2017 20:09  
Operator : SJ/JU  
Sample : I4312-05  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
72-11987

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.74  | 8.4     | ng    | 259538   | 1                     | 6.57  | 621143 | 20.0 |
| unknown6.35          | 6.35  | 77.5    | ng    | 2407020  | 1                     | 6.57  | 621143 | 20.0 |
| Hexanoic acid, 2-... | 7.27  | 2.3     | ng    | 79445    | 2                     | 7.85  | 702586 | 20.0 |
| 1-Iodo-2-methylun... | 10.52 | 2.0     | ng    | 46442    | 4                     | 11.07 | 457725 | 20.0 |
| 9-Octadecene, (E)-   | 11.32 | 2.4     | ng    | 55338    | 4                     | 11.07 | 457725 | 20.0 |
| 1-Heneicosanol       | 13.57 | 3.6     | ng    | 82637    | 5                     | 13.70 | 455984 | 20.0 |