

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
 Data File : BF098583.D  
 Acq On : 15 Sep 2017 21:57  
 Operator : SJ/JU  
 Sample : I5236-01 5X  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TR-04-091317

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-BF091117.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         | 5.245       | 534           | 537         | 547          | rBV      | 602749         | 655620        | 5.68%           | 1.530%        |
| 2         | 6.292       | 712           | 715         | 732          | rBV      | 652412         | 706913        | 6.13%           | 1.650%        |
| 3         | 6.416       | 732           | 736         | 742          | rVB      | 744435         | 717109        | 6.22%           | 1.674%        |
| 4         | 6.639       | 770           | 774         | 781          | rBV      | 912891         | 817226        | 7.08%           | 1.907%        |
| 5         | 6.798       | 797           | 801         | 808          | rVB      | 546552         | 518586        | 4.50%           | 1.210%        |
| 6         | 7.210       | 868           | 871         | 875          | rBV      | 464849         | 403379        | 3.50%           | 0.941%        |
| 7         | 7.928       | 987           | 993         | 996          | rBV      | 1312434        | 1141171       | 9.89%           | 2.663%        |
| 8         | 9.004       | 1172          | 1176        | 1179         | rBV      | 819647         | 685803        | 5.95%           | 1.601%        |
| 9         | 9.086       | 1185          | 1190        | 1194         | rBV      | 115863         | 117484        | 1.02%           | 0.274%        |
| 10        | 9.275       | 1217          | 1222        | 1226         | rBV4     | 111805         | 120891        | 1.05%           | 0.282%        |
| 11        | 9.404       | 1241          | 1244        | 1247         | rVB      | 167156         | 132671        | 1.15%           | 0.310%        |
| 12        | 9.680       | 1284          | 1291        | 1294         | rBV      | 1226271        | 1126725       | 9.77%           | 2.630%        |
| 13        | 10.469      | 1422          | 1425        | 1431         | rVB      | 424525         | 353197        | 3.06%           | 0.824%        |
| 14        | 10.586      | 1442          | 1445        | 1452         | rVB      | 165112         | 198749        | 1.72%           | 0.464%        |
| 15        | 10.692      | 1460          | 1463        | 1466         | rVB      | 196562         | 149777        | 1.30%           | 0.350%        |
| 16        | 11.163      | 1539          | 1543        | 1545         | rBV      | 1268161        | 1033866       | 8.96%           | 2.413%        |
| 17        | 11.186      | 1545          | 1547        | 1550         | rVB      | 284898         | 220878        | 1.91%           | 0.515%        |
| 18        | 11.486      | 1595          | 1598        | 1601         | rVB2     | 321857         | 290987        | 2.52%           | 0.679%        |
| 19        | 11.569      | 1606          | 1612        | 1615         | rVB      | 6117974        | 5533935       | 47.98%          | 12.915%       |
| 20        | 11.774      | 1643          | 1647        | 1649         | rBV      | 115911         | 122297        | 1.06%           | 0.285%        |
| 21        | 12.257      | 1723          | 1729        | 1730         | rBV      | 5990322        | 7571670       | 65.64%          | 17.671%       |
| 22        | 12.280      | 1730          | 1733        | 1739         | rVB      | 8669500        | 11534972      | 100.00%         | 26.920%       |
| 23        | 12.357      | 1742          | 1746        | 1748         | rBV      | 3532798        | 3002720       | 26.03%          | 7.008%        |
| 24        | 12.604      | 1782          | 1788        | 1790         | rBV      | 492895         | 539348        | 4.68%           | 1.259%        |
| 25        | 12.745      | 1809          | 1812        | 1815         | rBV      | 707254         | 578514        | 5.02%           | 1.350%        |
| 26        | 12.927      | 1837          | 1843        | 1848         | rBV2     | 133052         | 266541        | 2.31%           | 0.622%        |
| 27        | 12.998      | 1849          | 1855        | 1859         | rVV2     | 345750         | 410292        | 3.56%           | 0.958%        |
| 28        | 13.074      | 1865          | 1868        | 1871         | rBV      | 252497         | 231068        | 2.00%           | 0.539%        |
| 29        | 13.186      | 1884          | 1887        | 1892         | rVB4     | 137716         | 152449        | 1.32%           | 0.356%        |
| 30        | 13.504      | 1937          | 1941        | 1944         | rVB2     | 351066         | 468338        | 4.06%           | 1.093%        |
| 31        | 13.739      | 1979          | 1981        | 1985         | rVB      | 184911         | 172116        | 1.49%           | 0.402%        |
| 32        | 13.798      | 1986          | 1991        | 1994         | rVV2     | 959606         | 1035872       | 8.98%           | 2.418%        |
| 33        | 13.827      | 1994          | 1996        | 2000         | rVB      | 173469         | 151267        | 1.31%           | 0.353%        |
| 34        | 14.810      | 2160          | 2163        | 2170         | rVB      | 190356         | 286054        | 2.48%           | 0.668%        |

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Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

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-BF091117.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 35 | 15.086 | 2207 | 2210 | 2215 | rVB  | 118121 | 138756 | 1.20% | 0.324% |
| 36 | 15.145 | 2217 | 2220 | 2225 | rVV  | 188029 | 233371 | 2.02% | 0.545% |
| 37 | 15.204 | 2225 | 2230 | 2240 | rVB  | 647692 | 878272 | 7.61% | 2.050% |
| 38 | 16.198 | 2396 | 2399 | 2405 | rVB2 | 90157  | 149577 | 1.30% | 0.349% |

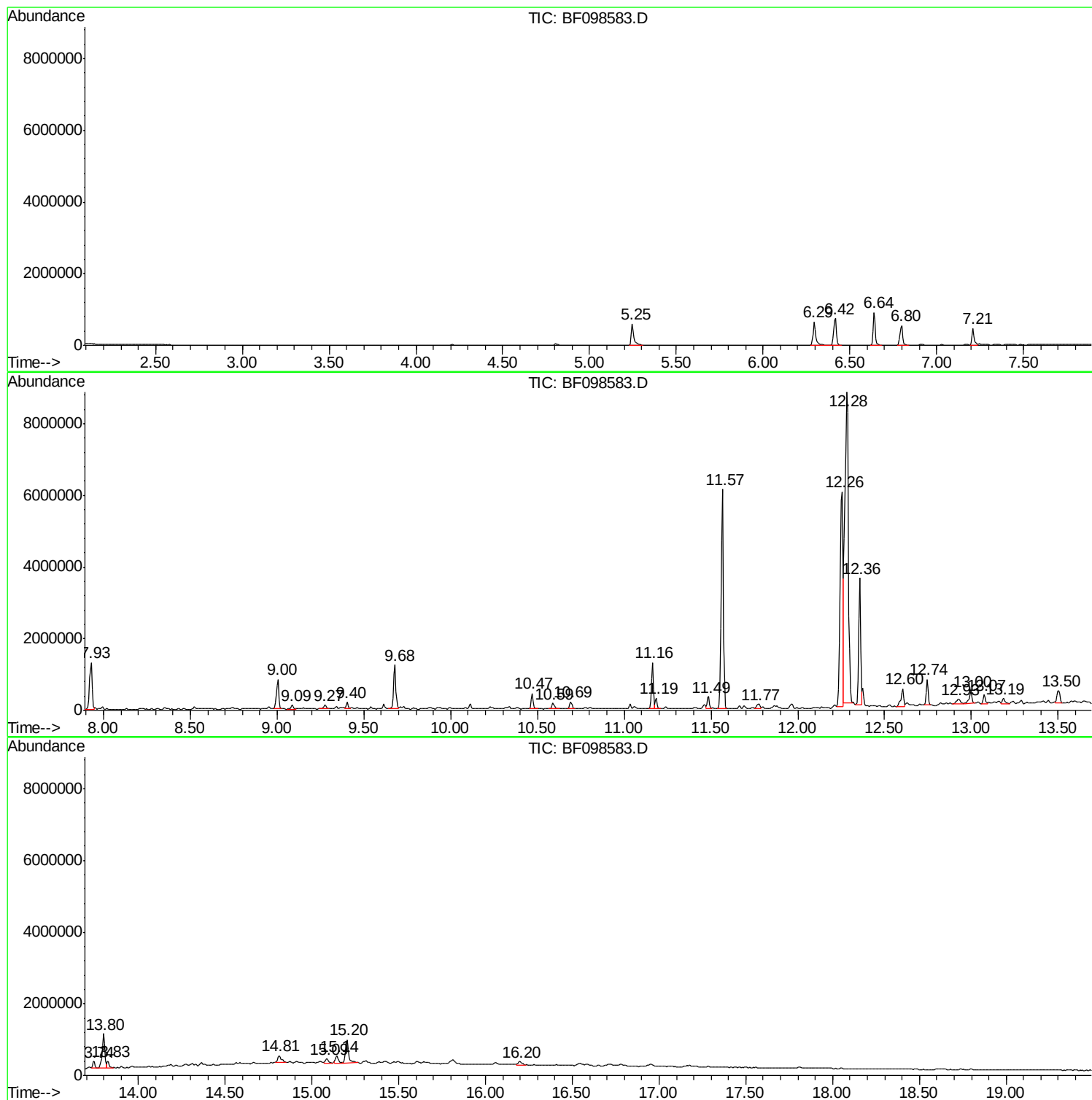
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

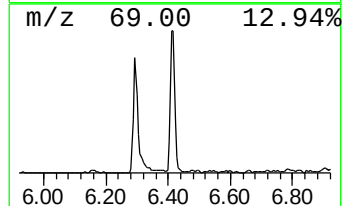
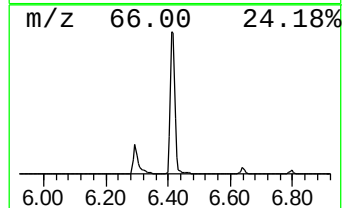
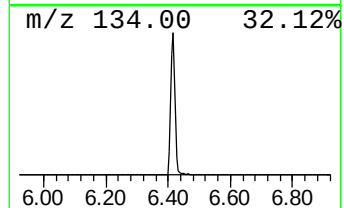
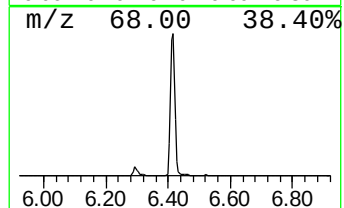
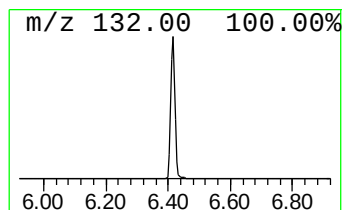
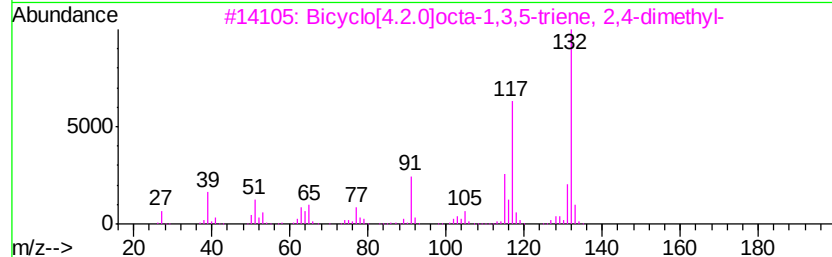
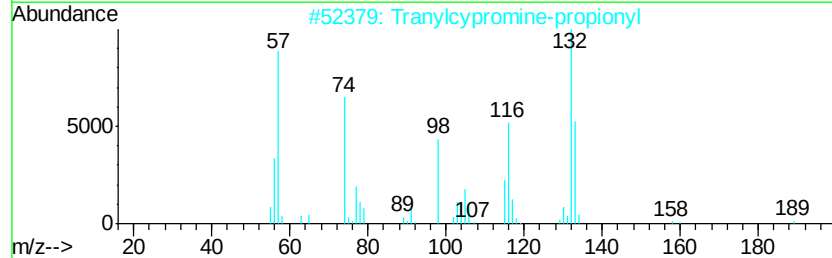
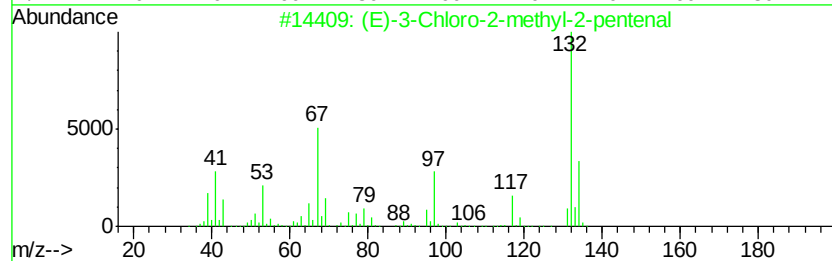
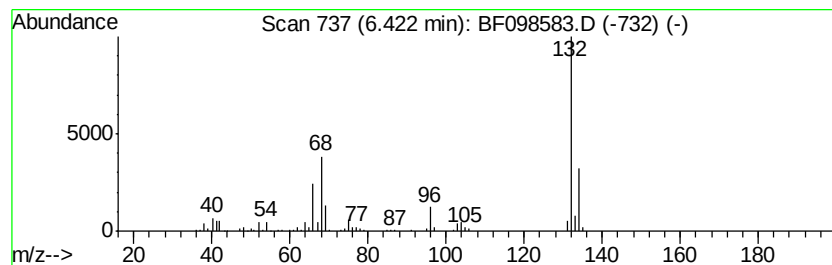
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown6.42 Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.42 | 17.55 ng | 717109 | 1,4-Dichlorobenzene-d4 | 6.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO  | 031357-76-3  | 17   |
| 2    |    | Tranylcypromine-propionyl           | 189 | C12H15NO | 1000123-86-3 | 10   |
| 3    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12   | 028749-81-7  | 9    |
| 4    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2   | 022291-85-6  | 9    |
| 5    |    | 1-Methylpyrrolo[1,2-a]pyrazine      | 132 | C8H8N2   | 064608-59-9  | 9    |



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BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

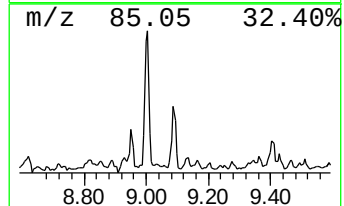
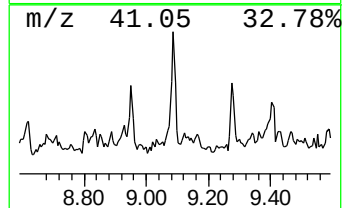
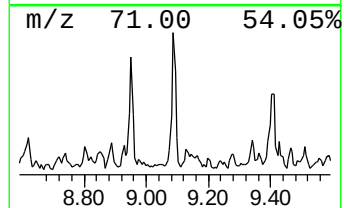
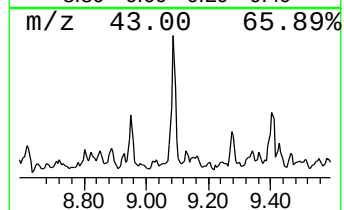
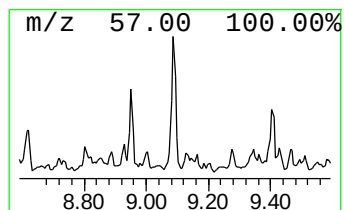
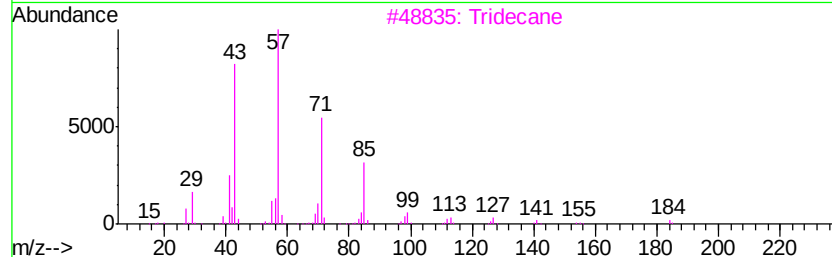
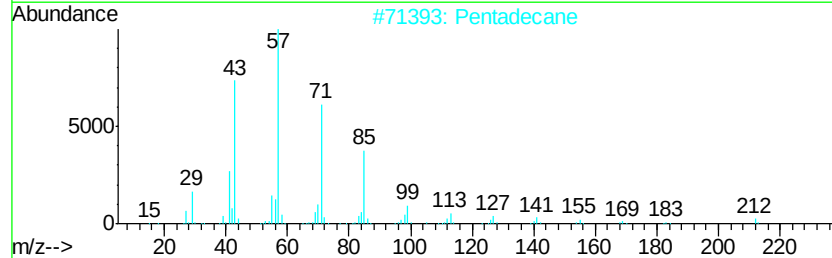
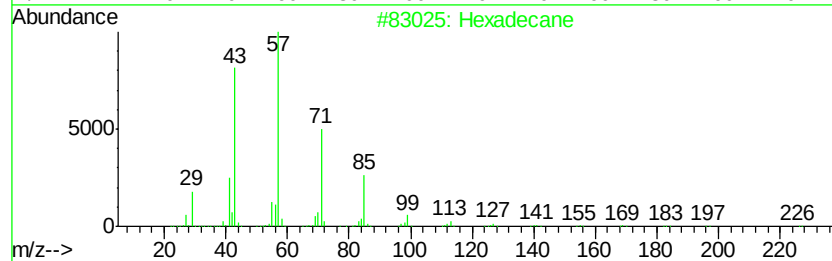
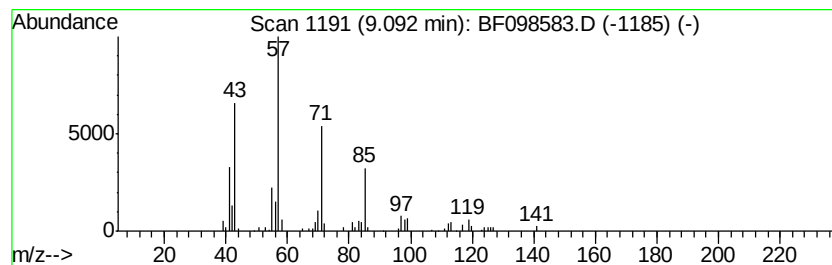
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Hexadecane Concentration Rank 18

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.09 | 2.09 ng | 117484 | Acenaphthene-d10 | 9.68 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane              | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    | Pentadecane             | 212 | C15H32  | 000629-62-9 | 86   |
| 3    |    | Tridecane               | 184 | C13H28  | 000629-50-5 | 86   |
| 4    |    | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 86   |
| 5    |    | Dodecane                | 170 | C12H26  | 000112-40-3 | 80   |



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ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

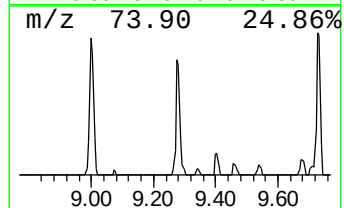
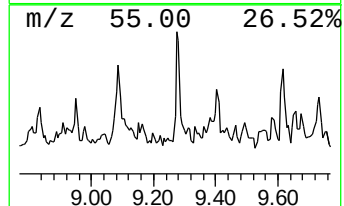
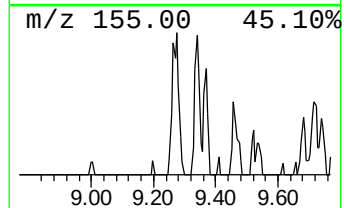
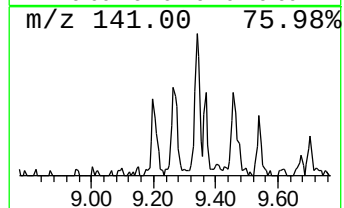
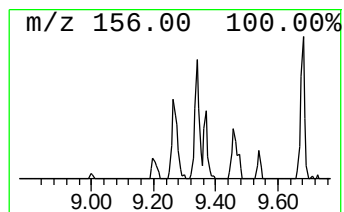
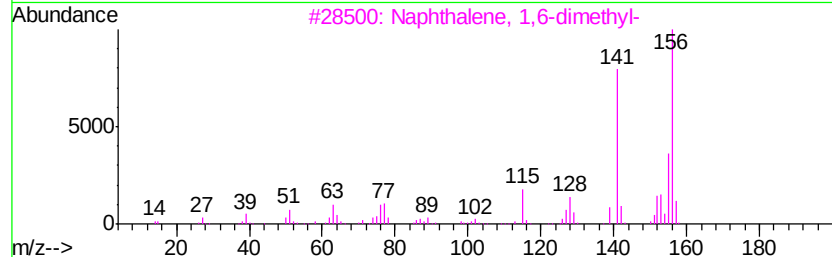
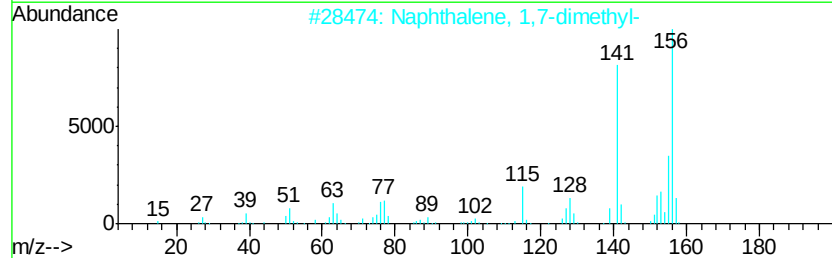
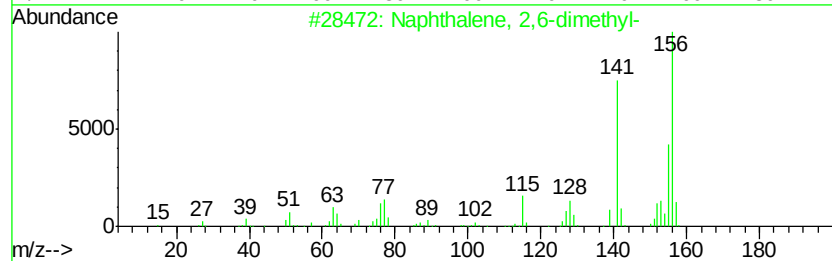
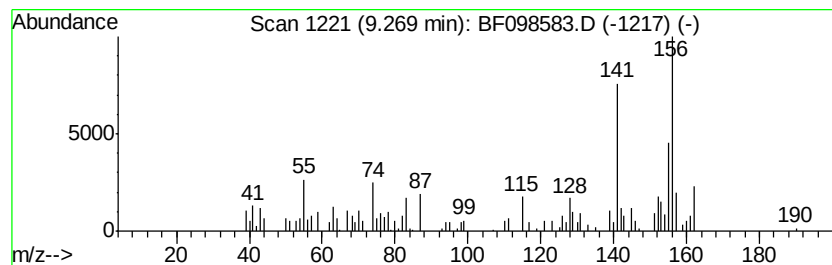
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Naphthalene, 2,6-dimethyl- Concentration Rank 17

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.27 | 2.15 ng | 120891 | Acenaphthene-d10 | 9.68 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 64   |
| 2    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 60   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 35   |
| 4    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 35   |
| 5    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 25   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

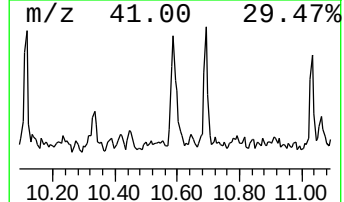
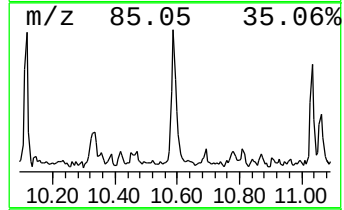
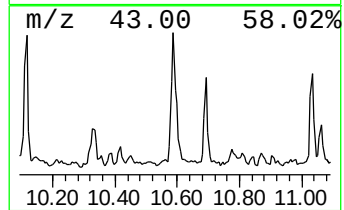
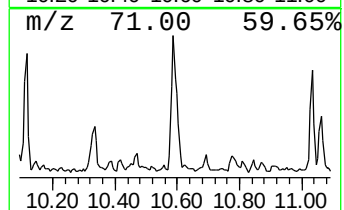
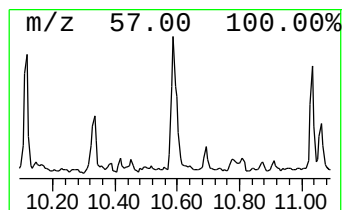
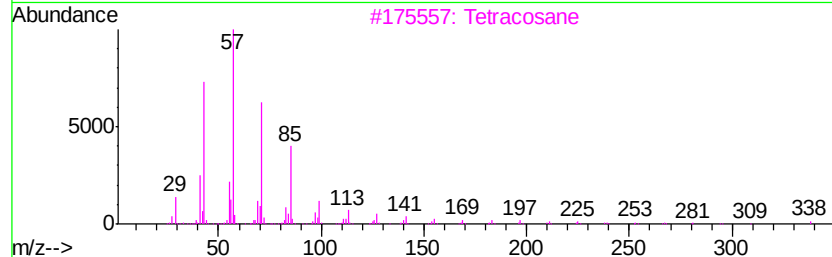
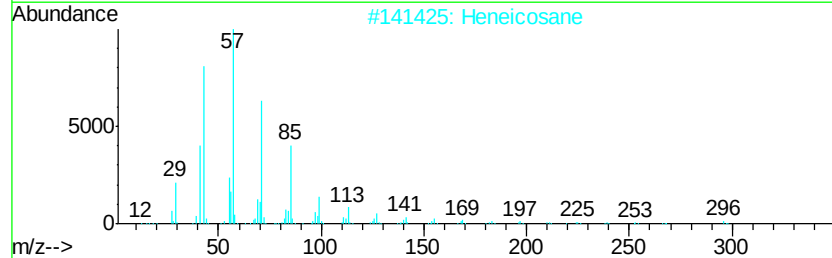
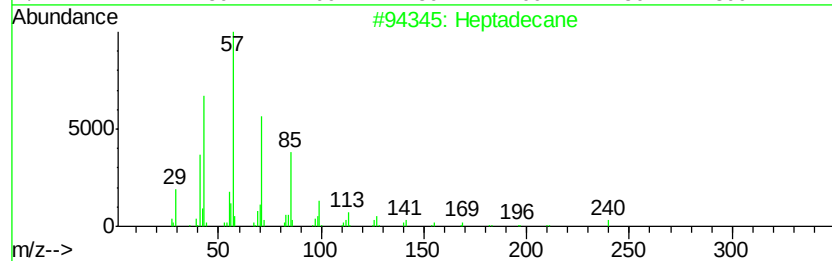
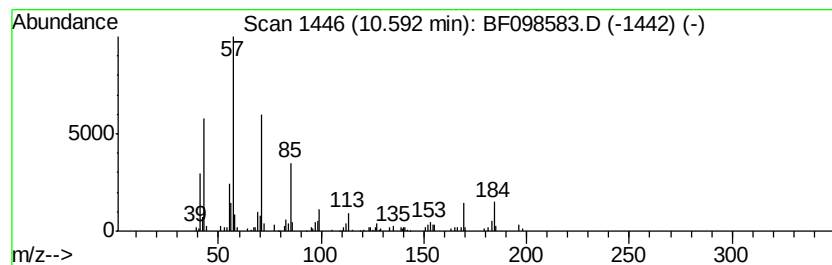
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Heptadecane Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.59 | 3.84 ng | 198749 | Phenanthrene-d10 | 11.16 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Heptadecane |              | 240 | C17H36  | 000629-78-7 | 90   |
| 2       | Heneicosane |              | 296 | C21H44  | 000629-94-7 | 87   |
| 3       | Tetracosane |              | 338 | C24H50  | 000646-31-1 | 86   |
| 4       | Docosane    |              | 310 | C22H46  | 000629-97-0 | 86   |
| 5       | Eicosane    |              | 282 | C20H42  | 000112-95-8 | 80   |



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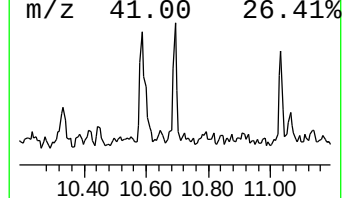
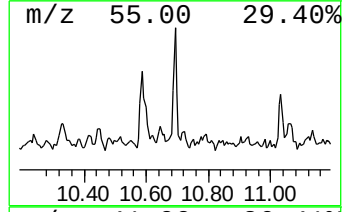
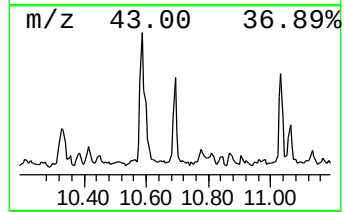
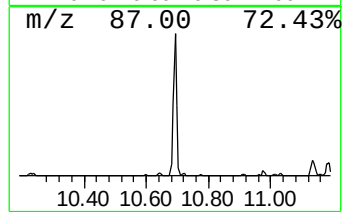
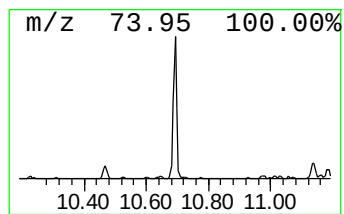
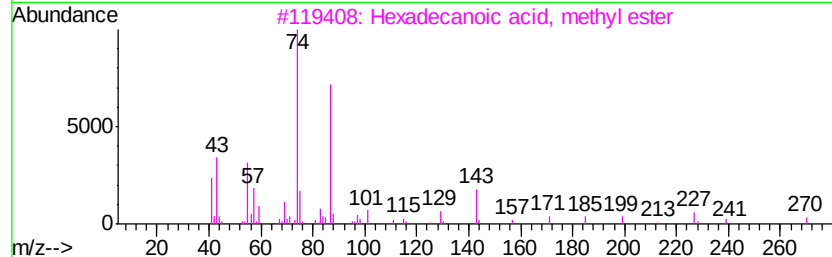
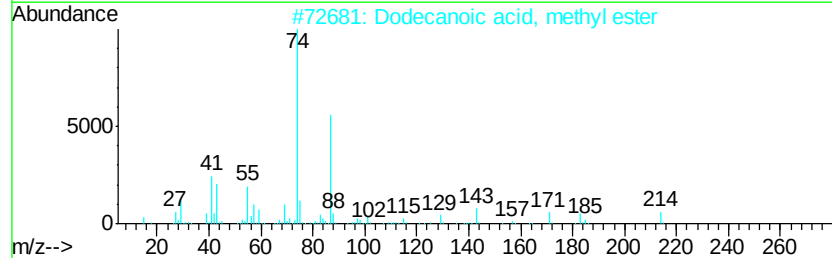
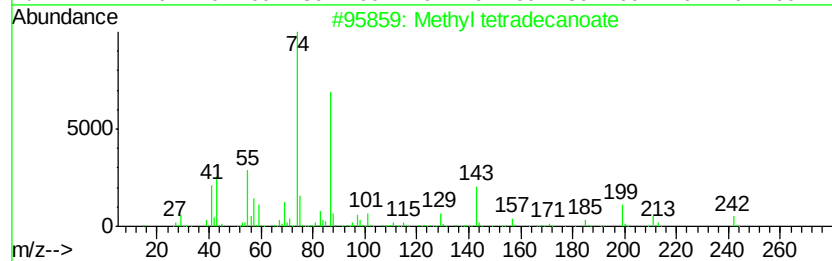
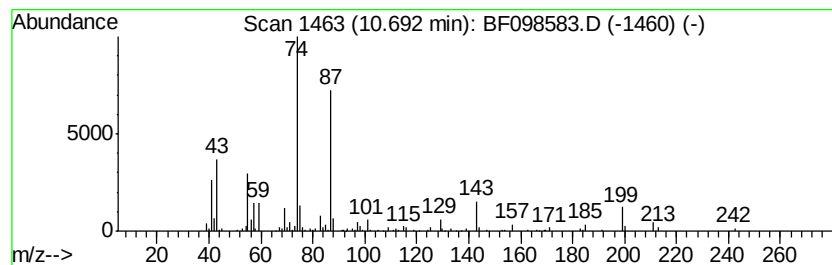
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BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Methyl tetradecanoate Concentration Rank 15

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------------|------------------|----------------|
| 10.69   | 2.90 ng | 149777                              | Phenanthrene-d10 | 11.16          |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |         | Methyl tetradecanoate               | 242 C15H30O2     | 000124-10-7 91 |
| 2       |         | Dodecanoic acid, methyl ester       | 214 C13H26O2     | 000111-82-0 86 |
| 3       |         | Hexadecanoic acid, methyl ester     | 270 C17H34O2     | 000112-39-0 86 |
| 4       |         | Decanoic acid, methyl ester         | 186 C11H22O2     | 000110-42-9 80 |
| 5       |         | Pentadecanoic acid, 14-methyl-, ... | 270 C17H34O2     | 005129-60-2 72 |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

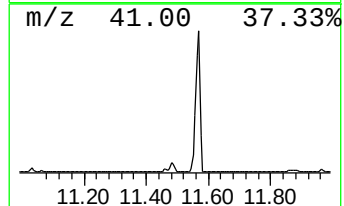
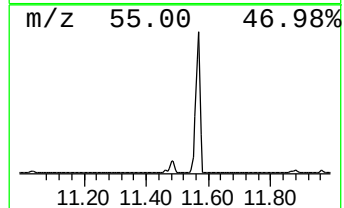
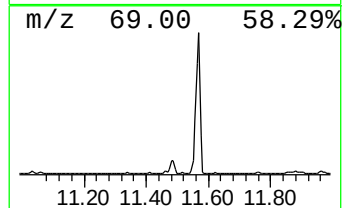
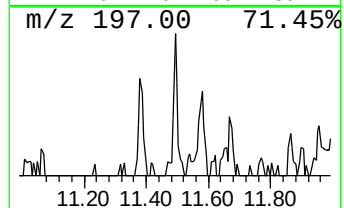
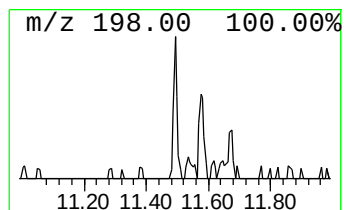
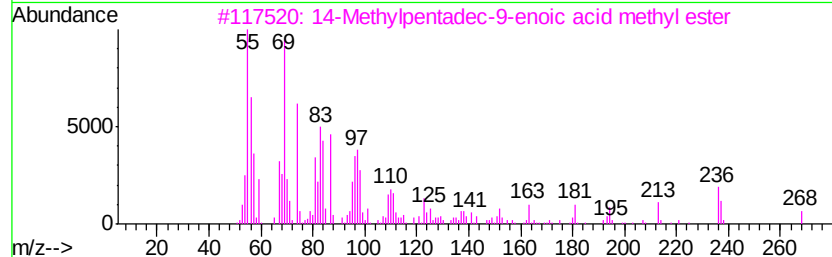
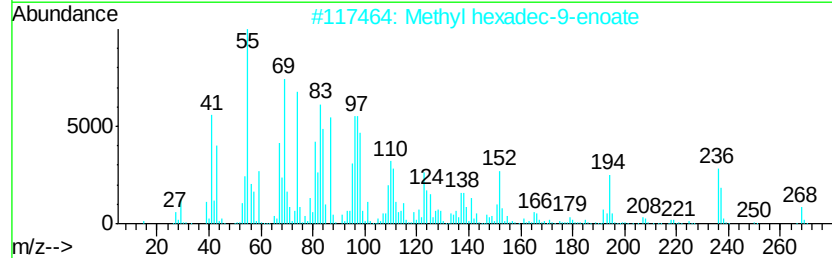
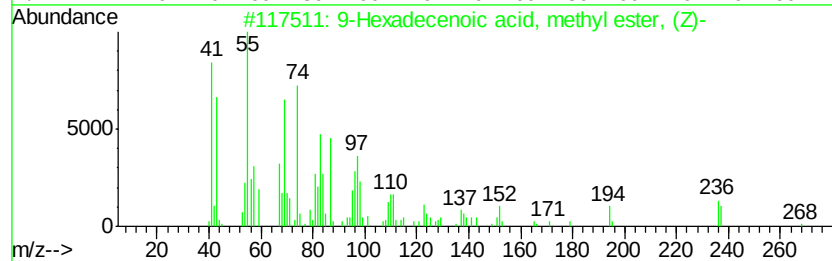
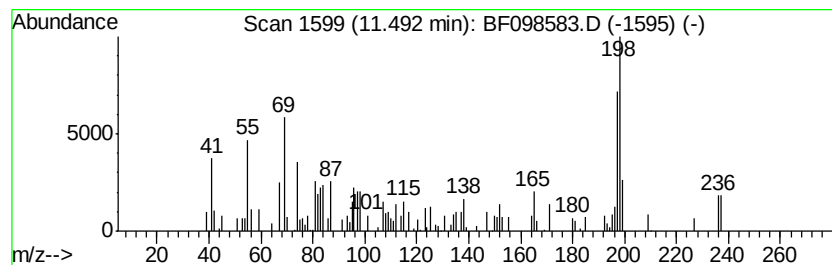
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ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 9-Hexadecenoic acid, methyl... Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 11.49     | 5.63 ng                             | 290987 | Phenanthrene-d10 | 11.16        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 9-Hexadecenoic acid, methyl este... | 268    | C17H32O2         | 001120-25-8  | 98   |
| 2         | Methyl hexadec-9-enoate             | 268    | C17H32O2         | 010030-74-7  | 90   |
| 3         | 14-Methylpentadec-9-enoic acid m... | 268    | C17H32O2         | 1000365-89-7 | 83   |
| 4         | Methyl myristoleate                 | 240    | C15H28O2         | 056219-06-8  | 76   |
| 5         | 7-Hexadecenoic acid, methyl este... | 268    | C17H32O2         | 056875-67-3  | 68   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

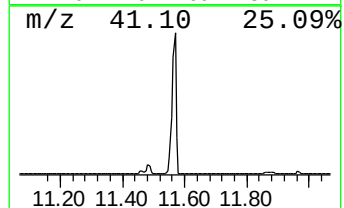
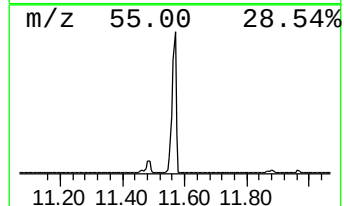
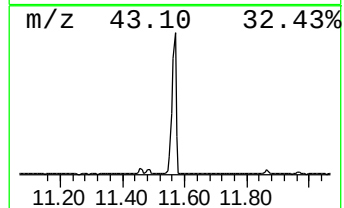
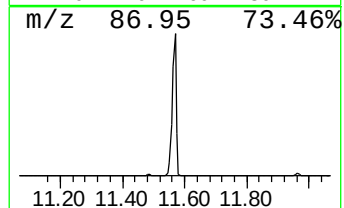
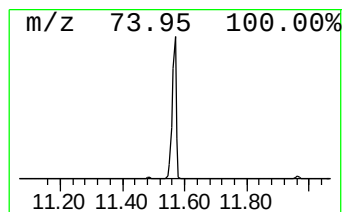
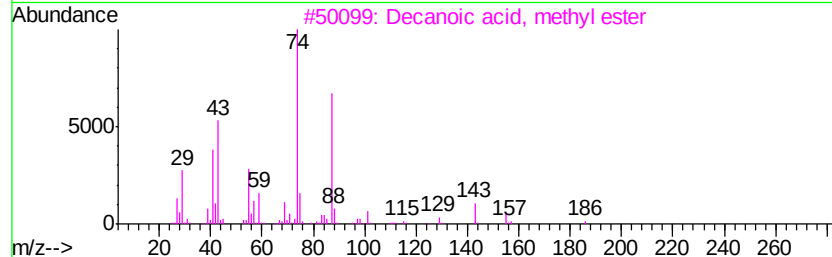
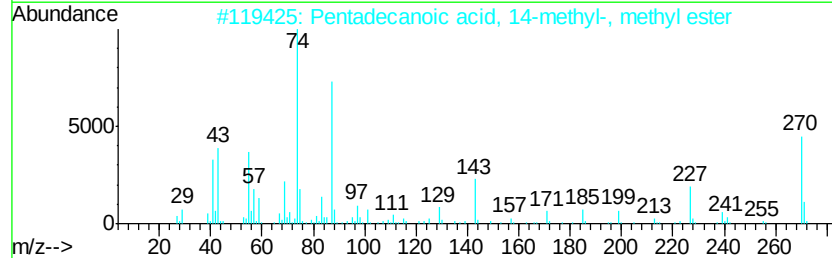
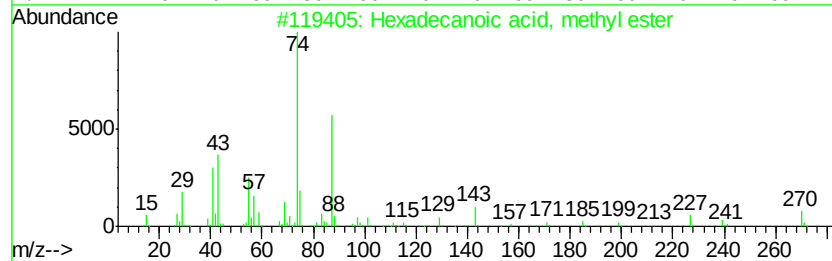
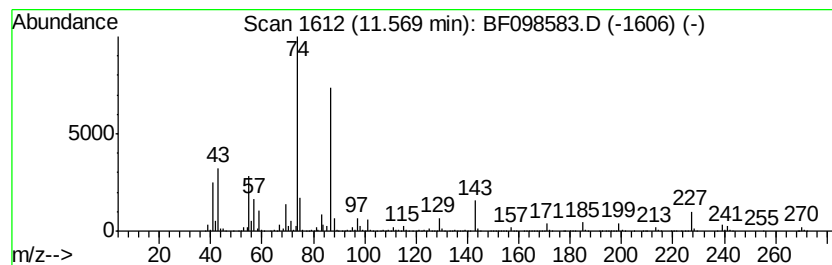
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 3

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 11.57 | 107.05 ng | 5533940 | Phenanthrene-d10 | 11.16 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 91   |
| 2    |    | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 90   |
| 3    |    | Decanoic acid, methyl ester         | 186 | C11H22O2 | 000110-42-9 | 87   |
| 4    |    | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 86   |
| 5    |    | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

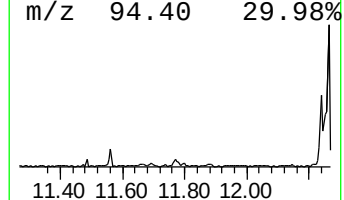
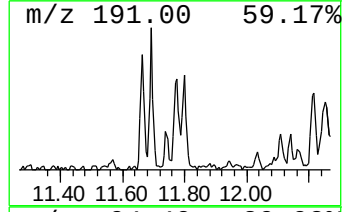
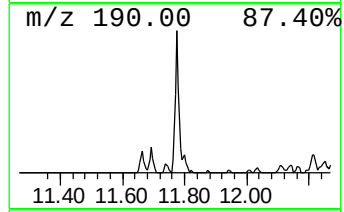
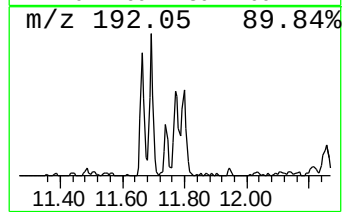
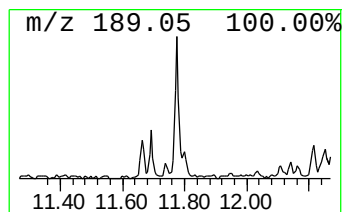
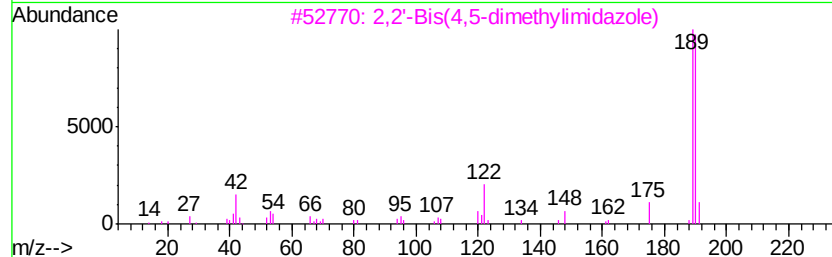
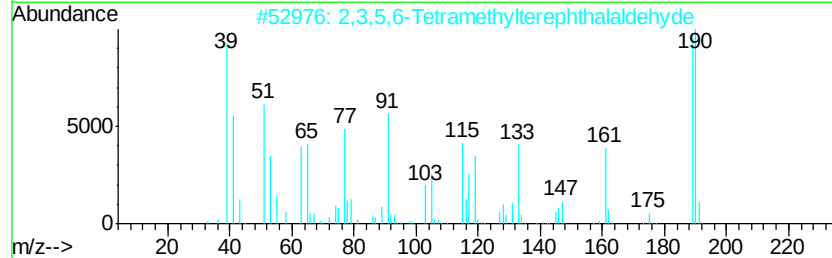
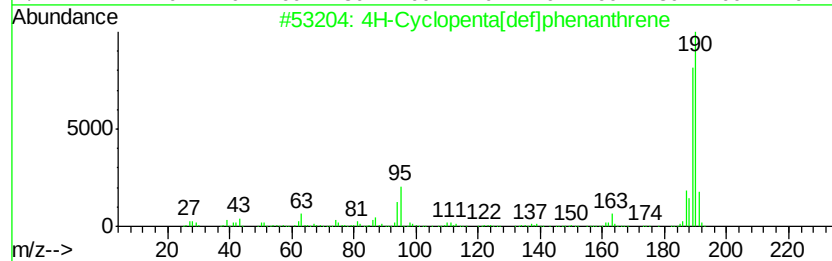
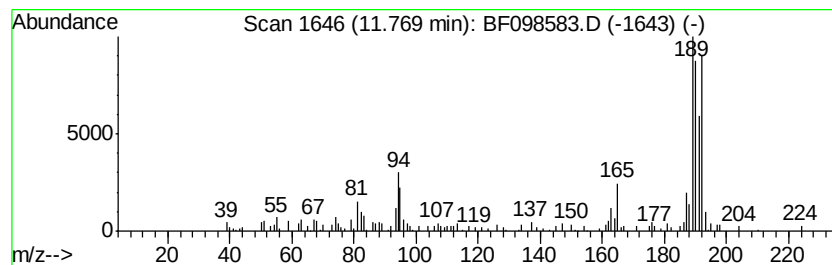
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ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 16

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.      |             |      |
|---------|---------|-------------------------------------|------------------|-----------|-------------|------|
| 11.77   | 2.37 ng | 122297                              | Phenanthrene-d10 | 11.16     |             |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm   | CAS#        | Qual |
| 1       |         | 4H-Cyclopenta[def]phenanthrene      | 190              | C15H10    | 000203-64-5 | 86   |
| 2       |         | 2,3,5,6-Tetramethylterephthalald... | 190              | C12H14O2  | 007072-01-7 | 49   |
| 3       |         | 2,2'-Bis(4,5-dimethylimidazole)     | 190              | C10H14N4  | 069286-06-2 | 47   |
| 4       |         | 6H-Cyclobuta[jk]phenanthrene        | 190              | C15H10    | 083469-43-6 | 47   |
| 5       |         | Pyrazole-4-carboxaldehyde, 3-(4-... | 190              | C10H7FN2O | 306936-57-2 | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

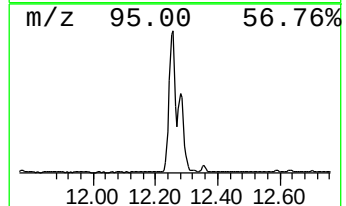
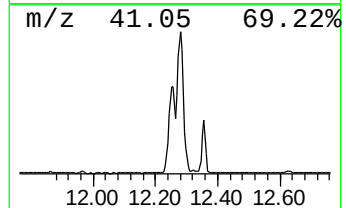
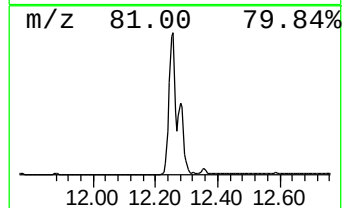
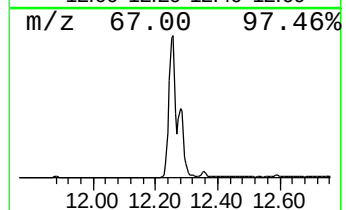
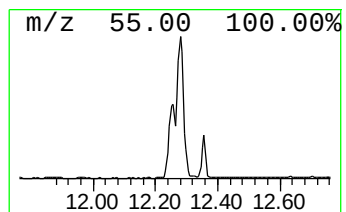
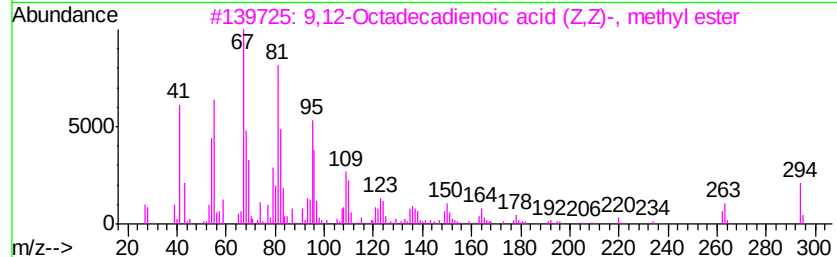
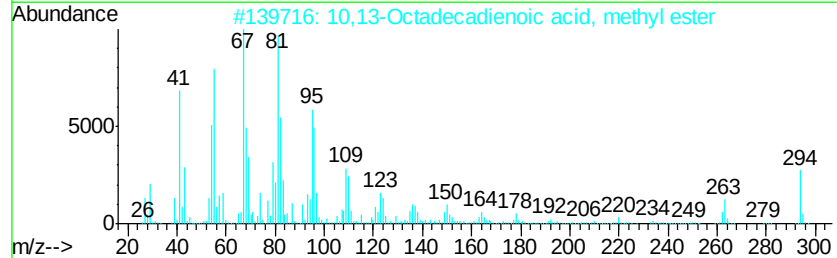
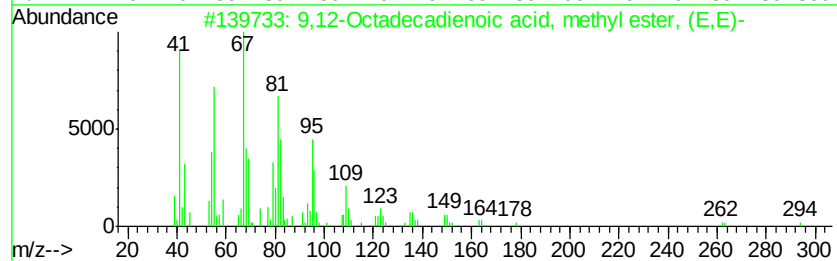
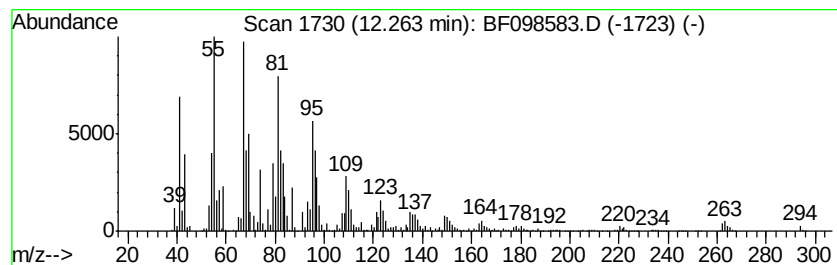
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 2

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.26 | 146.47 ng | 7571670 | Phenanthrene-d10 | 11.16 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4  | 99   |
| 2    |    | 10,13-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-62-2  | 99   |
| 3    |    | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0  | 99   |
| 4    |    | 9,15-Octadecadienoic acid, methy... | 294 | C19H34O2 | 017309-05-6  | 99   |
| 5    |    | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 1000336-44-0 | 99   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

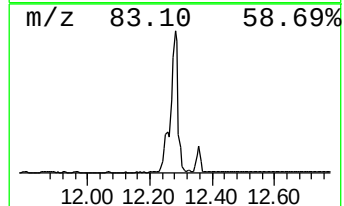
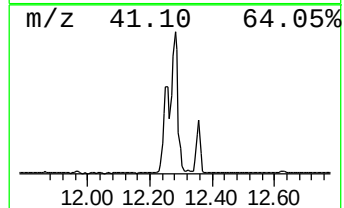
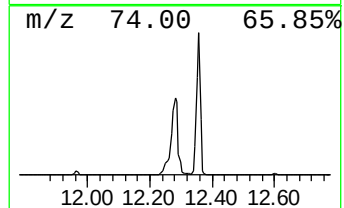
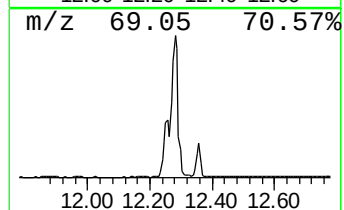
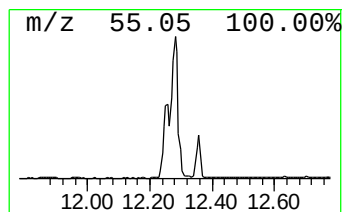
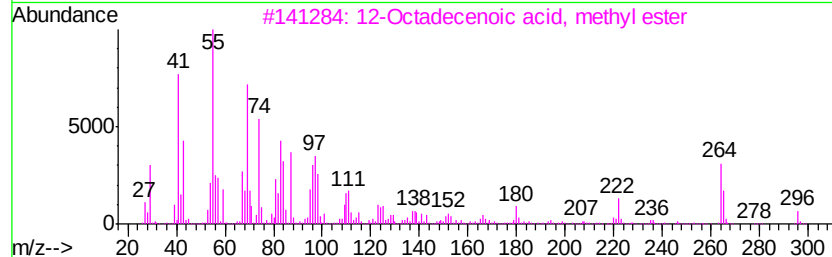
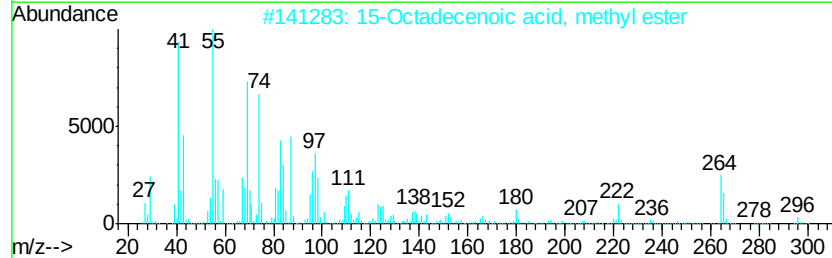
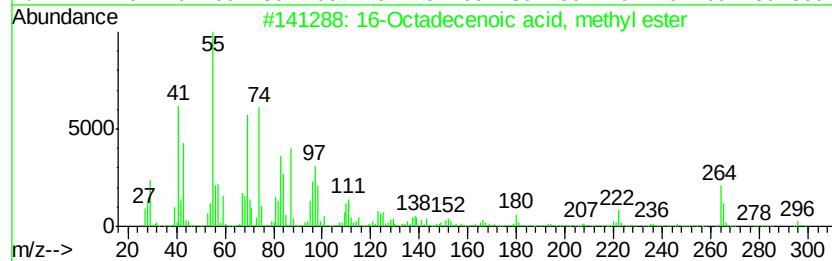
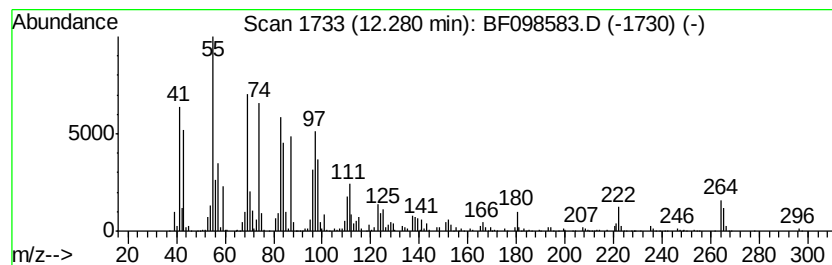
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 16-Octadecenoic acid, methyl ester Concentration Rank 1

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T.  |
|-------|-----------|----------|------------------|-------|
| 12.28 | 223.14 ng | 11535000 | Phenanthrene-d10 | 11.16 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 16-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-49-5 | 96   |
| 2    |    | 15-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 004764-72-1 | 96   |
| 3    |    | 12-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-46-2 | 96   |
| 4    |    | 13-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-47-3 | 96   |
| 5    |    | 14-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-48-4 | 94   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

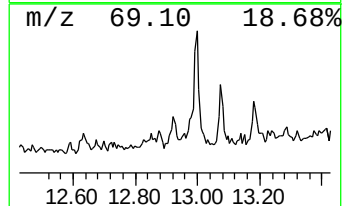
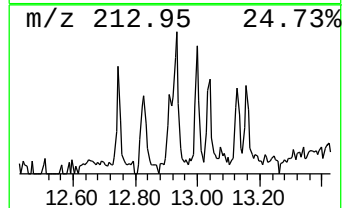
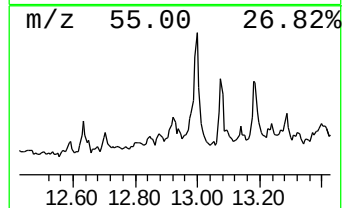
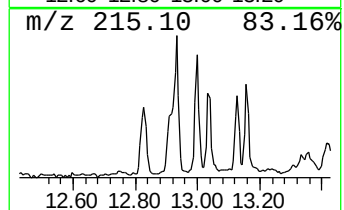
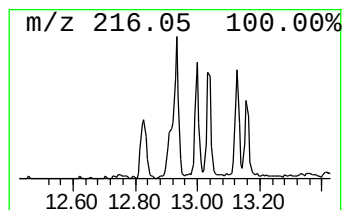
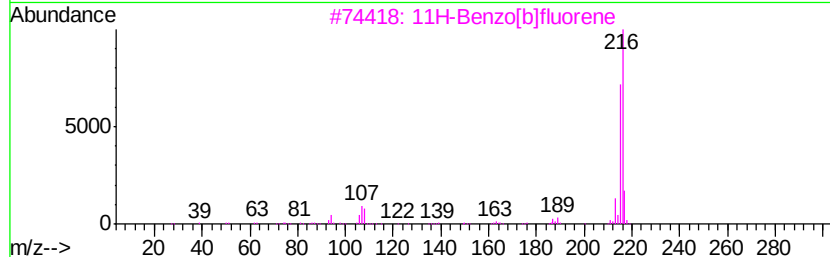
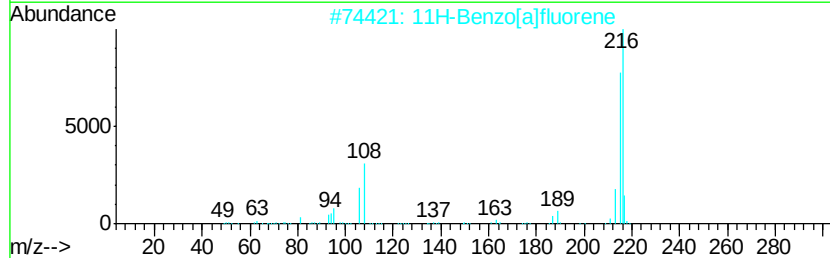
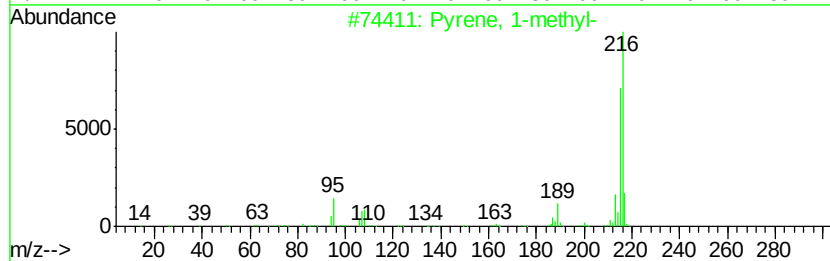
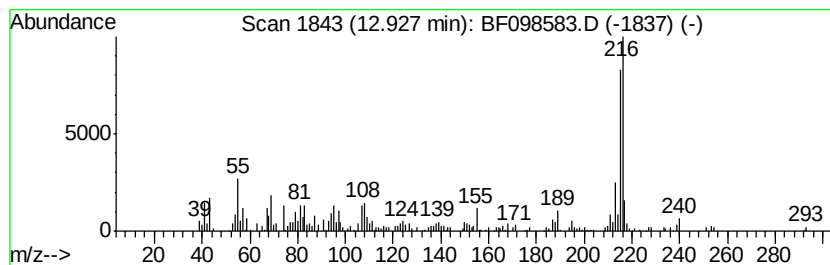
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BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 Pyrene, 1-methyl- Concentration Rank 9

| R.T.      | EstConc                 | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------|--------|------------------|-------------|-------|
| 12.93     | 5.15 ng                 | 266541 | Chrysene-d12     |             | 13.80 |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual  |
| 1         | Pyrene, 1-methyl-       | 216    | C17H12           | 002381-21-7 | 93    |
| 2         | 11H-Benzo[a]fluorene    | 216    | C17H12           | 000238-84-6 | 91    |
| 3         | 11H-Benzo[b]fluorene    | 216    | C17H12           | 000243-17-4 | 83    |
| 4         | Fluoranthene, 2-methyl- | 216    | C17H12           | 033543-31-6 | 81    |
| 5         | 7H-Benzo[c]fluorene     | 216    | C17H12           | 000205-12-9 | 81    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

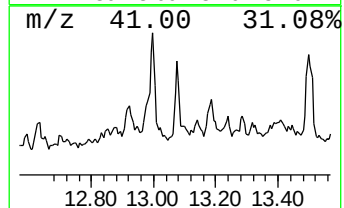
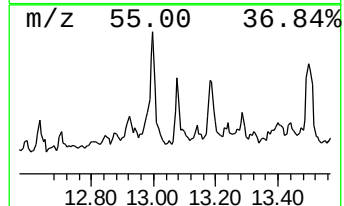
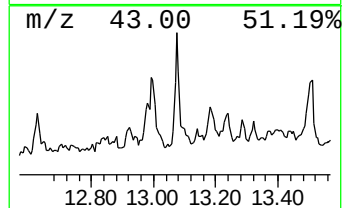
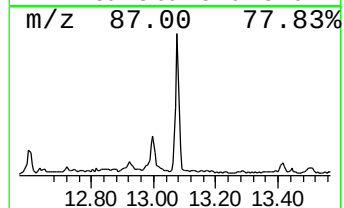
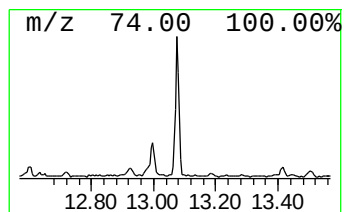
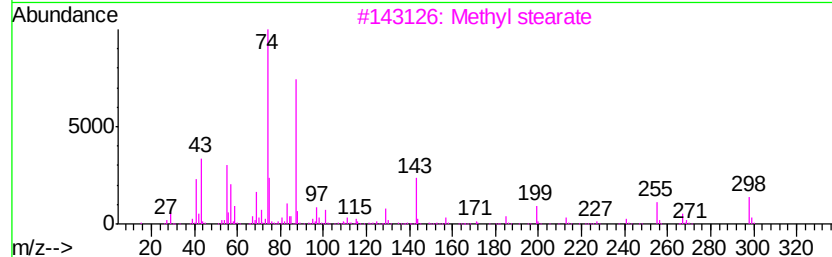
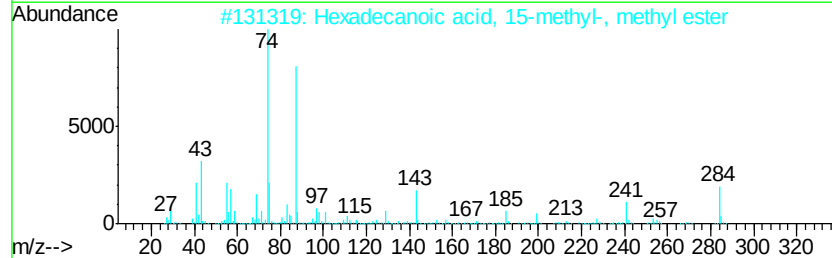
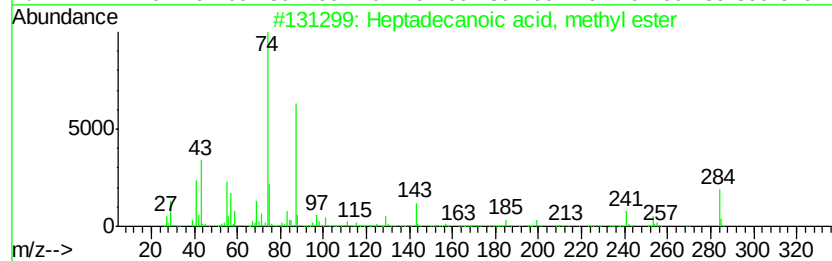
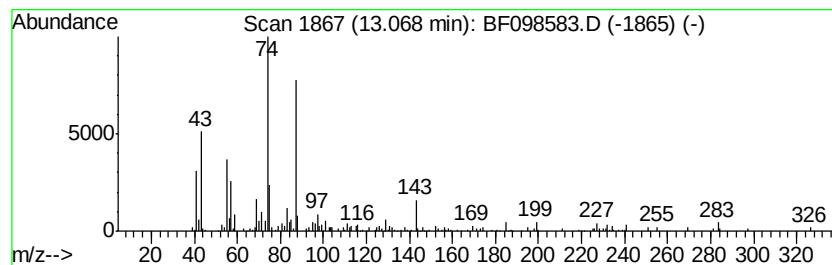
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Heptadecanoic acid, methyl ... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.07 | 4.46 ng | 231068 | Chrysene-d12     | 13.80 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecanoic acid, methyl ester    | 284 | C18H36O2 | 001731-92-6 | 94   |
| 2    |    | Hexadecanoic acid, 15-methyl-, m... | 284 | C18H36O2 | 006929-04-0 | 93   |
| 3    |    | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 91   |
| 4    |    | Nonadecanoic acid, methyl ester     | 312 | C20H40O2 | 001731-94-8 | 91   |
| 5    |    | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

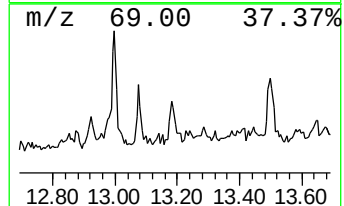
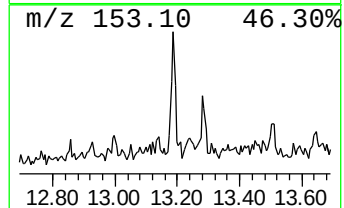
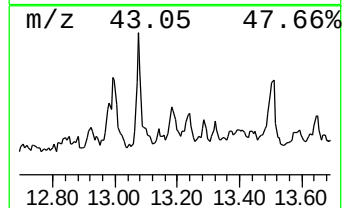
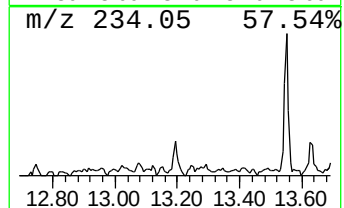
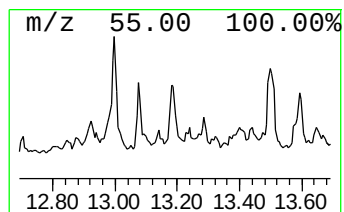
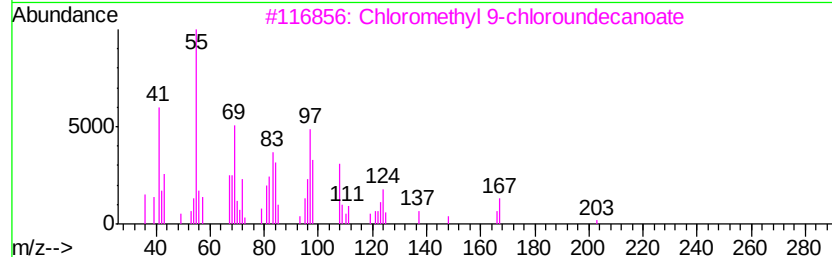
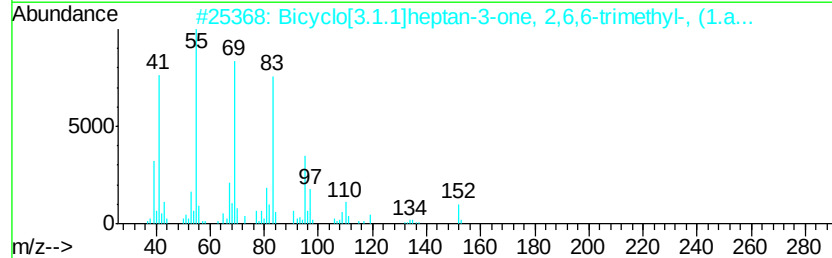
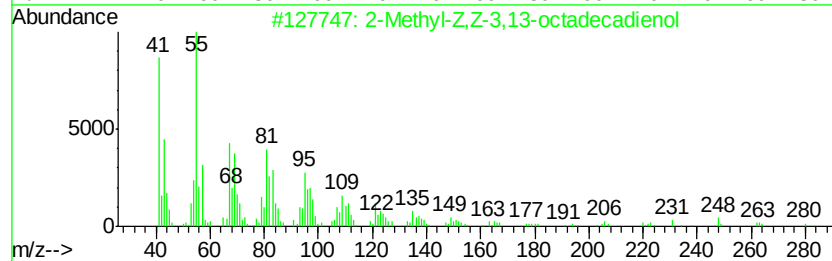
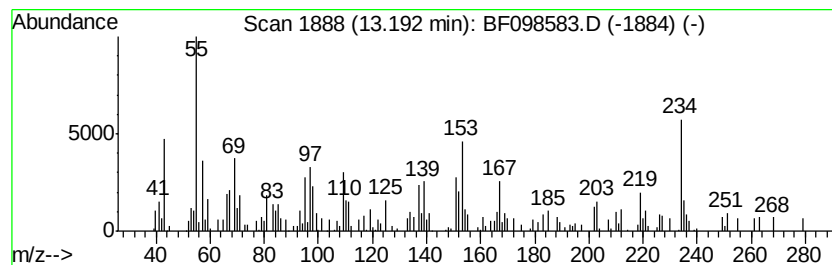
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 unknown13.19 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.19 | 2.94 ng | 152449 | Chrysene-d12     | 13.80 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 2-Methyl-Z,Z-3,13-octadecadienol    | 280 | C19H36O     | 1000130-90-5 | 38   |
| 2    |    |   | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O     | 015358-88-0  | 25   |
| 3    |    |   | Chloromethyl 9-chloroundecanoate    | 268 | C12H22Cl2O2 | 080418-95-7  | 14   |
| 4    |    |   | 9-Tetradecenal, (Z)-                | 210 | C14H26O     | 053939-27-8  | 12   |
| 5    |    |   | 10-Methyldodec-2-en-4-olide         | 210 | C13H22O2    | 1000370-41-0 | 10   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

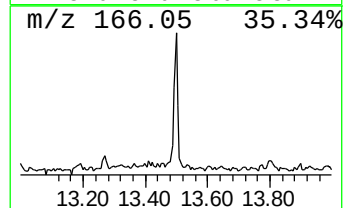
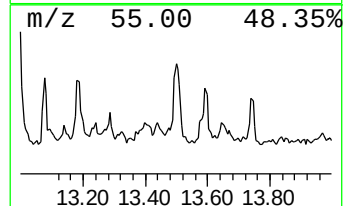
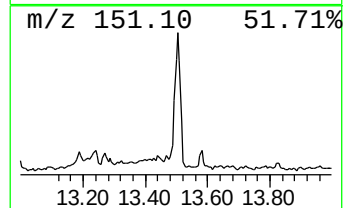
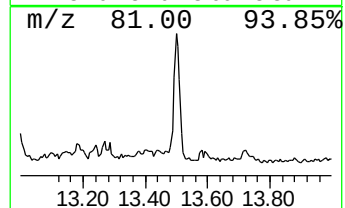
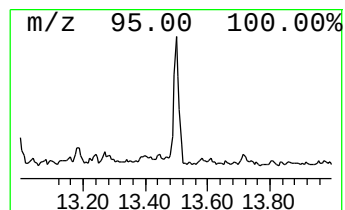
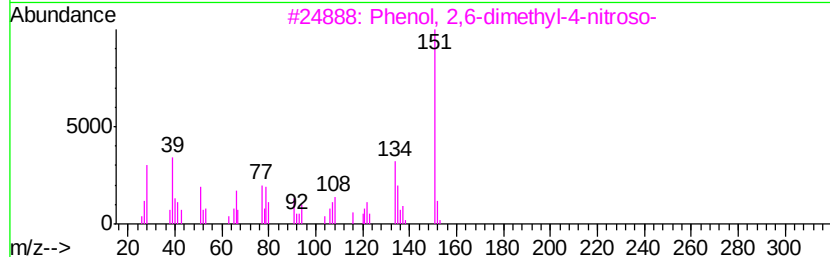
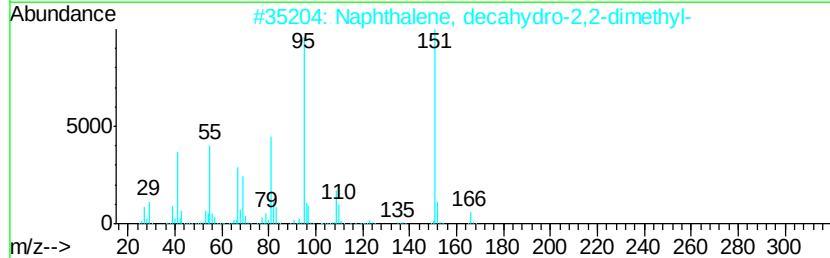
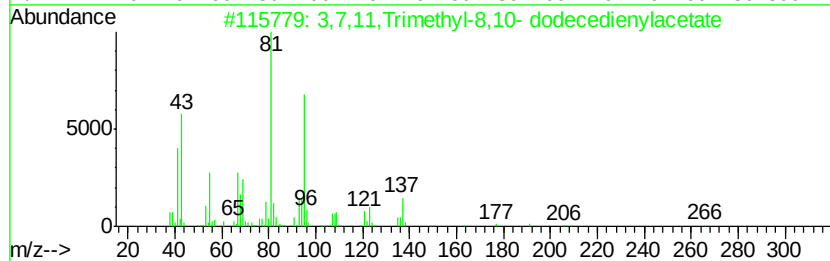
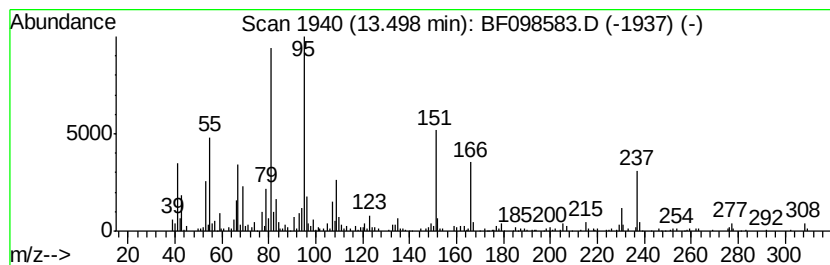
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ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 3,7,11,Trimethyl-8,10- dode... Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.50     | 9.04 ng                             | 468338 | Chrysene-d12     | 13.80        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 3,7,11,Trimethyl-8,10- dodecedie... | 266    | C17H30O2         | 1000374-10-3 | 55   |
| 2         | Naphthalene, decahydro-2,2-dimet... | 166    | C12H22           | 031124-85-3  | 27   |
| 3         | Phenol, 2,6-dimethyl-4-nitroso-     | 151    | C8H9NO2          | 013331-93-6  | 25   |
| 4         | 2,8-Bornanediol, stereoisomer       | 170    | C10H18O2         | 013429-50-0  | 22   |
| 5         | 4-Isoxazolecarboxamide, 5-amino-... | 275    | C13H13N3O4       | 1000320-20-3 | 14   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
Data File : BF098583.D  
Acq On : 15 Sep 2017 21:57  
Operator : SJ/JU  
Sample : I5236-01 5X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

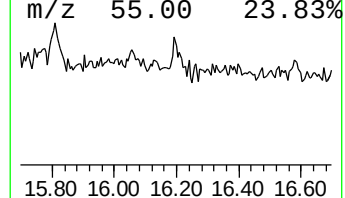
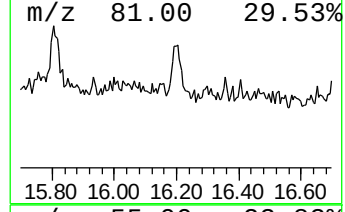
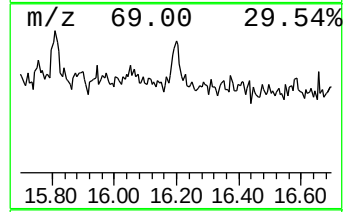
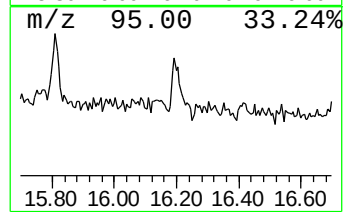
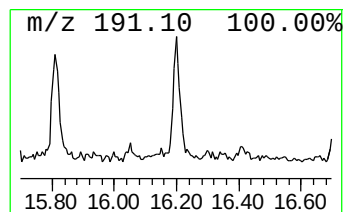
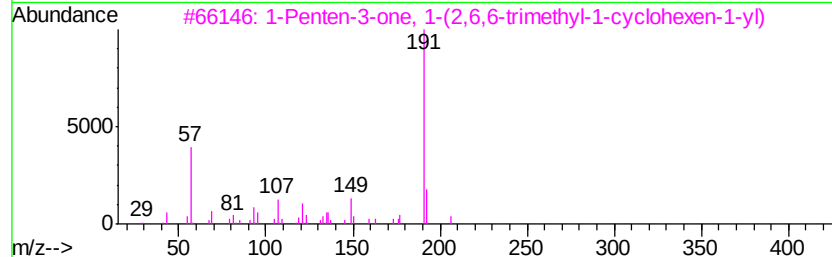
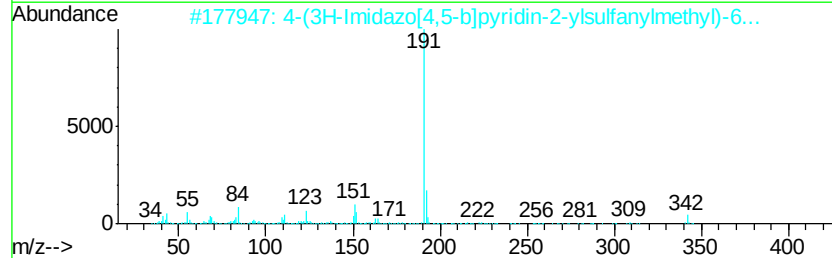
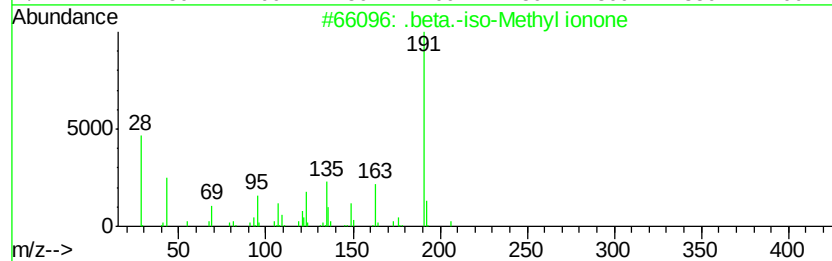
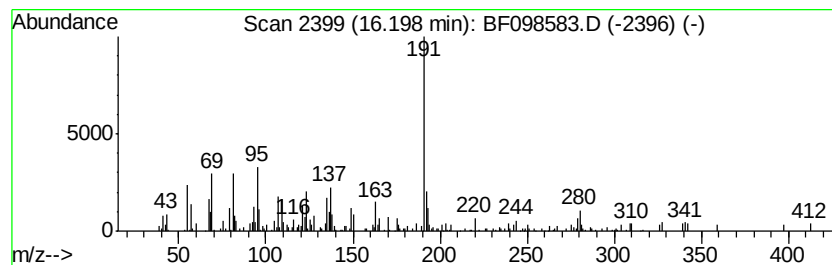
Instrument :  
BNA\_F  
ClientSampleId :  
TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 .beta.-iso-Methyl ionone Concentration Rank 12

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 16.20   | 3.41 ng | 149577                              | Perylene-d12     | 15.20           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | .beta.-iso-Methyl ionone            | 206 C14H22O      | 1000285-40-2 91 |
| 2       |         | 4-(3H-Imidazo[4,5-b]pyridin-2-yl... | 342 C15H18N8S    | 1000276-08-2 47 |
| 3       |         | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 C14H22O      | 000127-43-5 46  |
| 4       |         | 1-(9-Phenanthrenyl)-2-propanone     | 234 C17H14O      | 1000326-08-9 43 |
| 5       |         | Silane, [4-(1,1-dimethylethyl)ph... | 206 C13H22Si     | 018412-68-5 38  |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF091517\  
 Data File : BF098583.D  
 Acq On : 15 Sep 2017 21:57  
 Operator : SJ/JU  
 Sample : I5236-01 5X  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TR-04-091317

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF091117.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 |
| unknown6.42          | 6.42  | 17.6    | ng    | 717109   | 1                     | 6.64  | 817226  | 20.0 |
| Hexadecane           | 9.09  | 2.1     | ng    | 117484   | 3                     | 9.68  | 1126730 | 20.0 |
| Naphthalene, 2,6-... | 9.27  | 2.1     | ng    | 120891   | 3                     | 9.68  | 1126730 | 20.0 |
| Heptadecane          | 10.59 | 3.8     | ng    | 198749   | 4                     | 11.16 | 1033870 | 20.0 |
| Methyl tetradecan... | 10.69 | 2.9     | ng    | 149777   | 4                     | 11.16 | 1033870 | 20.0 |
| 9-Hexadecenoic ac... | 11.49 | 5.6     | ng    | 290987   | 4                     | 11.16 | 1033870 | 20.0 |
| Hexadecanoic acid... | 11.57 | 107.1   | ng    | 5533940  | 4                     | 11.16 | 1033870 | 20.0 |
| 4H-Cyclopenta[def... | 11.77 | 2.4     | ng    | 122297   | 4                     | 11.16 | 1033870 | 20.0 |
| 9,12-Octadecadien... | 12.26 | 146.5   | ng    | 7571670  | 4                     | 11.16 | 1033870 | 20.0 |
| 16-Octadecenoic a... | 12.28 | 223.1   | ng    | 11535000 | 4                     | 11.16 | 1033870 | 20.0 |
| Pyrene, 1-methyl-    | 12.93 | 5.2     | ng    | 266541   | 5                     | 13.80 | 1035870 | 20.0 |
| Heptadecanoic aci... | 13.07 | 4.5     | ng    | 231068   | 5                     | 13.80 | 1035870 | 20.0 |
| unknown13.19         | 13.19 | 2.9     | ng    | 152449   | 5                     | 13.80 | 1035870 | 20.0 |
| 3,7,11,Trimethyl-... | 13.50 | 9.0     | ng    | 468338   | 5                     | 13.80 | 1035870 | 20.0 |
| .beta.-iso-Methyl... | 16.20 | 3.4     | ng    | 149577   | 6                     | 15.20 | 878272  | 20.0 |