

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
 Data File : BF097446.D  
 Acq On : 7 Aug 2017 23:13  
 Operator : SJ/JU  
 Sample : I4639-04 5X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SA-06-080317-B

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-BF072517.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         | 5.004       | 544           | 547         | 568          | rBV      | 264868         | 449807        | 8.54%           | 1.345%        |
| 2         | 6.087       | 728           | 731         | 745          | rBV      | 317140         | 483711        | 9.18%           | 1.447%        |
| 3         | 6.187       | 745           | 748         | 757          | rVB      | 493205         | 512195        | 9.72%           | 1.532%        |
| 4         | 6.410       | 783           | 786         | 795          | rBV      | 601429         | 532160        | 10.10%          | 1.592%        |
| 5         | 6.563       | 809           | 812         | 820          | rBV      | 322316         | 338424        | 6.42%           | 1.012%        |
| 6         | 6.981       | 881           | 883         | 890          | rBV      | 236473         | 262780        | 4.99%           | 0.786%        |
| 7         | 7.692       | 1001          | 1004        | 1012         | rBV      | 848476         | 769513        | 14.60%          | 2.302%        |
| 8         | 8.310       | 1104          | 1109        | 1114         | rBV      | 190668         | 178965        | 3.40%           | 0.535%        |
| 9         | 8.516       | 1139          | 1144        | 1147         | rBV3     | 59222          | 73848         | 1.40%           | 0.221%        |
| 10        | 8.734       | 1179          | 1181        | 1184         | rVV2     | 80805          | 66301         | 1.26%           | 0.198%        |
| 11        | 8.769       | 1184          | 1187        | 1196         | rVB      | 568601         | 489636        | 9.29%           | 1.464%        |
| 12        | 8.869       | 1200          | 1204        | 1207         | rBV      | 265664         | 244871        | 4.65%           | 0.732%        |
| 13        | 9.128       | 1244          | 1248        | 1249         | rBV2     | 45839          | 56059         | 1.06%           | 0.168%        |
| 14        | 9.175       | 1253          | 1256        | 1257         | rBV      | 105813         | 97348         | 1.85%           | 0.291%        |
| 15        | 9.398       | 1290          | 1294        | 1297         | rBV      | 411625         | 345017        | 6.55%           | 1.032%        |
| 16        | 9.434       | 1297          | 1300        | 1305         | rVV2     | 760904         | 686774        | 13.03%          | 2.054%        |
| 17        | 9.628       | 1327          | 1333        | 1335         | rBV2     | 64107          | 104288        | 1.98%           | 0.312%        |
| 18        | 9.651       | 1335          | 1337        | 1341         | rVV3     | 68318          | 78813         | 1.50%           | 0.236%        |
| 19        | 9.716       | 1345          | 1348        | 1351         | rVV      | 119025         | 110526        | 2.10%           | 0.331%        |
| 20        | 9.751       | 1351          | 1354        | 1361         | rVV4     | 62097          | 93376         | 1.77%           | 0.279%        |
| 21        | 9.892       | 1375          | 1378        | 1381         | rVB      | 442250         | 340477        | 6.46%           | 1.018%        |
| 22        | 10.110      | 1410          | 1415        | 1418         | rBV      | 147865         | 186802        | 3.55%           | 0.559%        |
| 23        | 10.192      | 1426          | 1429        | 1432         | rBV      | 60501          | 56764         | 1.08%           | 0.170%        |
| 24        | 10.228      | 1432          | 1435        | 1439         | rVV3     | 225685         | 235550        | 4.47%           | 0.704%        |
| 25        | 10.363      | 1455          | 1458        | 1464         | rBV      | 422282         | 456157        | 8.66%           | 1.364%        |
| 26        | 10.557      | 1487          | 1491        | 1494         | rBV2     | 60654          | 74610         | 1.42%           | 0.223%        |
| 27        | 10.680      | 1509          | 1512        | 1516         | rVB3     | 47548          | 53772         | 1.02%           | 0.161%        |
| 28        | 10.804      | 1530          | 1533        | 1536         | rBV      | 331900         | 308809        | 5.86%           | 0.924%        |
| 29        | 10.839      | 1536          | 1539        | 1544         | rVB      | 116669         | 131192        | 2.49%           | 0.392%        |
| 30        | 10.910      | 1547          | 1551        | 1554         | rBV      | 616794         | 524655        | 9.96%           | 1.569%        |
| 31        | 10.992      | 1561          | 1565        | 1568         | rBV4     | 45535          | 68518         | 1.30%           | 0.205%        |
| 32        | 11.233      | 1602          | 1606        | 1608         | rBV      | 281768         | 262641        | 4.98%           | 0.786%        |
| 33        | 11.333      | 1619          | 1623        | 1626         | rBV      | 2824368        | 2626727       | 49.85%          | 7.856%        |
| 34        | 11.480      | 1643          | 1648        | 1653         | rBV      | 754880         | 799735        | 15.18%          | 2.392%        |

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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-BF072517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |         |
|----|--------|------|------|------|------|---------|---------|---------|---------|
| 35 | 11.522 | 1653 | 1655 | 1666 | rVB5 | 62308   | 129691  | 2.46%   | 0.388%  |
| 36 | 11.633 | 1671 | 1674 | 1682 | rVB  | 232949  | 286960  | 5.45%   | 0.858%  |
| 37 | 12.016 | 1734 | 1739 | 1741 | rBV  | 3619789 | 5011780 | 95.12%  | 14.990% |
| 38 | 12.045 | 1741 | 1744 | 1749 | rVV  | 4316213 | 5269068 | 100.00% | 15.759% |
| 39 | 12.122 | 1753 | 1757 | 1760 | rBV2 | 1800403 | 1642896 | 31.18%  | 4.914%  |
| 40 | 12.192 | 1760 | 1769 | 1778 | rBV2 | 1812166 | 3686784 | 69.97%  | 11.027% |
| 41 | 12.263 | 1778 | 1781 | 1785 | rVB  | 215681  | 206446  | 3.92%   | 0.617%  |
| 42 | 12.351 | 1791 | 1796 | 1800 | rVB2 | 319651  | 403694  | 7.66%   | 1.207%  |
| 43 | 12.392 | 1800 | 1803 | 1808 | rBV  | 139569  | 127971  | 2.43%   | 0.383%  |
| 44 | 12.492 | 1817 | 1820 | 1823 | rVB  | 426559  | 349181  | 6.63%   | 1.044%  |
| 45 | 12.557 | 1828 | 1831 | 1833 | rBV2 | 70857   | 77382   | 1.47%   | 0.231%  |
| 46 | 12.604 | 1835 | 1839 | 1843 | rVV3 | 87053   | 139535  | 2.65%   | 0.417%  |
| 47 | 12.669 | 1847 | 1850 | 1852 | rVV2 | 198957  | 191171  | 3.63%   | 0.572%  |
| 48 | 12.698 | 1852 | 1855 | 1860 | rVV3 | 189524  | 325502  | 6.18%   | 0.974%  |
| 49 | 12.757 | 1860 | 1865 | 1868 | rVV3 | 128494  | 201302  | 3.82%   | 0.602%  |
| 50 | 12.839 | 1876 | 1879 | 1882 | rVB  | 222828  | 192647  | 3.66%   | 0.576%  |
| 51 | 12.945 | 1894 | 1897 | 1900 | rBV  | 115496  | 116460  | 2.21%   | 0.348%  |
| 52 | 12.998 | 1903 | 1906 | 1909 | rVV2 | 124033  | 143229  | 2.72%   | 0.428%  |
| 53 | 13.027 | 1909 | 1911 | 1918 | rVB2 | 69082   | 74739   | 1.42%   | 0.224%  |
| 54 | 13.086 | 1918 | 1921 | 1925 | rVB  | 71794   | 75597   | 1.43%   | 0.226%  |
| 55 | 13.263 | 1947 | 1951 | 1954 | rVB2 | 236645  | 294202  | 5.58%   | 0.880%  |
| 56 | 13.504 | 1989 | 1992 | 1994 | rBV  | 179418  | 164640  | 3.12%   | 0.492%  |
| 57 | 13.533 | 1994 | 1997 | 2000 | rVV  | 570644  | 590871  | 11.21%  | 1.767%  |
| 58 | 13.563 | 2000 | 2002 | 2005 | rVB  | 64756   | 57766   | 1.10%   | 0.173%  |
| 59 | 13.886 | 2053 | 2057 | 2059 | rBV  | 66945   | 95340   | 1.81%   | 0.285%  |
| 60 | 14.033 | 2078 | 2082 | 2085 | rBV2 | 111892  | 156058  | 2.96%   | 0.467%  |
| 61 | 14.121 | 2095 | 2097 | 2103 | rVB3 | 82734   | 84542   | 1.60%   | 0.253%  |
| 62 | 14.327 | 2129 | 2132 | 2136 | rBV  | 87912   | 95145   | 1.81%   | 0.285%  |
| 63 | 14.421 | 2144 | 2148 | 2154 | rBV  | 235789  | 290477  | 5.51%   | 0.869%  |
| 64 | 14.874 | 2222 | 2225 | 2229 | rBV  | 281221  | 288171  | 5.47%   | 0.862%  |
| 65 | 15.257 | 2288 | 2290 | 2301 | rVB2 | 114933  | 242946  | 4.61%   | 0.727%  |
| 66 | 16.110 | 2430 | 2435 | 2441 | rBV3 | 66924   | 144493  | 2.74%   | 0.432%  |
| 67 | 17.286 | 2628 | 2635 | 2636 | rBV6 | 49170   | 102447  | 1.94%   | 0.306%  |
| 68 | 17.686 | 2699 | 2703 | 2708 | rBV7 | 32954   | 75123   | 1.43%   | 0.225%  |

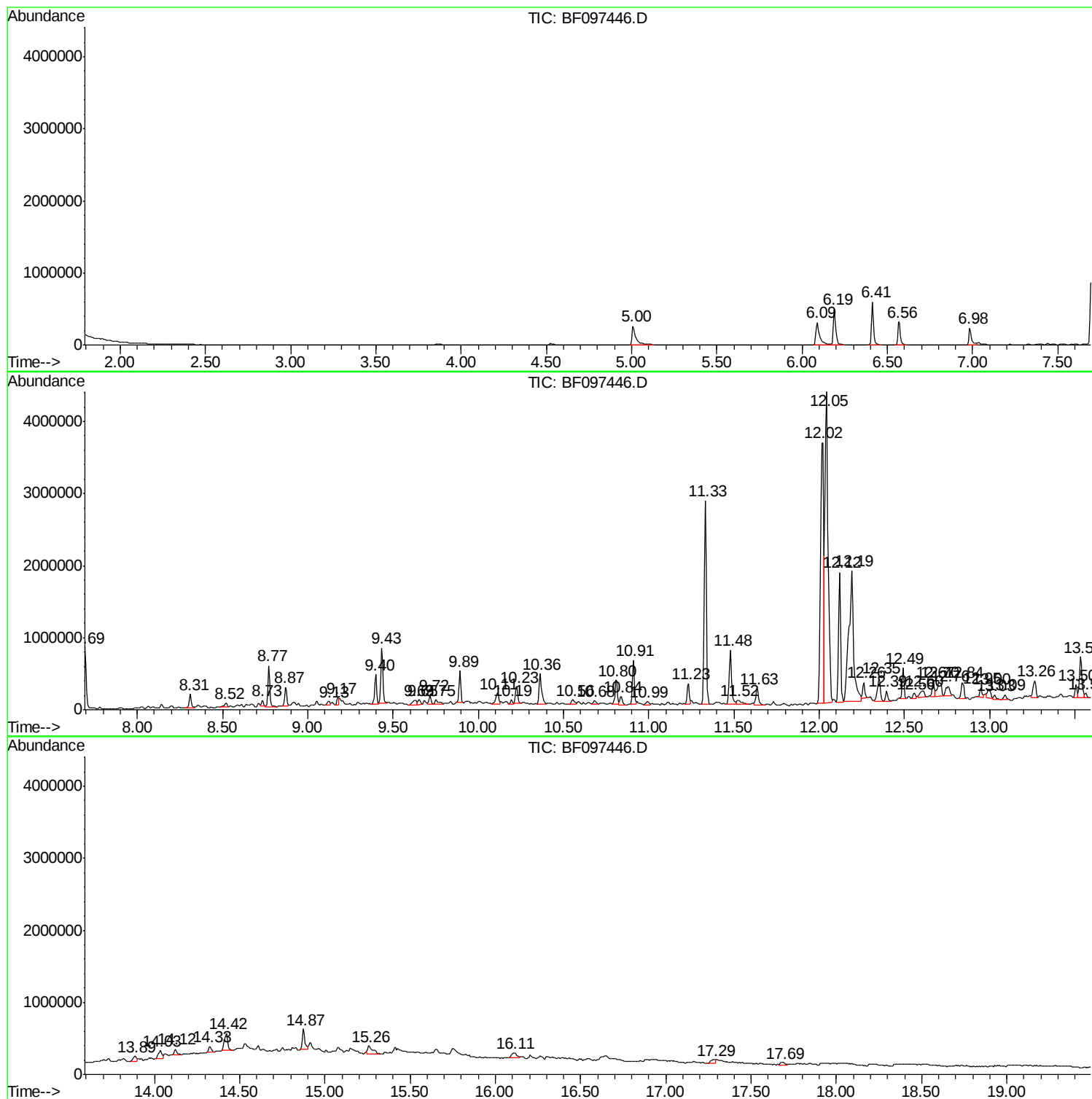
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Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.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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Instrument :  
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SA-06-080317-B

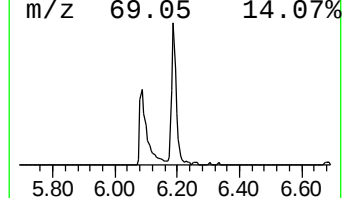
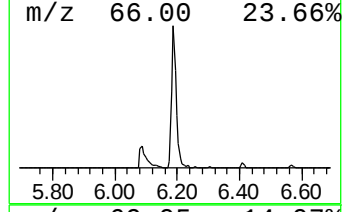
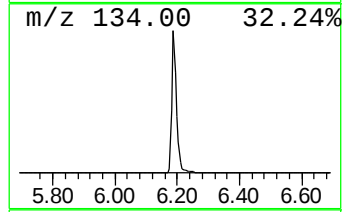
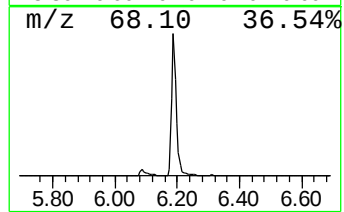
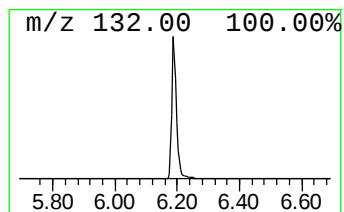
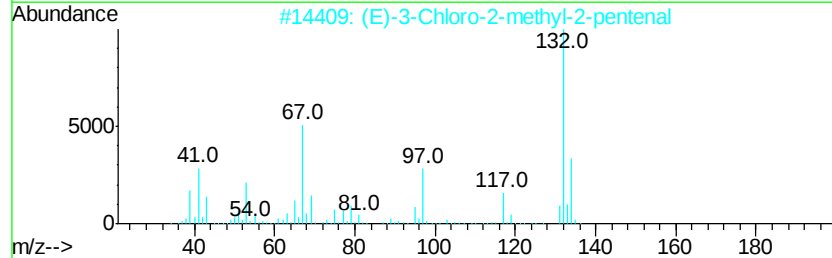
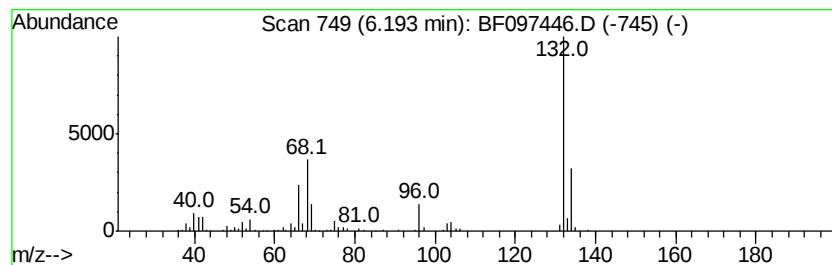
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown6.19 Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.19 | 19.25 ng | 512195 | 1,4-Dichlorobenzene-d4 | 6.41 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 25   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 1-Methylpyrrolo[1,2-a]pyrazine      | 132 | C8H8N2    | 064608-59-9  | 9    |



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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

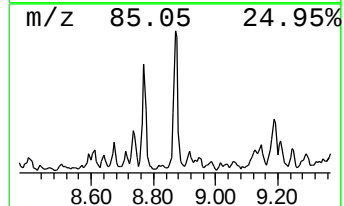
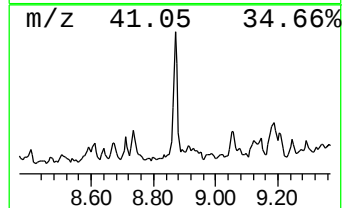
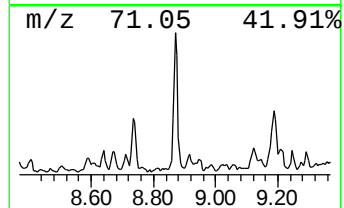
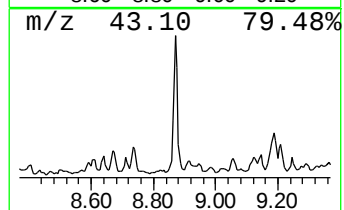
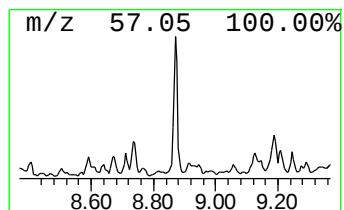
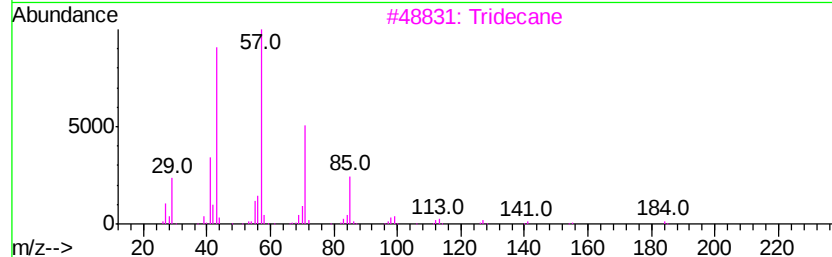
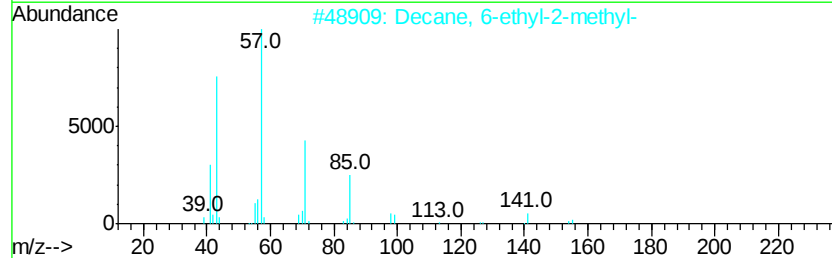
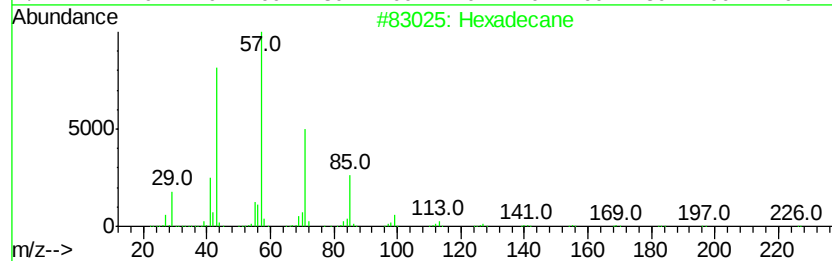
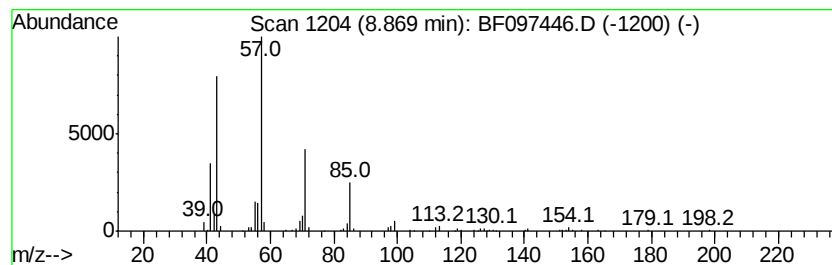
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Hexadecane Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 8.87 | 7.13 ng | 244871 | Acenaphthene-d10 | 9.43 |

| Hit# of | 5                         | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|---------------------------|--------------|-----|---------|-------------|------|
| 1       | Hexadecane                |              | 226 | C16H34  | 000544-76-3 | 90   |
| 2       | Decane, 6-ethyl-2-methyl- |              | 184 | C13H28  | 062108-21-8 | 90   |
| 3       | Tridecane                 |              | 184 | C13H28  | 000629-50-5 | 90   |
| 4       | Tridecane, 7-propyl-      |              | 226 | C16H34  | 055045-09-5 | 80   |
| 5       | Nonane, 5-butyl-          |              | 184 | C13H28  | 017312-63-9 | 64   |



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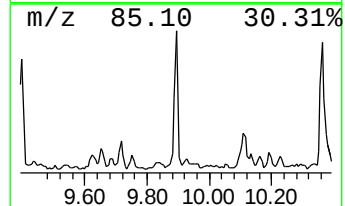
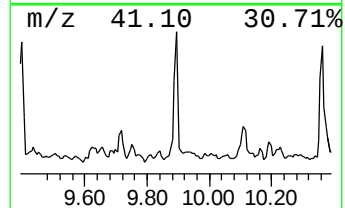
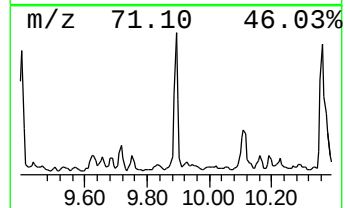
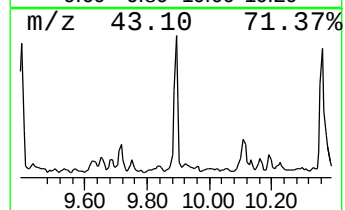
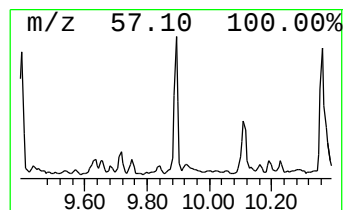
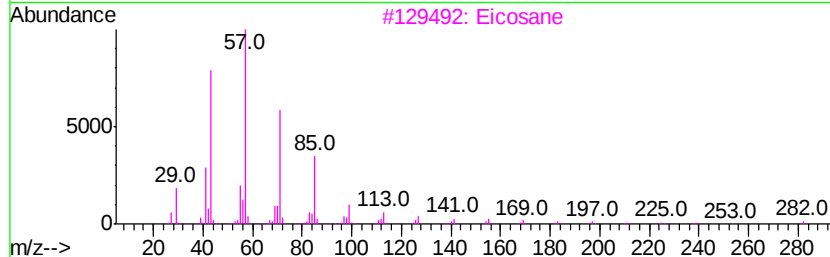
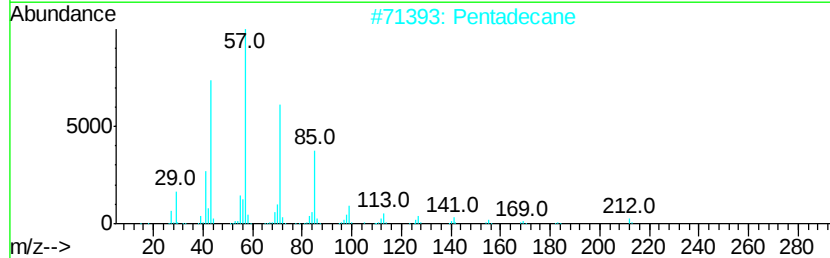
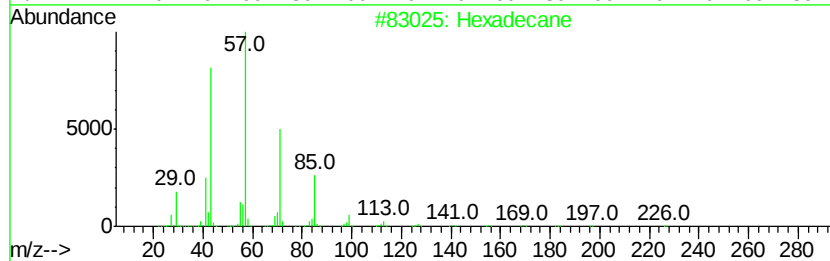
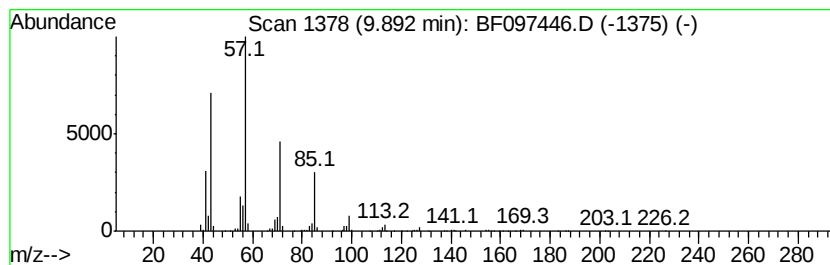
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Pentadecane Concentration Rank 18

| R.T.    | EstConc                             | Area          | Relative to ISTD | R.T.      |
|---------|-------------------------------------|---------------|------------------|-----------|
| 9.89    | 9.92 ng                             | 340477        | Acenaphthene-d10 | 9.43      |
| Hit# of | 5                                   | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | Hexadecane                          | 226 C16H34    | 000544-76-3      | 97        |
| 2       | Pentadecane                         | 212 C15H32    | 000629-62-9      | 90        |
| 3       | Eicosane                            | 282 C20H42    | 000112-95-8      | 83        |
| 4       | Octane, 3,5-dimethyl-               | 142 C10H22    | 015869-93-9      | 81        |
| 5       | Sulfurous acid, 2-ethylhexyl hex... | 278 C14H30O3S | 1000309-20-2     | 80        |



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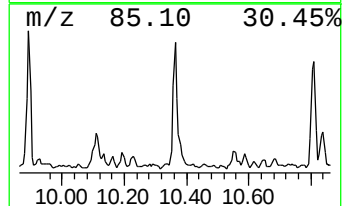
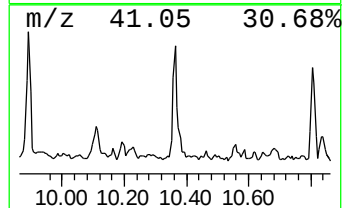
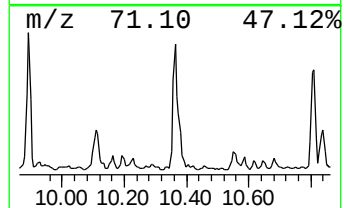
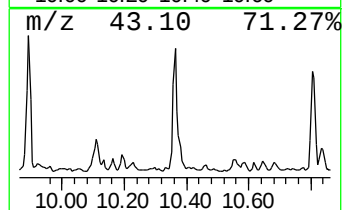
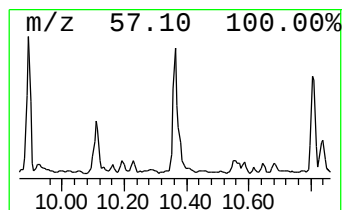
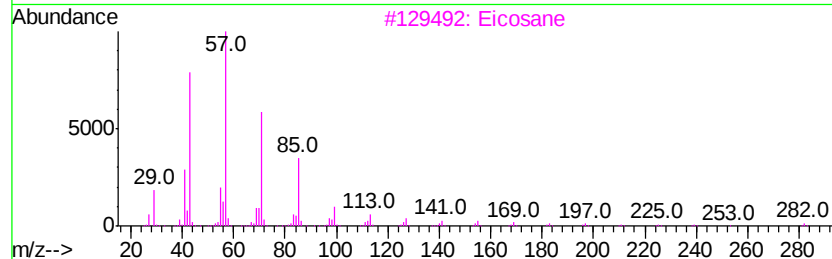
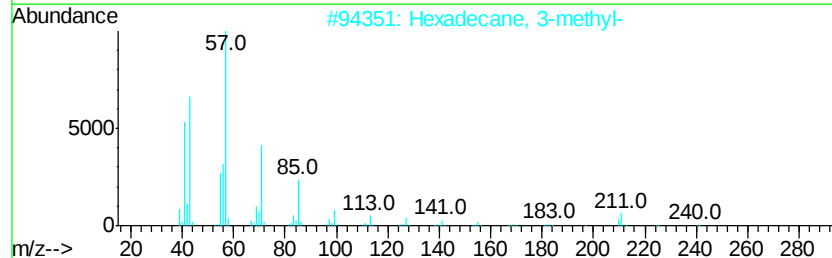
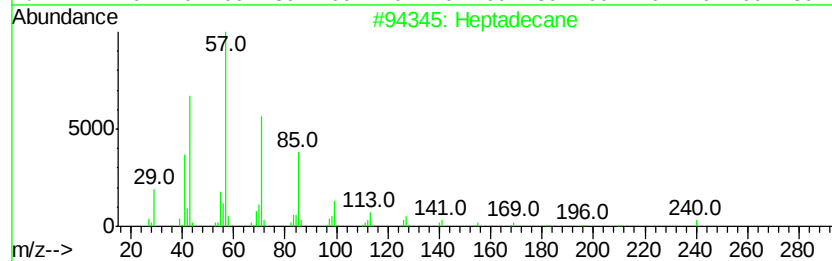
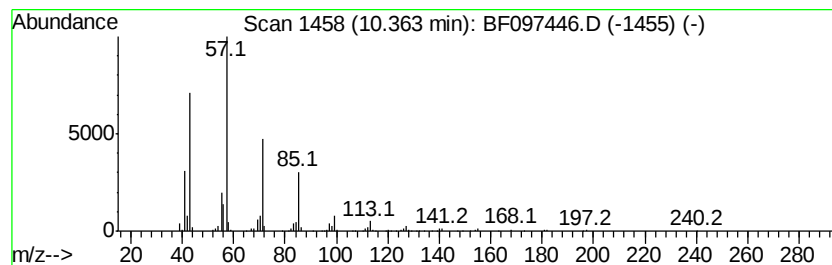
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Heptadecane Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.36 | 17.39 ng | 456157 | Phenanthrene-d10 | 10.91 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Heptadecane           |              | 240 | C17H36  | 000629-78-7 | 95   |
| 2       | Hexadecane, 3-methyl- |              | 240 | C17H36  | 006418-43-5 | 90   |
| 3       | Eicosane              |              | 282 | C20H42  | 000112-95-8 | 90   |
| 4       | Hexadecane            |              | 226 | C16H34  | 000544-76-3 | 87   |
| 5       | Pentadecane           |              | 212 | C15H32  | 000629-62-9 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

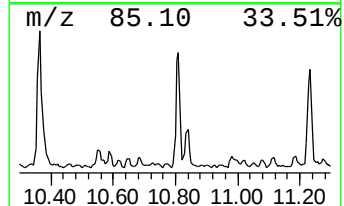
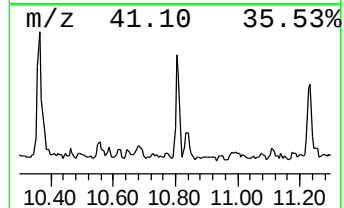
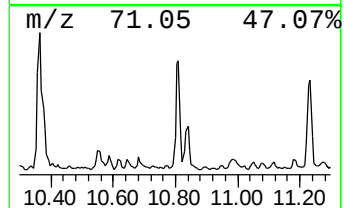
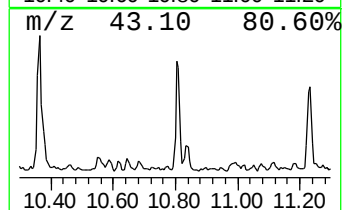
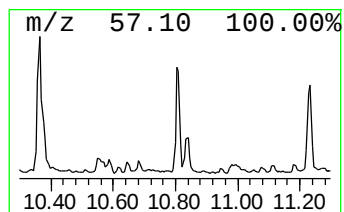
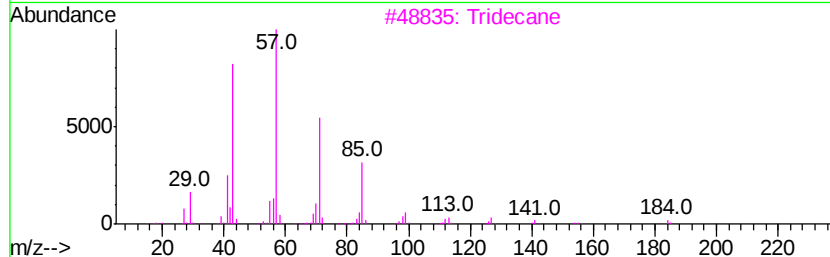
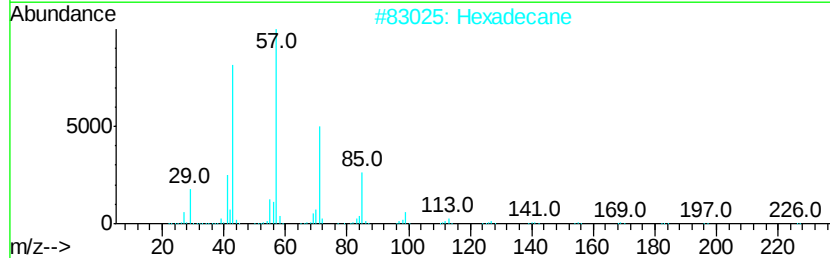
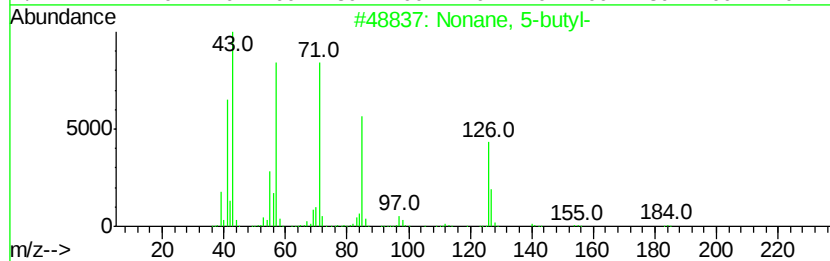
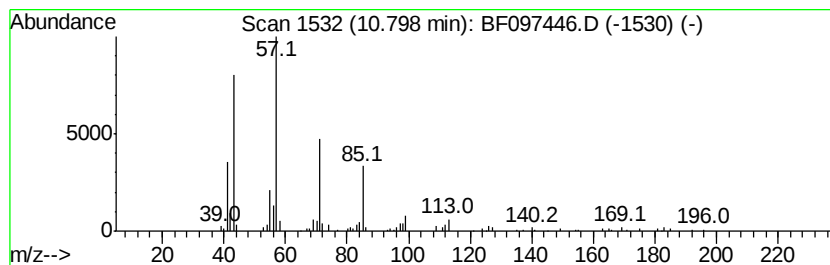
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Nonane, 5-butyl- Concentration Rank 12

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.80 | 11.77 ng | 308809 | Phenanthrene-d10 | 10.91 |

| Hit# of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|---------|---|------------------|-----|---------|-------------|------|
| 1       |   | Nonane, 5-butyl- | 184 | C13H28  | 017312-63-9 | 93   |
| 2       |   | Hexadecane       | 226 | C16H34  | 000544-76-3 | 91   |
| 3       |   | Tridecane        | 184 | C13H28  | 000629-50-5 | 90   |
| 4       |   | Eicosane         | 282 | C20H42  | 000112-95-8 | 90   |
| 5       |   | Tetradecane      | 198 | C14H30  | 000629-59-4 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

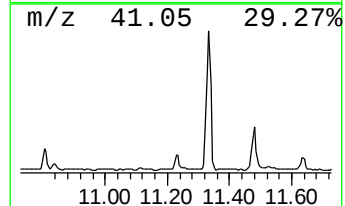
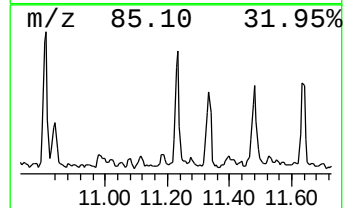
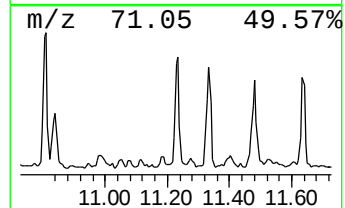
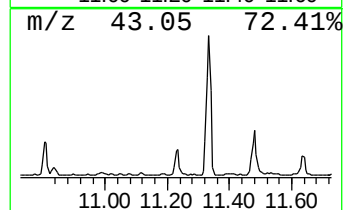
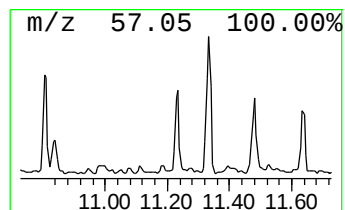
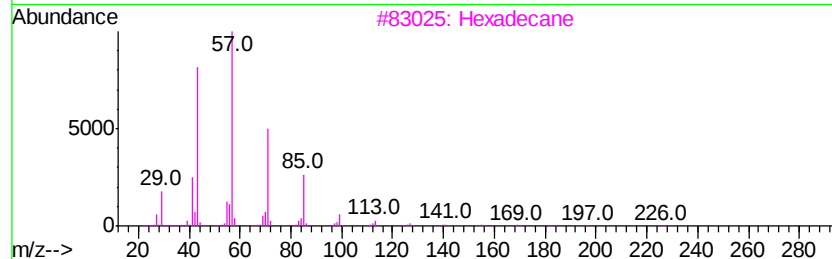
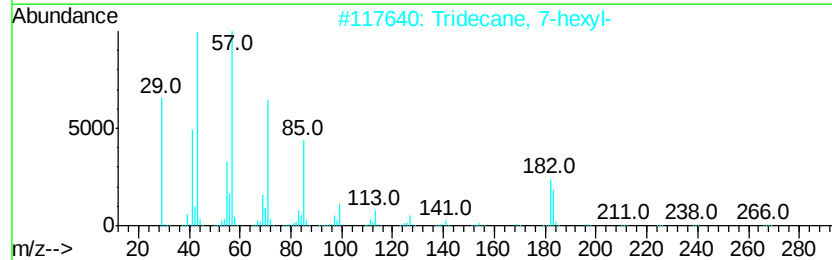
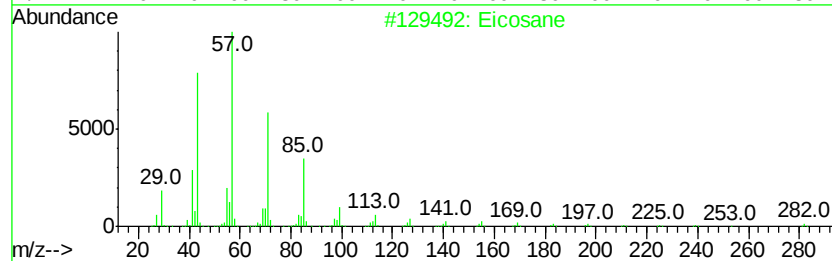
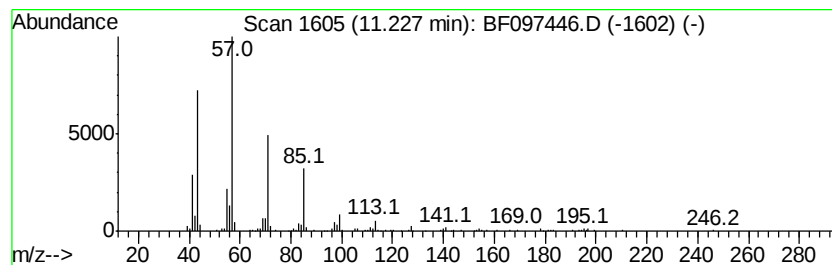
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Eicosane Concentration Rank 16

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.23 | 10.01 ng | 262641 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane            | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    | Tridecane, 7-hexyl- | 268 | C19H40  | 007225-66-3 | 90   |
| 3    |    | Hexadecane          | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    | Tetracosane         | 338 | C24H50  | 000646-31-1 | 86   |
| 5    |    | Heneicosane         | 296 | C21H44  | 000629-94-7 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

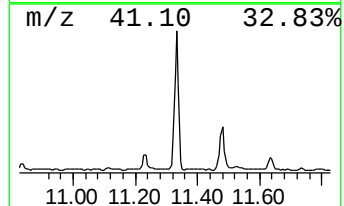
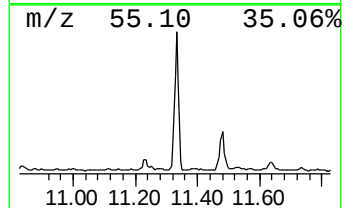
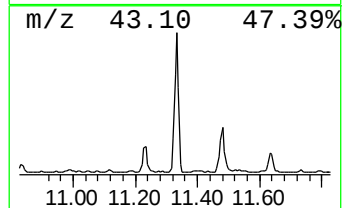
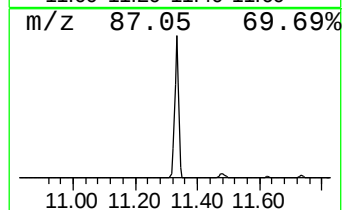
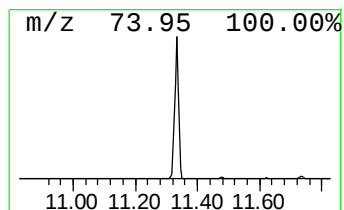
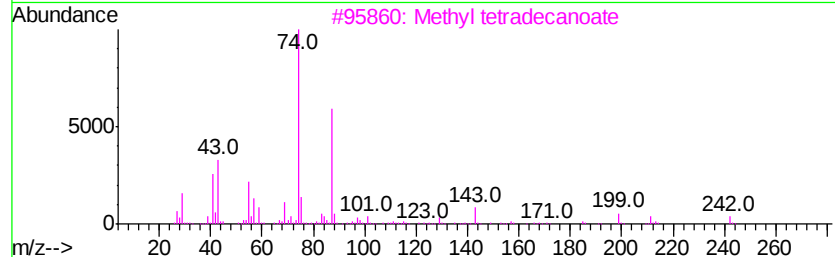
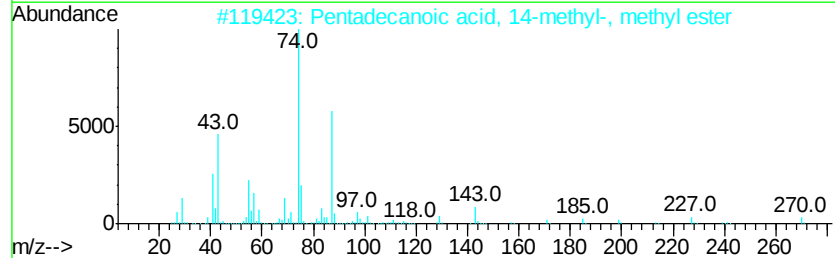
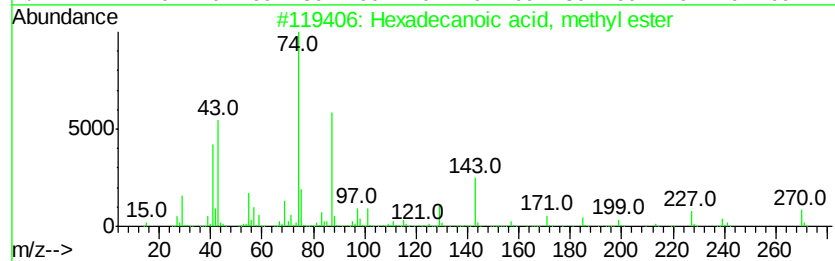
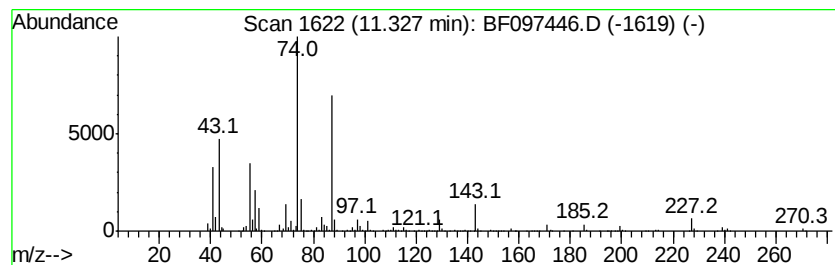
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 4

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 11.33 | 100.13 ng | 2626730 | Phenanthrene-d10 | 10.91 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------------|-----|----------|-------------|------|
| 1       |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 94   |
| 2       |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 91   |
| 3       |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 90   |
| 4       |   | Dodecanoic acid, methyl ester       | 214 | C13H26O2 | 000111-82-0 | 90   |
| 5       |   | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

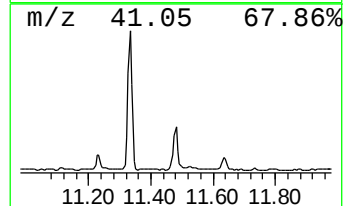
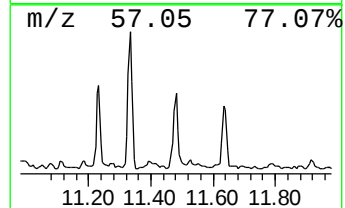
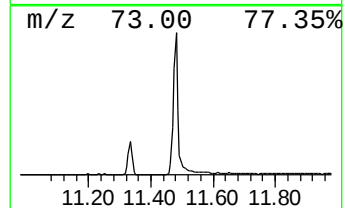
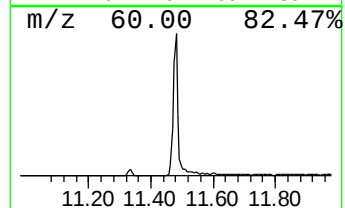
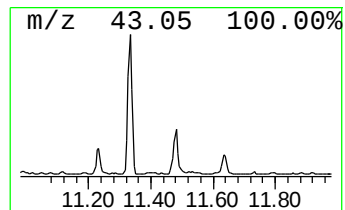
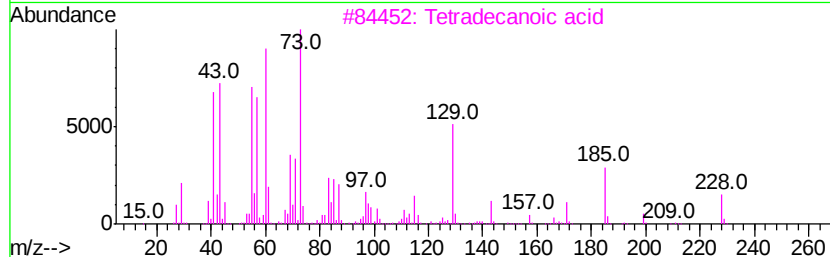
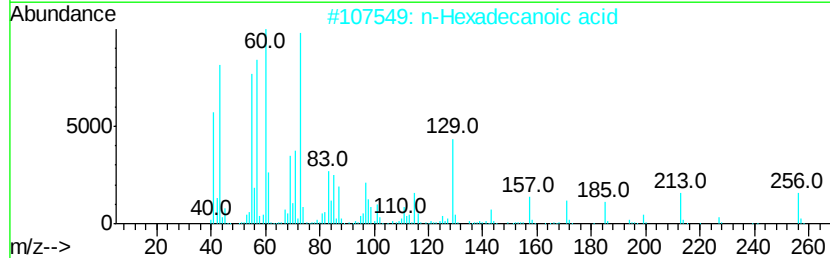
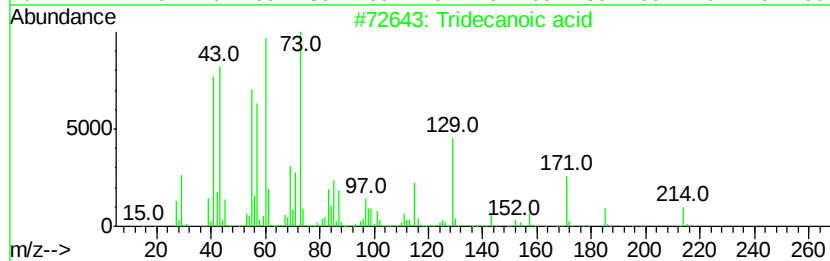
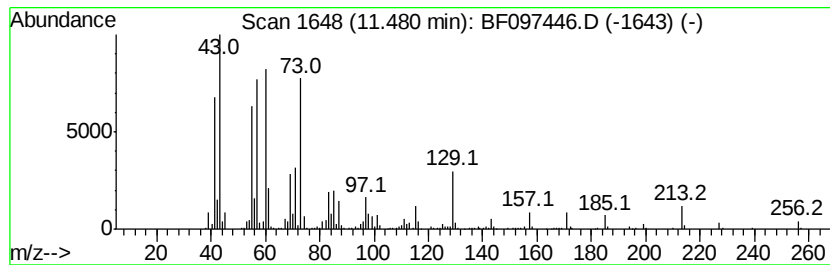
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Tridecanoic acid Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.48 | 30.49 ng | 799735 | Phenanthrene-d10 | 10.91 |

| Hit# of 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|-----------|---------------------|-----|----------|-------------|------|
| 1         | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 95   |
| 2         | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 94   |
| 3         | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 86   |
| 4         | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 74   |
| 5         | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 62   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

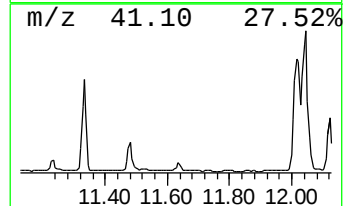
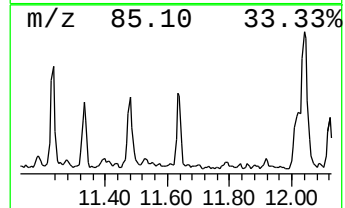
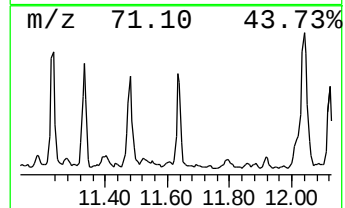
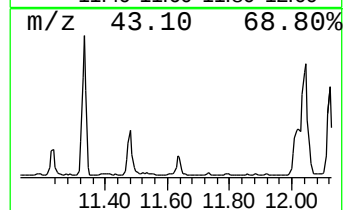
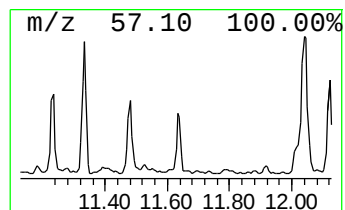
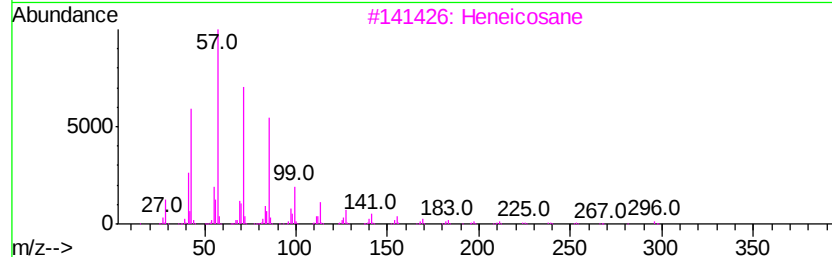
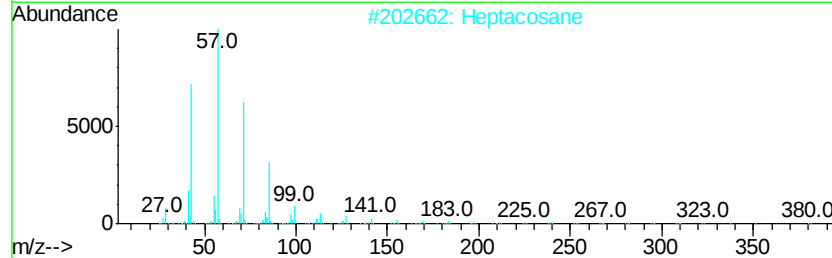
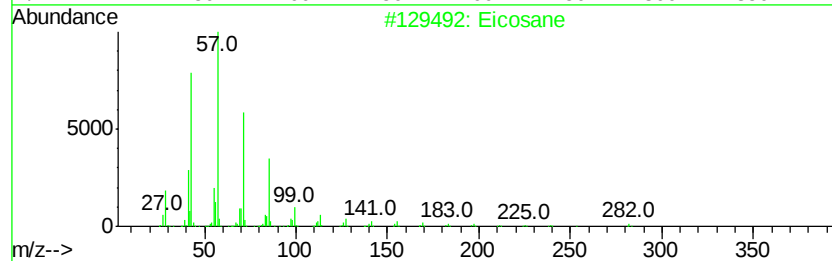
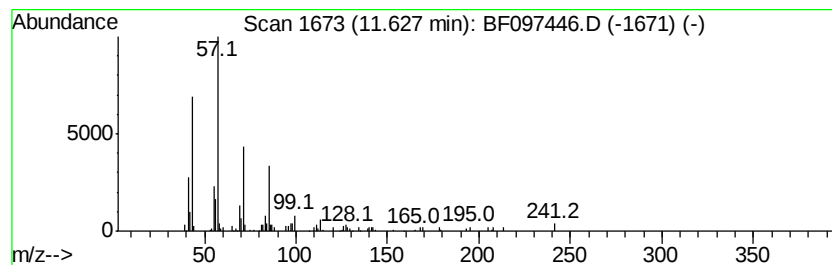
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Heneicosane Concentration Rank 14

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.63 | 10.94 ng | 286960 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    | Heptacosane  | 380 | C27H56  | 000593-49-7 | 90   |
| 3    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 90   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 5    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

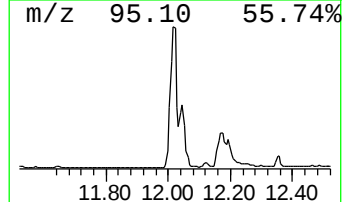
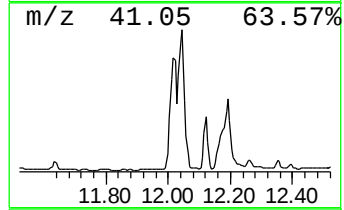
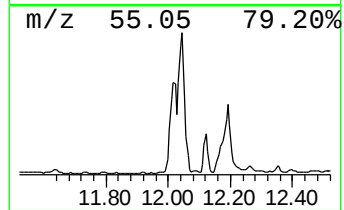
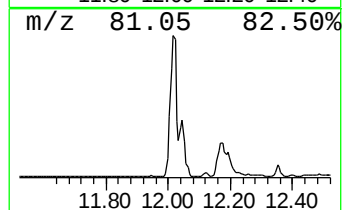
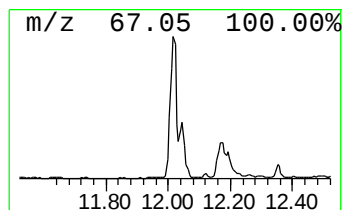
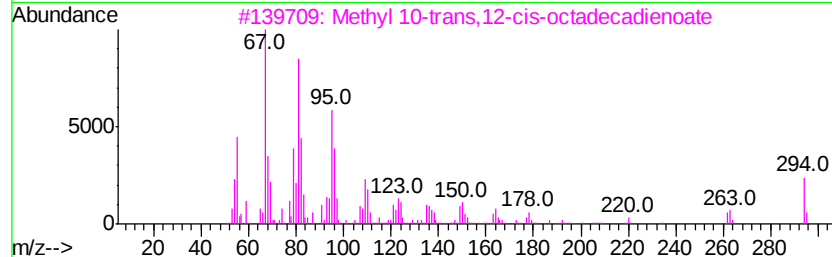
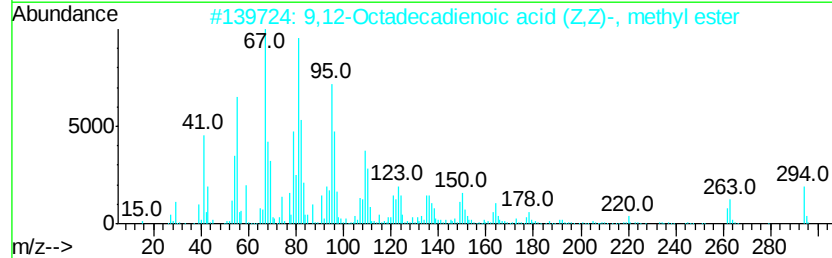
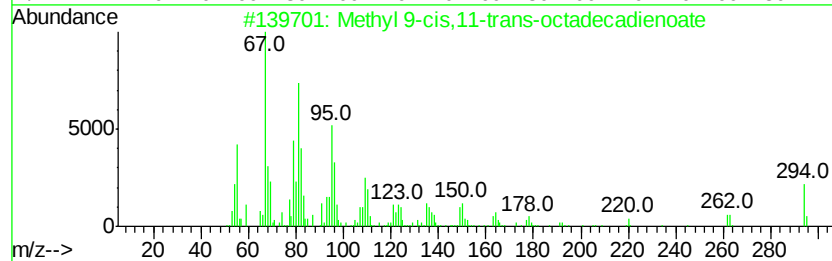
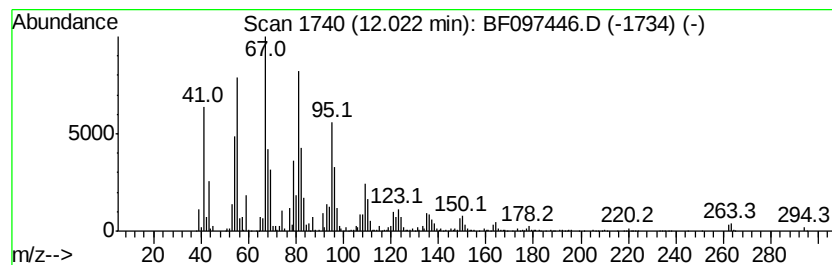
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Methyl 9-cis,11-trans-octad... Concentration Rank 2

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.02 | 191.05 ng | 5011780 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 1000336-44-0 | 99   |
| 2    |    | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0  | 99   |
| 3    |    | Methyl 10-trans,12-cis-octadecad... | 294 | C19H34O2 | 1000336-44-2 | 99   |
| 4    |    | 10,13-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-62-2  | 99   |
| 5    |    | 9,15-Octadecadienoic acid, methy... | 294 | C19H34O2 | 017309-05-6  | 99   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

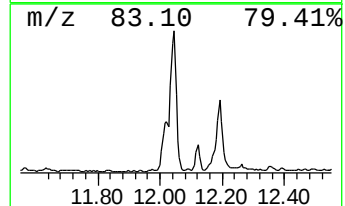
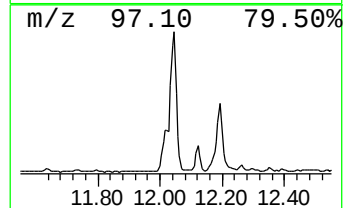
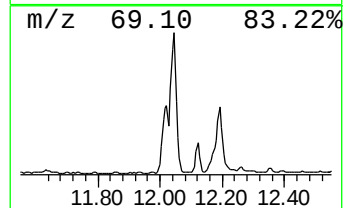
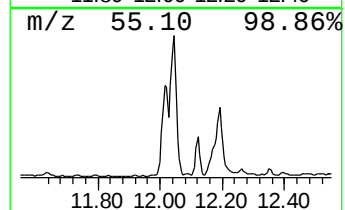
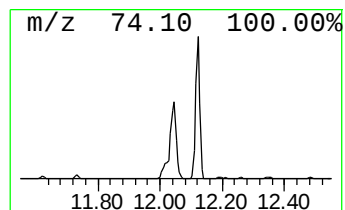
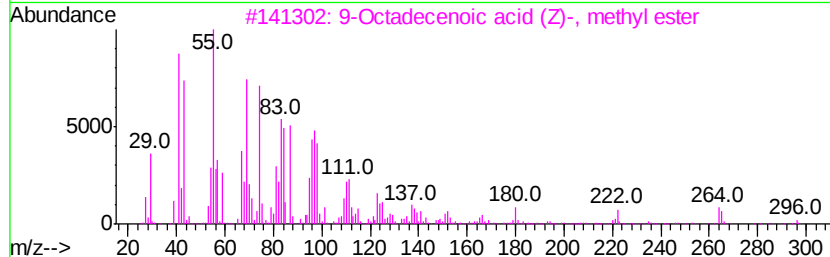
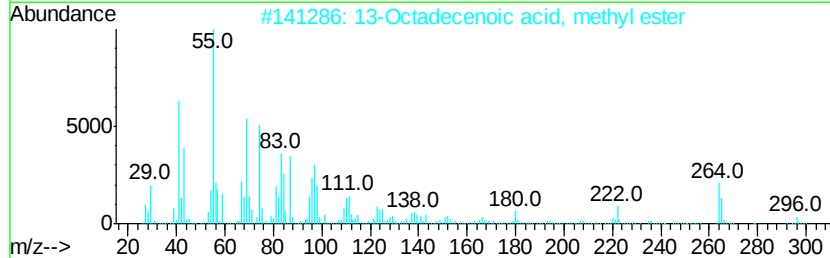
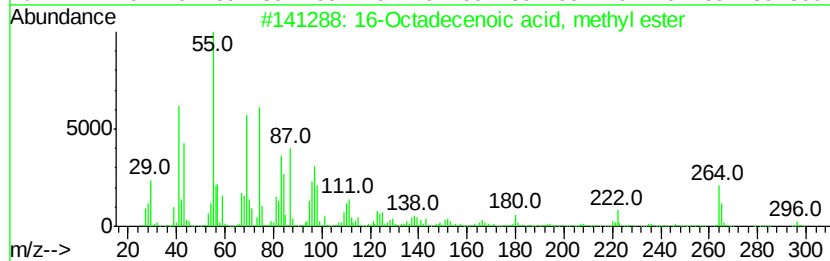
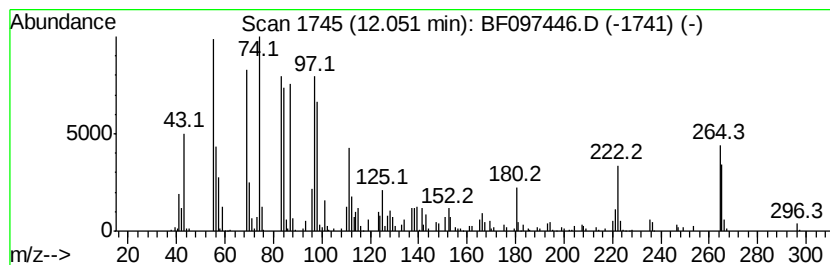
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 16-Octadecenoic acid, methyl... Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.05 | 200.86 ng | 5269070 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 16-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 056554-49-5 | 96   |
| 2    |    | 13-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 056554-47-3 | 95   |
| 3    |    | 9-Octadecenoic acid (Z)-, methyl...  | 296 | C19H36O2 | 000112-62-9 | 94   |
| 4    |    | 15-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 004764-72-1 | 93   |
| 5    |    | 6-Octadecenoic acid, methyl ester... | 296 | C19H36O2 | 002777-58-4 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

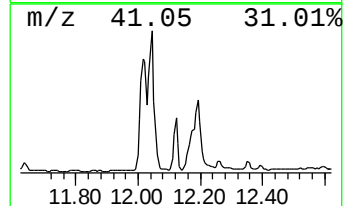
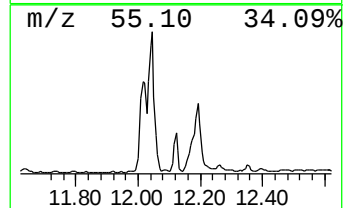
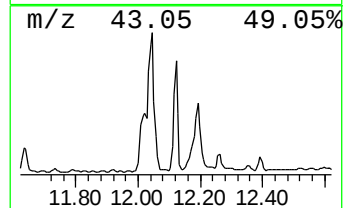
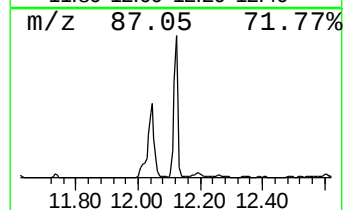
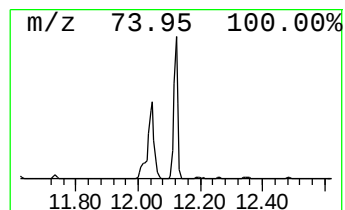
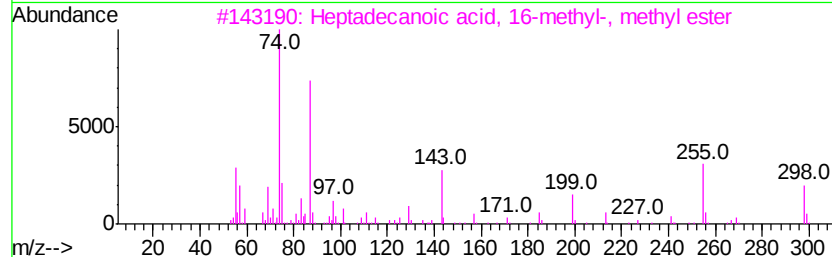
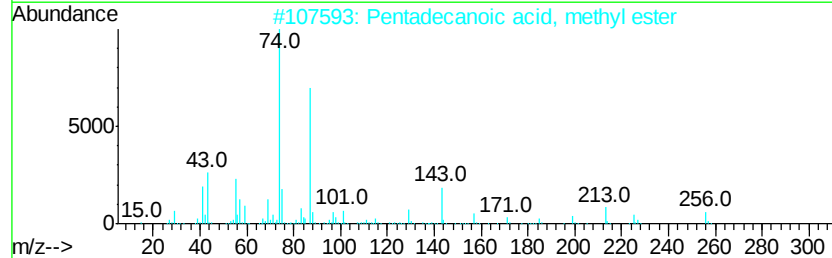
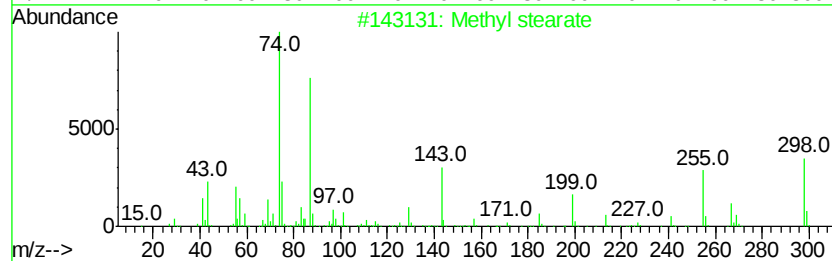
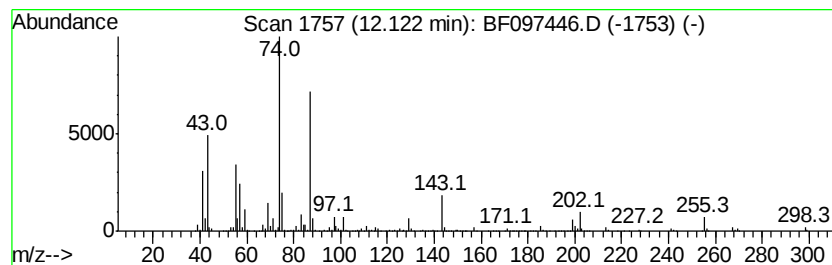
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Methyl stearate Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.12 | 62.63 ng | 1642900 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 97   |
| 2    |    | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 93   |
| 3    |    | Heptadecanoic acid, 16-methyl-, ... | 298 | C19H38O2 | 005129-61-3 | 90   |
| 4    |    | Heptadecanoic acid, methyl ester    | 284 | C18H36O2 | 001731-92-6 | 87   |
| 5    |    | Heptadecanoic acid, 15-methyl-, ... | 298 | C19H38O2 | 054833-55-5 | 87   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

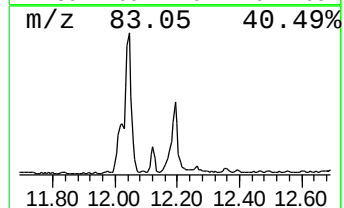
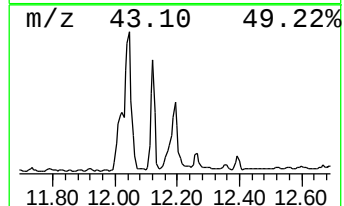
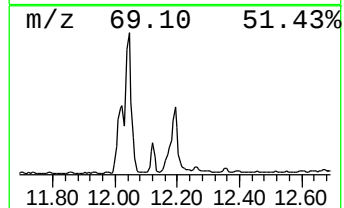
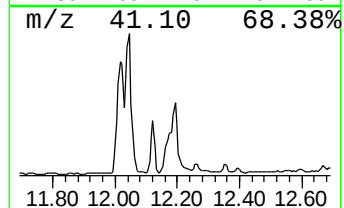
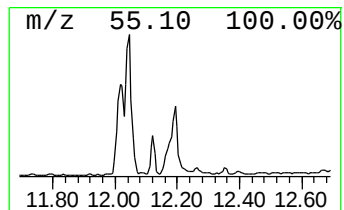
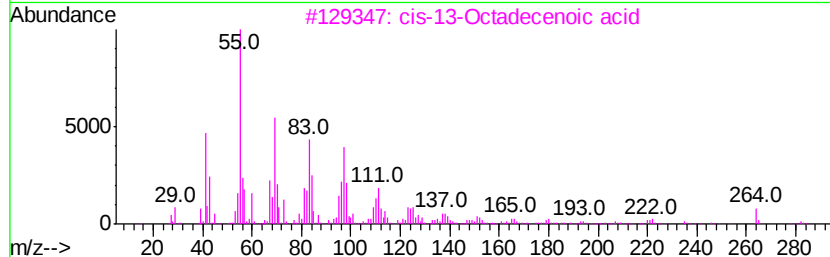
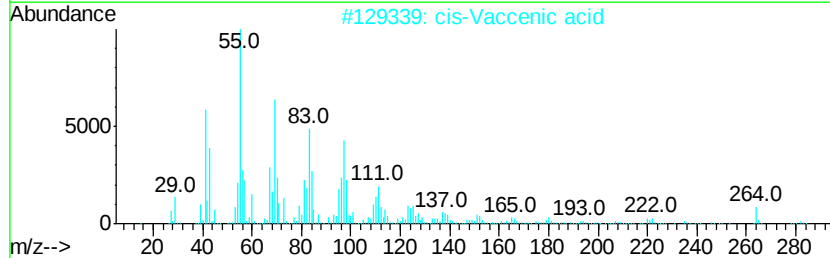
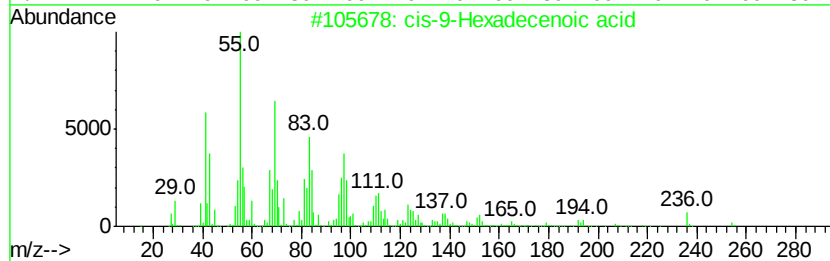
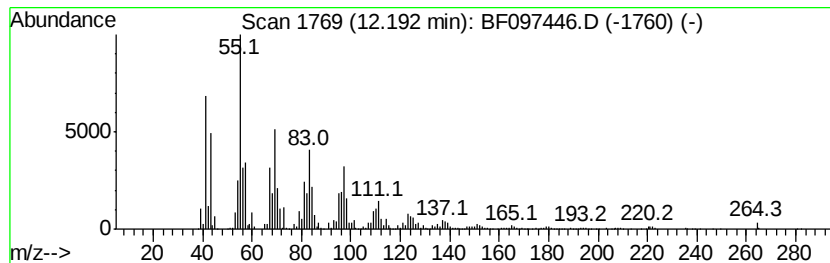
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 cis-9-Hexadecenoic acid Concentration Rank 3

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.19 | 140.54 ng | 3686780 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | cis-9-Hexadecenoic acid             | 254 | C16H30O2   | 1000333-19-5 | 91   |
| 2    |    | cis-Vaccenic acid                   | 282 | C18H34O2   | 000506-17-2  | 91   |
| 3    |    | cis-13-Octadecenoic acid            | 282 | C18H34O2   | 013126-39-1  | 91   |
| 4    |    | Palmitoleic acid                    | 254 | C16H30O2   | 000373-49-9  | 87   |
| 5    |    | 9-octadecanoic acid, 2,2,3,3,4,4... | 464 | C22H35F7O2 | 1000376-61-4 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

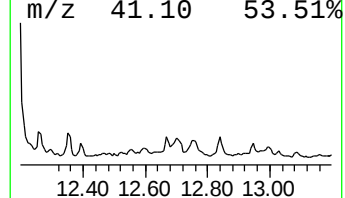
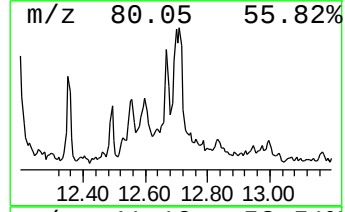
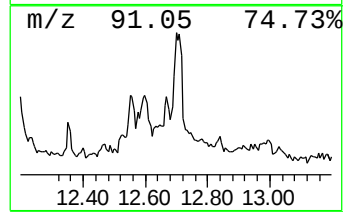
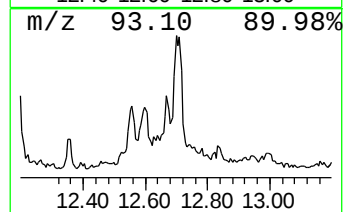
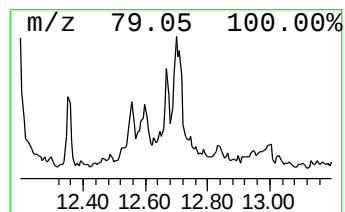
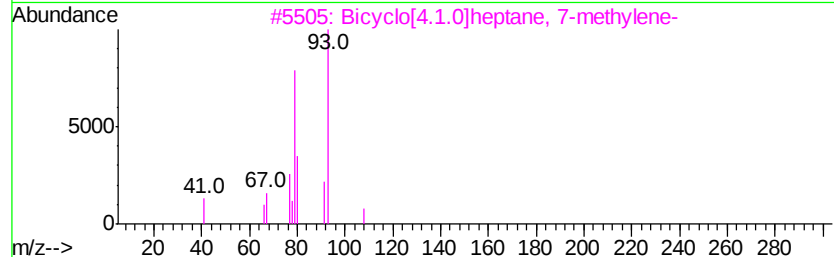
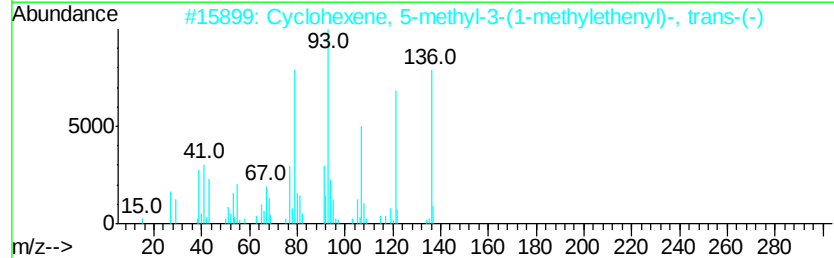
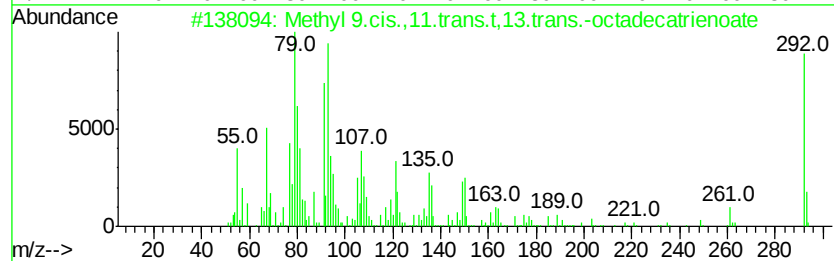
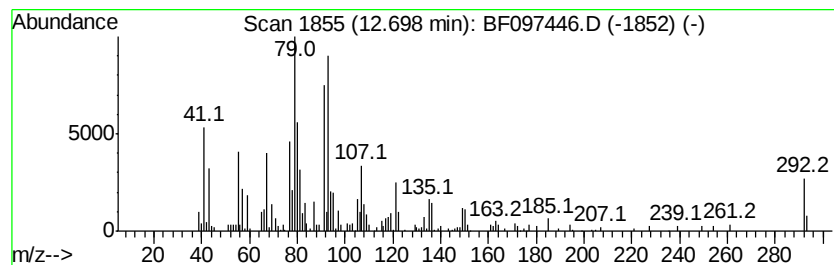
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Methyl 9.cis.,11.trans.t,13... Concentration Rank 13

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.70 | 11.02 ng | 325502 | Chrysene-d12     | 13.53 |

| Hit# | of | Tentative ID                          | MW  | MolForm  | CAS#         | Qual |
|------|----|---------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Methyl 9.cis.,11.trans.t,13.trans...  | 292 | C19H32O2 | 1000336-42-6 | 97   |
| 2    |    | Cyclohexene, 5-methyl-3-(1-methyl-... | 136 | C10H16   | 056816-08-1  | 72   |
| 3    |    | Bicyclo[4.1.0]heptane, 7-methylene-   | 108 | C8H12    | 054211-14-2  | 68   |
| 4    |    | 6-Butyl-1,4-cycloheptadiene           | 150 | C11H18   | 022735-58-6  | 64   |
| 5    |    | Bicyclo[3.3.1]non-2-en-9-one          | 136 | C9H12O   | 004844-11-5  | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

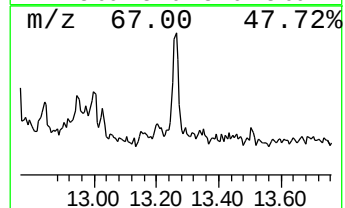
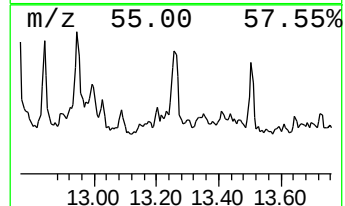
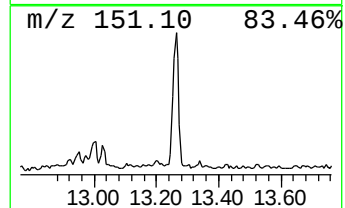
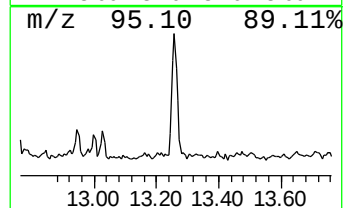
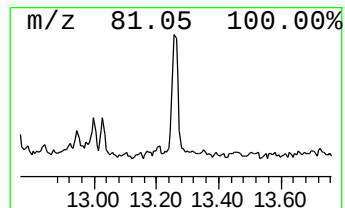
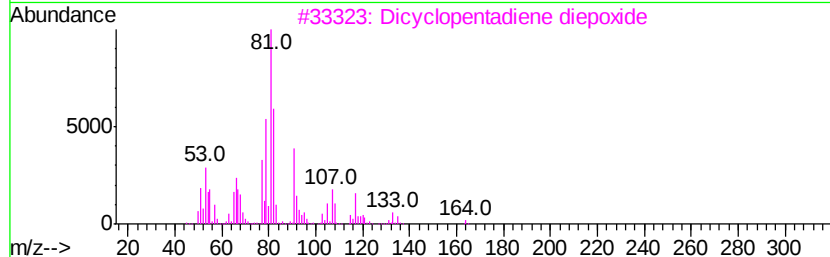
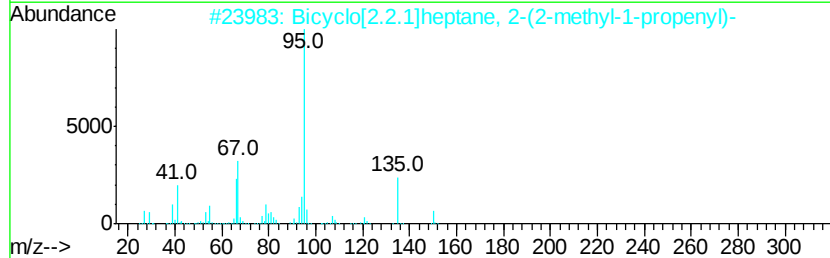
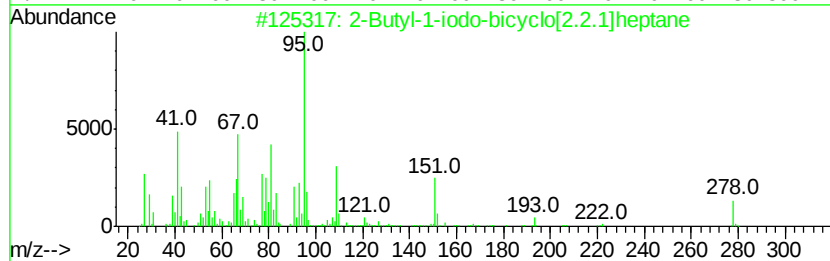
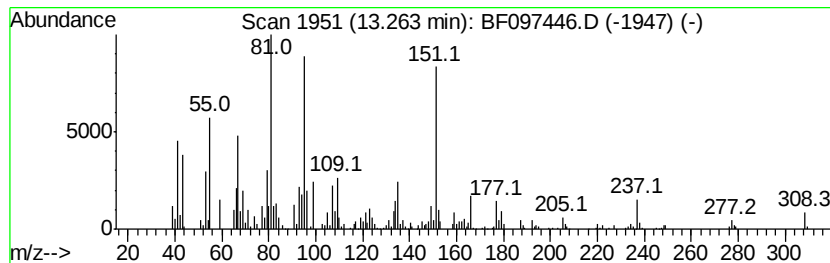
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown13.26 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.26 | 9.96 ng | 294202 | Chrysene-d12     | 13.53 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Butyl-1-iodo-bicyclo[2.2.1]hep... | 278 | C11H19I  | 1000190-59-7 | 45   |
| 2    |    |   | Bicyclo[2.2.1]heptane, 2-(2-meth... | 150 | C11H18   | 061142-27-6  | 35   |
| 3    |    |   | Dicyclopentadiene diepoxide         | 164 | C10H12O2 | 000081-21-0  | 30   |
| 4    |    |   | 2(5H)-Furanone, 4-methyl-5,5-bis... | 206 | C13H18O2 | 099202-98-9  | 30   |
| 5    |    |   | Cis-3-ethyl-endo-tricyclo[5.2.1.... | 164 | C12H20   | 1000215-29-1 | 15   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

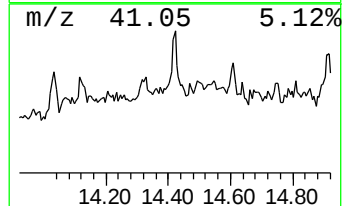
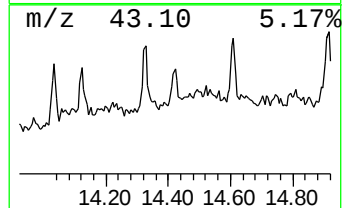
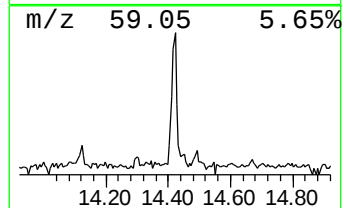
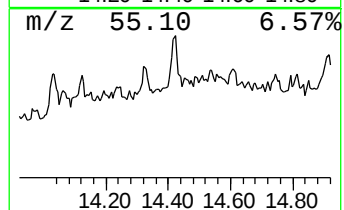
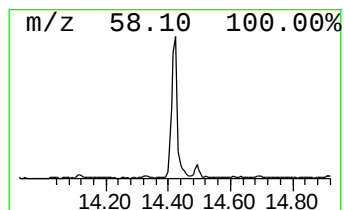
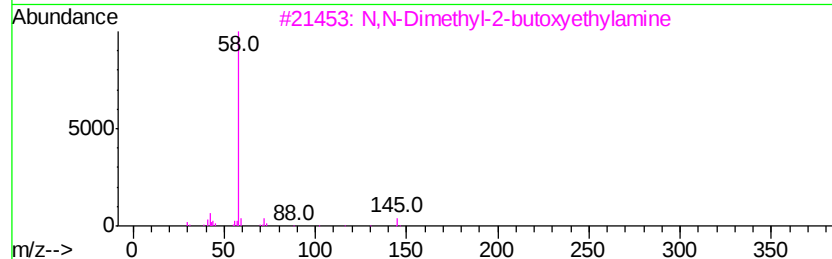
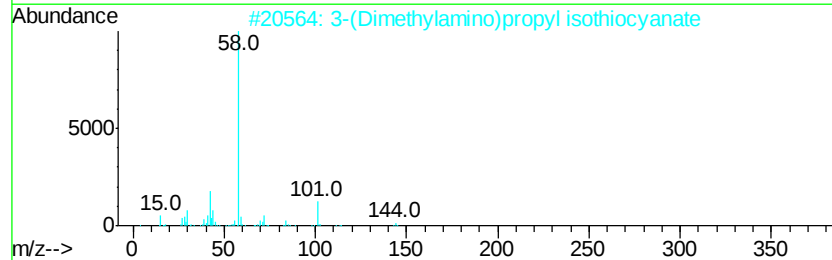
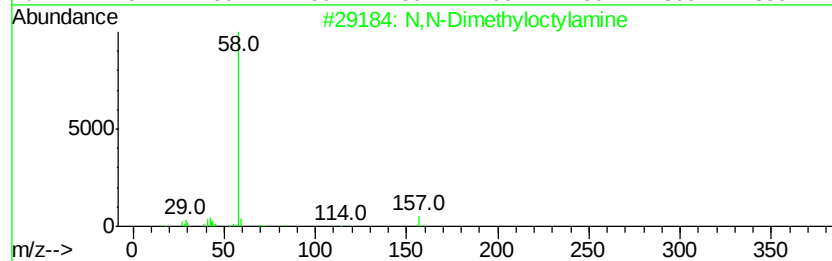
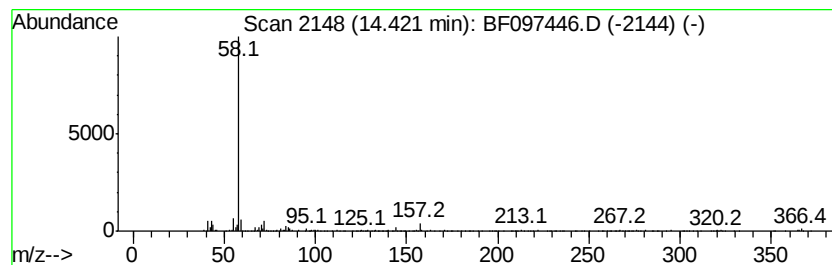
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 N,N-Dimethyloctylamine Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.42 | 20.16 ng | 290477 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | N,N-Dimethyloctylamine              | 157 | C10H23N   | 007378-99-6 | 64   |
| 2    |    | 3-(Dimethylamino)propyl isothioc... | 144 | C6H12N2S  | 027421-70-1 | 64   |
| 3    |    | N,N-Dimethyl-2-butoxyethylamine     | 145 | C8H19NO   | 071126-58-4 | 64   |
| 4    |    | n-Octyl trimethyl ammonium bromide  | 251 | C11H26BrN | 002083-68-3 | 64   |
| 5    |    | Acetone, 1-[4-(dimethylaminoetho... | 221 | C13H19NO2 | 086608-11-9 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

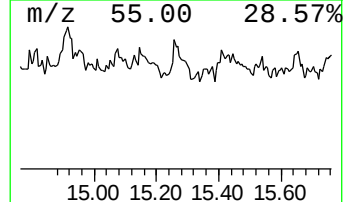
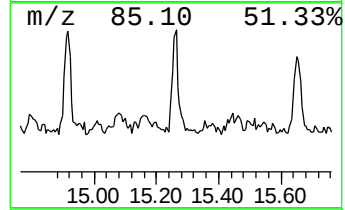
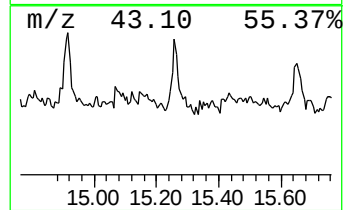
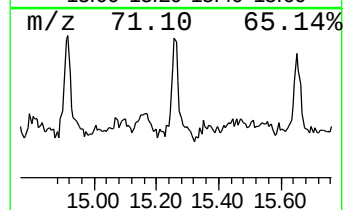
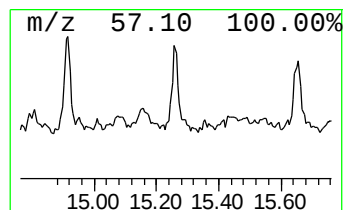
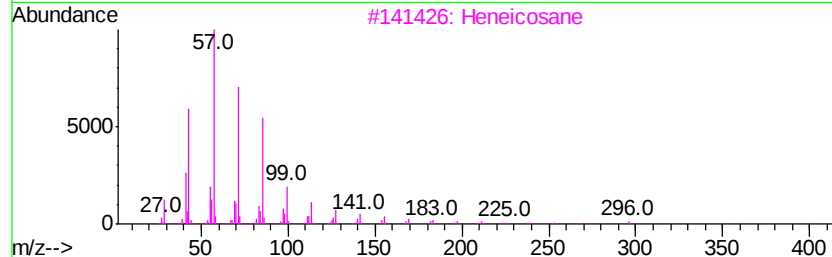
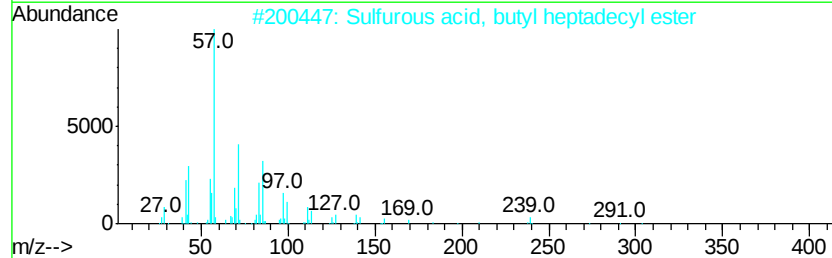
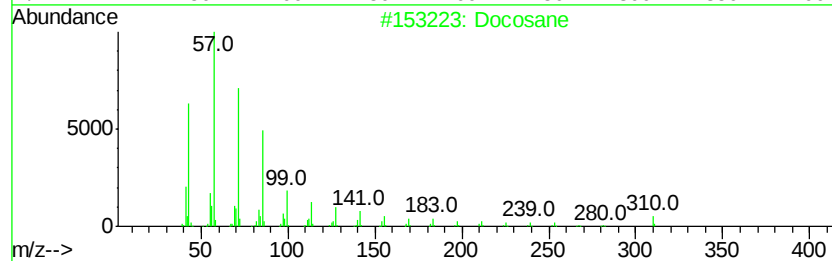
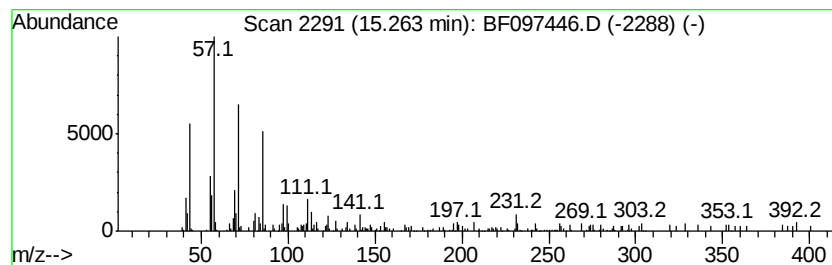
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Docosane Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.26 | 16.86 ng | 242946 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Docosane                            | 310 | C22H46    | 000629-97-0  | 94   |
| 2    |    | Sulfurous acid, butyl heptadecyl... | 376 | C21H44O3S | 1000309-18-4 | 86   |
| 3    |    | Heneicosane                         | 296 | C21H44    | 000629-94-7  | 76   |
| 4    |    | Heptacosane                         | 380 | C27H56    | 000593-49-7  | 74   |
| 5    |    | Eicosane, 2-methyl-                 | 296 | C21H44    | 001560-84-5  | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

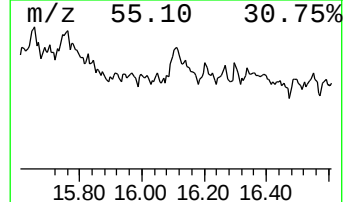
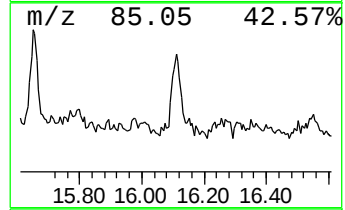
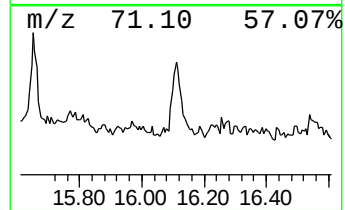
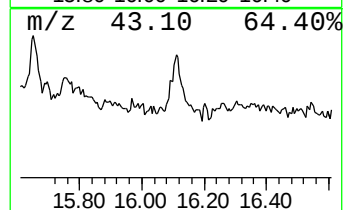
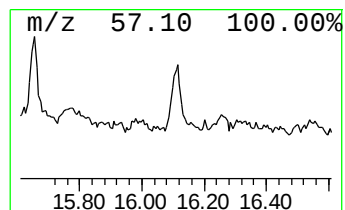
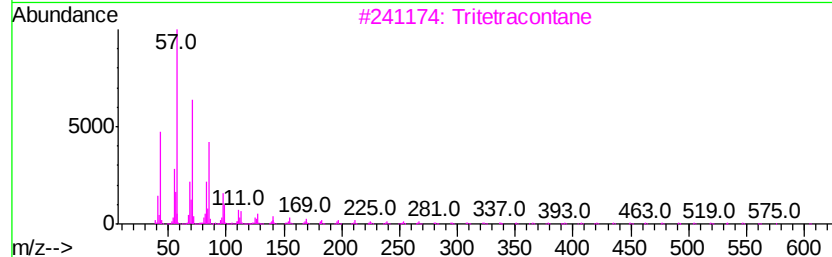
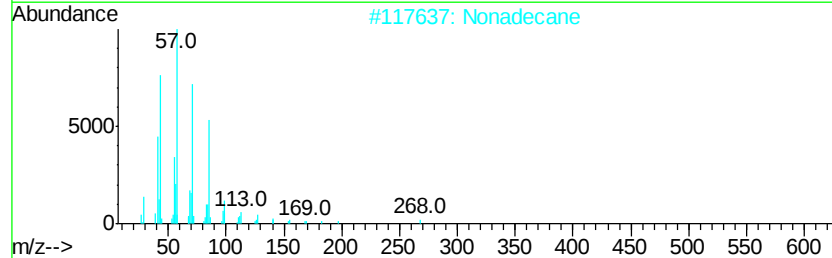
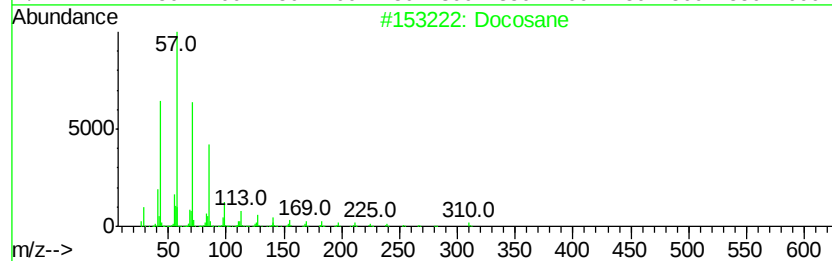
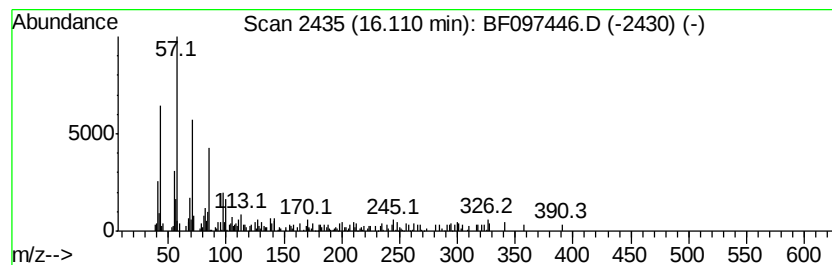
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Nonadecane Concentration Rank 15

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.11 | 10.03 ng | 144493 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Docosane                            | 310 | C22H46    | 000629-97-0  | 90   |
| 2    |    | Nonadecane                          | 268 | C19H40    | 000629-92-5  | 78   |
| 3    |    | Tritetracontane                     | 605 | C43H88    | 007098-21-7  | 64   |
| 4    |    | Sulfurous acid, butyl octadecyl ... | 390 | C22H46O3S | 1000309-18-5 | 64   |
| 5    |    | Tetradecane                         | 198 | C14H30    | 000629-59-4  | 60   |



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Data File : BF097446.D  
Acq On : 7 Aug 2017 23:13  
Operator : SJ/JU  
Sample : I4639-04 5X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

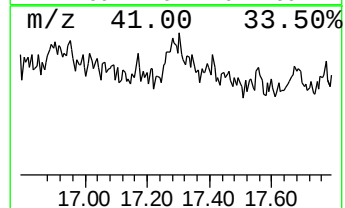
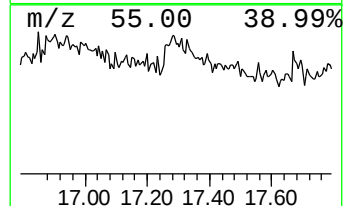
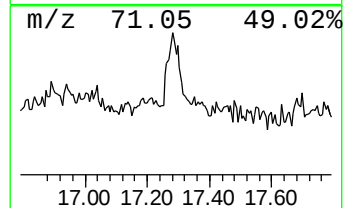
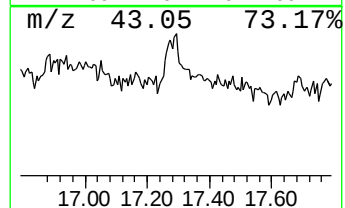
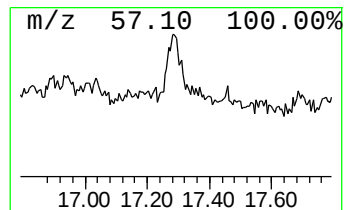
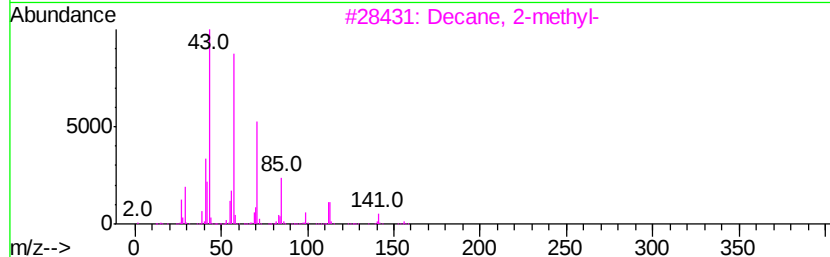
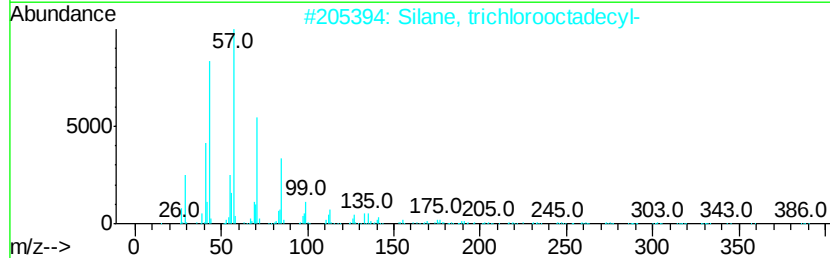
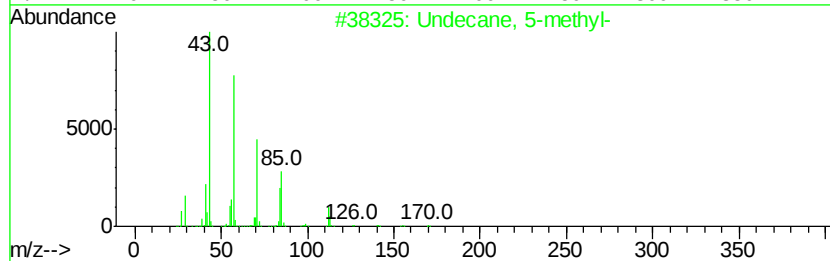
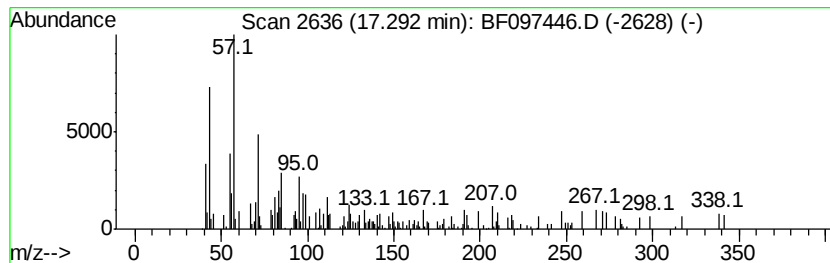
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 unknown17.29 Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.29 | 7.11 ng | 102447 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                | MW  | MolForm     | CAS#        | Qual |
|------|----|-----------------------------|-----|-------------|-------------|------|
| 1    | 5  | Undecane, 5-methyl-         | 170 | C12H26      | 001632-70-8 | 49   |
| 2    |    | Silane, trichlorooctadecyl- | 386 | C18H37Cl3Si | 000112-04-9 | 49   |
| 3    |    | Decane, 2-methyl-           | 156 | C11H24      | 006975-98-0 | 49   |
| 4    |    | Ethanol, 2-(dodecyloxy)-    | 230 | C14H30O2    | 004536-30-5 | 49   |
| 5    |    | Eicosane                    | 282 | C20H42      | 000112-95-8 | 49   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
 Data File : BF097446.D  
 Acq On : 7 Aug 2017 23:13  
 Operator : SJ/JU  
 Sample : I4639-04 5X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SA-06-080317-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.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 |
| unknown6.19          | 6.19  | 19.3    | ng    | 512195   | 1                     | 6.41  | 532160 | 20.0 |
| Hexadecane           | 8.87  | 7.1     | ng    | 244871   | 3                     | 9.43  | 686774 | 20.0 |
| Pentadecane          | 9.89  | 9.9     | ng    | 340477   | 3                     | 9.43  | 686774 | 20.0 |
| Heptadecane          | 10.36 | 17.4    | ng    | 456157   | 4                     | 10.91 | 524655 | 20.0 |
| Nonane, 5-butyl-     | 10.80 | 11.8    | ng    | 308809   | 4                     | 10.91 | 524655 | 20.0 |
| Eicosane             | 11.23 | 10.0    | ng    | 262641   | 4                     | 10.91 | 524655 | 20.0 |
| Hexadecanoic acid... | 11.33 | 100.1   | ng    | 2626730  | 4                     | 10.91 | 524655 | 20.0 |
| Tridecanoic acid     | 11.48 | 30.5    | ng    | 799735   | 4                     | 10.91 | 524655 | 20.0 |
| Heneicosane          | 11.63 | 10.9    | ng    | 286960   | 4                     | 10.91 | 524655 | 20.0 |
| Methyl 9-cis,11-t... | 12.02 | 191.1   | ng    | 5011780  | 4                     | 10.91 | 524655 | 20.0 |
| 16-Octadecenoic a... | 12.05 | 200.9   | ng    | 5269070  | 4                     | 10.91 | 524655 | 20.0 |
| Methyl stearate      | 12.12 | 62.6    | ng    | 1642900  | 4                     | 10.91 | 524655 | 20.0 |
| cis-9-Hexadeceno...  | 12.19 | 140.5   | ng    | 3686780  | 4                     | 10.91 | 524655 | 20.0 |
| Methyl 9.cis.,11...  | 12.70 | 11.0    | ng    | 325502   | 5                     | 13.53 | 590871 | 20.0 |
| unknown13.26         | 13.26 | 10.0    | ng    | 294202   | 5                     | 13.53 | 590871 | 20.0 |
| N,N-Dimethyloctyl... | 14.42 | 20.2    | ng    | 290477   | 6                     | 14.87 | 288171 | 20.0 |
| Docosane             | 15.26 | 16.9    | ng    | 242946   | 6                     | 14.87 | 288171 | 20.0 |
| Nonadecane           | 16.11 | 10.0    | ng    | 144493   | 6                     | 14.87 | 288171 | 20.0 |
| unknown17.29         | 17.29 | 7.1     | ng    | 102447   | 6                     | 14.87 | 288171 | 20.0 |