

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\  
 Data File : BF096914.D  
 Acq On : 20 Jul 2017 19:12  
 Operator : SJ/JU  
 Sample : I4308-01  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NB-08-071917

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-BF071317.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.740       | 464           | 468         | 480          | rBV      | 227170         | 341033        | 14.43%          | 1.384%        |
| 2         | 5.181       | 538           | 543         | 546          | rBV      | 2205912        | 2269511       | 96.01%          | 9.208%        |
| 3         | 6.228       | 715           | 721         | 736          | rBV      | 2211404        | 2347016       | 99.28%          | 9.523%        |
| 4         | 6.340       | 736           | 740         | 743          | rBV      | 2321522        | 2280159       | 96.46%          | 9.251%        |
| 5         | 6.569       | 775           | 779         | 782          | rBV      | 721950         | 614281        | 25.99%          | 2.492%        |
| 6         | 6.722       | 801           | 805         | 809          | rBV      | 1770287        | 1752153       | 74.12%          | 7.109%        |
| 7         | 7.134       | 870           | 875         | 878          | rBV      | 1466104        | 1475106       | 62.40%          | 5.985%        |
| 8         | 7.263       | 893           | 897         | 900          | rVB      | 26668          | 26282         | 1.11%           | 0.107%        |
| 9         | 7.851       | 992           | 997         | 1000         | rBV      | 1141740        | 962884        | 40.73%          | 3.907%        |
| 10        | 8.928       | 1175          | 1180        | 1184         | rBV      | 2270757        | 2363930       | 100.00%         | 9.591%        |
| 11        | 9.028       | 1193          | 1197        | 1202         | rVB      | 26440          | 28128         | 1.19%           | 0.114%        |
| 12        | 9.328       | 1243          | 1248        | 1251         | rBV      | 406680         | 383759        | 16.23%          | 1.557%        |
| 13        | 9.551       | 1283          | 1286        | 1289         | rVB      | 48449          | 38470         | 1.63%           | 0.156%        |
| 14        | 9.598       | 1289          | 1294        | 1298         | rBV      | 952737         | 862248        | 36.48%          | 3.498%        |
| 15        | 10.022      | 1360          | 1366        | 1368         | rBV3     | 21147          | 28825         | 1.22%           | 0.117%        |
| 16        | 10.051      | 1368          | 1371        | 1374         | rVB      | 86944          | 73811         | 3.12%           | 0.299%        |
| 17        | 10.269      | 1403          | 1408        | 1411         | rBV2     | 43460          | 60559         | 2.56%           | 0.246%        |
| 18        | 10.386      | 1423          | 1428        | 1431         | rBV      | 1407813        | 1348868       | 57.06%          | 5.473%        |
| 19        | 10.522      | 1448          | 1451        | 1458         | rVB2     | 150933         | 169925        | 7.19%           | 0.689%        |
| 20        | 10.580      | 1458          | 1461        | 1463         | rBV      | 30102          | 28869         | 1.22%           | 0.117%        |
| 21        | 10.610      | 1463          | 1466        | 1469         | rBV2     | 38745          | 41389         | 1.75%           | 0.168%        |
| 22        | 10.645      | 1469          | 1472        | 1474         | rBV2     | 32021          | 28958         | 1.22%           | 0.117%        |
| 23        | 10.751      | 1487          | 1490        | 1494         | rBV3     | 26810          | 31785         | 1.34%           | 0.129%        |
| 24        | 10.833      | 1502          | 1504        | 1510         | rVB4     | 21827          | 31070         | 1.31%           | 0.126%        |
| 25        | 10.969      | 1523          | 1527        | 1530         | rBV      | 109528         | 98673         | 4.17%           | 0.400%        |
| 26        | 11.033      | 1535          | 1538        | 1541         | rVV      | 149034         | 131065        | 5.54%           | 0.532%        |
| 27        | 11.075      | 1541          | 1545        | 1547         | rVV      | 929335         | 798466        | 33.78%          | 3.240%        |
| 28        | 11.098      | 1547          | 1549        | 1552         | rVB      | 54880          | 48329         | 2.04%           | 0.196%        |
| 29        | 11.145      | 1552          | 1557        | 1561         | rVB2     | 33746          | 38439         | 1.63%           | 0.156%        |
| 30        | 11.192      | 1561          | 1565        | 1567         | rBV      | 86892          | 73088         | 3.09%           | 0.297%        |
| 31        | 11.298      | 1580          | 1583        | 1588         | rVB2     | 27827          | 33566         | 1.42%           | 0.136%        |
| 32        | 11.392      | 1596          | 1599        | 1606         | rVB      | 56683          | 54322         | 2.30%           | 0.220%        |
| 33        | 11.492      | 1612          | 1616        | 1619         | rVB      | 348686         | 309666        | 13.10%          | 1.256%        |
| 34        | 11.627      | 1636          | 1639        | 1641         | rBV      | 80045          | 66652         | 2.82%           | 0.270%        |

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Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON Filtering: 5  
 Sampling : 1 Min Area: 3 % of largest Peak  
 Start Thrs: 0.2 Max Peaks: 100  
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

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

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.798 | 1665 | 1668 | 1672 | rVB  | 36119   | 31827   | 1.35%  | 0.129% |
| 36 | 12.169 | 1727 | 1731 | 1733 | rBV  | 847283  | 778963  | 32.95% | 3.161% |
| 37 | 12.192 | 1733 | 1735 | 1742 | rVB  | 918987  | 841111  | 35.58% | 3.413% |
| 38 | 12.280 | 1746 | 1750 | 1753 | rBV  | 316331  | 267382  | 11.31% | 1.085% |
| 39 | 12.504 | 1783 | 1788 | 1792 | rBV  | 136119  | 137614  | 5.82%  | 0.558% |
| 40 | 12.551 | 1792 | 1796 | 1803 | rVB3 | 34859   | 59322   | 2.51%  | 0.241% |
| 41 | 12.663 | 1810 | 1815 | 1818 | rBV  | 1508332 | 1534635 | 64.92% | 6.227% |
| 42 | 12.833 | 1839 | 1844 | 1847 | rBV4 | 29253   | 31200   | 1.32%  | 0.127% |
| 43 | 12.910 | 1854 | 1857 | 1860 | rBV3 | 21681   | 26155   | 1.11%  | 0.106% |
| 44 | 13.427 | 1940 | 1945 | 1947 | rBV6 | 14675   | 27647   | 1.17%  | 0.112% |
| 45 | 13.568 | 1964 | 1969 | 1974 | rBV3 | 118742  | 147514  | 6.24%  | 0.599% |
| 46 | 13.698 | 1987 | 1991 | 1994 | rBV  | 614320  | 625963  | 26.48% | 2.540% |
| 47 | 13.721 | 1994 | 1995 | 1999 | rVB2 | 71225   | 55038   | 2.33%  | 0.223% |
| 48 | 14.198 | 2074 | 2076 | 2080 | rBV2 | 33737   | 37710   | 1.60%  | 0.153% |
| 49 | 14.698 | 2158 | 2161 | 2163 | rBV  | 87350   | 108067  | 4.57%  | 0.438% |
| 50 | 14.957 | 2203 | 2205 | 2210 | rVB  | 62015   | 64887   | 2.74%  | 0.263% |
| 51 | 15.010 | 2211 | 2214 | 2220 | rBV  | 70151   | 96369   | 4.08%  | 0.391% |
| 52 | 15.068 | 2220 | 2224 | 2230 | rVV  | 461157  | 534133  | 22.60% | 2.167% |

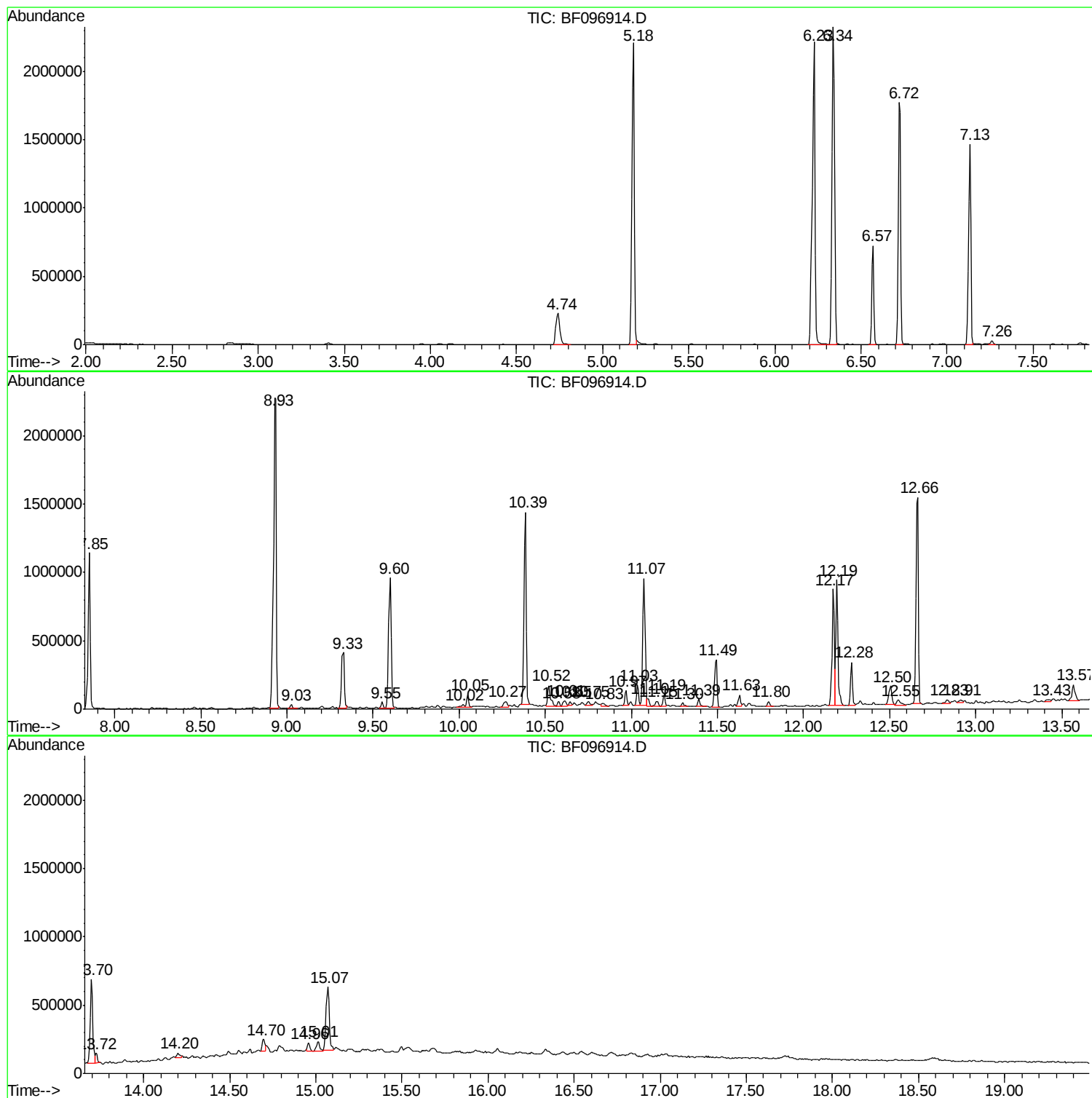
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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

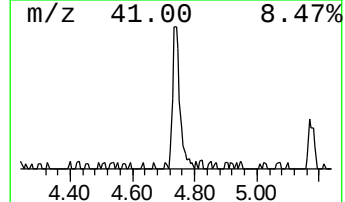
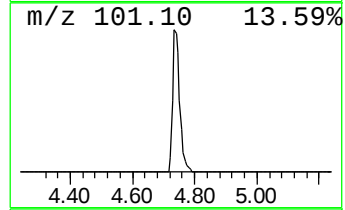
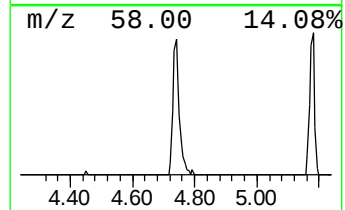
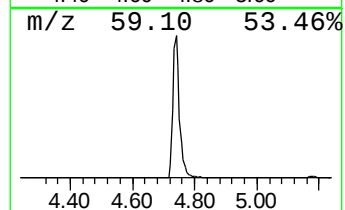
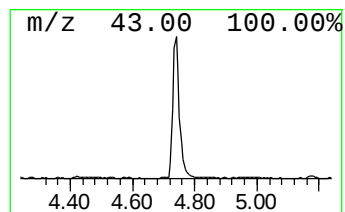
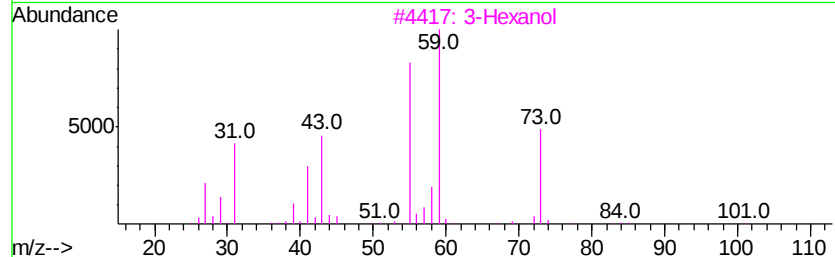
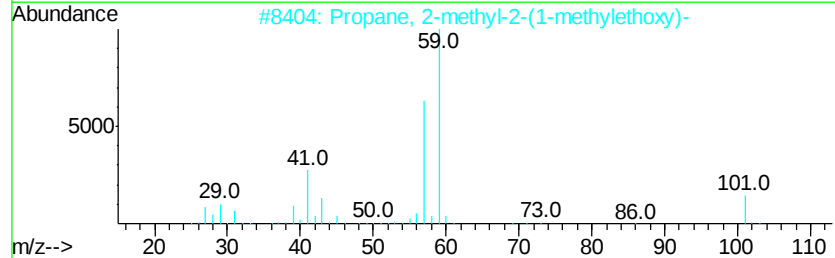
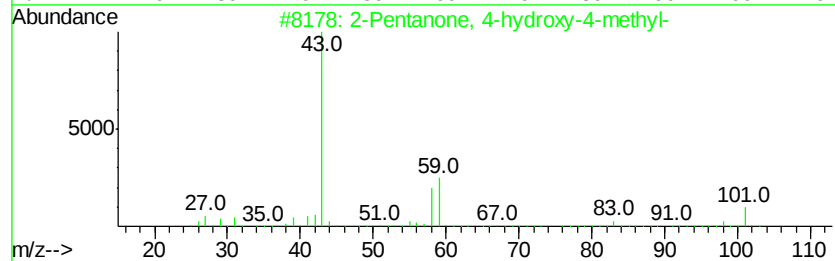
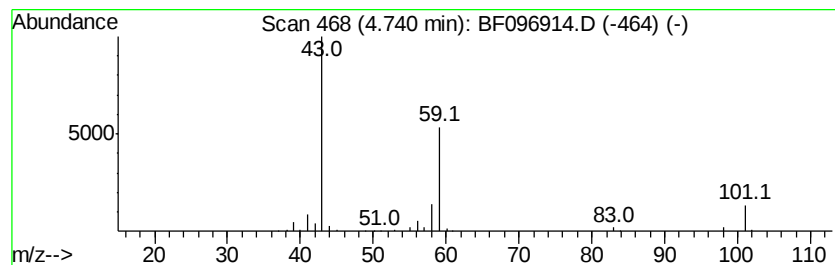
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.74 | 11.10 ng | 341033 | 1,4-Dichlorobenzene-d4 | 6.57 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-      | 116 | C6H12O2 | 000123-42-2 | 59   |
| 2    |    |   | Propane, 2-methyl-2-(1-methylethoxy)- | 116 | C7H16O  | 017348-59-3 | 36   |
| 3    |    |   | 3-Hexanol                             | 102 | C6H14O  | 000623-37-0 | 33   |
| 4    |    |   | 3-Hexanol, 4-methyl-                  | 116 | C7H16O  | 000615-29-2 | 33   |
| 5    |    |   | 2-Hexanol, 2-methyl-                  | 116 | C7H16O  | 000625-23-0 | 28   |



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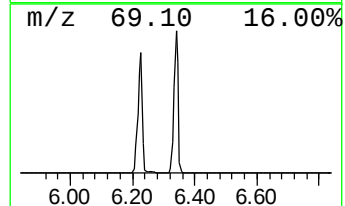
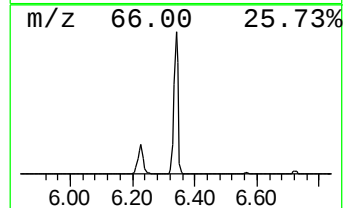
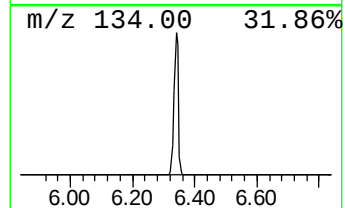
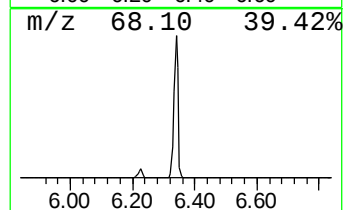
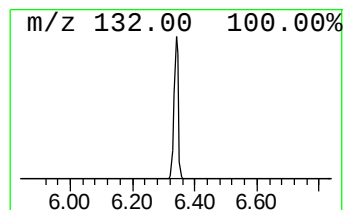
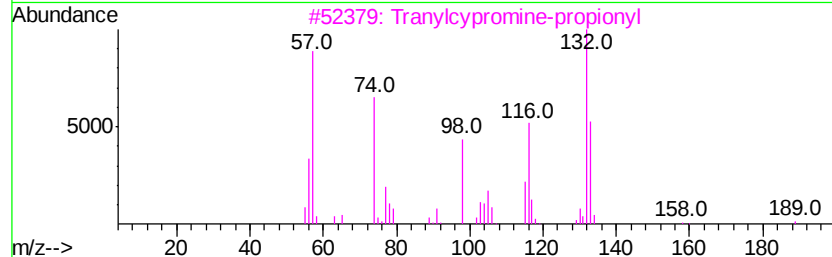
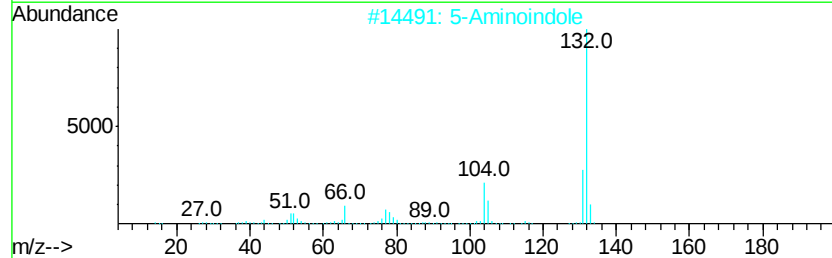
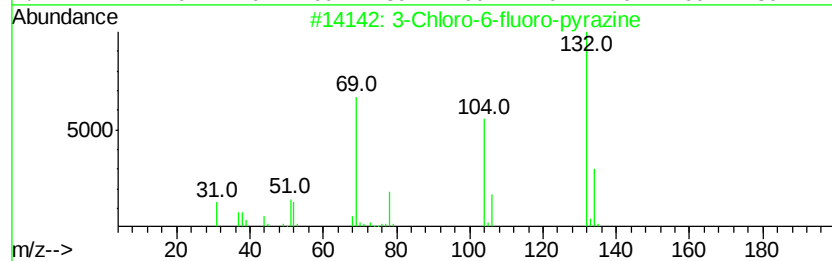
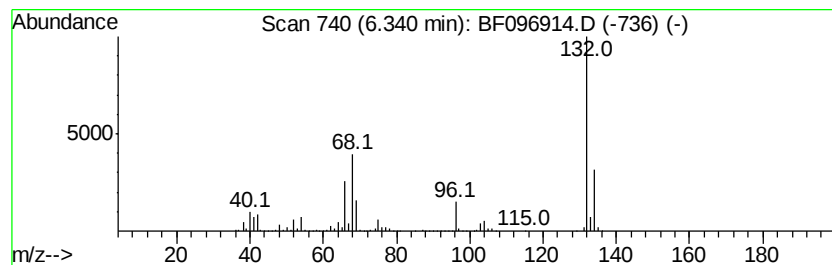
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF071317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 unknown6.34 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.34 | 74.24 ng | 2280160 | 1,4-Dichlorobenzene-d4 | 6.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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

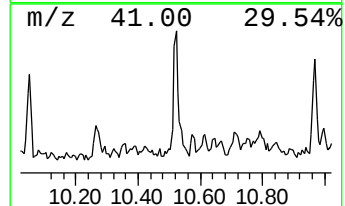
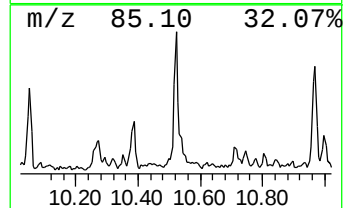
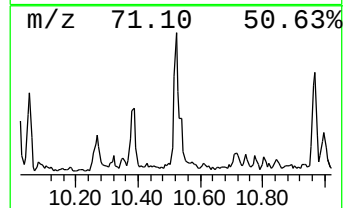
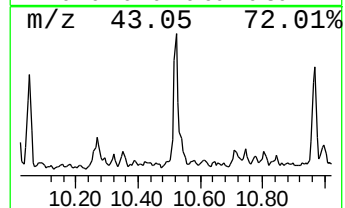
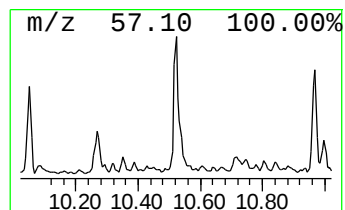
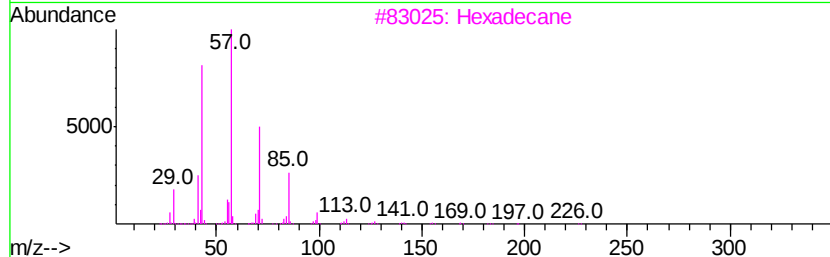
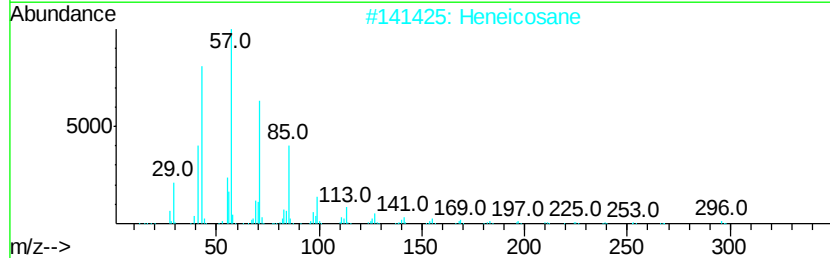
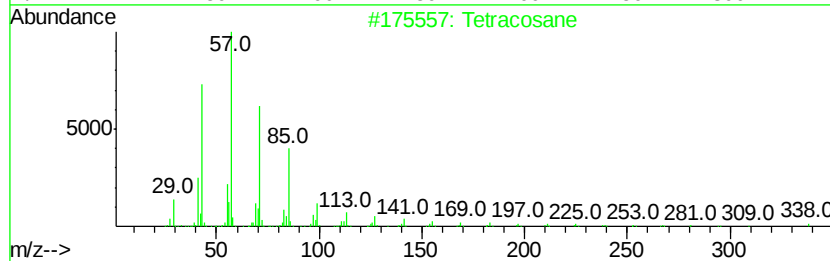
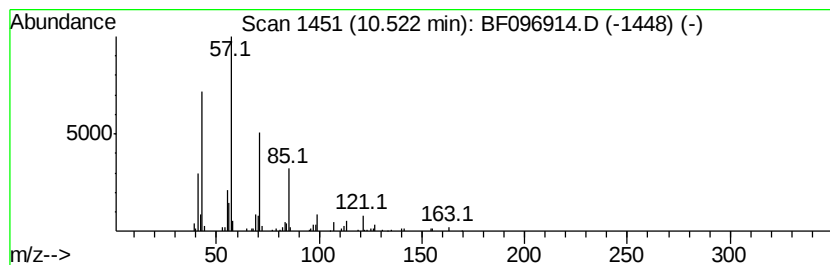
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Tetracosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.52 | 4.26 ng | 169925 | Phenanthrene-d10 | 11.07 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane  | 338 | C24H50  | 000646-31-1 | 86   |
| 2    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 86   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 80   |
| 5    |    | Pentadecane  | 212 | C15H32  | 000629-62-9 | 80   |



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

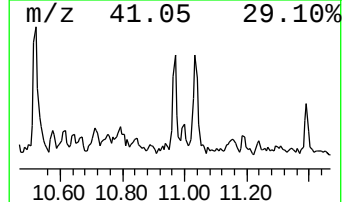
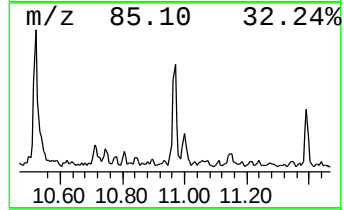
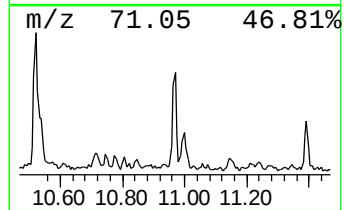
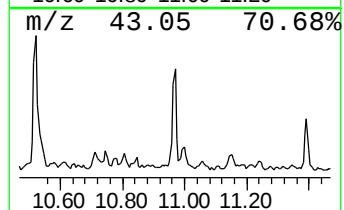
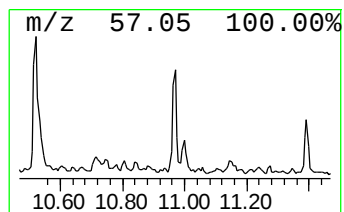
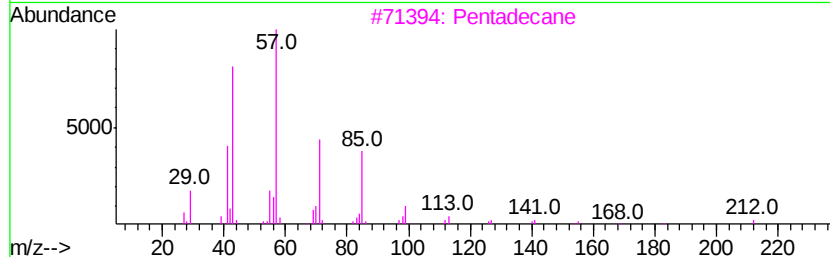
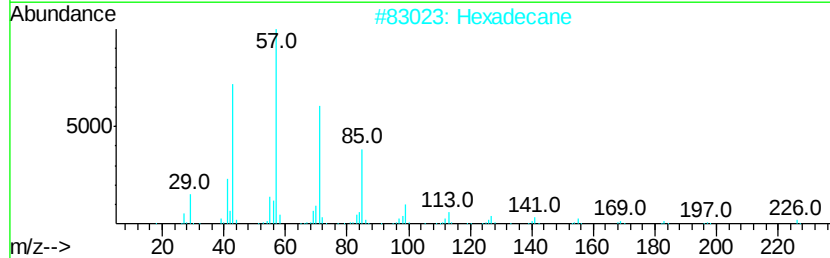
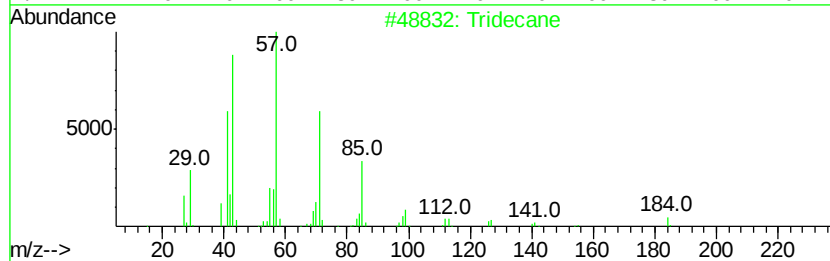
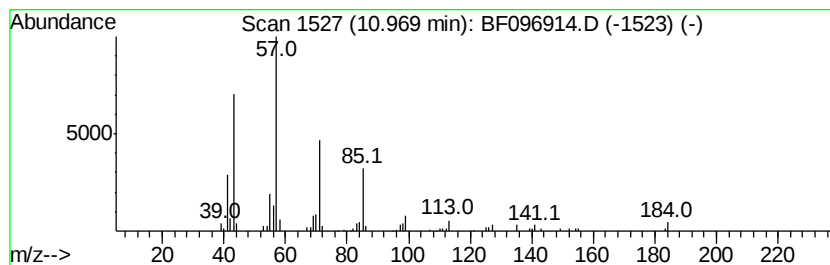
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 10.97 | 2.47 ng | 98673 | Phenanthrene-d10 | 11.07 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|---------|-------------------------------------|--------------|-----|-----------|--------------|------|
| 1       | Tridecane                           |              | 184 | C13H28    | 000629-50-5  | 90   |
| 2       | Hexadecane                          |              | 226 | C16H34    | 000544-76-3  | 90   |
| 3       | Pentadecane                         |              | 212 | C15H32    | 000629-62-9  | 90   |
| 4       | Sulfurous acid, butyl dodecyl ester |              | 306 | C16H34O3S | 1000309-17-9 | 90   |
| 5       | Eicosane                            |              | 282 | C20H42    | 000112-95-8  | 87   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

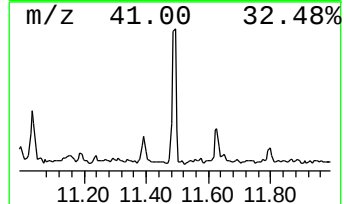
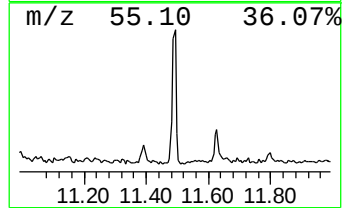
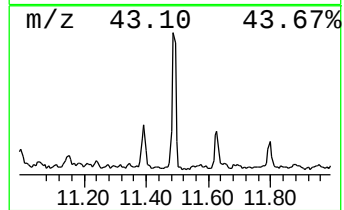
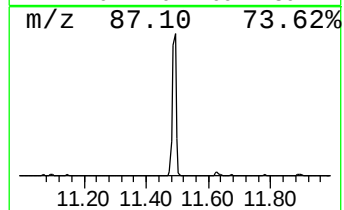
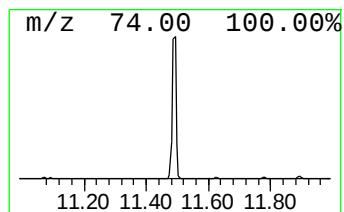
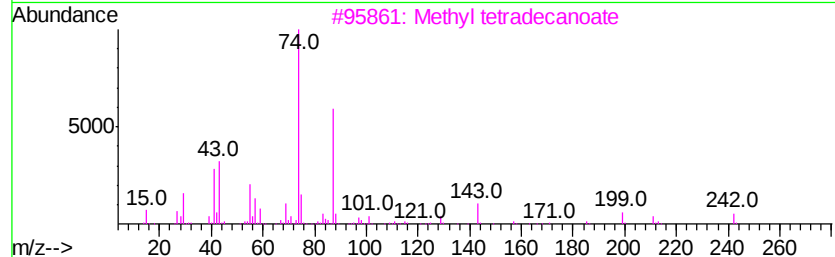
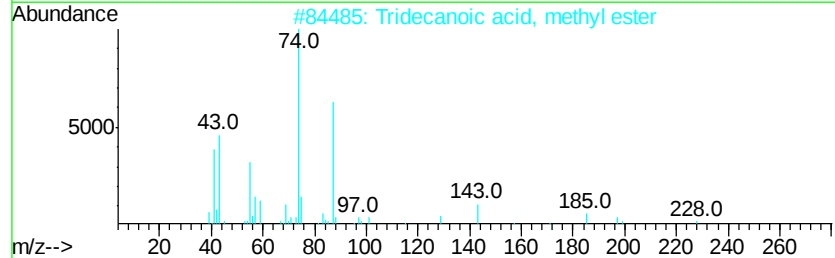
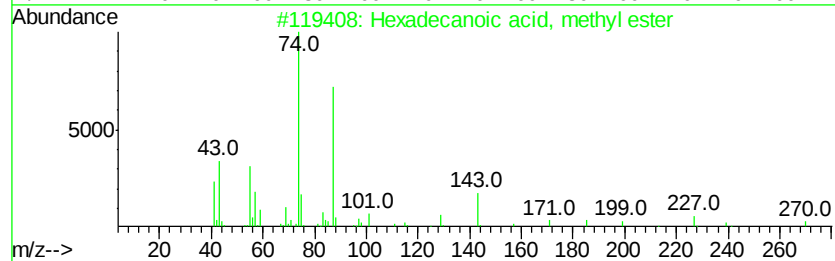
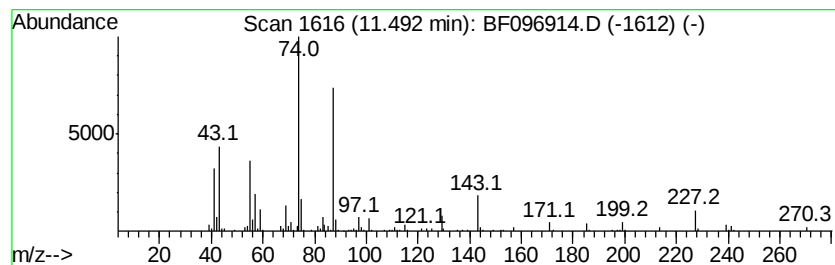
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.49 | 7.76 ng | 309666 | Phenanthrene-d10 | 11.07 |

| Hit# of | 5 | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|---------|---|----------------------------------|-----|----------|-------------|------|
| 1       |   | Hexadecanoic acid, methyl ester  | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2       |   | Tridecanoic acid, methyl ester   | 228 | C14H28O2 | 001731-88-0 | 87   |
| 3       |   | Methyl tetradecanoate            | 242 | C15H30O2 | 000124-10-7 | 87   |
| 4       |   | Pentadecanoic acid, methyl ester | 256 | C16H32O2 | 007132-64-1 | 86   |
| 5       |   | Dodecanoic acid, methyl ester    | 214 | C13H26O2 | 000111-82-0 | 80   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

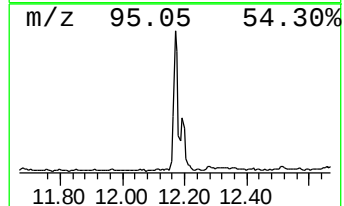
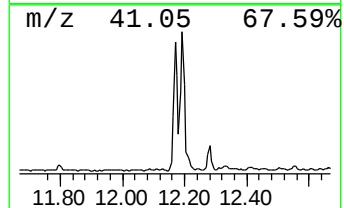
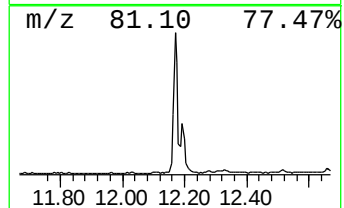
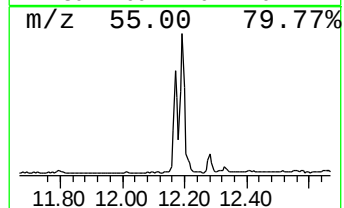
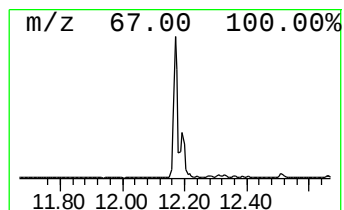
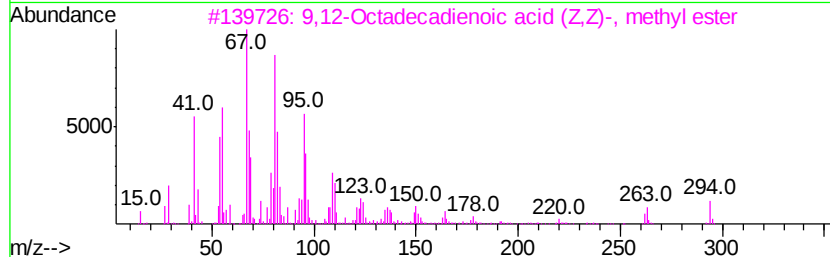
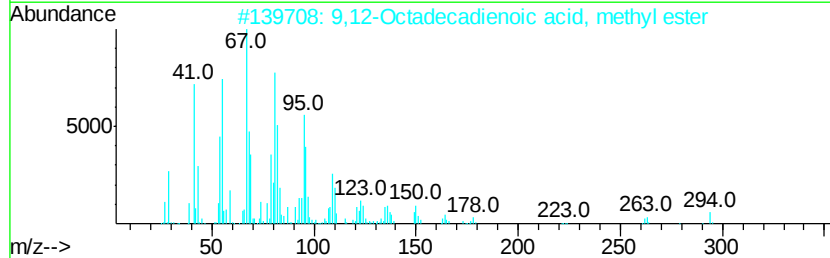
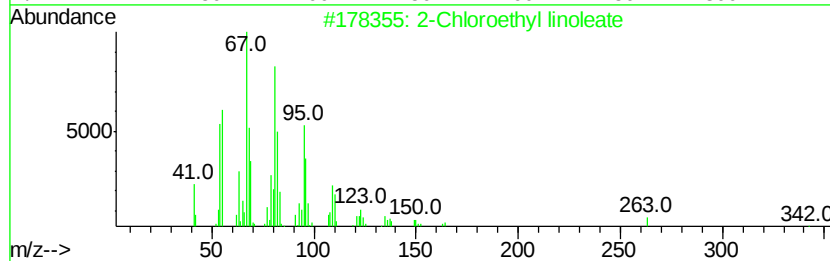
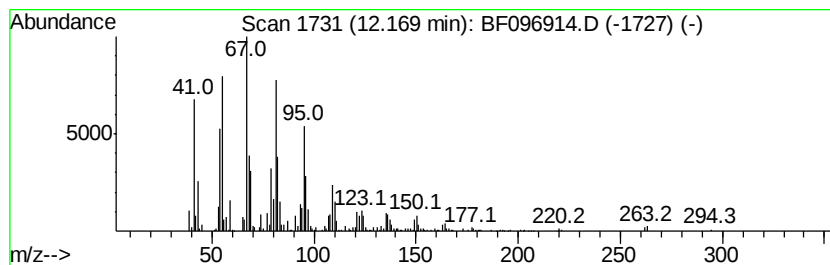
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2-Chloroethyl linoleate Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.17 | 19.51 ng | 778963 | Phenanthrene-d10 | 11.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 2-Chloroethyl linoleate             | 342 | C20H35ClO2 | 025525-76-2 | 94   |
| 2    |    | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2   | 002462-85-3 | 93   |
| 3    |    | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2   | 000112-63-0 | 91   |
| 4    |    | 1-Tridecyne                         | 180 | C13H24     | 026186-02-7 | 91   |
| 5    |    | 9,12-Octadecadienoic acid (Z,Z)-    | 280 | C18H32O2   | 000060-33-3 | 90   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

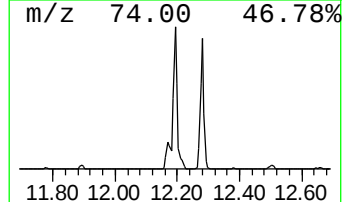
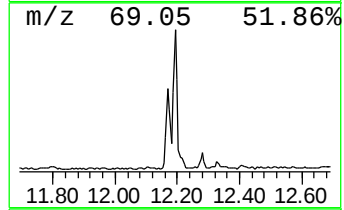
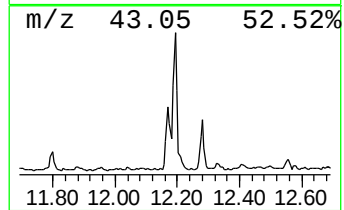
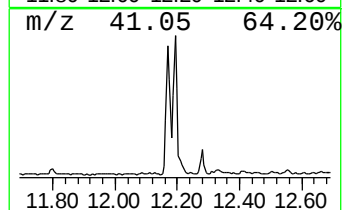
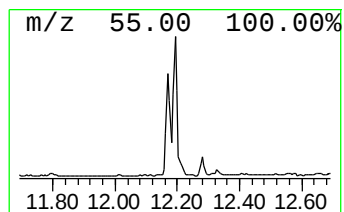
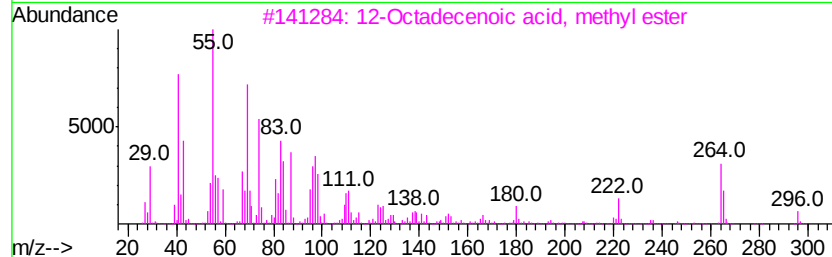
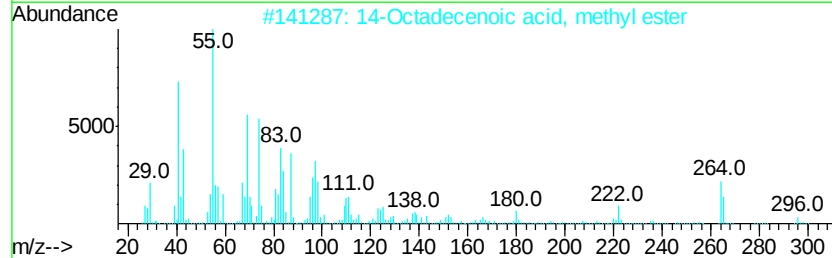
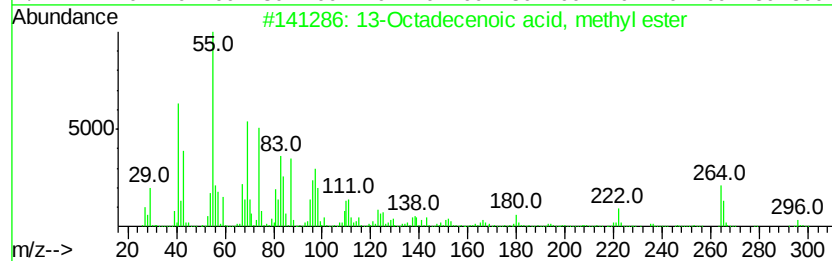
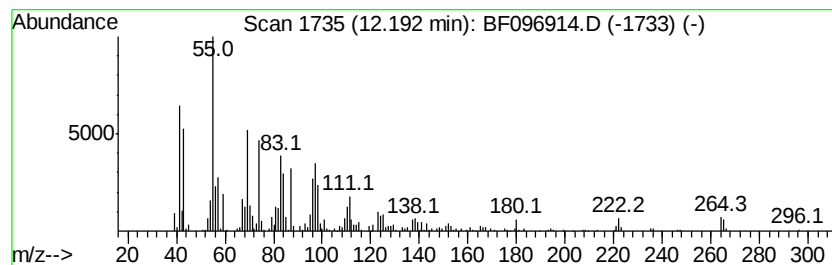
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 13-Octadecenoic acid, methyl... Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.19 | 21.07 ng | 841111 | Phenanthrene-d10 | 11.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 13-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-47-3 | 95   |
| 2    |    | 14-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-48-4 | 95   |
| 3    |    | 12-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-46-2 | 94   |
| 4    |    | 10-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 013481-95-3 | 94   |
| 5    |    | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9 | 94   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

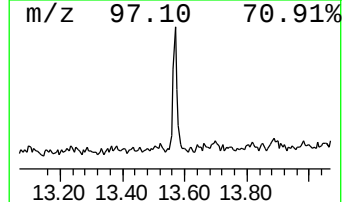
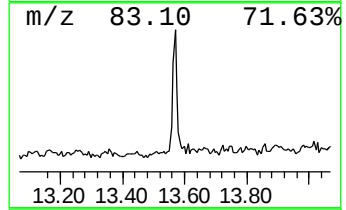
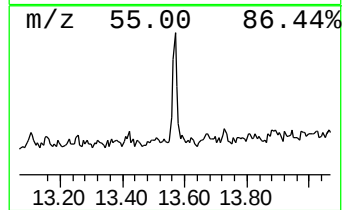
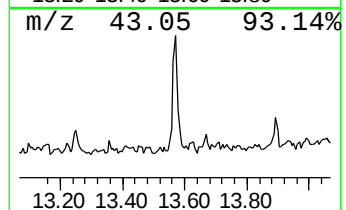
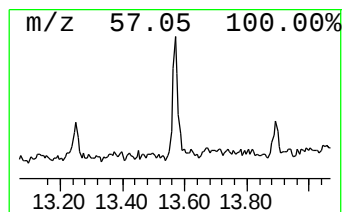
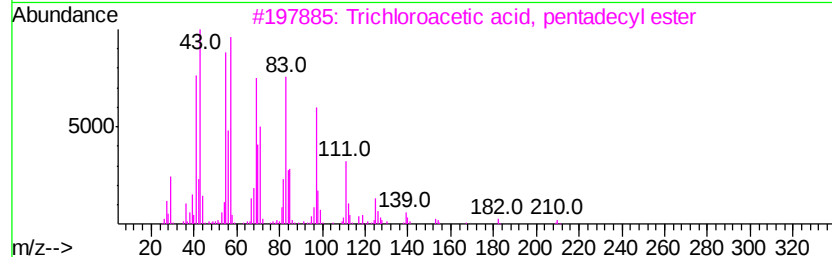
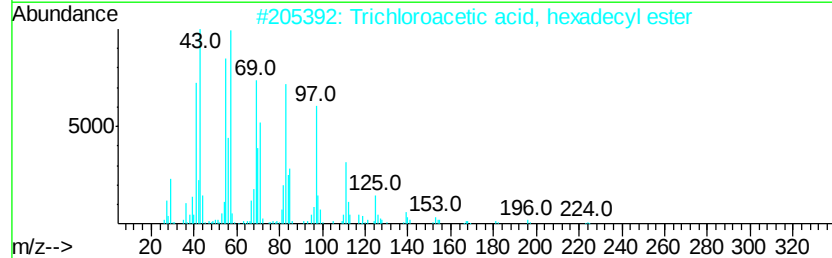
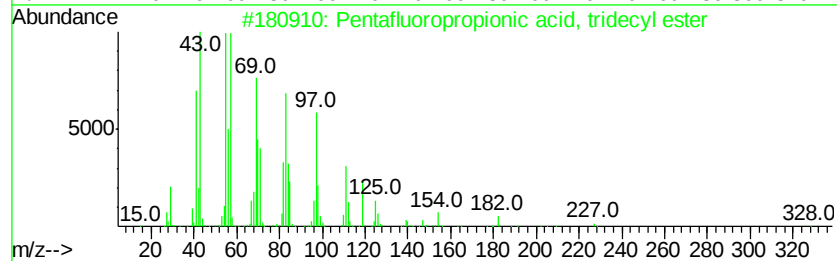
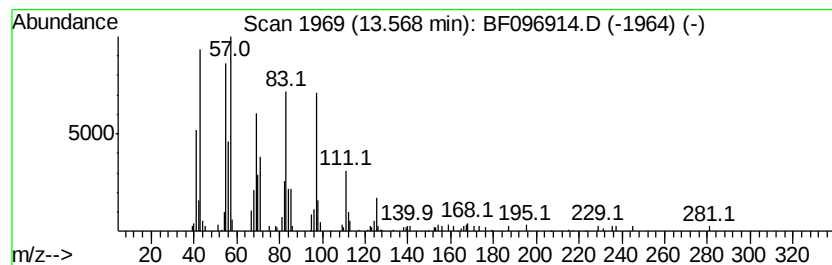
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Pentafluoropropionic acid, ... Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.57 | 4.71 ng | 147514 | Chrysene-d12     | 13.70 |

| Hit# of | 5 | Tentative ID                           | MW  | MolForm     | CAS#        | Qual |
|---------|---|----------------------------------------|-----|-------------|-------------|------|
| 1       |   | Pentafluoropropionic acid, tridecyl... | 346 | C16H27F5O2  | 959261-22-4 | 91   |
| 2       |   | Trichloroacetic acid, hexadecyl ...    | 386 | C18H33Cl3O2 | 074339-54-1 | 91   |
| 3       |   | Trichloroacetic acid, pentadecyl...    | 372 | C17H31Cl3O2 | 074339-53-0 | 91   |
| 4       |   | 11-Tricosene                           | 322 | C23H46      | 052078-56-5 | 91   |
| 5       |   | Trichloroacetic acid, tetradecyl...    | 358 | C16H29Cl3O2 | 074339-52-9 | 90   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

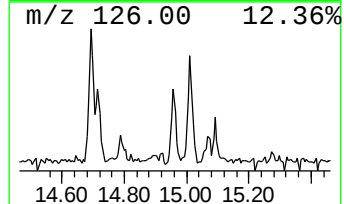
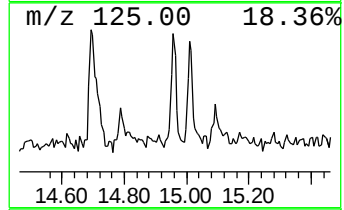
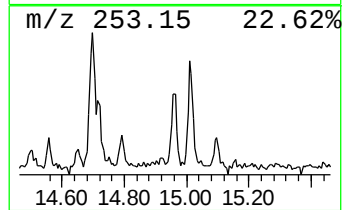
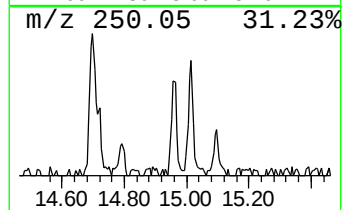
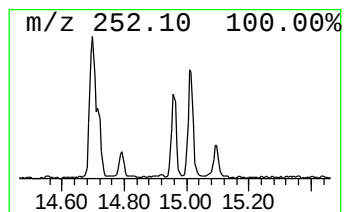
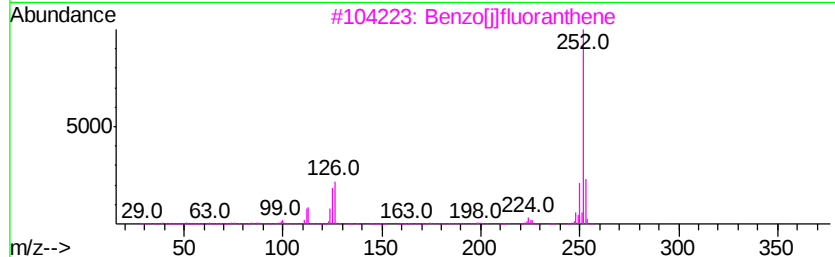
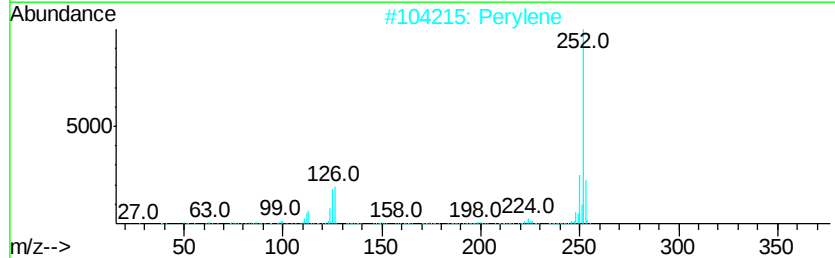
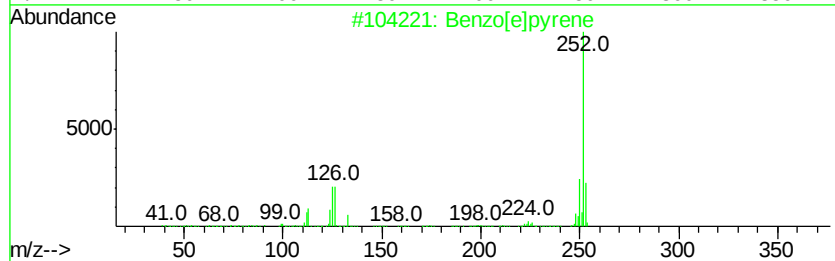
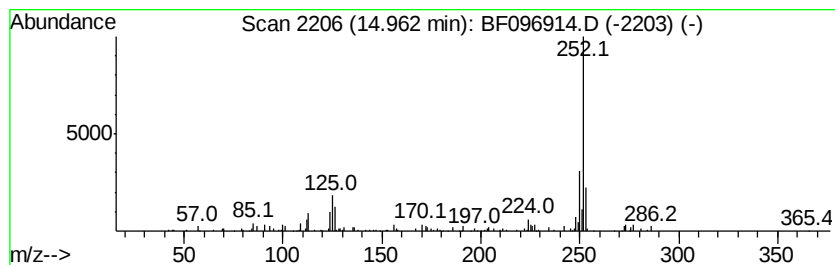
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzo[e]pyrene Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.96 | 2.43 ng | 64887 | Perylene-d12     | 15.07 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene            | 252 | C20H12  | 000192-97-2 | 93   |
| 2    |    |   | Perylene                  | 252 | C20H12  | 000198-55-0 | 93   |
| 3    |    |   | Benzo[il]fluoranthene     | 252 | C20H12  | 000205-82-3 | 86   |
| 4    |    |   | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 76   |
| 5    |    |   | Benzo[a]pyrene            | 252 | C20H12  | 000050-32-8 | 70   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF072017\  
Data File : BF096914.D  
Acq On : 20 Jul 2017 19:12  
Operator : SJ/JU  
Sample : I4308-01  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NB-08-071917

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

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.74  | 11.1 ng |       | 341033   | 1                     | 6.57  | 614281 | 20.0 |
| unknown6.34          | 6.34  | 74.2 ng |       | 2280160  | 1                     | 6.57  | 614281 | 20.0 |
| Tetracosane          | 10.52 | 4.3 ng  |       | 169925   | 4                     | 11.07 | 798466 | 20.0 |
| Tridecane            | 10.97 | 2.5 ng  |       | 98673    | 4                     | 11.07 | 798466 | 20.0 |
| Hexadecanoic acid... | 11.49 | 7.8 ng  |       | 309666   | 4                     | 11.07 | 798466 | 20.0 |
| 2-Chloroethyl lin... | 12.17 | 19.5 ng |       | 778963   | 4                     | 11.07 | 798466 | 20.0 |
| 13-Octadecenoic a... | 12.19 | 21.1 ng |       | 841111   | 4                     | 11.07 | 798466 | 20.0 |
| Pentafluoropropio... | 13.57 | 4.7 ng  |       | 147514   | 5                     | 13.70 | 625963 | 20.0 |
| Benzo[e]pyrene       | 14.96 | 2.4 ng  |       | 64887    | 6                     | 15.07 | 534133 | 20.0 |