

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\  
 Data File : BF090419.D  
 Acq On : 6 Oct 2016 9:35  
 Operator : UM/SJ  
 Sample : H5065-02 5X  
 Misc :  
 ALS Vial : 36 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-BC-2-10032016

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-BF100516.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         | 4.863       | 218           | 221         | 225          | rBV      | 25356          | 30346         | 1.12%           | 0.105%        |
| 2         | 5.183       | 247           | 249         | 252          | rBV      | 330553         | 320498        | 11.78%          | 1.113%        |
| 3         | 5.594       | 283           | 285         | 288          | rBV      | 1742583        | 1498093       | 55.07%          | 5.203%        |
| 4         | 5.811       | 301           | 304         | 314          | rBV      | 77791          | 126791        | 4.66%           | 0.440%        |
| 5         | 6.612       | 371           | 374         | 376          | rBV      | 1970396        | 1640144       | 60.29%          | 5.696%        |
| 6         | 6.760       | 385           | 387         | 390          | rBV      | 1857054        | 1568104       | 57.64%          | 5.446%        |
| 7         | 7.000       | 405           | 408         | 410          | rBV      | 2150711        | 1787934       | 65.73%          | 6.209%        |
| 8         | 7.160       | 419           | 422         | 424          | rBV      | 938455         | 1107850       | 40.73%          | 3.848%        |
| 9         | 7.560       | 454           | 457         | 459          | rBV      | 919094         | 962018        | 35.36%          | 3.341%        |
| 10        | 8.292       | 518           | 521         | 523          | rBV      | 2530831        | 2441555       | 89.75%          | 8.480%        |
| 11        | 9.366       | 612           | 615         | 617          | rBV      | 2049804        | 1698224       | 62.43%          | 5.898%        |
| 12        | 9.755       | 647           | 649         | 652          | rBV      | 450752         | 402128        | 14.78%          | 1.397%        |
| 13        | 10.052      | 672           | 675         | 677          | rBV      | 3120726        | 2720319       | 100.00%         | 9.448%        |
| 14        | 10.475      | 709           | 712         | 714          | rBV2     | 17605          | 34808         | 1.28%           | 0.121%        |
| 15        | 10.829      | 741           | 743         | 746          | rBV2     | 633010         | 894086        | 32.87%          | 3.105%        |
| 16        | 10.955      | 752           | 754         | 758          | rBV2     | 44506          | 60945         | 2.24%           | 0.212%        |
| 17        | 11.149      | 769           | 771         | 774          | rBV3     | 18430          | 32278         | 1.19%           | 0.112%        |
| 18        | 11.195      | 774           | 775         | 777          | rVB2     | 24396          | 32537         | 1.20%           | 0.113%        |
| 19        | 11.412      | 791           | 794         | 798          | rBV2     | 27043          | 52329         | 1.92%           | 0.182%        |
| 20        | 11.481      | 798           | 800         | 803          | rBV      | 24645          | 46712         | 1.72%           | 0.162%        |
| 21        | 11.538      | 803           | 805         | 808          | rBV      | 2777310        | 2433407       | 89.45%          | 8.451%        |
| 22        | 11.721      | 818           | 821         | 823          | rBV3     | 33065          | 52649         | 1.94%           | 0.183%        |
| 23        | 11.778      | 823           | 826         | 830          | rBV5     | 40199          | 100232        | 3.68%           | 0.348%        |
| 24        | 11.995      | 842           | 845         | 846          | rBV      | 21958          | 35710         | 1.31%           | 0.124%        |
| 25        | 12.063      | 848           | 851         | 852          | rBV      | 173480         | 179202        | 6.59%           | 0.622%        |
| 26        | 12.246      | 865           | 867         | 869          | rBV2     | 18217          | 31403         | 1.15%           | 0.109%        |
| 27        | 12.601      | 896           | 898         | 903          | rBV2     | 131902         | 339825        | 12.49%          | 1.180%        |
| 28        | 12.749      | 909           | 911         | 914          | rBV2     | 51657          | 122598        | 4.51%           | 0.426%        |
| 29        | 12.795      | 914           | 915         | 918          | rVV3     | 58864          | 91503         | 3.36%           | 0.318%        |
| 30        | 12.852      | 918           | 920         | 924          | rVV      | 88848          | 119896        | 4.41%           | 0.416%        |
| 31        | 13.058      | 936           | 938         | 940          | rVB      | 55751          | 50023         | 1.84%           | 0.174%        |
| 32        | 13.126      | 942           | 944         | 947          | rBV      | 1656621        | 1430683       | 52.59%          | 4.969%        |
| 33        | 13.367      | 963           | 965         | 967          | rBV      | 41514          | 41142         | 1.51%           | 0.143%        |
| 34        | 13.709      | 993           | 995         | 1002         | rVB      | 45088          | 69296         | 2.55%           | 0.241%        |

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 Max Peaks: 100  
 Peak Location: TOP

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 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 14.041 | 1022 | 1024 | 1031 | rVB2 | 74270   | 178217  | 6.55%  | 0.619% |
| 36 | 14.189 | 1034 | 1037 | 1039 | rBV  | 2041114 | 2000512 | 73.54% | 6.948% |
| 37 | 14.875 | 1095 | 1097 | 1100 | rVB  | 46112   | 52023   | 1.91%  | 0.181% |
| 38 | 14.955 | 1102 | 1104 | 1108 | rVV2 | 82695   | 134668  | 4.95%  | 0.468% |
| 39 | 15.058 | 1111 | 1113 | 1116 | rVV  | 134402  | 167107  | 6.14%  | 0.580% |
| 40 | 15.127 | 1117 | 1119 | 1122 | rVB  | 69793   | 91852   | 3.38%  | 0.319% |
| 41 | 15.332 | 1135 | 1137 | 1139 | rBV  | 81889   | 104197  | 3.83%  | 0.362% |
| 42 | 15.378 | 1139 | 1141 | 1149 | rVB2 | 85296   | 267095  | 9.82%  | 0.928% |
| 43 | 15.687 | 1165 | 1168 | 1171 | rBV  | 963080  | 1553974 | 57.12% | 5.397% |
| 44 | 15.732 | 1171 | 1172 | 1175 | rVB  | 40684   | 47872   | 1.76%  | 0.166% |
| 45 | 15.938 | 1188 | 1190 | 1197 | rVB  | 43621   | 81611   | 3.00%  | 0.283% |
| 46 | 16.190 | 1209 | 1212 | 1215 | rVB  | 86333   | 128945  | 4.74%  | 0.448% |
| 47 | 16.453 | 1232 | 1235 | 1245 | rVB2 | 158397  | 443640  | 16.31% | 1.541% |
| 48 | 16.647 | 1249 | 1252 | 1254 | rVV2 | 79632   | 154976  | 5.70%  | 0.538% |
| 49 | 16.715 | 1256 | 1258 | 1268 | rVB3 | 48467   | 146254  | 5.38%  | 0.508% |
| 50 | 17.275 | 1304 | 1307 | 1310 | rVB2 | 73470   | 147453  | 5.42%  | 0.512% |
| 51 | 17.618 | 1334 | 1337 | 1347 | rVB8 | 51832   | 216659  | 7.96%  | 0.752% |
| 52 | 17.881 | 1355 | 1360 | 1366 | rBV4 | 76573   | 281704  | 10.36% | 0.978% |
| 53 | 18.338 | 1397 | 1400 | 1401 | rBV3 | 20002   | 39569   | 1.45%  | 0.137% |

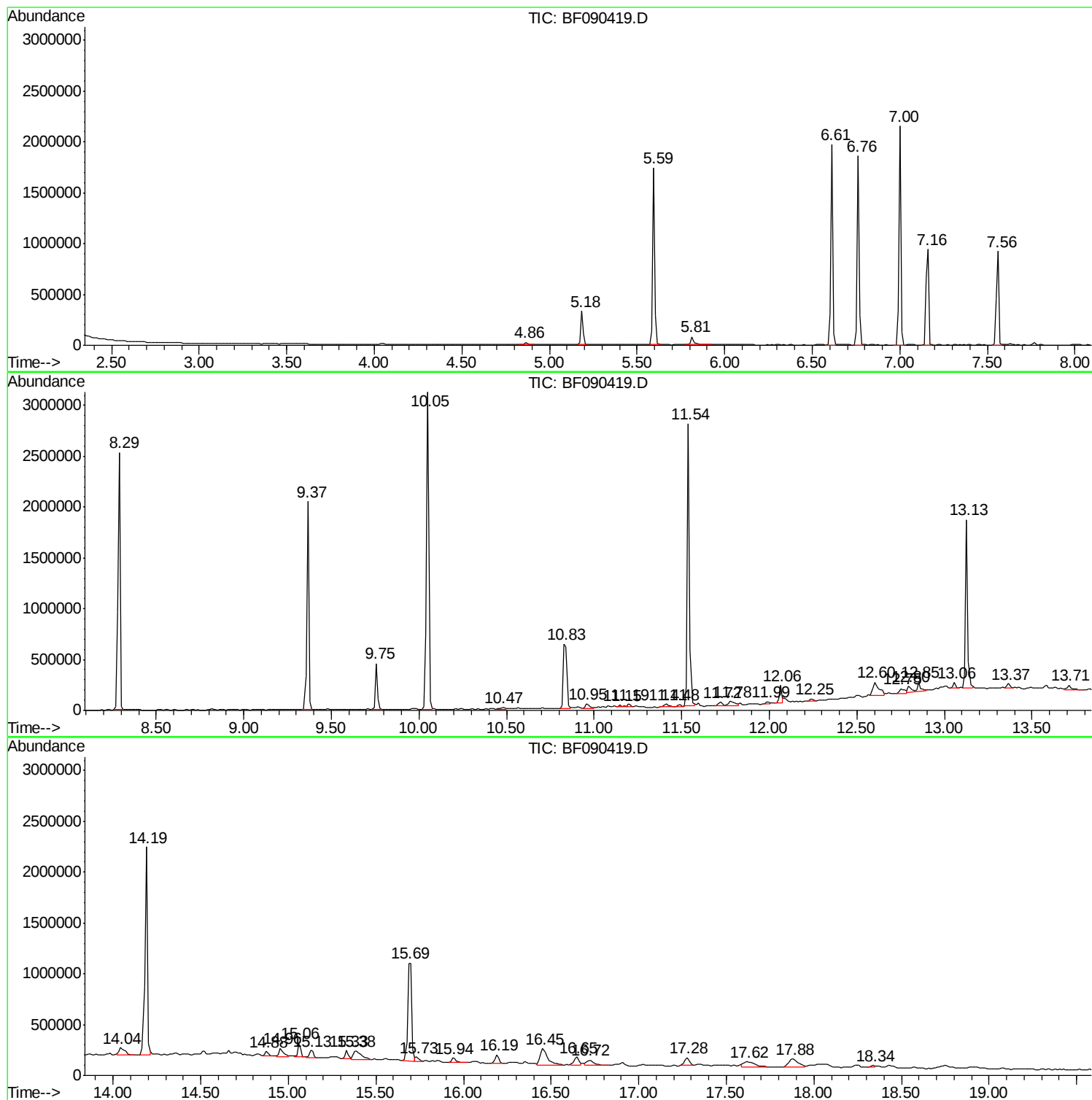
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF100516\  
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Acq On : 6 Oct 2016 9:35  
Operator : UM/SJ  
Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

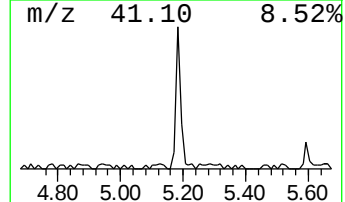
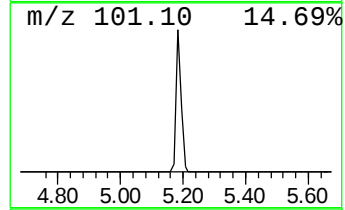
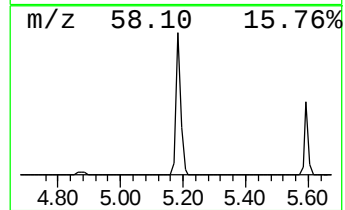
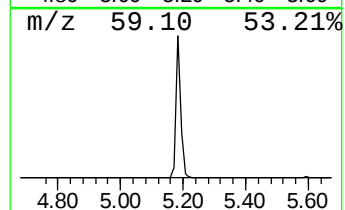
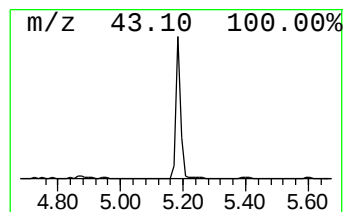
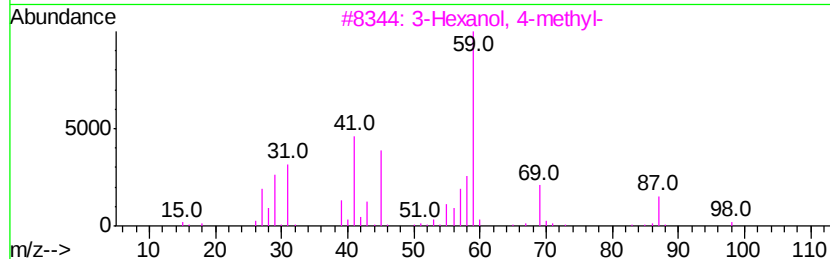
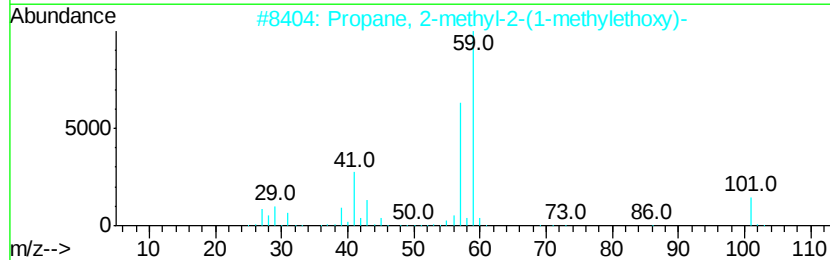
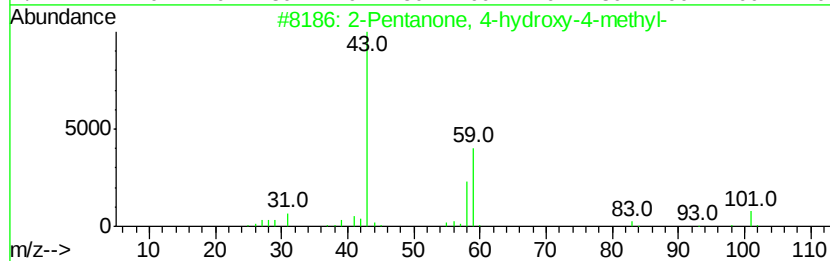
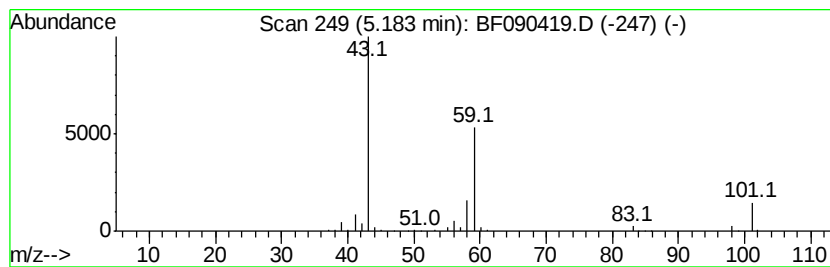
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.18 | 3.59 ng | 320498 | 1,4-Dichlorobenzene-d4 | 7.00 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 50   |
| 2    |    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 36   |
| 3    |    | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 4    |    | 3-Octanol                           | 130 | C8H18O   | 000589-98-0 | 33   |
| 5    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 32   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

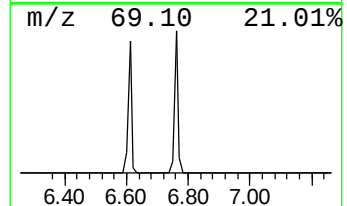
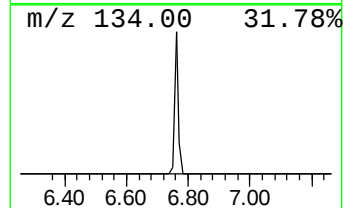
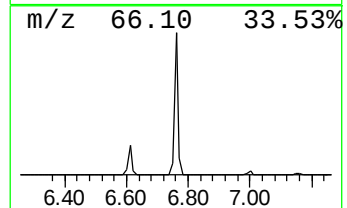
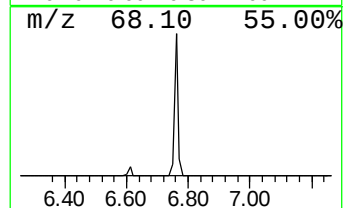
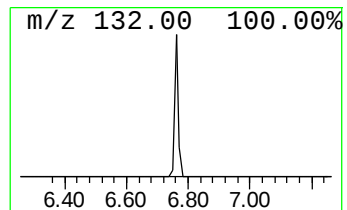
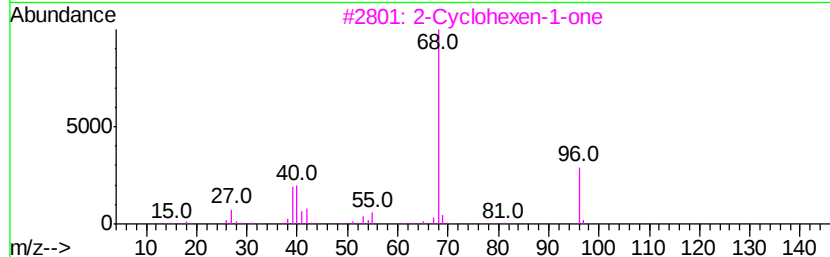
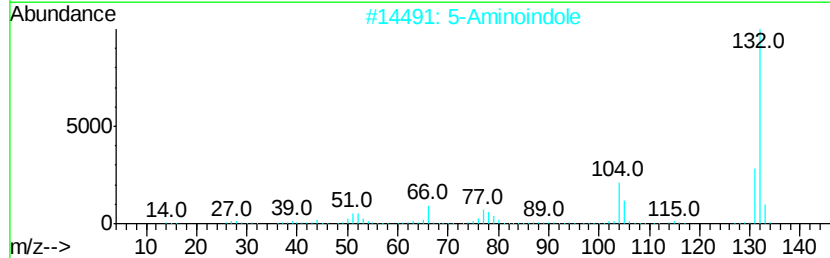
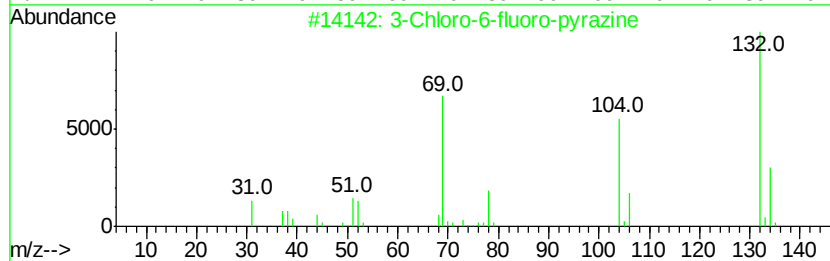
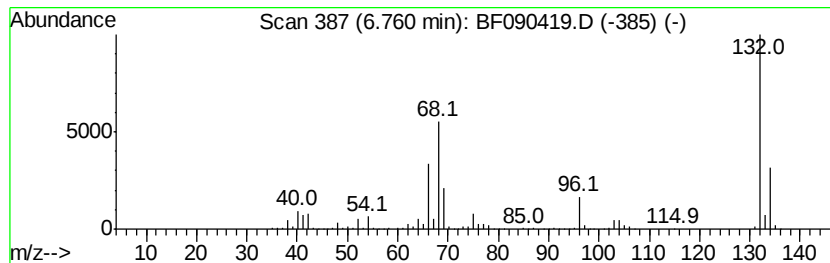
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.76 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.76 | 17.54 ng | 1568100 | 1,4-Dichlorobenzene-d4 | 7.00 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 22   |
| 3    |    | 2-Cyclohexen-1-one               | 96  | C6H8O     | 000930-68-7  | 11   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | Benzene, 1-ethenyl-3,5-dimethyl- | 132 | C10H12    | 005379-20-4  | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

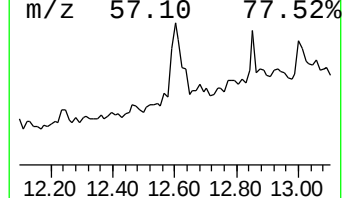
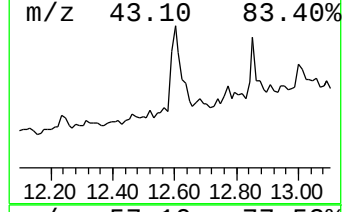
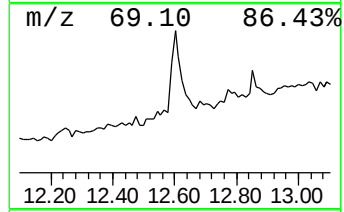
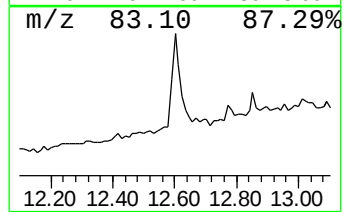
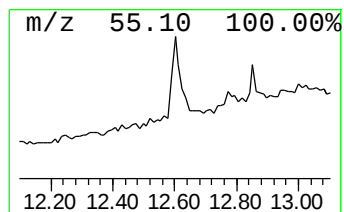
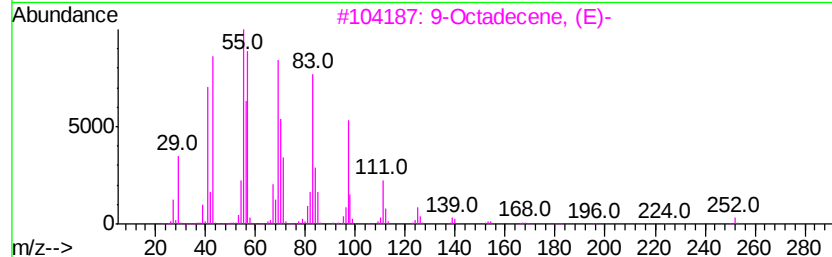
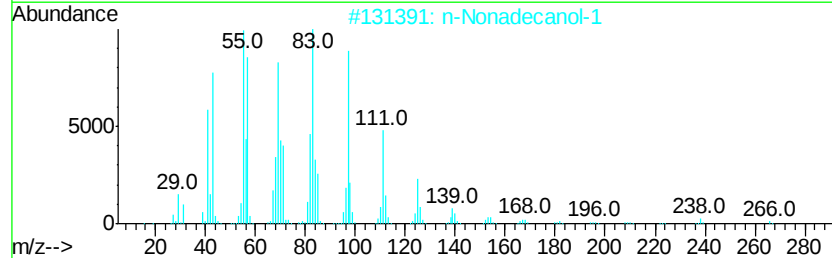
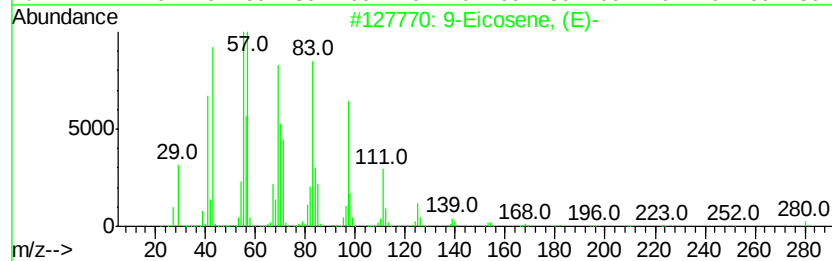
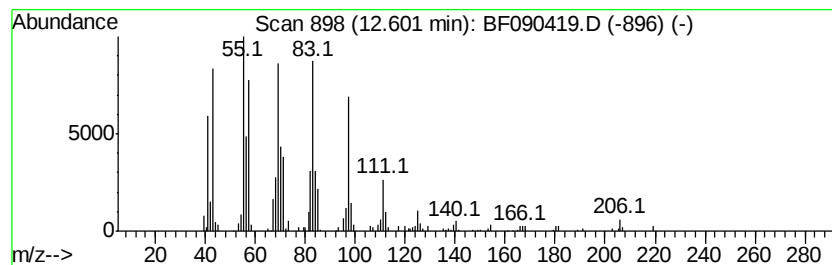
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 9-Eicosene, (E)- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.60 | 2.79 ng | 339825 | Phenanthrene-d10 | 11.54 |

| Hit# of | 5                  | Tentative ID | MW      | MolForm     | CAS# | Qual |
|---------|--------------------|--------------|---------|-------------|------|------|
| 1       | 9-Eicosene, (E)-   | 280          | C20H40  | 074685-29-3 | 94   |      |
| 2       | n-Nonadecanol-1    | 284          | C19H40O | 001454-84-8 | 91   |      |
| 3       | 9-Octadecene, (E)- | 252          | C18H36  | 007206-25-9 | 91   |      |
| 4       | n-Heptadecanol-1   | 256          | C17H36O | 001454-85-9 | 91   |      |
| 5       | 1-Nonadecene       | 266          | C19H38  | 018435-45-5 | 91   |      |



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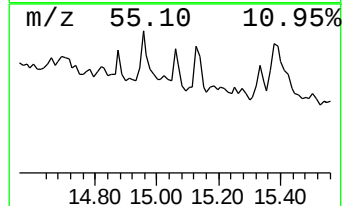
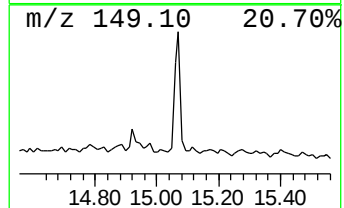
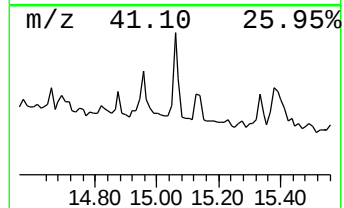
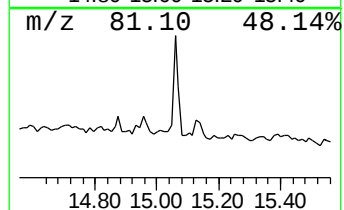
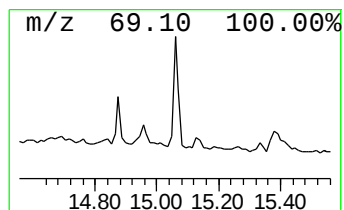
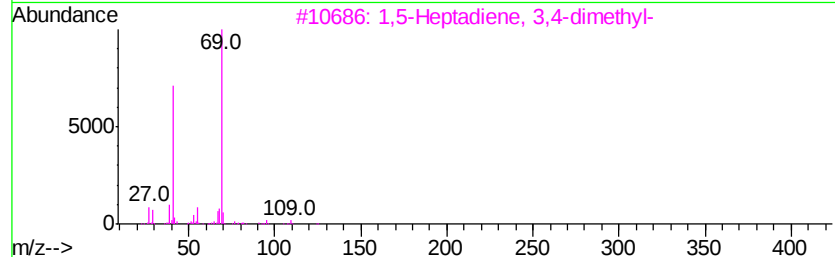
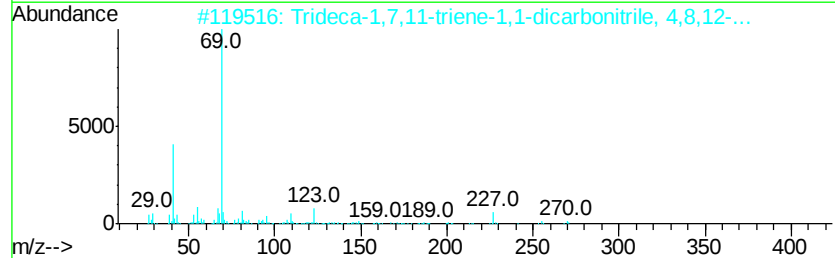
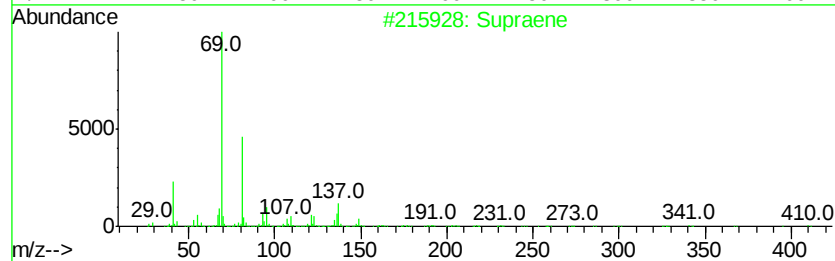
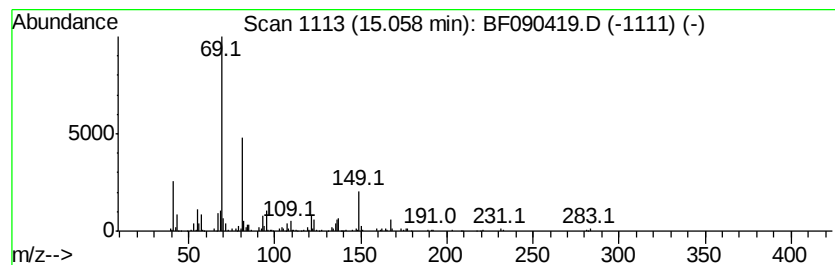
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BNA\_F  
ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Supraene Concentration Rank 9

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.            |
|---------|-------------------------------------|--------------|------------------|-----------------|
| 15.06   | 2.15 ng                             | 167107       | Perylene-d12     | 15.70           |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual       |
| 1       | Supraene                            |              | 410 C30H50       | 007683-64-9 74  |
| 2       | Trideca-1,7,11-triene-1,1-dicarb... |              | 270 C18H26N2     | 1000161-46-0 47 |
| 3       | 1,5-Heptadiene, 3,4-dimethyl-       |              | 124 C9H16        | 1000061-77-9 47 |
| 4       | 2,6-Octadiene, 4,5-dimethyl-        |              | 138 C10H18       | 018476-57-8 43  |
| 5       | 2,6-Octadiene, 4-methyl-            |              | 124 C9H16        | 074498-94-5 43  |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF100516\  
Data File : BF090419.D  
Acq On : 6 Oct 2016 9:35  
Operator : UM/SJ  
Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

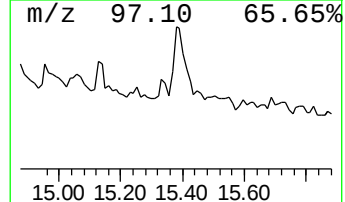
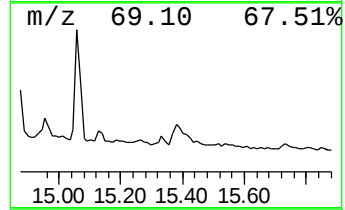
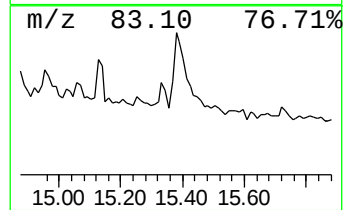
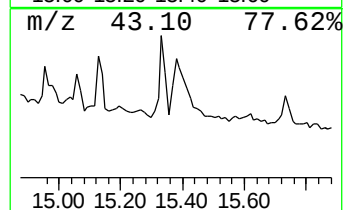
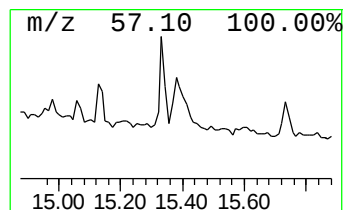
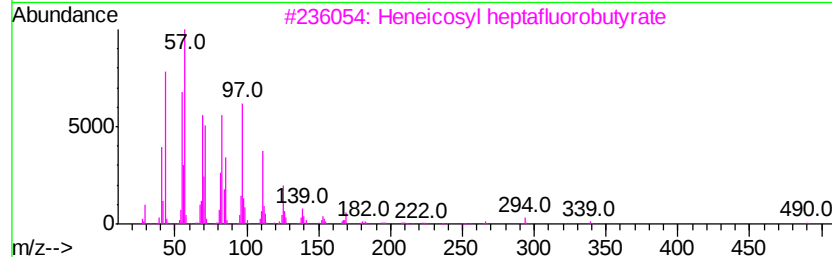
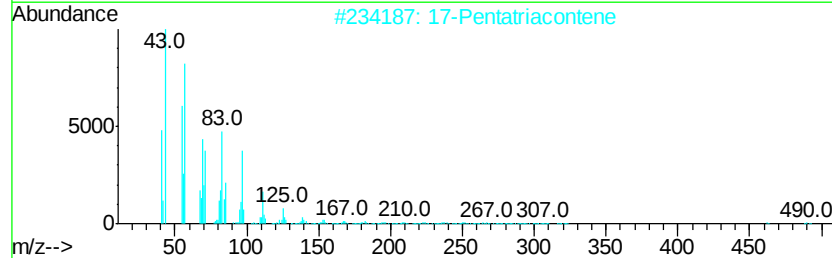
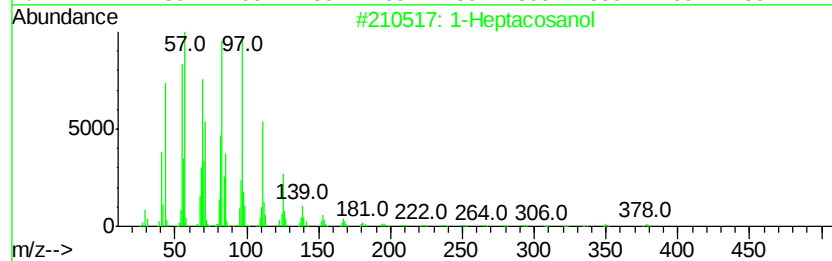
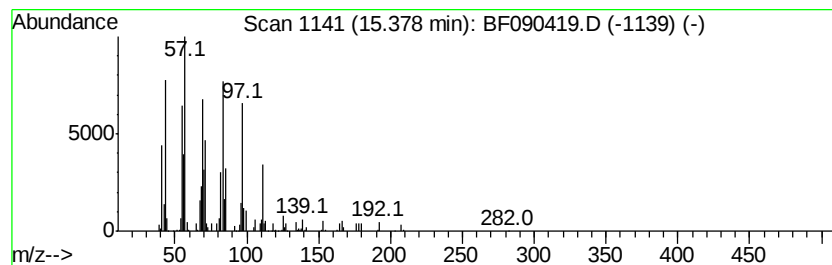
Instrument :  
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ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 1-Heptacosanol Concentration Rank 6

| R.T.      | EstConc                        | Area   | Relative to ISTD | R.T.         |      |
|-----------|--------------------------------|--------|------------------|--------------|------|
| 15.38     | 3.44 ng                        | 267095 | Perylene-d12     | 15.70        |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Heptacosanol                 | 396    | C27H56O          | 002004-39-9  | 91   |
| 2         | 17-Pentatriacontene            | 491    | C35H70           | 006971-40-0  | 90   |
| 3         | Heneicosyl heptafluorobutvrate | 508    | C25H43F7O2       | 1000351-83-8 | 90   |
| 4         | 1-Docosene                     | 308    | C22H44           | 001599-67-3  | 90   |
| 5         | 1-Heneicosanol                 | 312    | C21H44O          | 015594-90-8  | 74   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF100516\  
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Operator : UM/SJ  
Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

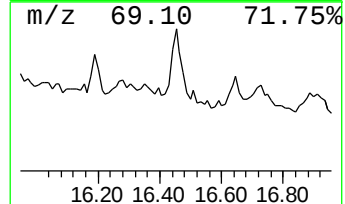
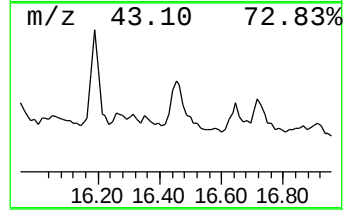
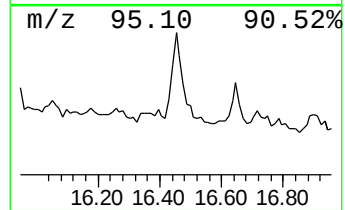
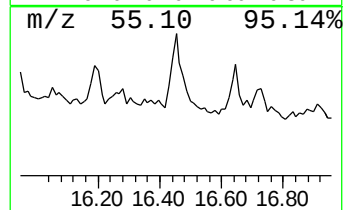
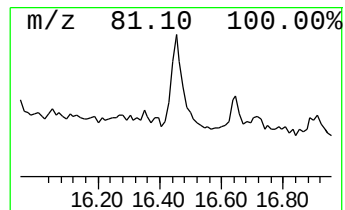
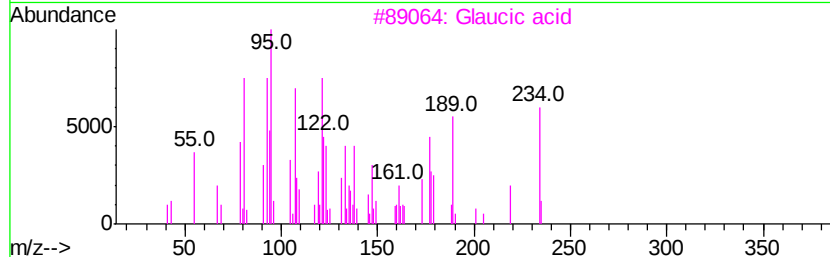
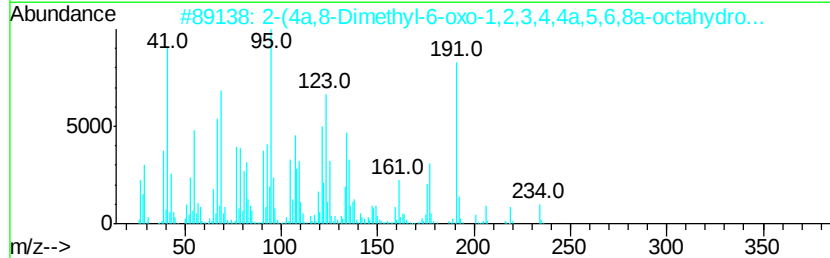
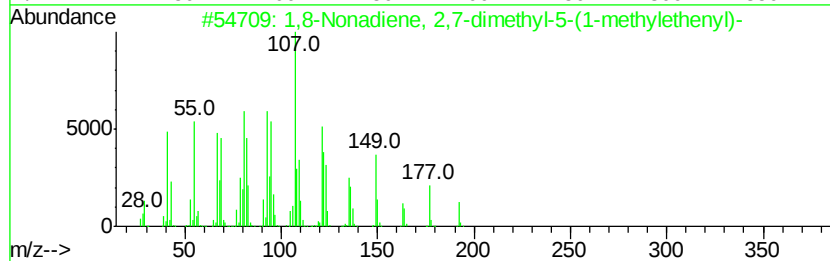
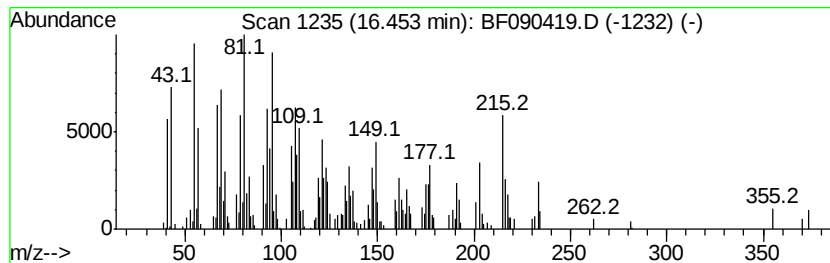
Instrument :  
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ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 1,8-Nonadiene, 2,7-dimethyl... Concentration Rank 3

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 16.45     | 5.71 ng                             | 443640 | Perylene-d12     | 15.70        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1,8-Nonadiene, 2,7-dimethyl-5-(1... | 192    | C14H24           | 068702-20-5  | 53   |
| 2         | 2-(4a,8-Dimethyl-6-oxo-1,2,3,4,4... | 234    | C15H22O2         | 1000190-21-8 | 50   |
| 3         | Glaucic acid                        | 234    | C15H22O2         | 087745-30-0  | 49   |
| 4         | 2-Methyl-2-bornene                  | 150    | C11H18           | 072540-93-3  | 44   |
| 5         | Bicyclo[2.2.2]octane, 2-methyl-     | 124    | C9H16            | 000766-53-0  | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF100516\  
Data File : BF090419.D  
Acq On : 6 Oct 2016 9:35  
Operator : UM/SJ  
Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

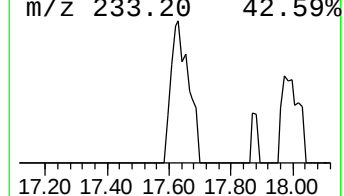
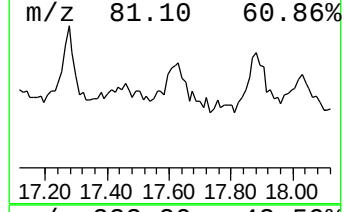
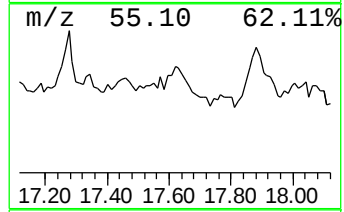
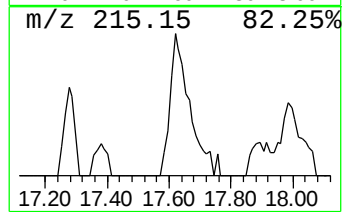
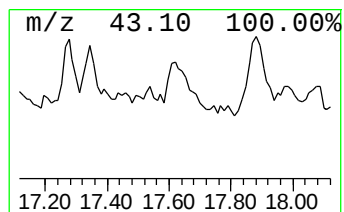
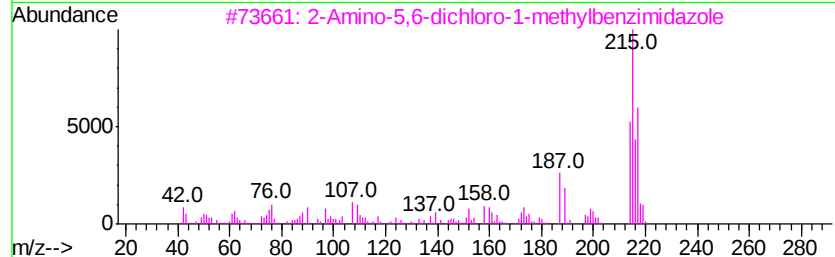
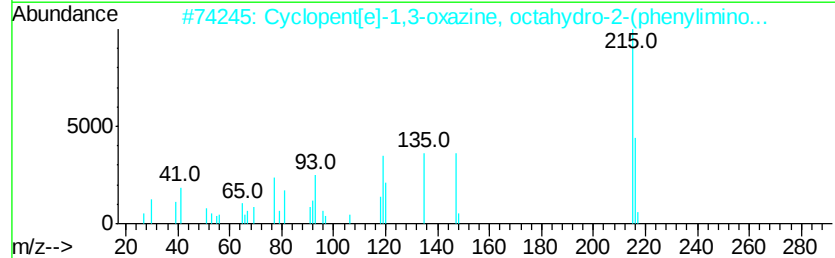
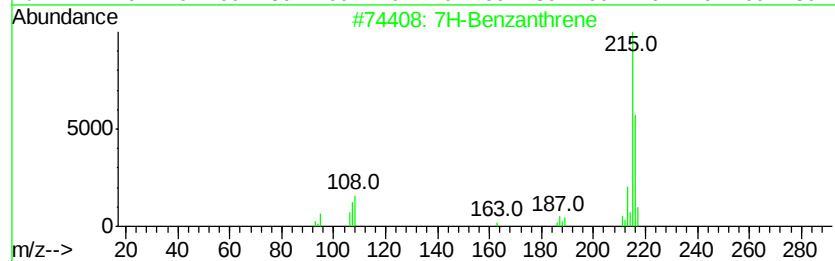
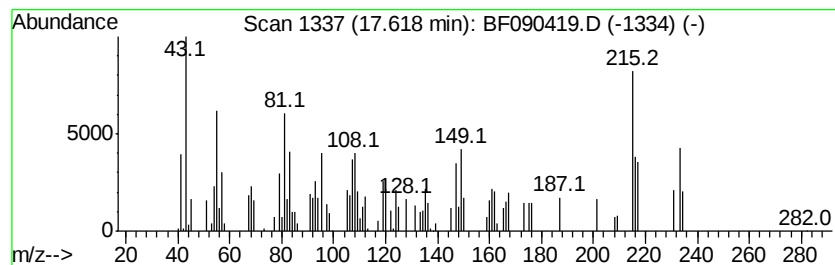
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown17.62 Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.62 | 2.79 ng | 216659 | Perylene-d12     | 15.70 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 7H-Benzanthrene                      | 216 | C17H12     | 000199-94-0  | 38   |
| 2    |    | Cyclopent[el]-1,3-oxazine, octahy... | 216 | C13H16N2O  | 1000138-92-2 | 22   |
| 3    |    | 2-Amino-5,6-dichloro-1-methylben...  | 215 | C8H7Cl2N3  | 030486-79-4  | 22   |
| 4    |    | 2,6-Dichloro-4-trifluoromethylpv...  | 215 | C6H2Cl2F3N | 039890-98-7  | 14   |
| 5    |    | 2-Chloro-7-phenyltropone             | 216 | C13H9ClO   | 090128-01-1  | 12   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF100516\  
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Operator : UM/SJ  
Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

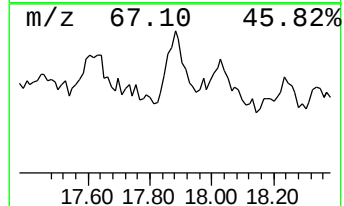
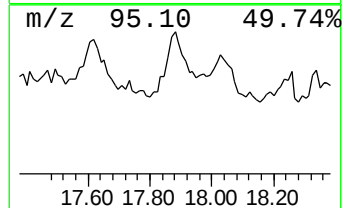
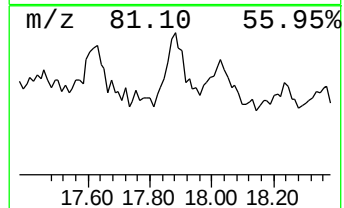
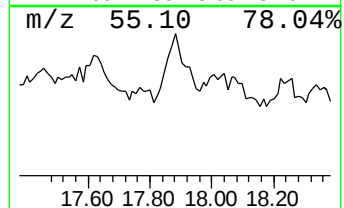
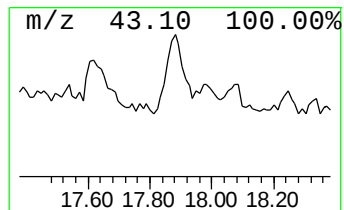
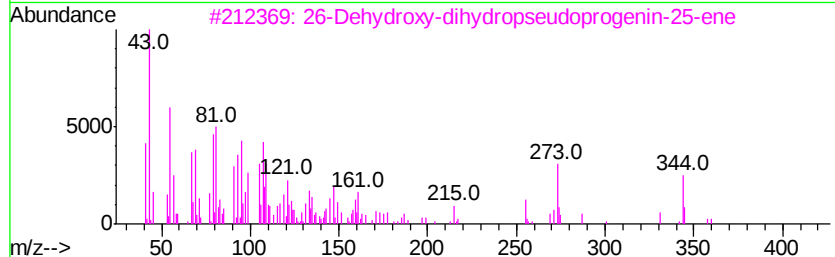
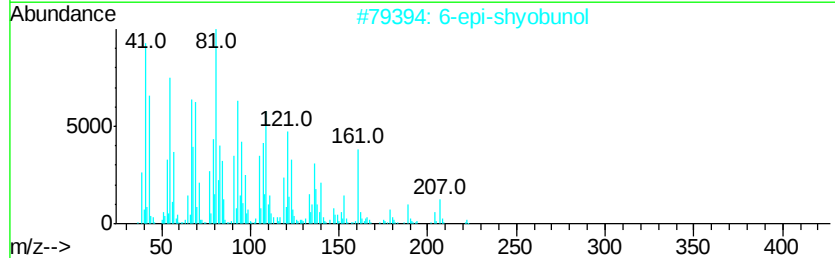
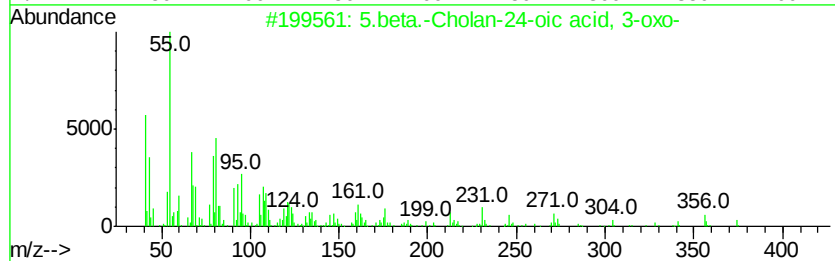
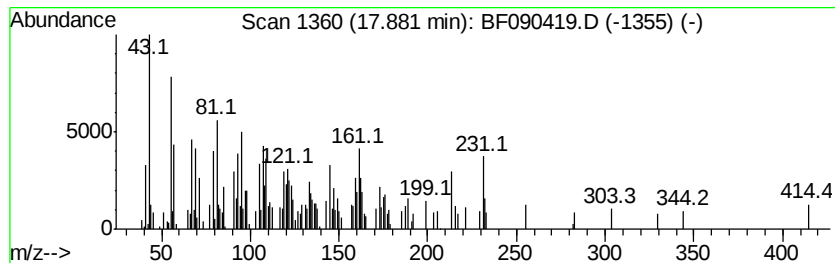
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown17.88 Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.88 | 3.63 ng | 281704 | Perylene-d12     | 15.70 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|--------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 5.beta.-Cholan-24-oic acid, 3-oxo-   | 374 | C24H38O3 | 001553-56-6  | 38   |
| 2    |    | 6-epi-shyobunol                      | 222 | C15H26O  | 1000374-18-3 | 37   |
| 3    |    | 26-Dehydroxy-dihdropseudoprogenin... | 400 | C27H44O2 | 1000255-22-7 | 27   |
| 4    |    | 9,19-Cyclolanostan-3-ol, acetate...  | 470 | C32H54O2 | 004575-74-0  | 25   |
| 5    |    | 4-Hydroxy-4-(4,6-dimethylcyclohe...  | 196 | C12H20O2 | 002316-79-2  | 22   |



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Sample : H5065-02 5X  
Misc :  
ALS Vial : 36 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-BC-2-10032016

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF100516.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... | 5.18  | 3.6 ng  |       | 320498   | 1                     | 7.00  | 1787930 | 20.0 |
| unknown6.76          | 6.76  | 17.5 ng |       | 1568100  | 1                     | 7.00  | 1787930 | 20.0 |
| 9-Eicosene, (E)-     | 12.60 | 2.8 ng  |       | 339825   | 4                     | 11.54 | 2433410 | 20.0 |
| Supraene             | 15.06 | 2.1 ng  |       | 167107   | 6                     | 15.70 | 1553970 | 20.0 |
| 1-Heptacosanol       | 15.38 | 3.4 ng  |       | 267095   | 6                     | 15.70 | 1553970 | 20.0 |
| 1,8-Nonadiene, 2,... | 16.45 | 5.7 ng  |       | 443640   | 6                     | 15.70 | 1553970 | 20.0 |
| unknown17.62         | 17.62 | 2.8 ng  |       | 216659   | 6                     | 15.70 | 1553970 | 20.0 |
| unknown17.88         | 17.88 | 3.6 ng  |       | 281704   | 6                     | 15.70 | 1553970 | 20.0 |