

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
 Data File : BF088913.D  
 Acq On : 11 Jul 2016 21:51  
 Operator : UM/SJ  
 Sample : H3921-08  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 N8-B6(0-2)C

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-BF063016.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         | 1.678       | 6             | 8           | 13           | rVV      | 148507         | 203510        | 10.12%          | 1.125%        |
| 2         | 1.793       | 17            | 18          | 28           | rVB      | 1091715        | 2010607       | 100.00%         | 11.112%       |
| 3         | 1.918       | 28            | 29          | 77           | rVB      | 73528          | 431623        | 21.47%          | 2.386%        |
| 4         | 4.867       | 285           | 287         | 292          | rBB      | 69802          | 94424         | 4.70%           | 0.522%        |
| 5         | 5.302       | 323           | 325         | 328          | rBV      | 1038507        | 1480246       | 73.62%          | 8.181%        |
| 6         | 6.353       | 414           | 417         | 420          | rBV      | 1205534        | 1344244       | 66.86%          | 7.430%        |
| 7         | 6.479       | 426           | 428         | 431          | rBV      | 1576575        | 1436006       | 71.42%          | 7.937%        |
| 8         | 6.707       | 446           | 448         | 450          | rBB      | 365096         | 366631        | 18.23%          | 2.026%        |
| 9         | 6.867       | 459           | 462         | 464          | rBB      | 1398184        | 1186738       | 59.02%          | 6.559%        |
| 10        | 7.279       | 495           | 498         | 500          | rBV      | 1035272        | 1001468       | 49.81%          | 5.535%        |
| 11        | 7.725       | 534           | 537         | 543          | rVB2     | 25359          | 39686         | 1.97%           | 0.219%        |
| 12        | 7.999       | 558           | 561         | 563          | rBV      | 543752         | 493947        | 24.57%          | 2.730%        |
| 13        | 9.073       | 652           | 655         | 657          | rBV      | 1787575        | 1553048       | 77.24%          | 8.584%        |
| 14        | 9.473       | 687           | 690         | 692          | rVB      | 123779         | 162153        | 8.06%           | 0.896%        |
| 15        | 9.748       | 711           | 714         | 716          | rBV      | 465426         | 448534        | 22.31%          | 2.479%        |
| 16        | 10.525      | 780           | 782         | 785          | rBV      | 761863         | 785053        | 39.05%          | 4.339%        |
| 17        | 10.571      | 785           | 786         | 789          | rVB      | 21571          | 24821         | 1.23%           | 0.137%        |
| 18        | 10.662      | 789           | 794         | 799          | rBV      | 25998          | 33764         | 1.68%           | 0.187%        |
| 19        | 11.222      | 841           | 843         | 844          | rBV      | 440373         | 398060        | 19.80%          | 2.200%        |
| 20        | 11.245      | 844           | 845         | 846          | rVV      | 30427          | 21069         | 1.05%           | 0.116%        |
| 21        | 12.422      | 946           | 948         | 950          | rVB      | 80691          | 65750         | 3.27%           | 0.363%        |
| 22        | 12.651      | 964           | 968         | 969          | rBV      | 62468          | 81849         | 4.07%           | 0.452%        |
| 23        | 12.685      | 969           | 971         | 977          | rVV      | 240609         | 302750        | 15.06%          | 1.673%        |
| 24        | 12.799      | 979           | 981         | 984          | rVB      | 1058375        | 1189131       | 59.14%          | 6.572%        |
| 25        | 13.702      | 1058          | 1060        | 1064         | rBV2     | 113161         | 156226        | 7.77%           | 0.863%        |
| 26        | 13.839      | 1069          | 1072        | 1079         | rBV      | 435718         | 478117        | 23.78%          | 2.643%        |
| 27        | 13.954      | 1079          | 1082        | 1085         | rBV2     | 12530          | 21647         | 1.08%           | 0.120%        |
| 28        | 14.045      | 1087          | 1090        | 1094         | rBV      | 28828          | 58606         | 2.91%           | 0.324%        |
| 29        | 14.354      | 1113          | 1117        | 1125         | rBV      | 170009         | 301214        | 14.98%          | 1.665%        |
| 30        | 14.651      | 1139          | 1143        | 1145         | rBV      | 101072         | 137909        | 6.86%           | 0.762%        |
| 31        | 14.754      | 1150          | 1152        | 1154         | rBV      | 113749         | 107386        | 5.34%           | 0.594%        |
| 32        | 14.834      | 1157          | 1159        | 1163         | rBV      | 47352          | 81500         | 4.05%           | 0.450%        |
| 33        | 14.948      | 1165          | 1169        | 1172         | rBV      | 179015         | 351680        | 17.49%          | 1.944%        |
| 34        | 15.097      | 1181          | 1182        | 1185         | rVB      | 20822          | 28103         | 1.40%           | 0.155%        |

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N8-B6(0-2)C

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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-BF063016.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 15.154 | 1185 | 1187 | 1190 | rBV  | 30056  | 38792  | 1.93%  | 0.214% |
| 36 | 15.211 | 1190 | 1192 | 1195 | rVV  | 260996 | 310239 | 15.43% | 1.715% |
| 37 | 15.303 | 1198 | 1200 | 1203 | rBV  | 22674  | 34783  | 1.73%  | 0.192% |
| 38 | 15.440 | 1210 | 1212 | 1214 | rBV  | 40096  | 46623  | 2.32%  | 0.258% |
| 39 | 15.645 | 1228 | 1230 | 1232 | rBV  | 64488  | 104183 | 5.18%  | 0.576% |
| 40 | 15.691 | 1232 | 1234 | 1237 | rVV  | 49162  | 84653  | 4.21%  | 0.468% |
| 41 | 15.748 | 1237 | 1239 | 1245 | rVB2 | 14634  | 34062  | 1.69%  | 0.188% |
| 42 | 16.343 | 1288 | 1291 | 1294 | rVB2 | 34533  | 59328  | 2.95%  | 0.328% |
| 43 | 16.503 | 1302 | 1305 | 1308 | rBV2 | 13113  | 38348  | 1.91%  | 0.212% |
| 44 | 16.605 | 1311 | 1314 | 1317 | rVV  | 21507  | 44913  | 2.23%  | 0.248% |
| 45 | 16.686 | 1317 | 1321 | 1325 | rVV5 | 18667  | 65355  | 3.25%  | 0.361% |
| 46 | 16.788 | 1325 | 1330 | 1336 | rVV4 | 37270  | 109923 | 5.47%  | 0.608% |
| 47 | 16.880 | 1336 | 1338 | 1344 | rVB  | 9485   | 22550  | 1.12%  | 0.125% |
| 48 | 17.017 | 1347 | 1350 | 1354 | rBV5 | 21188  | 50644  | 2.52%  | 0.280% |
| 49 | 17.086 | 1354 | 1356 | 1362 | rVB6 | 17810  | 33141  | 1.65%  | 0.183% |
| 50 | 18.320 | 1460 | 1464 | 1469 | rBV  | 51581  | 138245 | 6.88%  | 0.764% |

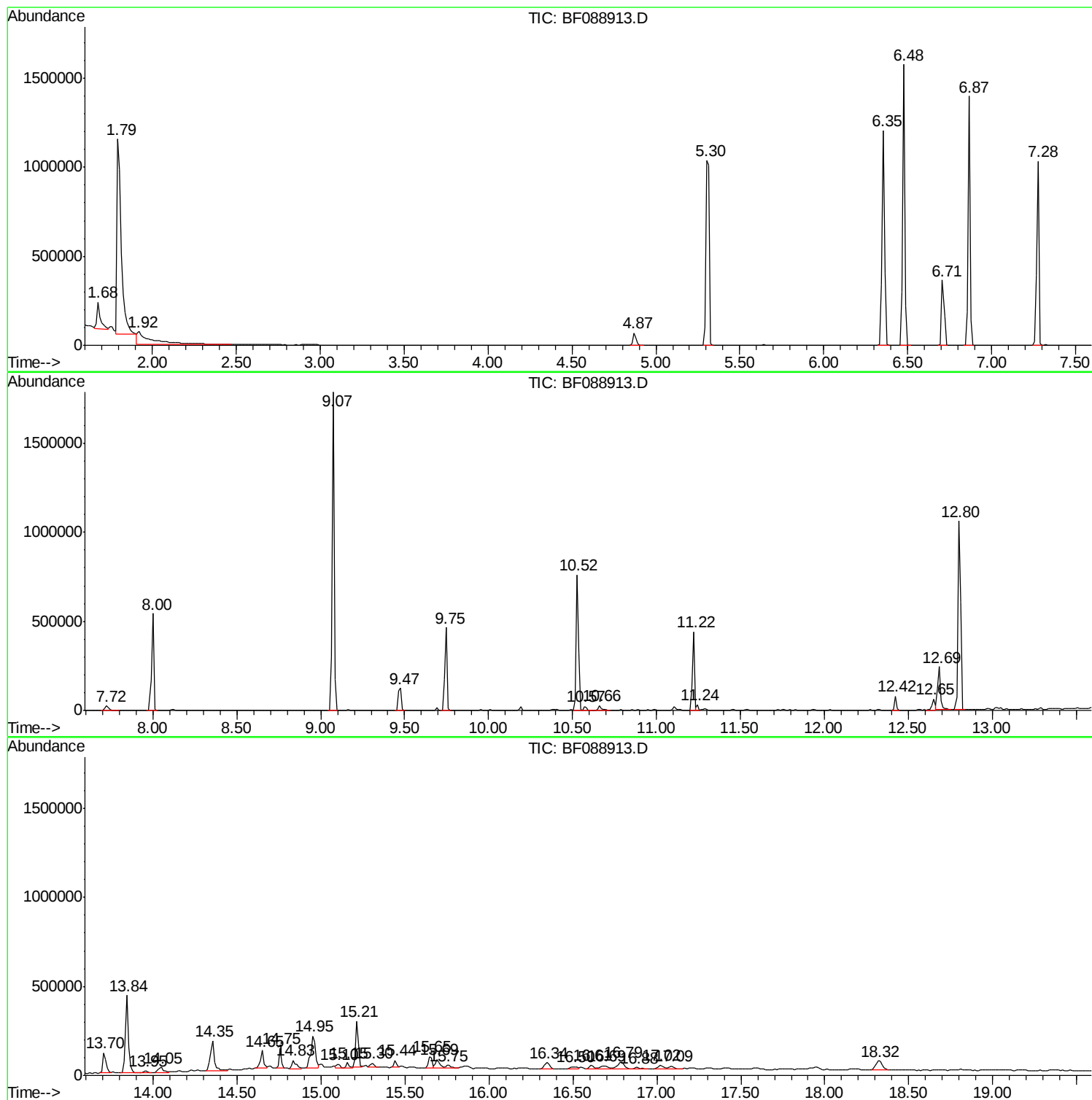
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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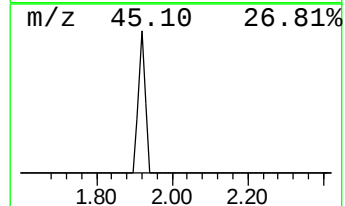
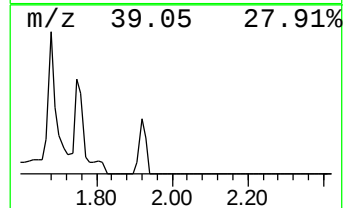
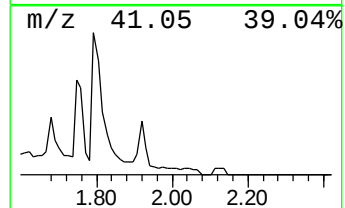
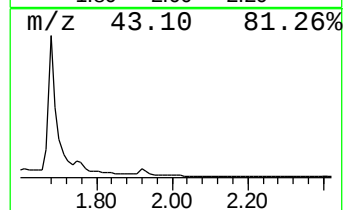
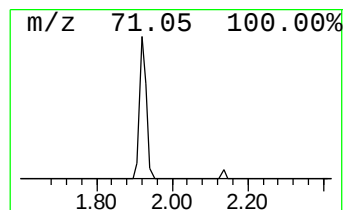
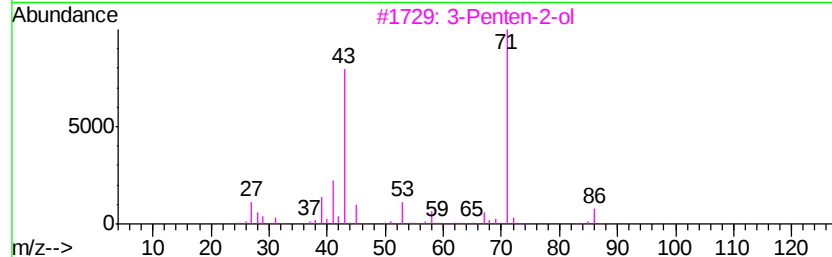
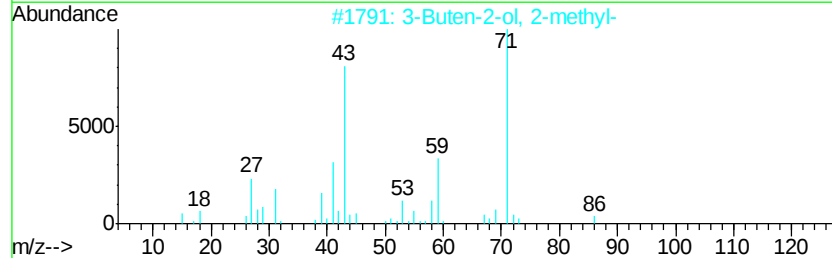
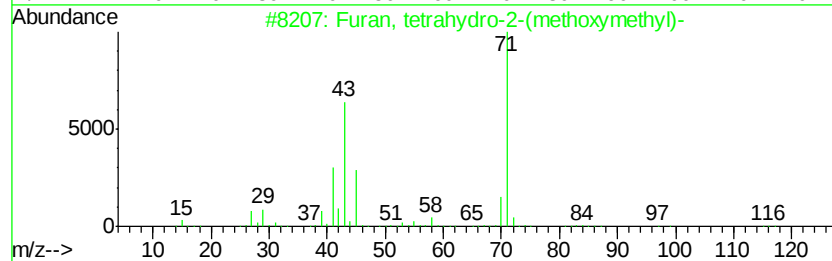
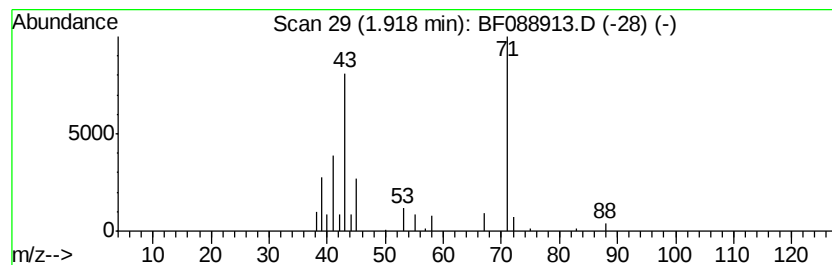
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Furan, tetrahydro-2-(methox... Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 1.92 | 23.55 ng | 431623 | 1,4-Dichlorobenzene-d4 | 6.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Furan, tetrahydro-2-(methoxymeth... | 116 | C6H12O2 | 019354-27-9 | 64   |
| 2    |    | 3-Buten-2-ol, 2-methyl-             | 86  | C5H10O  | 000115-18-4 | 50   |
| 3    |    | 3-Penten-2-ol                       | 86  | C5H10O  | 001569-50-2 | 50   |
| 4    |    | Cyclopropanemethanol, .alpha.-bu... | 128 | C8H16O  | 004379-16-2 | 45   |
| 5    |    | 1,6-Heptadien-4-ol                  | 112 | C7H12O  | 002883-45-6 | 42   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

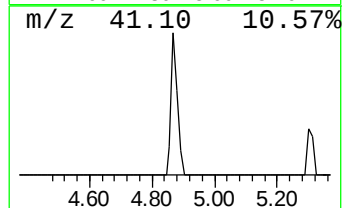
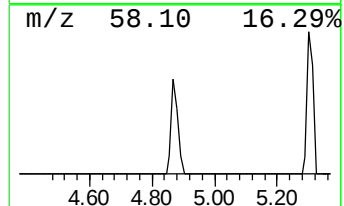
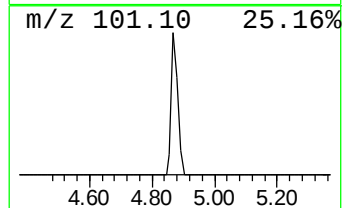
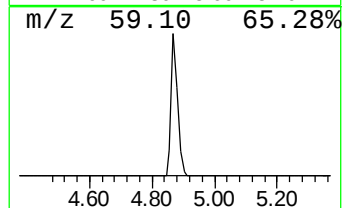
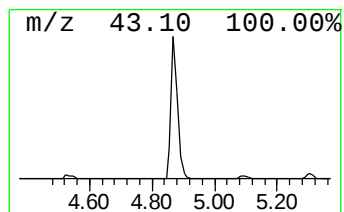
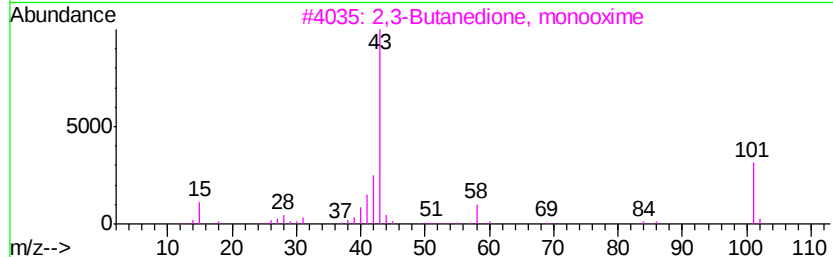
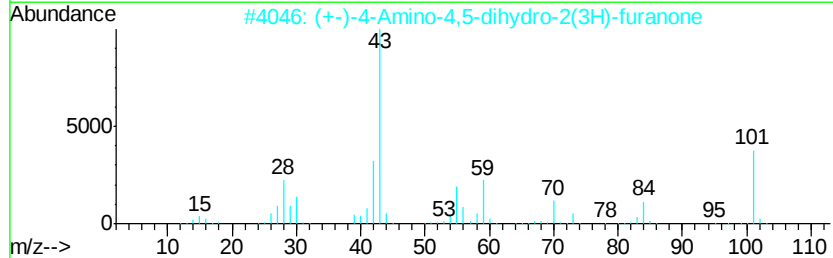
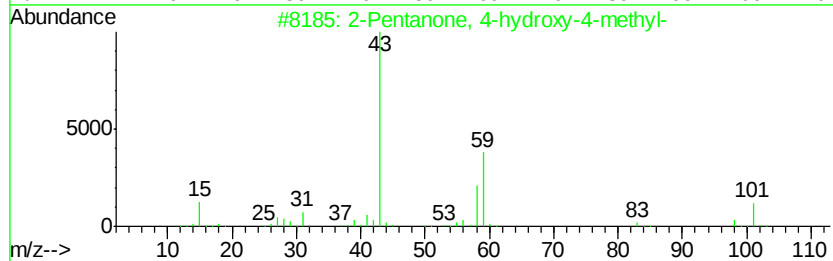
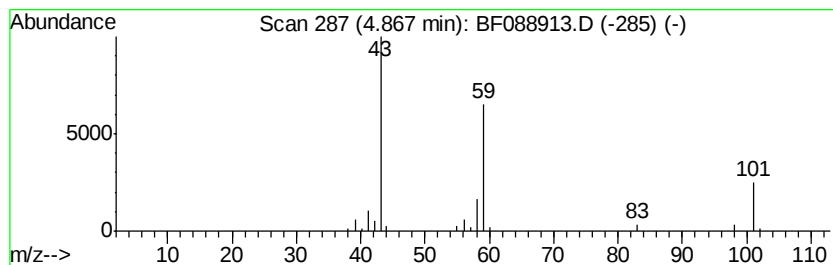
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 4.87 | 5.15 ng | 94424 | 1,4-Dichlorobenzene-d4 | 6.71 |

| Hit# | of | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-       | 116 | C6H12O2 | 000123-42-2 | 56   |
| 2    |    | (+)-4-Amino-4,5-dihydro-2(3H)-furanone | 101 | C4H7NO2 | 016504-58-8 | 9    |
| 3    |    | 2,3-Butanedione, monooxime             | 101 | C4H7NO2 | 000057-71-6 | 9    |
| 4    |    | Morpholine, 4-methyl-                  | 101 | C5H11NO | 000109-02-4 | 9    |
| 5    |    | Butane, 1-ethoxy-                      | 102 | C6H14O  | 000628-81-9 | 9    |



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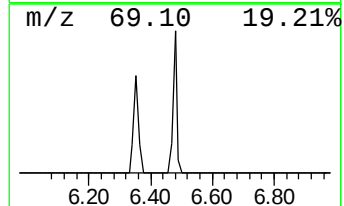
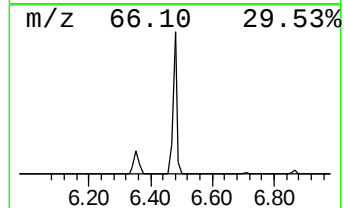
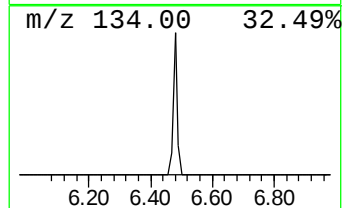
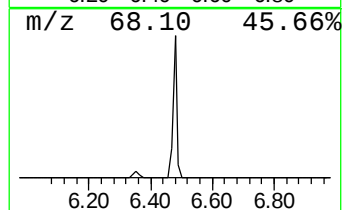
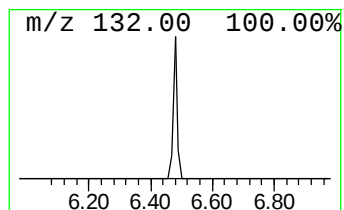
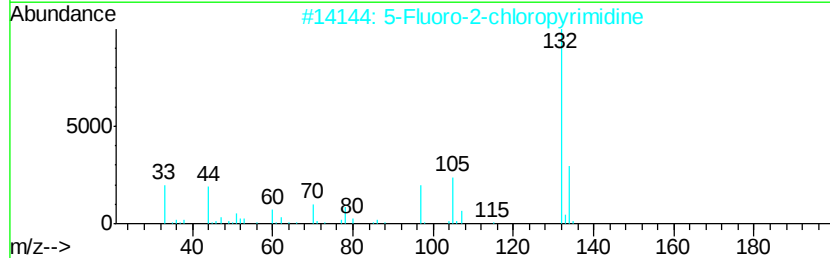
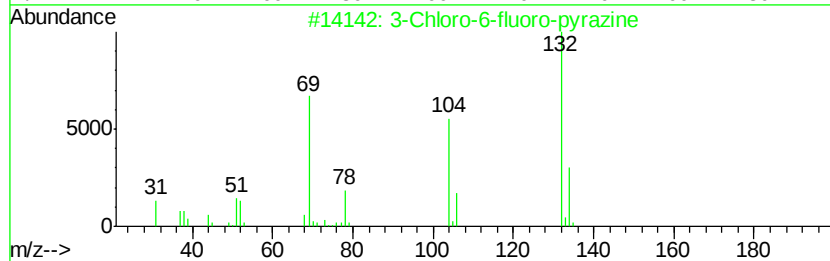
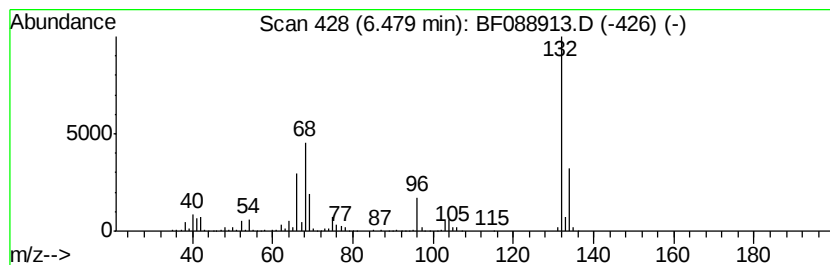
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 unknown6.48 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.48 | 78.34 ng | 1436010 | 1,4-Dichlorobenzene-d4 | 6.71 |

| Hit# | of                                  | Tentative ID | MW        | MolForm      | CAS# | Qual |
|------|-------------------------------------|--------------|-----------|--------------|------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine          | 132          | C4H2ClFN2 | 1000146-10-7 | 27   |      |
| 2    | 5-Fluoro-2-chloropyrimidine         | 132          | C4H2ClFN2 | 062802-42-0  | 16   |      |
| 3    | Tranvlcyproline-propionyl           | 189          | C12H15NO  | 1000123-86-3 | 10   |      |
| 4    | Cyclopropanamine, 1-phenyl-         | 133          | C9H11N    | 1000351-26-2 | 10   |      |
| 5    | N-(3,4-Methylenedioxyphenyl)isop... | 267          | C17H17NO2 | 1000378-96-6 | 10   |      |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

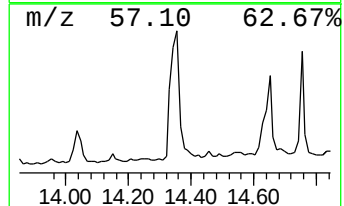
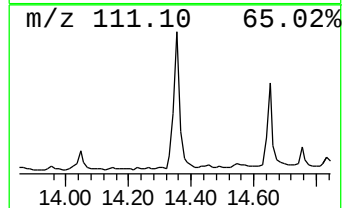
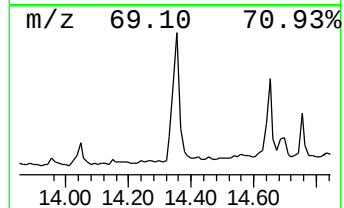
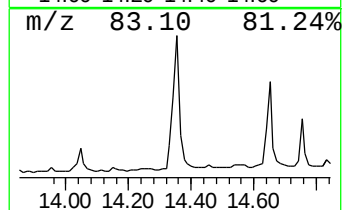
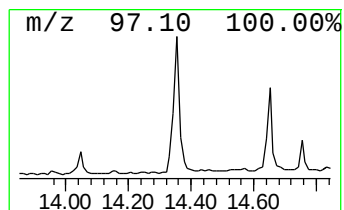
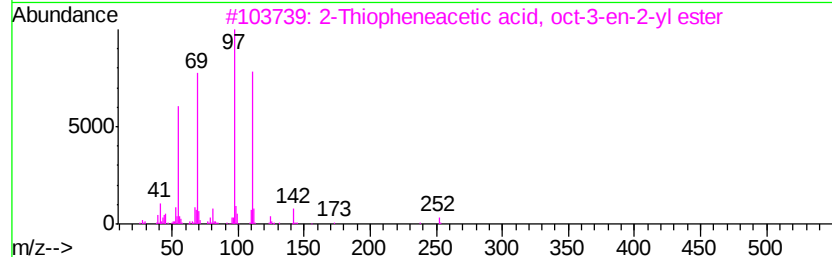
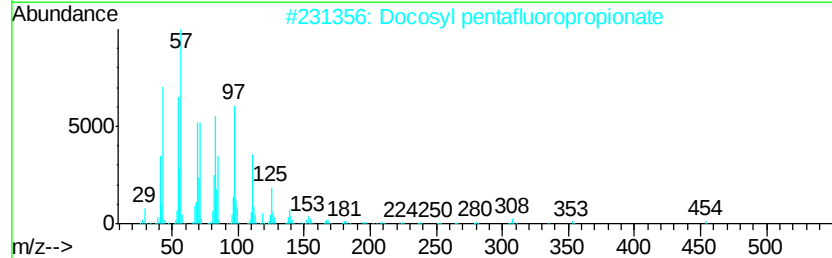
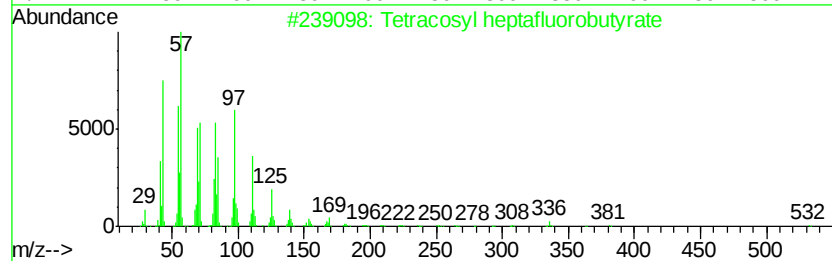
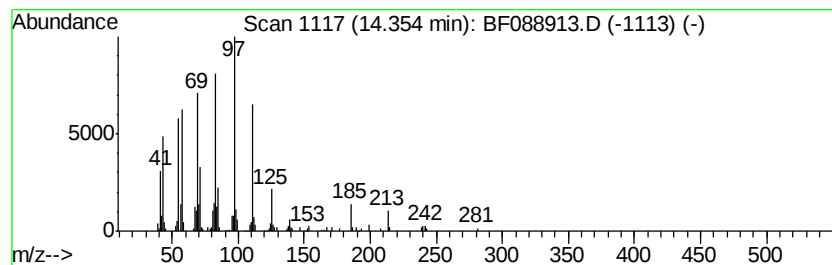
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown14.35 Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.35 | 12.60 ng | 301214 | Chrysene-d12     | 13.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Tetracosyl heptafluorobutyrate      | 550 | C28H49F7O2 | 1000351-83-7 | 38   |
| 2    |    | Docosyl pentafluoropropionate       | 472 | C25H45F5O2 | 1000351-80-9 | 38   |
| 3    |    | 2-Thiopheneacetic acid, oct-3-en... | 252 | C14H20O2S  | 1000299-38-6 | 38   |
| 4    |    | Tricosyl trifluoroacetate           | 436 | C25H47F3O2 | 1000351-75-1 | 38   |
| 5    |    | Tricosyl pentafluoropropionate      | 486 | C26H47F5O2 | 1000351-81-0 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

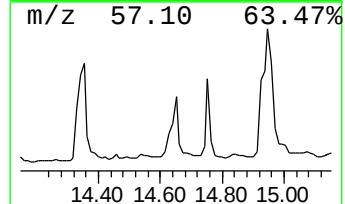
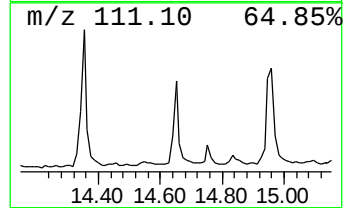
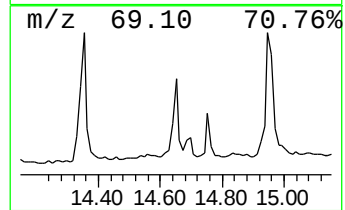
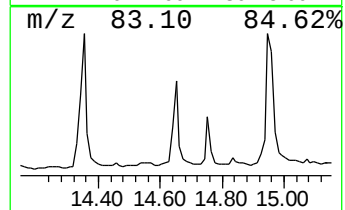
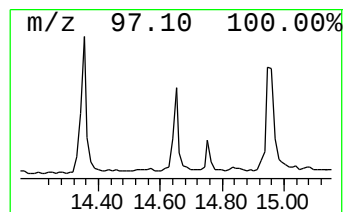
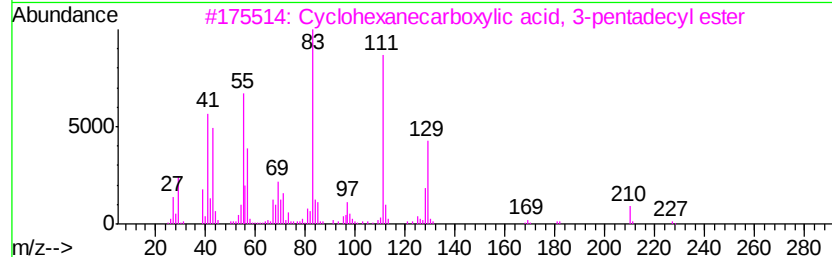
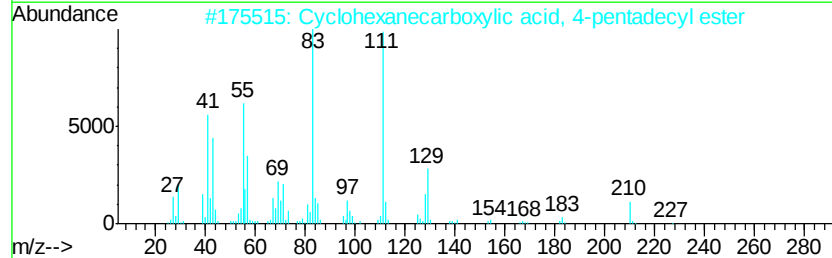
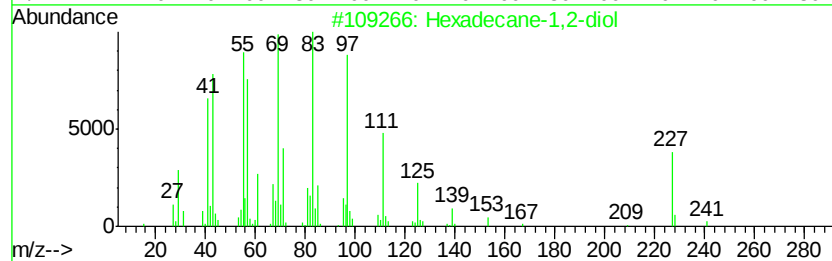
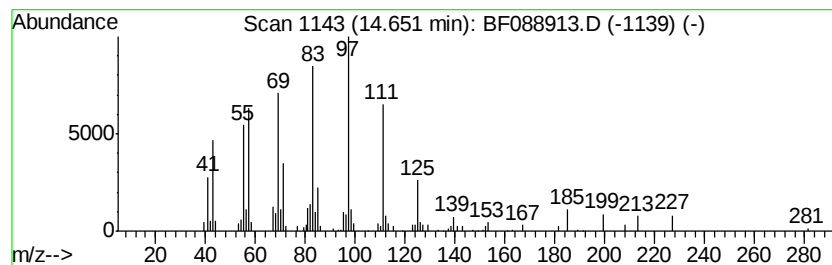
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 unknown14.65 Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.65 | 8.89 ng | 137909 | Perylene-d12     | 15.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Hexadecane-1,2-diol                 | 258 | C16H34O2   | 006920-24-7  | 47   |
| 2    |    | Cyclohexanecarboxylic acid, 4-pe... | 338 | C22H42O2   | 1000280-60-1 | 38   |
| 3    |    | Cyclohexanecarboxylic acid, 3-pe... | 338 | C22H42O2   | 1000280-60-0 | 38   |
| 4    |    | 2-Thiopheneacetic acid, oct-3-en... | 252 | C14H20O2S  | 1000299-38-6 | 35   |
| 5    |    | Dotriacontyl heptafluorobutyrate    | 662 | C36H65F7O2 | 1000351-84-2 | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

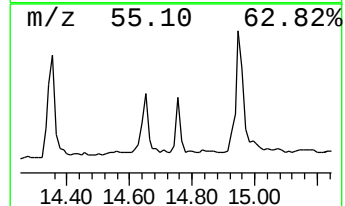
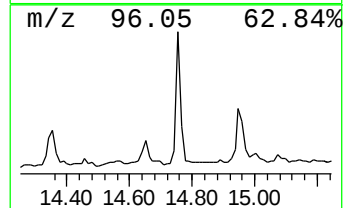
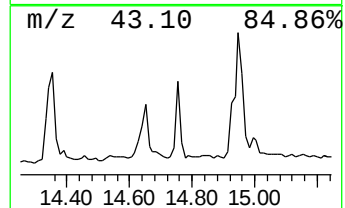
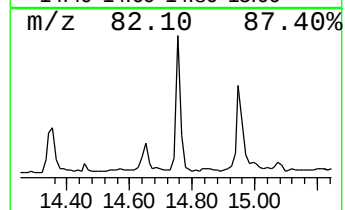
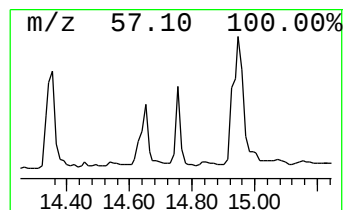
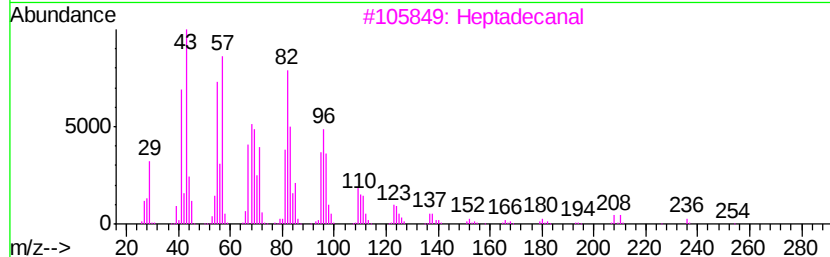
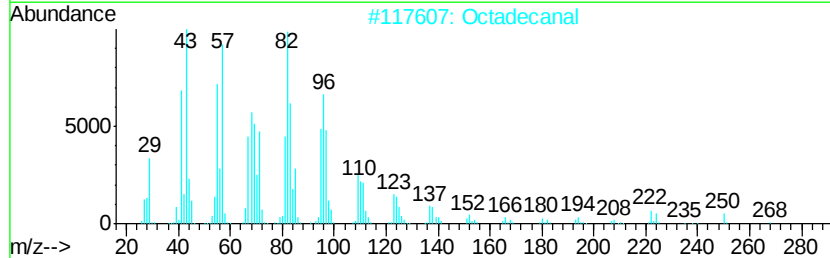
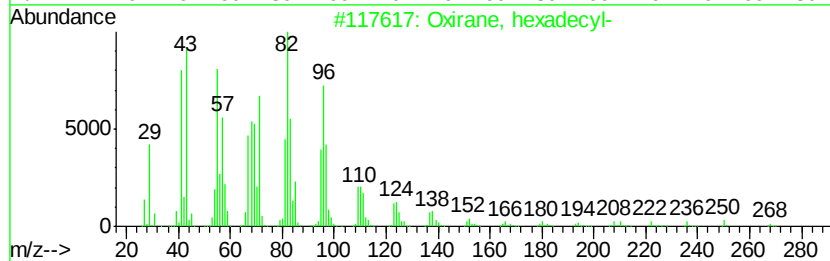
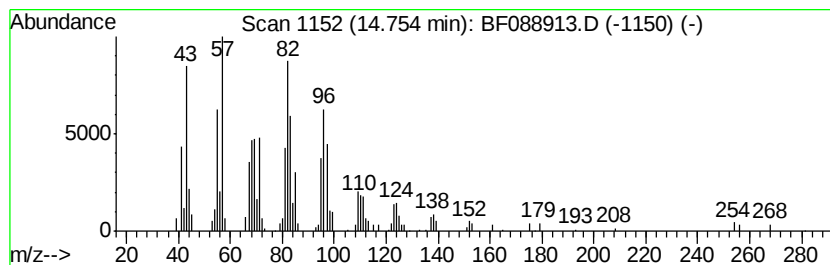
Instrument :  
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ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.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 10 Oxirane, hexadecyl- Concentration Rank 11

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.         |      |
|-----------|----------------------|--------|------------------|--------------|------|
| 14.75     | 6.92 ng              | 107386 | Perylene-d12     | 15.21        |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#         | Qual |
| 1         | Oxirane, hexadecyl-  | 268    | C18H36O          | 007390-81-0  | 91   |
| 2         | Octadecanal          | 268    | C18H36O          | 000638-66-4  | 91   |
| 3         | Heptadecanal         | 254    | C17H34O          | 1000376-70-0 | 91   |
| 4         | Oxirane, heptadecyl- | 282    | C19H38O          | 067860-04-2  | 91   |
| 5         | Hexadecanal          | 240    | C16H32O          | 000629-80-1  | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

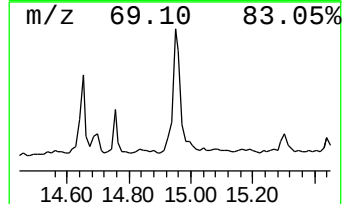
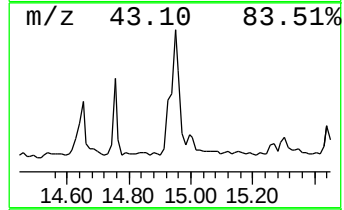
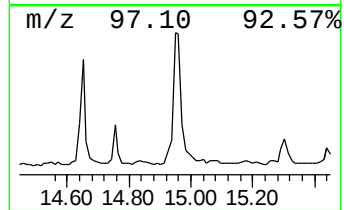
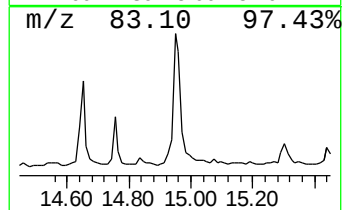
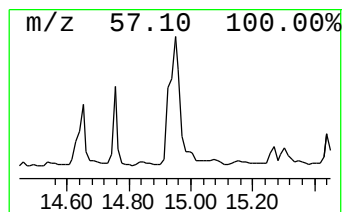
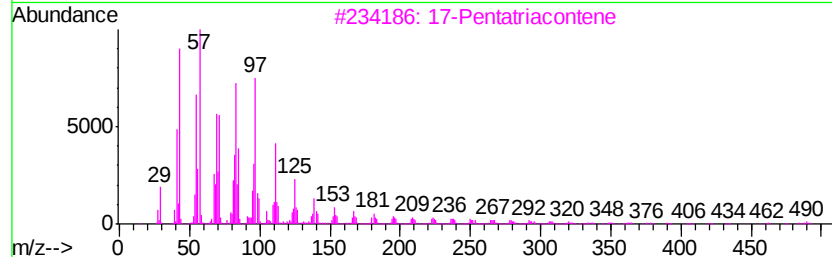
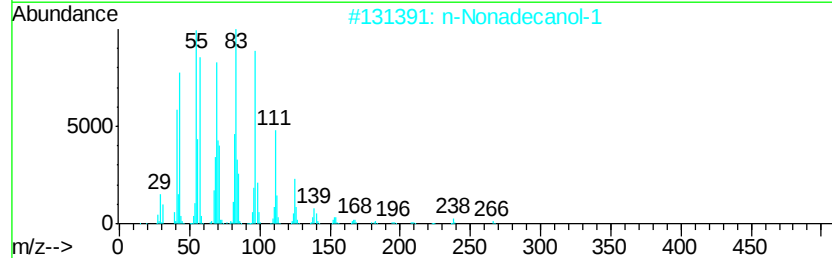
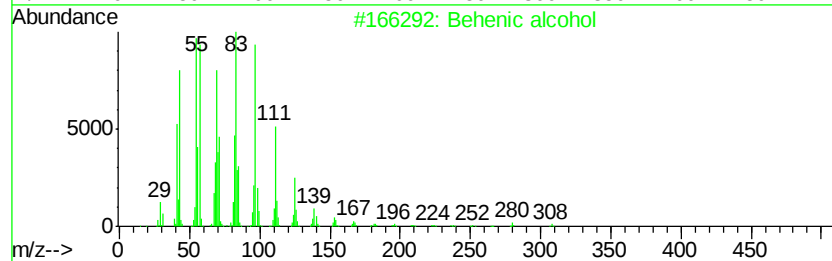
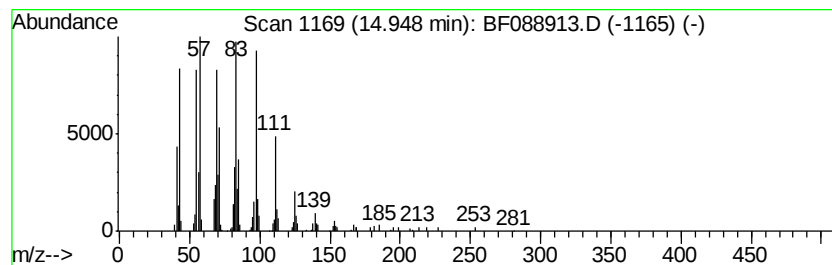
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Behenic alcohol Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.95 | 22.67 ng | 351680 | Perylene-d12     | 15.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Behenic alcohol                     | 326 | C22H46O    | 000661-19-8 | 94   |
| 2    |    | n-Nonadecanol-1                     | 284 | C19H40O    | 001454-84-8 | 76   |
| 3    |    | 17-Pentatriacontene                 | 491 | C35H70     | 006971-40-0 | 76   |
| 4    |    | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2 | 000400-57-7 | 70   |
| 5    |    | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 70   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

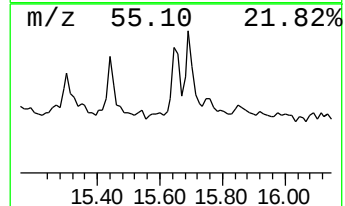
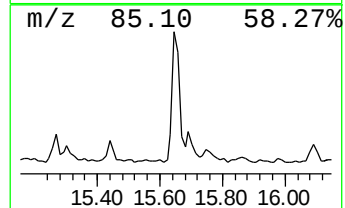
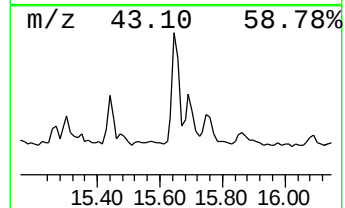
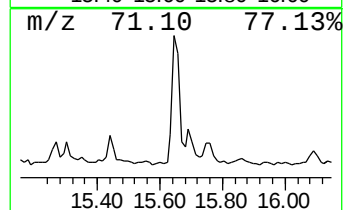
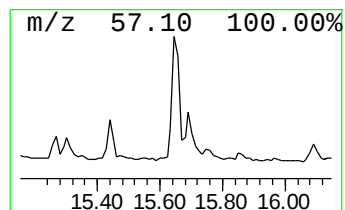
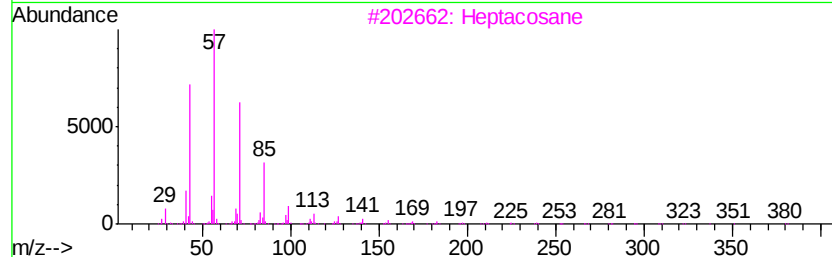
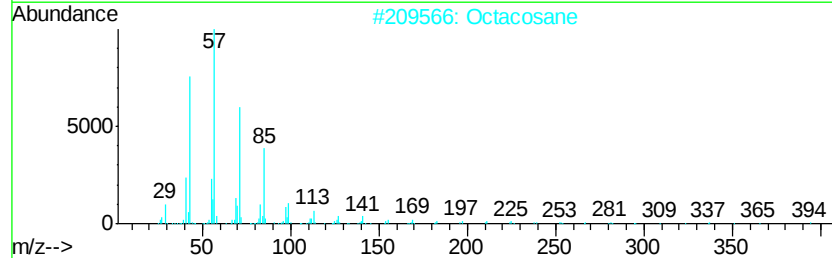
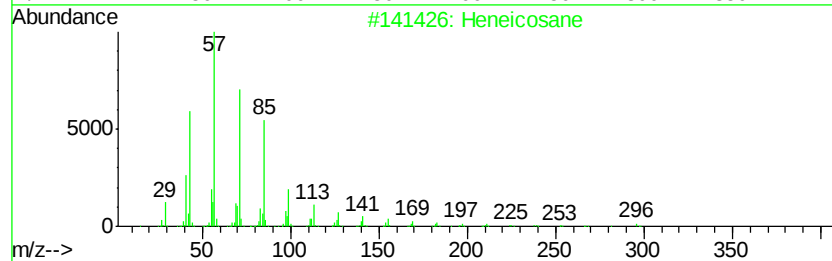
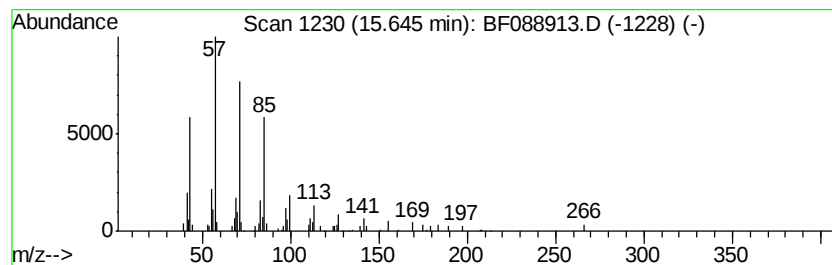
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Heneicosane Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.65 | 6.72 ng | 104183 | Perylene-d12     | 15.21 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 91   |
| 3    |    |   | Heptacosane           | 380 | C27H56  | 000593-49-7 | 91   |
| 4    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 5    |    |   | Octadecane            | 254 | C18H38  | 000593-45-3 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

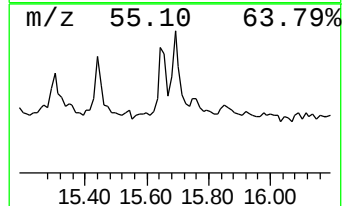
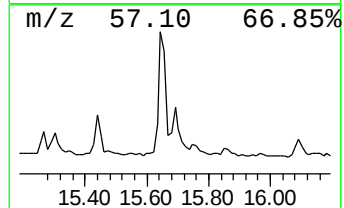
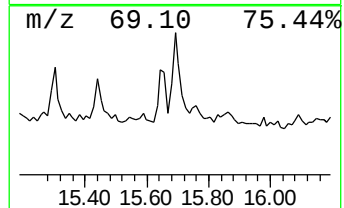
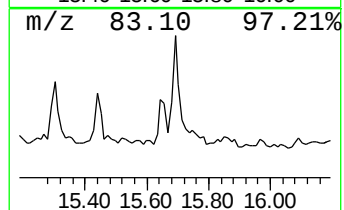
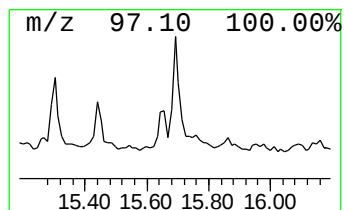
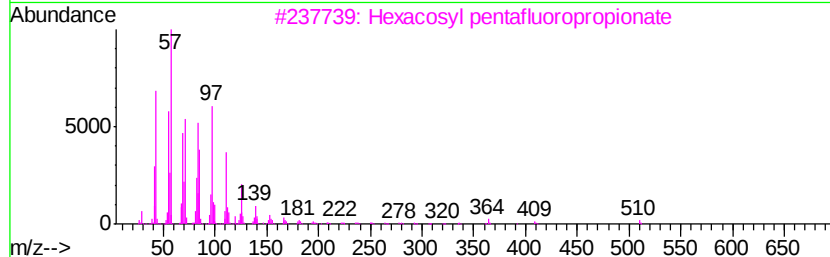
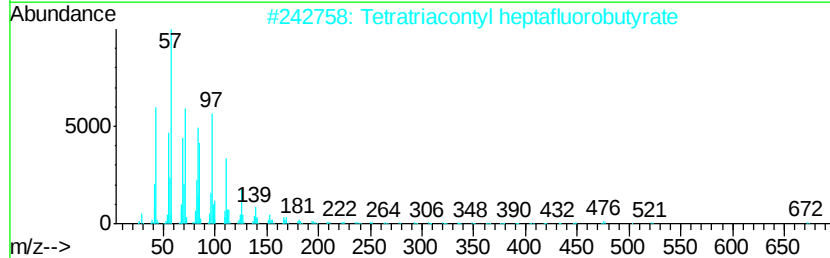
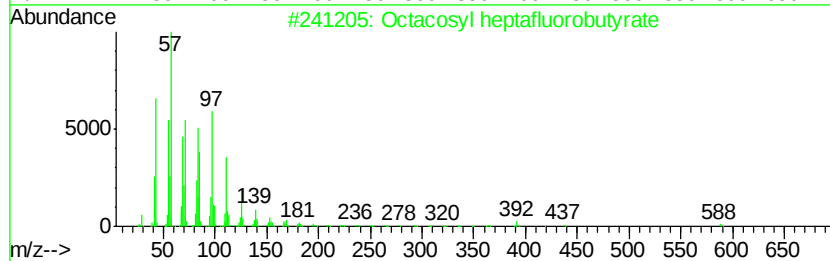
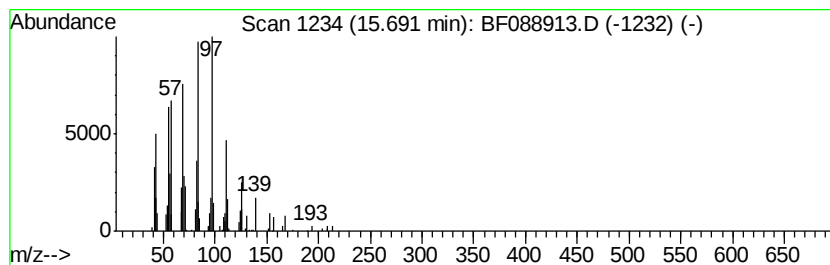
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown15.69 Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.69 | 5.46 ng | 84653 | Perylene-d12     | 15.21 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Octacosyl heptafluorobutyrate       | 606 | C32H57F7O2 | 1000351-83-6 | 43   |
| 2    |    |   | Tetratriacontyl heptafluorobutyrate | 691 | C38H69F7O2 | 1000351-84-1 | 43   |
| 3    |    |   | Hexacosyl pentafluoropropionate     | 528 | C29H53F5O2 | 1000351-81-2 | 43   |
| 4    |    |   | Hexatriacontyl pentafluoropropio... | 669 | C39H73F5O2 | 1000351-89-0 | 43   |
| 5    |    |   | Dotriacontyl heptafluorobutyrate    | 662 | C36H65F7O2 | 1000351-84-2 | 43   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

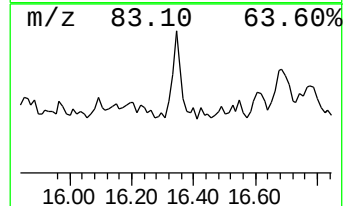
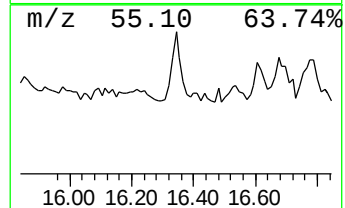
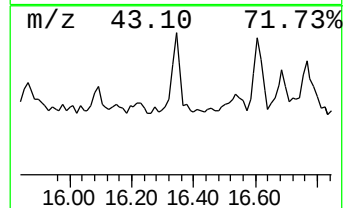
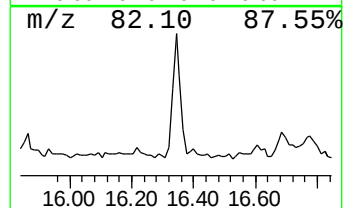
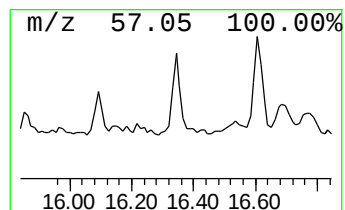
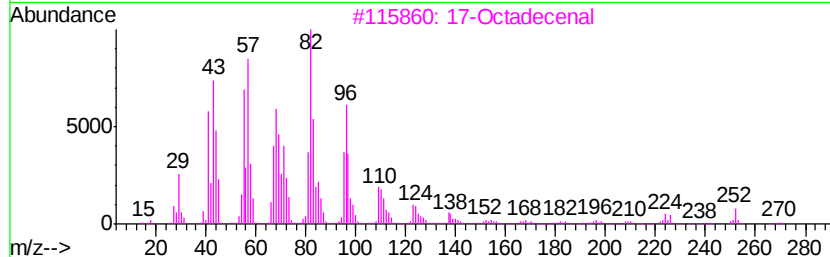
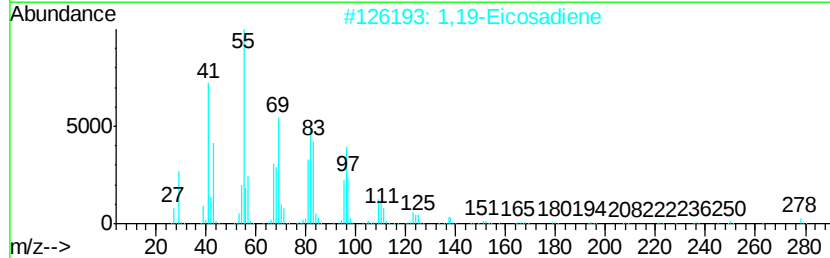
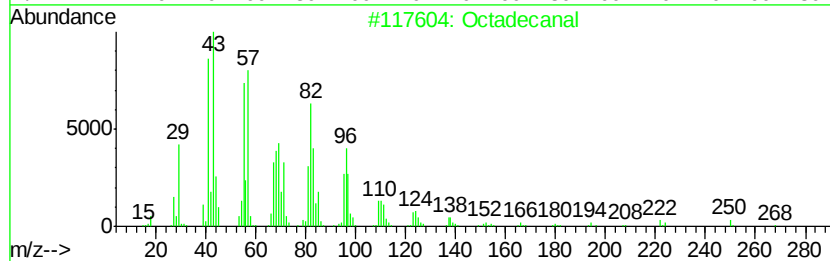
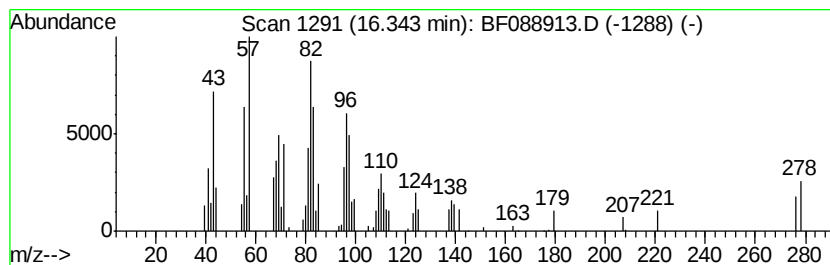
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ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.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 15 Octadecanal Concentration Rank 17

| R.T.      | EstConc                 | Area  | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|-------|------------------|-------------|------|
| 16.34     | 3.82 ng                 | 59328 | Perylene-d12     | 15.21       |      |
| Hit# of 5 | Tentative ID            | MW    | MolForm          | CAS#        | Qual |
| 1         | Octadecanal             | 268   | C18H36O          | 000638-66-4 | 72   |
| 2         | 1,19-Eicosadiene        | 278   | C20H38           | 014811-95-1 | 58   |
| 3         | 17-Octadecenal          | 266   | C18H34O          | 056554-86-0 | 58   |
| 4         | Oxirane, hexadecyl-     | 268   | C18H36O          | 007390-81-0 | 58   |
| 5         | 1-Dodecen-1-ol, acetate | 226   | C14H26O2         | 056438-08-5 | 52   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

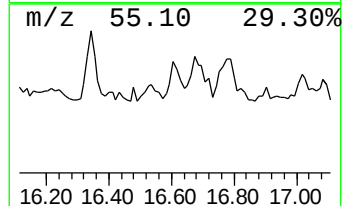
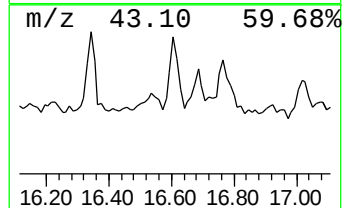
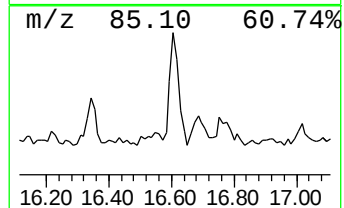
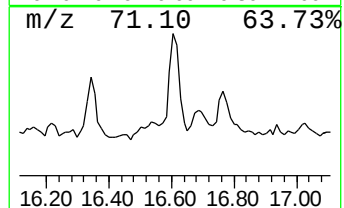
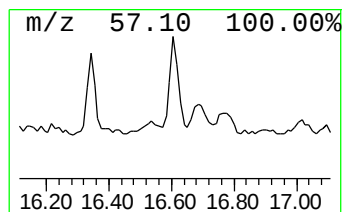
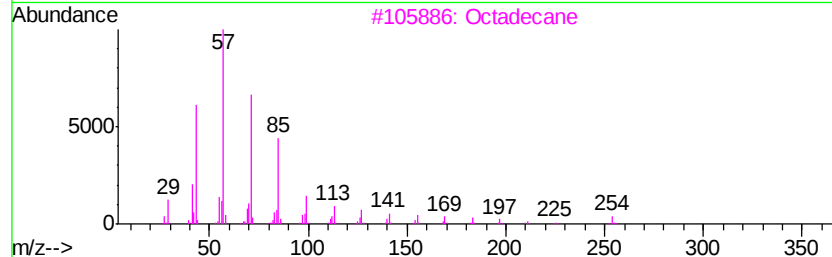
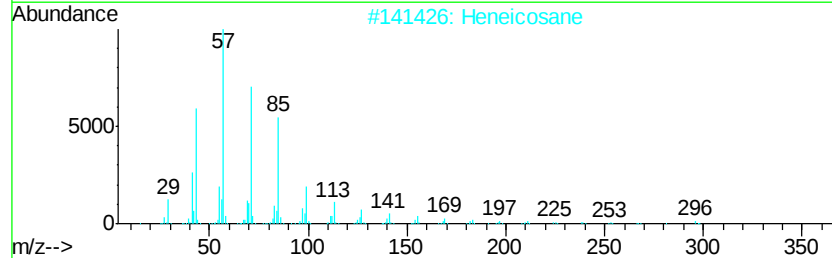
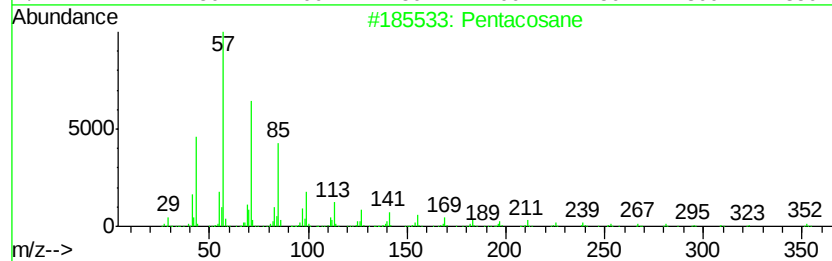
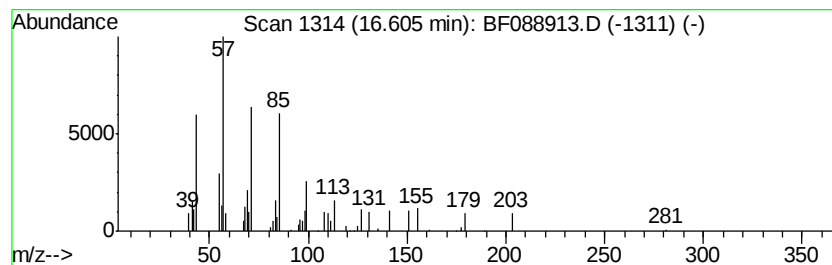
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Pentacosane Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.61 | 2.90 ng | 44913 | Perylene-d12     | 15.21 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Pentacosane         | 352 | C25H52  | 000629-99-2 | 59   |
| 2    |    | Heneicosane         | 296 | C21H44  | 000629-94-7 | 53   |
| 3    |    | Octadecane          | 254 | C18H38  | 000593-45-3 | 52   |
| 4    |    | Eicosane            | 282 | C20H42  | 000112-95-8 | 50   |
| 5    |    | Octadecane, 1-iodo- | 380 | C18H37I | 000629-93-6 | 50   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

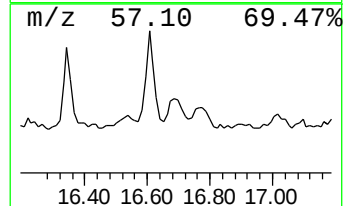
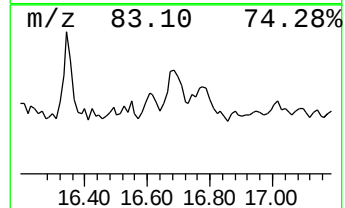
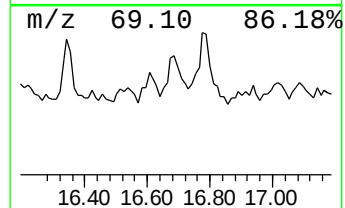
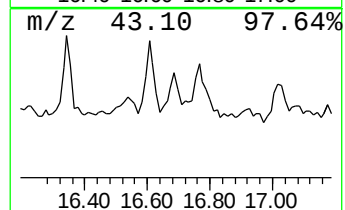
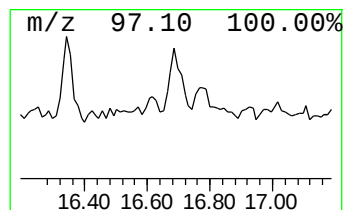
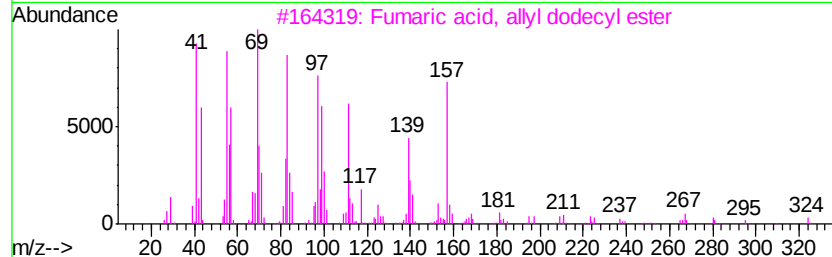
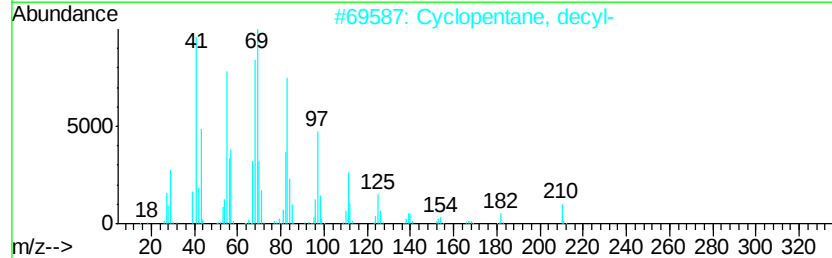
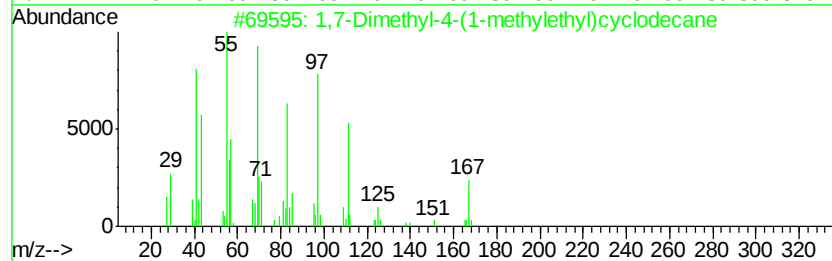
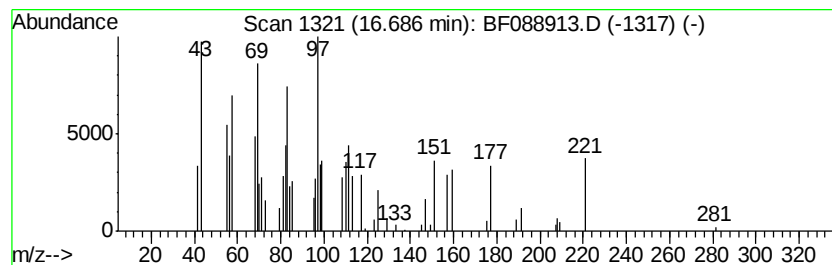
Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.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 17 1,7-Dimethyl-4-(1-methyleth... Concentration Rank 16

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 16.69     | 4.21 ng                             | 65355 | Perylene-d12     | 15.21        |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | 1,7-Dimethyl-4-(1-methylethyl)cy... | 210   | C15H30           | 000645-10-3  | 53   |
| 2         | Cyclopentane, decyl-                | 210   | C15H30           | 001795-21-7  | 47   |
| 3         | Fumaric acid, allyl dodecyl ester   | 324   | C19H32O4         | 1000330-12-8 | 47   |
| 4         | Acetic acid, chloro-, hexadecyl ... | 318   | C18H35ClO2       | 052132-58-8  | 47   |
| 5         | 1-Heneicosanol                      | 312   | C21H44O          | 015594-90-8  | 43   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

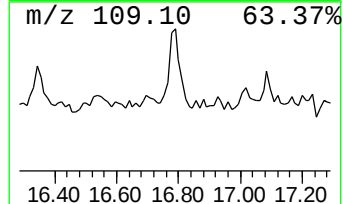
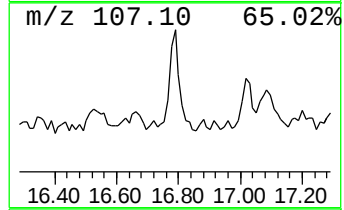
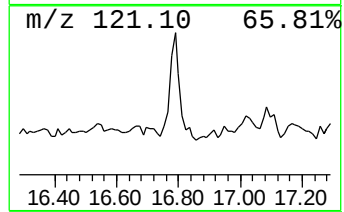
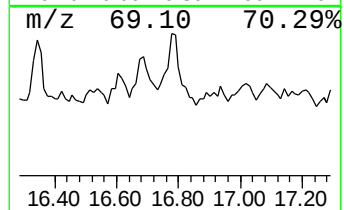
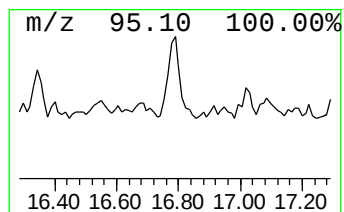
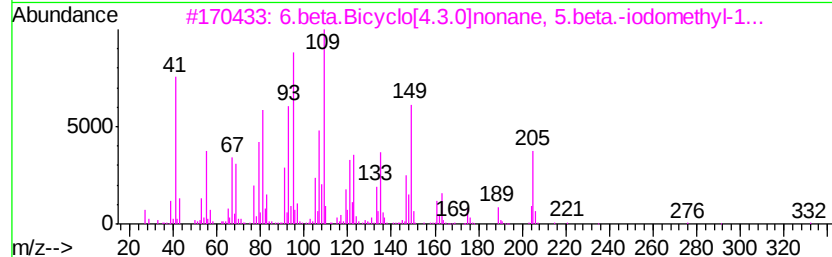
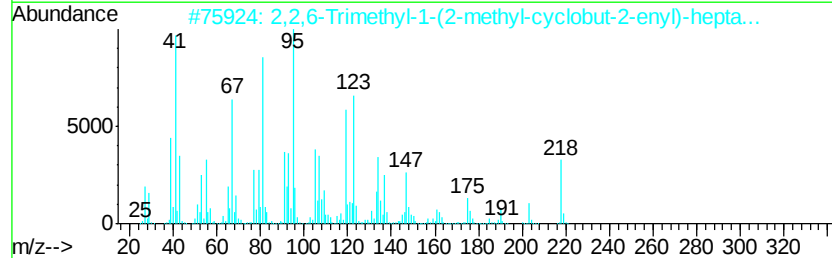
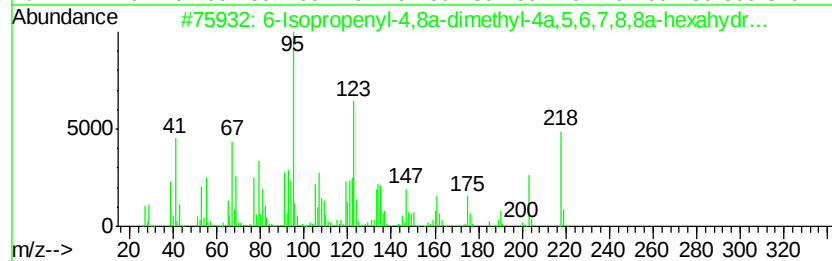
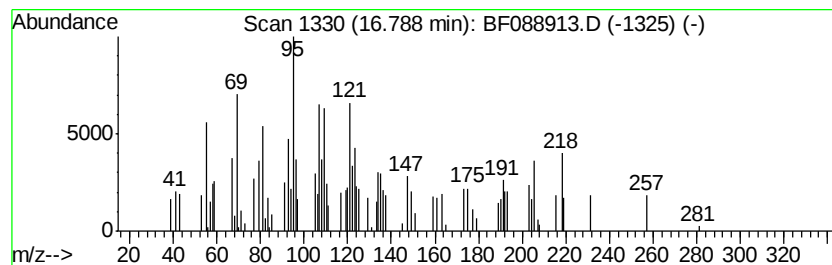
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 6-Isopropenyl-4,8a-dimethyl... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.79 | 7.09 ng | 109923 | Perylene-d12     | 15.21 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 6-Isopropenyl-4,8a-dimethyl-4a,5... | 218 | C15H22O  | 086917-79-5  | 55   |
| 2    |    |   | 2,2,6-Trimethyl-1-(2-methyl-cycl... | 218 | C15H22O  | 1000188-72-8 | 46   |
| 3    |    |   | 6.beta.Bicyclo[4.3.0]nonane, 5.b... | 332 | C15H25I  | 1000195-85-9 | 35   |
| 4    |    |   | Caparratriene                       | 206 | C15H26   | 1000374-08-7 | 27   |
| 5    |    |   | Isobornyl acetate                   | 196 | C12H20O2 | 000125-12-2  | 18   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

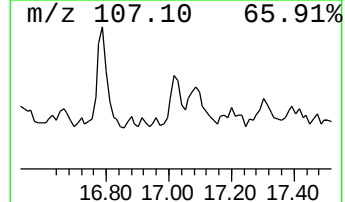
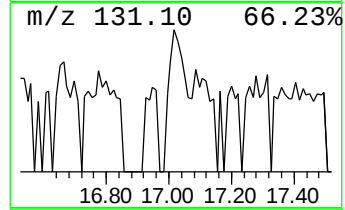
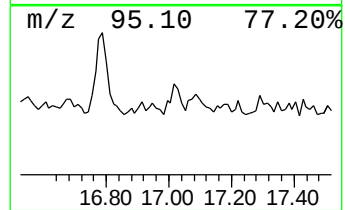
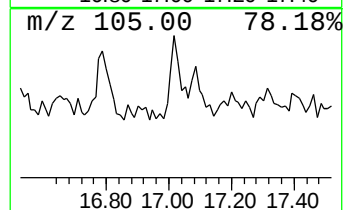
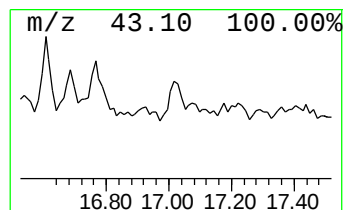
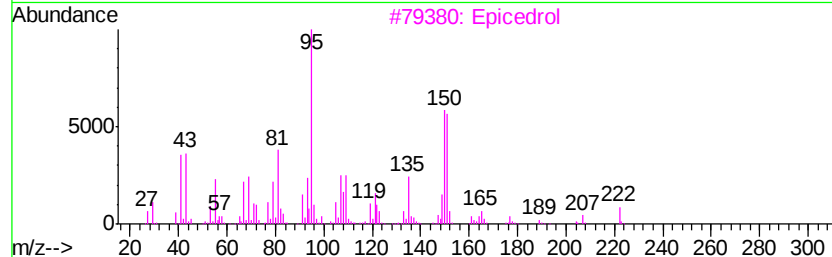
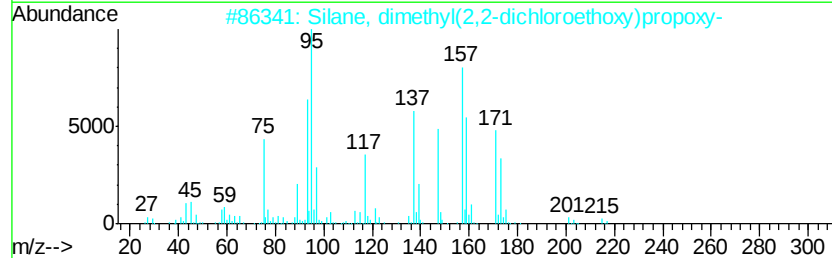
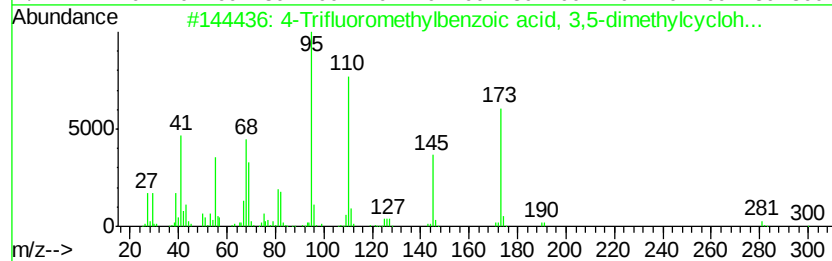
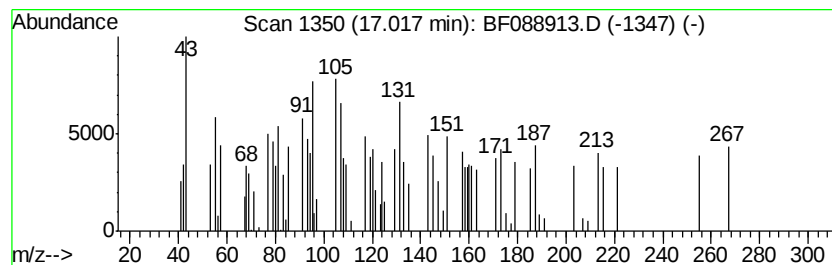
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 unknown17.02 Concentration Rank 18

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 17.02 | 3.26 ng | 50644 | Perylene-d12     | 15.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 4-Trifluoromethylbenzoic acid, 3... | 300 | C16H19F3O2   | 1000280-03-3 | 10   |
| 2    |    | Silane, dimethyl(2,2-dichloroeth... | 230 | C7H16Cl2O2Si | 1000347-60-9 | 10   |
| 3    |    | Epicedrol                           | 222 | C15H26O      | 1000156-22-8 | 10   |
| 4    |    | 2,4-Hexadienoanilide                | 187 | C12H13NO     | 001483-88-1  | 9    |
| 5    |    | Acetic acid, 2-(2-acetyl-3-oxois... | 229 | C10H15NO5    | 1000185-79-0 | 9    |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\  
Data File : BF088913.D  
Acq On : 11 Jul 2016 21:51  
Operator : UM/SJ  
Sample : H3921-08  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

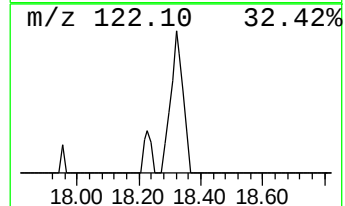
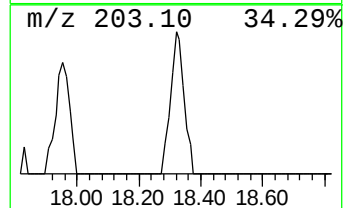
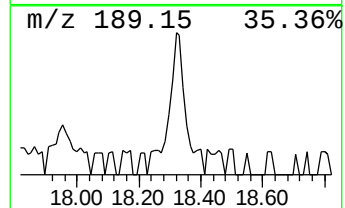
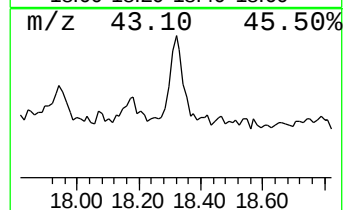
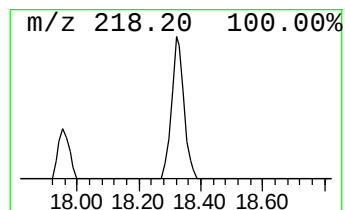
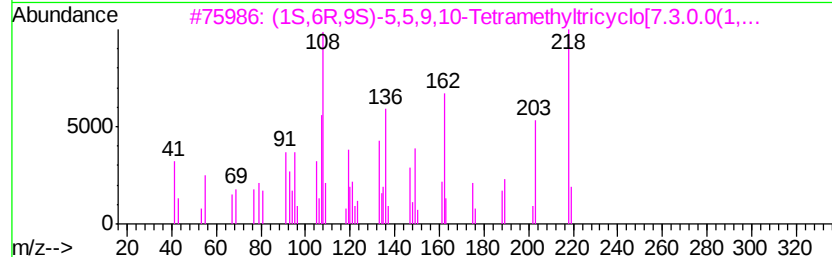
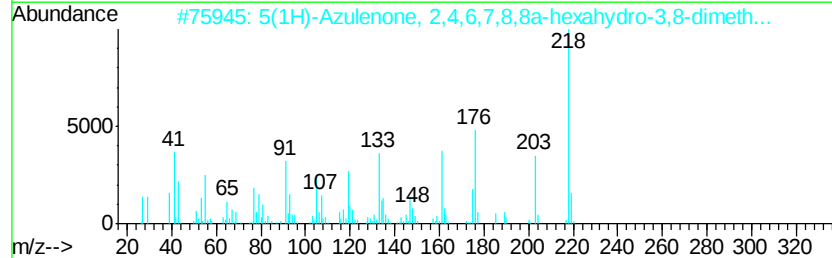
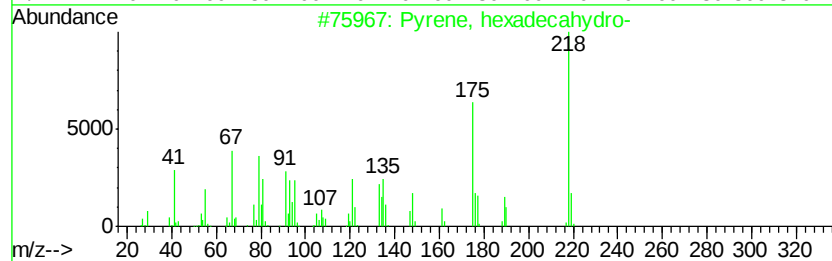
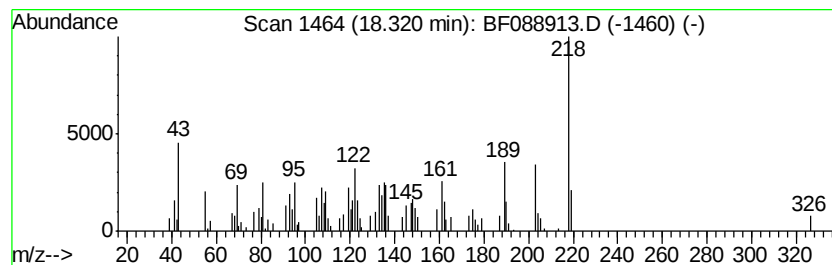
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Pyrene, hexadecahydro- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.32 | 8.91 ng | 138245 | Perylene-d12     | 15.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Pyrene, hexadecahydro-              | 218 | C16H26  | 002435-85-0  | 83   |
| 2    |    | 5(1H)-Azulenone, 2,4,6,7,8,8a-he... | 218 | C15H22O | 006754-66-1  | 70   |
| 3    |    | (1S,6R,9S)-5,5,9,10-Tetramethylt... | 218 | C16H26  | 1000298-97-8 | 64   |
| 4    |    | 2(1H)Naphthalenone, 3,5,6,7,8,8a... | 218 | C15H22O | 1000188-66-5 | 53   |
| 5    |    | 2(3H)-Naphthalenone, 4,4a,5,6,7,... | 218 | C15H22O | 019598-45-9  | 46   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF071116\  
 Data File : BF088913.D  
 Acq On : 11 Jul 2016 21:51  
 Operator : UM/SJ  
 Sample : H3921-08  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 N8-B6(0-2)C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF063016.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 |
| Furan, tetrahydro... | 1.92  | 23.6    | ng    | 431623   | 1                     | 6.71  | 366631 | 20.0 |
| 2-Pentanone, 4-hy... | 4.87  | 5.2     | ng    | 94424    | 1                     | 6.71  | 366631 | 20.0 |
| unknown6.48          | 6.48  | 78.3    | ng    | 1436010  | 1                     | 6.71  | 366631 | 20.0 |
| unknown14.35         | 14.35 | 12.6    | ng    | 301214   | 5                     | 13.84 | 478117 | 20.0 |
| unknown14.65         | 14.65 | 8.9     | ng    | 137909   | 6                     | 15.21 | 310239 | 20.0 |
| Oxirane, hexadecyl-  | 14.75 | 6.9     | ng    | 107386   | 6                     | 15.21 | 310239 | 20.0 |
| Behenic alcohol      | 14.95 | 22.7    | ng    | 351680   | 6                     | 15.21 | 310239 | 20.0 |
| Heneicosane          | 15.65 | 6.7     | ng    | 104183   | 6                     | 15.21 | 310239 | 20.0 |
| unknown15.69         | 15.69 | 5.5     | ng    | 84653    | 6                     | 15.21 | 310239 | 20.0 |
| Octadecanal          | 16.34 | 3.8     | ng    | 59328    | 6                     | 15.21 | 310239 | 20.0 |
| Pentacosane          | 16.61 | 2.9     | ng    | 44913    | 6                     | 15.21 | 310239 | 20.0 |
| 1,7-Dimethyl-4-(1... | 16.69 | 4.2     | ng    | 65355    | 6                     | 15.21 | 310239 | 20.0 |
| 6-Isopropenyl-4,8... | 16.79 | 7.1     | ng    | 109923   | 6                     | 15.21 | 310239 | 20.0 |
| unknown17.02         | 17.02 | 3.3     | ng    | 50644    | 6                     | 15.21 | 310239 | 20.0 |
| Pyrene, hexadecah... | 18.32 | 8.9     | ng    | 138245   | 6                     | 15.21 | 310239 | 20.0 |