

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
 Data File : BF101172.D  
 Acq On : 6 Dec 2017 17:18  
 Operator : SJ/JU  
 Sample : I6683-03 10X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EO-01-120517-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-BF113017.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         | 5.457       | 570           | 573         | 585          | rBV      | 84674          | 89735         | 3.43%           | 1.120%        |
| 2         | 6.475       | 743           | 746         | 753          | rBV      | 85794          | 84776         | 3.24%           | 1.058%        |
| 3         | 6.616       | 767           | 770         | 776          | rBV      | 120111         | 102187        | 3.91%           | 1.275%        |
| 4         | 6.845       | 806           | 809         | 813          | rBV      | 331587         | 259818        | 9.93%           | 3.243%        |
| 5         | 6.998       | 832           | 835         | 842          | rVB      | 96667          | 81463         | 3.11%           | 1.017%        |
| 6         | 7.410       | 902           | 905         | 907          | rBV      | 58710          | 54069         | 2.07%           | 0.675%        |
| 7         | 8.128       | 1024          | 1027        | 1033         | rVB      | 406849         | 331529        | 12.67%          | 4.138%        |
| 8         | 9.198       | 1206          | 1209        | 1214         | rBV      | 143298         | 119231        | 4.56%           | 1.488%        |
| 9         | 9.275       | 1218          | 1222        | 1225         | rBV      | 36688          | 29083         | 1.11%           | 0.363%        |
| 10        | 9.592       | 1273          | 1276        | 1284         | rVB4     | 19586          | 32304         | 1.23%           | 0.403%        |
| 11        | 9.804       | 1305          | 1312        | 1317         | rBV      | 41650          | 38312         | 1.46%           | 0.478%        |
| 12        | 9.886       | 1322          | 1326        | 1329         | rVV      | 414460         | 347760        | 13.29%          | 4.341%        |
| 13        | 10.304      | 1391          | 1397        | 1400         | rBV2     | 44811          | 41492         | 1.59%           | 0.518%        |
| 14        | 10.675      | 1457          | 1460        | 1464         | rBV      | 55442          | 47364         | 1.81%           | 0.591%        |
| 15        | 10.775      | 1475          | 1477        | 1488         | rBV      | 39705          | 47239         | 1.81%           | 0.590%        |
| 16        | 11.375      | 1575          | 1579        | 1581         | rBV      | 381053         | 319138        | 12.20%          | 3.983%        |
| 17        | 11.398      | 1581          | 1583        | 1589         | rVB      | 69938          | 57937         | 2.21%           | 0.723%        |
| 18        | 11.757      | 1640          | 1644        | 1647         | rBV      | 843282         | 640143        | 24.46%          | 7.990%        |
| 19        | 12.451      | 1754          | 1762        | 1764         | rBV      | 2341595        | 2616683       | 100.00%         | 32.660%       |
| 20        | 12.469      | 1764          | 1765        | 1770         | rVB2     | 1743483        | 1182456       | 45.19%          | 14.759%       |
| 21        | 12.545      | 1775          | 1778        | 1782         | rVB      | 270804         | 238425        | 9.11%           | 2.976%        |
| 22        | 12.592      | 1782          | 1786        | 1789         | rBV      | 72948          | 73765         | 2.82%           | 0.921%        |
| 23        | 12.822      | 1822          | 1825        | 1831         | rVB      | 112983         | 105003        | 4.01%           | 1.311%        |
| 24        | 12.957      | 1844          | 1848        | 1851         | rBV      | 109885         | 86560         | 3.31%           | 1.080%        |
| 25        | 13.033      | 1856          | 1861        | 1869         | rBV5     | 14025          | 35178         | 1.34%           | 0.439%        |
| 26        | 13.145      | 1872          | 1880        | 1883         | rBV4     | 18628          | 33128         | 1.27%           | 0.413%        |
| 27        | 14.021      | 2024          | 2029        | 2031         | rBV2     | 303146         | 322323        | 12.32%          | 4.023%        |
| 28        | 14.045      | 2031          | 2033        | 2036         | rVB      | 44377          | 35161         | 1.34%           | 0.439%        |
| 29        | 14.404      | 2090          | 2094        | 2098         | rBV2     | 21759          | 30070         | 1.15%           | 0.375%        |
| 30        | 14.898      | 2173          | 2178        | 2182         | rVB      | 27636          | 40241         | 1.54%           | 0.502%        |
| 31        | 15.063      | 2203          | 2206        | 2209         | rBV      | 35311          | 50469         | 1.93%           | 0.630%        |
| 32        | 15.368      | 2255          | 2258        | 2262         | rVB      | 28678          | 35228         | 1.35%           | 0.440%        |
| 33        | 15.433      | 2265          | 2269        | 2275         | rBV      | 45956          | 53764         | 2.05%           | 0.671%        |
| 34        | 15.498      | 2275          | 2280        | 2284         | rBV      | 216579         | 287422        | 10.98%          | 3.587%        |

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EO-01-120517-B

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-BF113017.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |       |       |       |       |        |
|----|--------|------|------|------|-------|-------|-------|-------|--------|
| 35 | 16.604 | 2465 | 2468 | 2478 | rVB10 | 12538 | 27619 | 1.06% | 0.345% |
| 36 | 17.427 | 2604 | 2608 | 2615 | rVB2  | 17362 | 34844 | 1.33% | 0.435% |

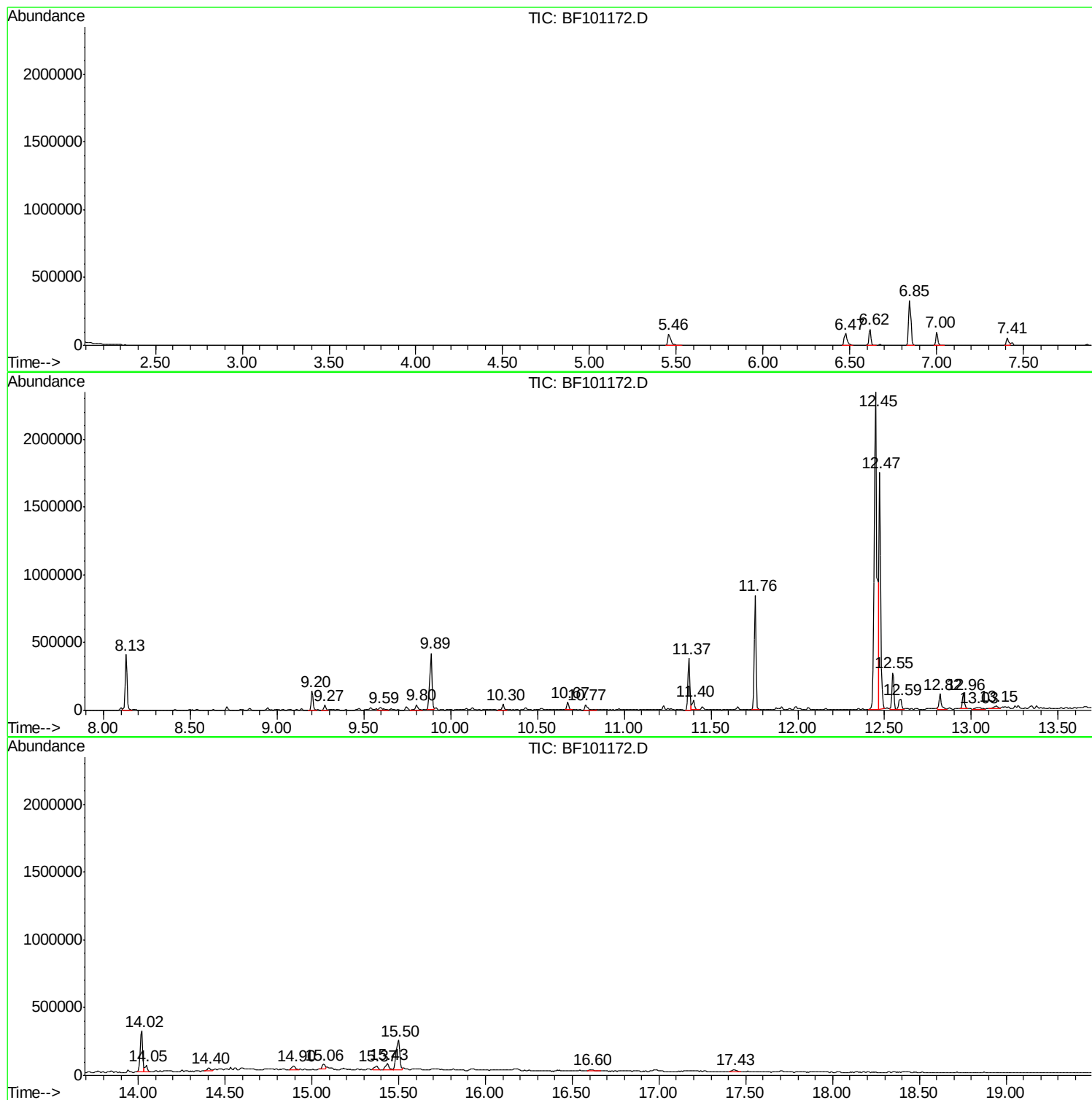
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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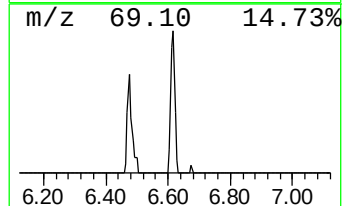
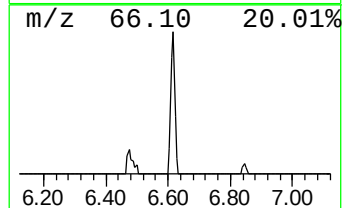
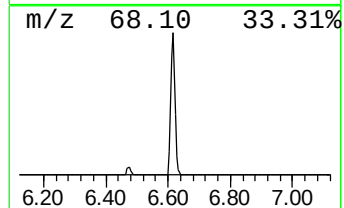
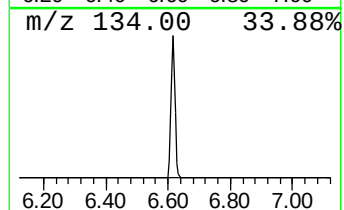
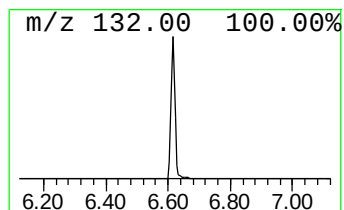
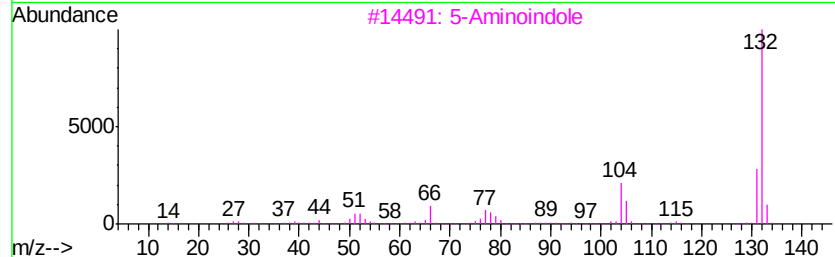
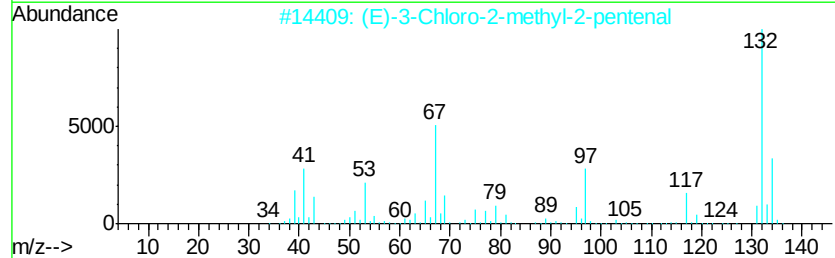
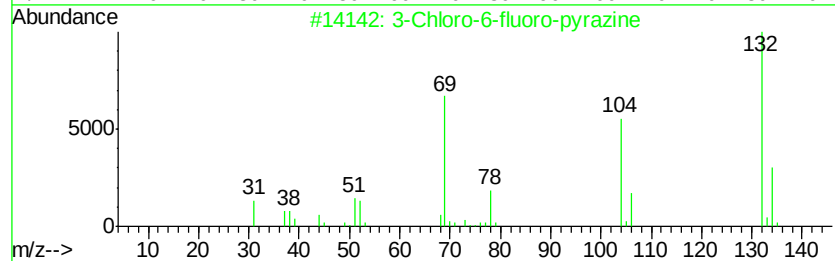
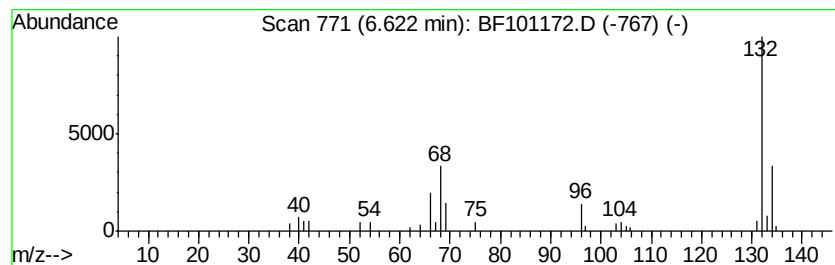
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown6.62 Concentration Rank 4

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.62 | 7.87 ng | 102187 | 1,4-Dichlorobenzene-d4 | 6.85 |

| 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  | 17   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | 4-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-60-2  | 9    |
| 5    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 9    |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

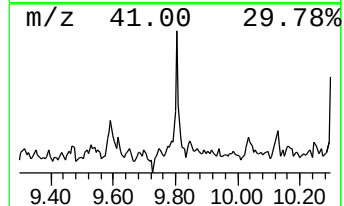
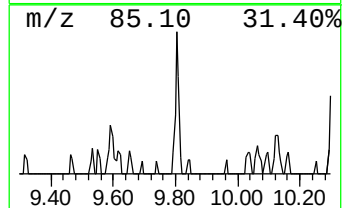
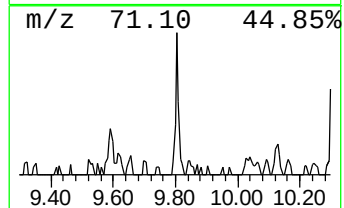
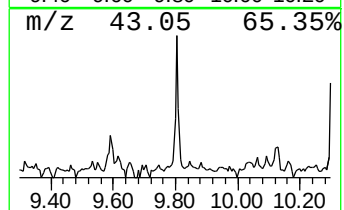
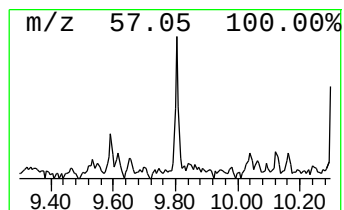
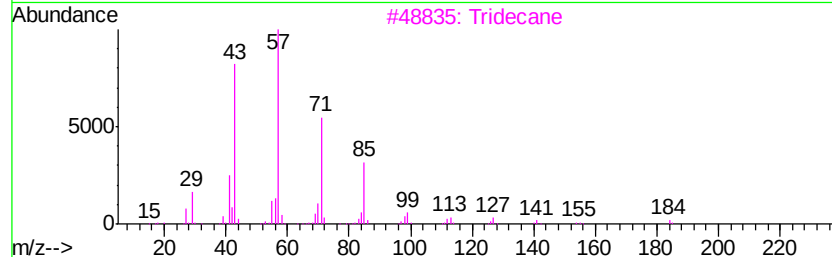
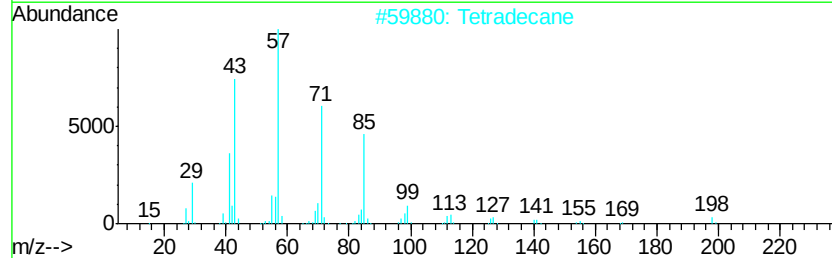
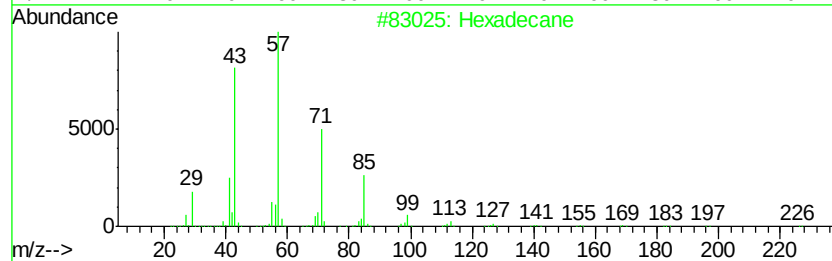
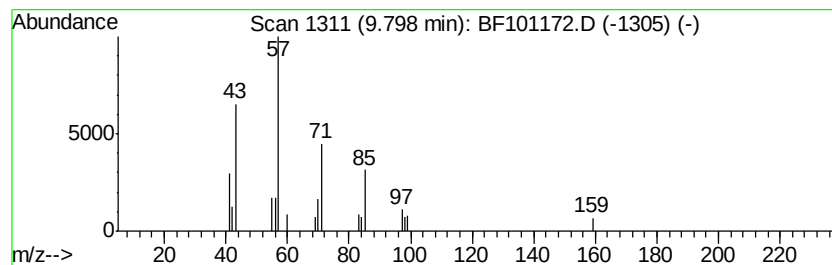
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Hexadecane Concentration Rank 10

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 9.80 | 2.20 ng | 38312 | Acenaphthene-d10 | 9.89 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane        | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    | Tetradecane       | 198 | C14H30  | 000629-59-4 | 86   |
| 3    |    | Tridecane         | 184 | C13H28  | 000629-50-5 | 78   |
| 4    |    | Decane, 2-methyl- | 156 | C11H24  | 006975-98-0 | 64   |
| 5    |    | Pentadecane       | 212 | C15H32  | 000629-62-9 | 59   |



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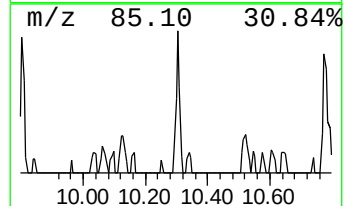
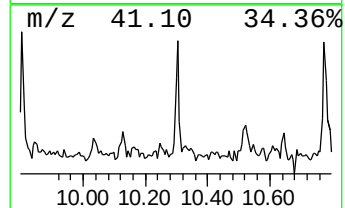
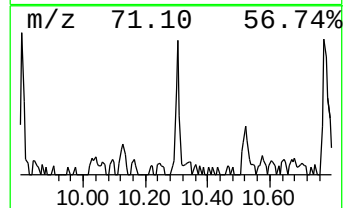
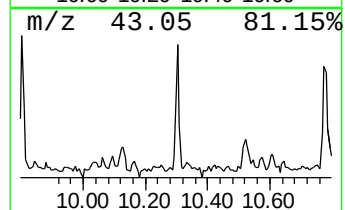
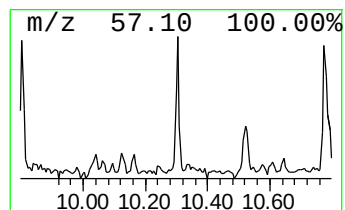
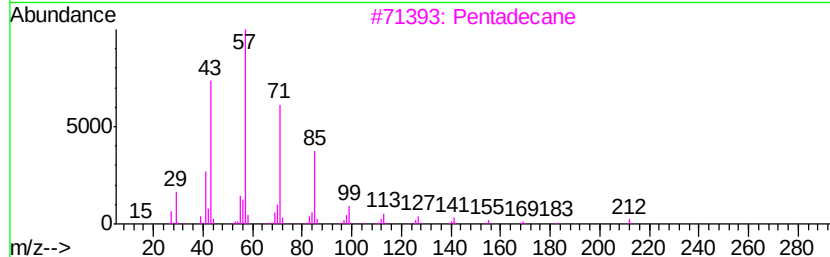
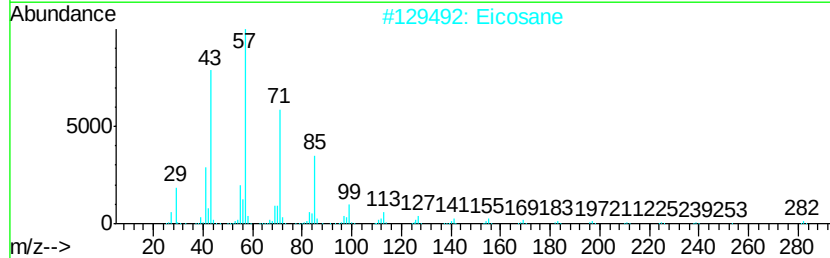
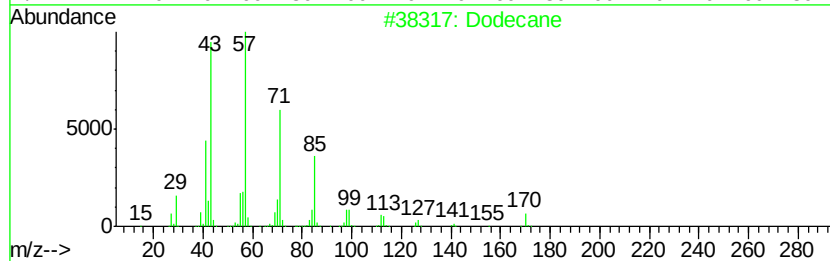
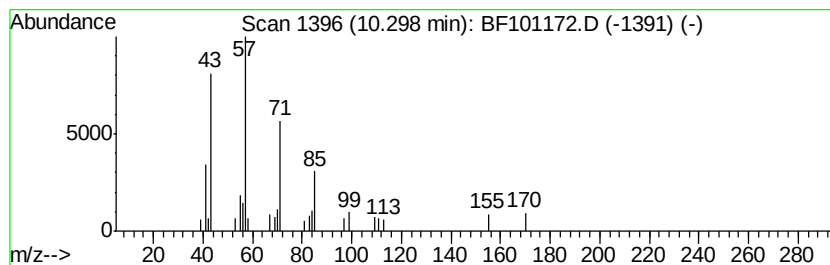
Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Dodecane Concentration Rank 9

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 10.30   | 2.39 ng                             | 41492        | Acenaphthene-d10 | 9.89      |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Dodecane                            | 170 C12H26   | 000112-40-3      | 90        |
| 2       | Eicosane                            | 282 C20H42   | 000112-95-8      | 86        |
| 3       | Pentadecane                         | 212 C15H32   | 000629-62-9      | 80        |
| 4       | Tetracosane                         | 338 C24H50   | 000646-31-1      | 80        |
| 5       | Heptadecane, 2,6,10,15-tetramethyl- | 296 C21H44   | 054833-48-6      | 78        |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

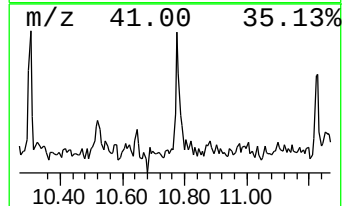
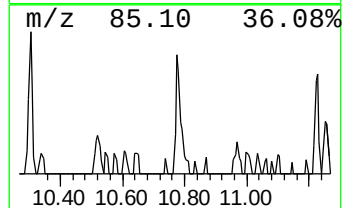
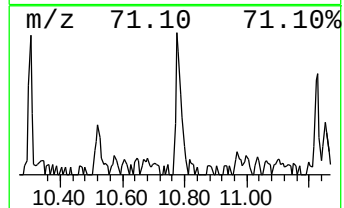
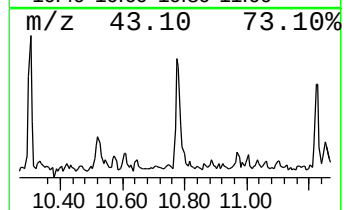
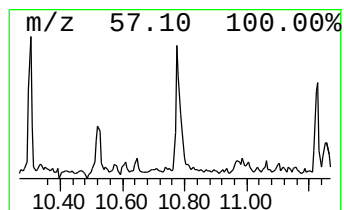
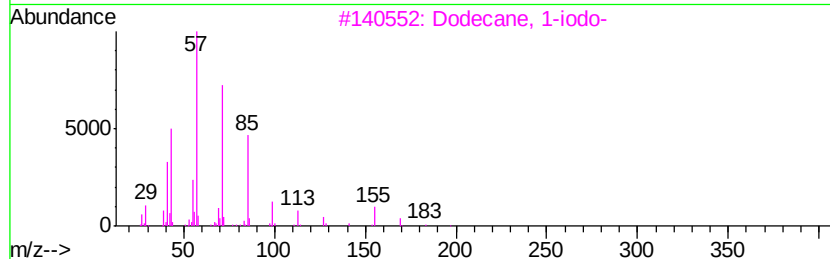
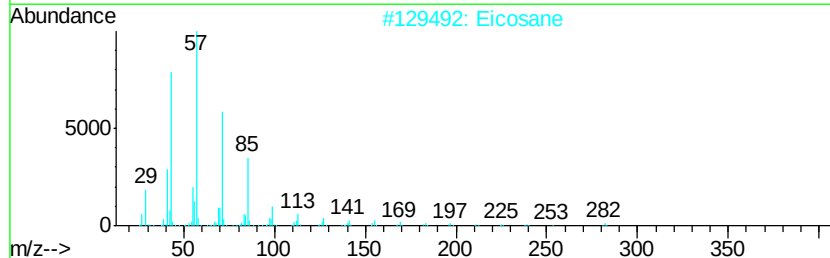
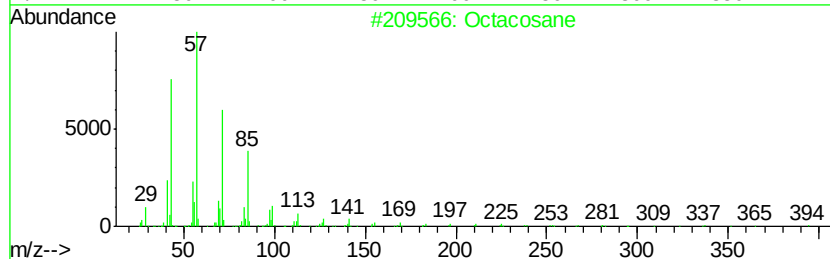
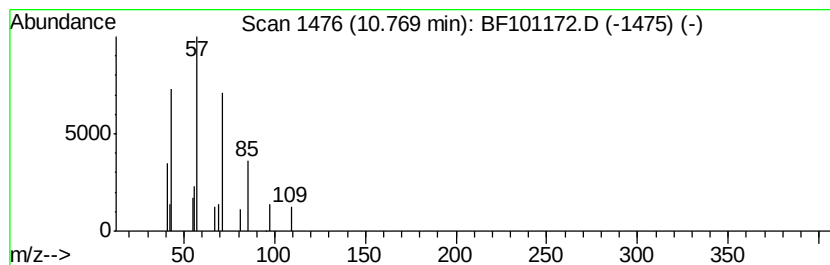
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Octacosane Concentration Rank 6

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 10.77 | 2.96 ng | 47239 | Phenanthrene-d10 | 11.37 |

| Hit# of | 5                        | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|--------------------------|--------------|-----|---------|-------------|------|
| 1       | Octacosane               |              | 394 | C28H58  | 000630-02-4 | 83   |
| 2       | Eicosane                 |              | 282 | C20H42  | 000112-95-8 | 72   |
| 3       | Dodecane, 1-iodo-        |              | 296 | C12H25I | 004292-19-7 | 72   |
| 4       | Tridecane                |              | 184 | C13H28  | 000629-50-5 | 72   |
| 5       | Decane, 2,3,5-trimethyl- |              | 184 | C13H28  | 062238-11-3 | 64   |



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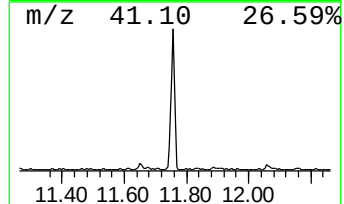
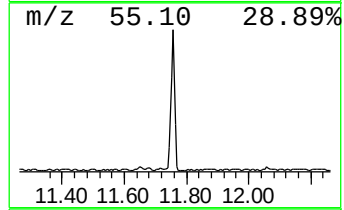
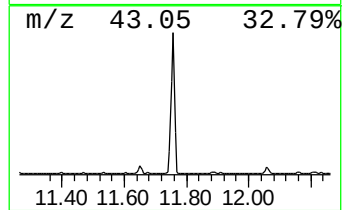
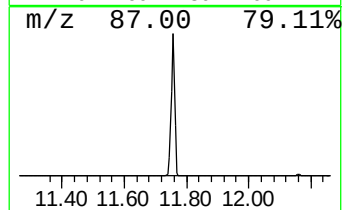
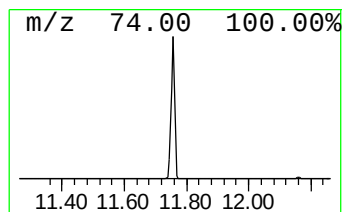
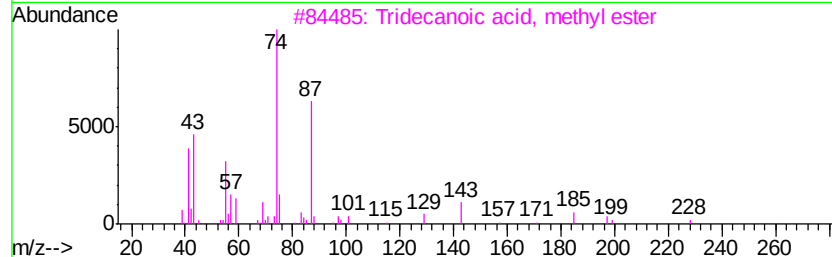
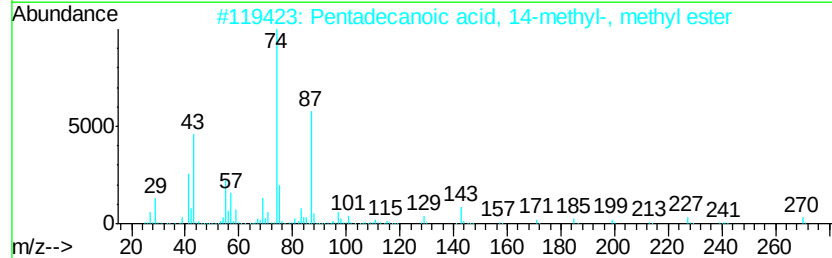
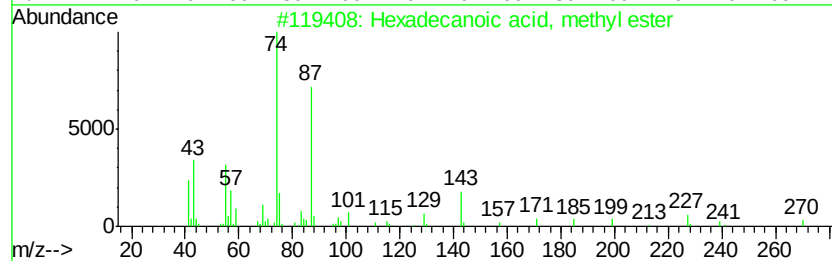
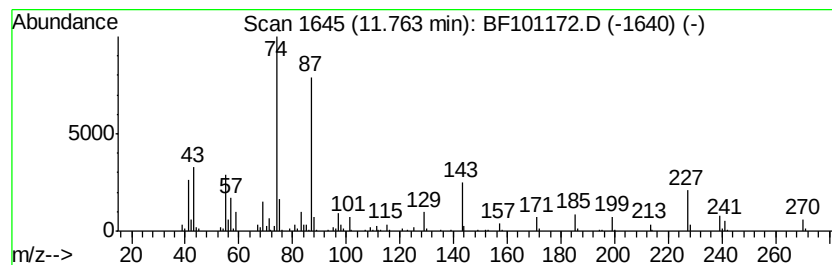
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.76 | 40.12 ng | 640143 | Phenanthrene-d10 | 11.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2    |    | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 92   |
| 3    |    | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 90   |
| 4    |    | Decanoic acid, methyl ester         | 186 | C11H22O2 | 000110-42-9 | 87   |
| 5    |    | Hexadecanoic acid, 15-methyl-, m... | 284 | C18H36O2 | 006929-04-0 | 83   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

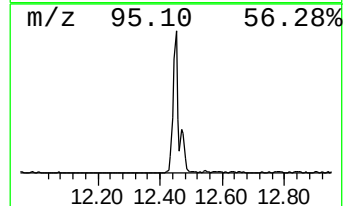
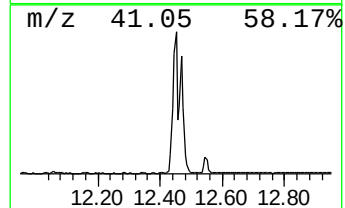
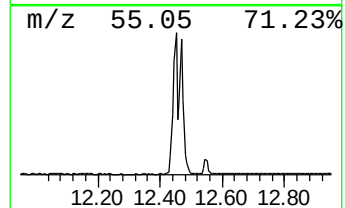
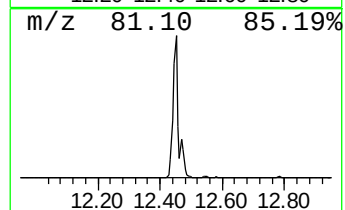
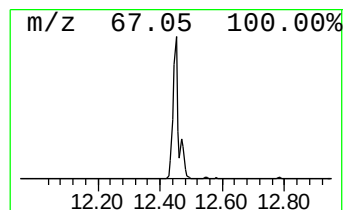
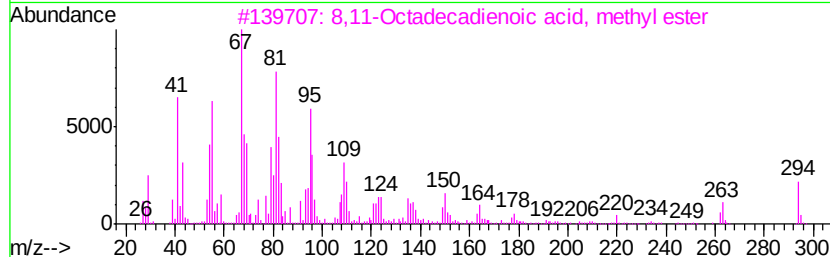
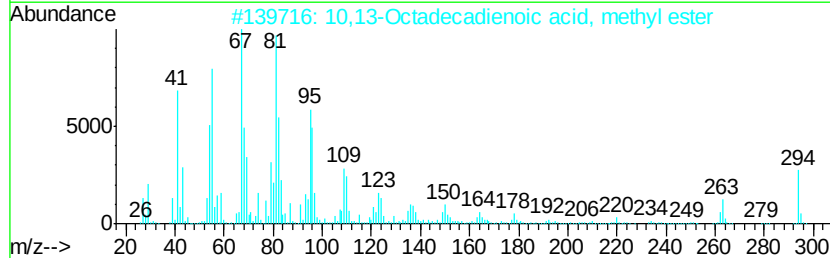
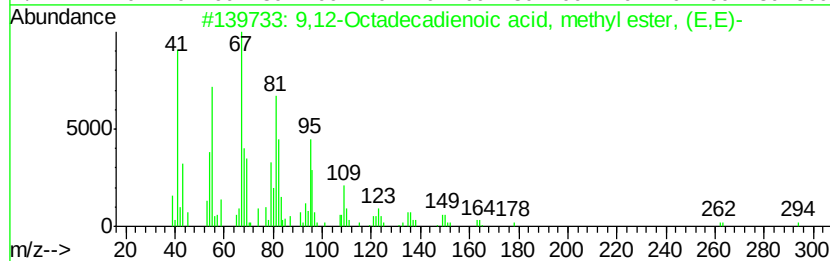
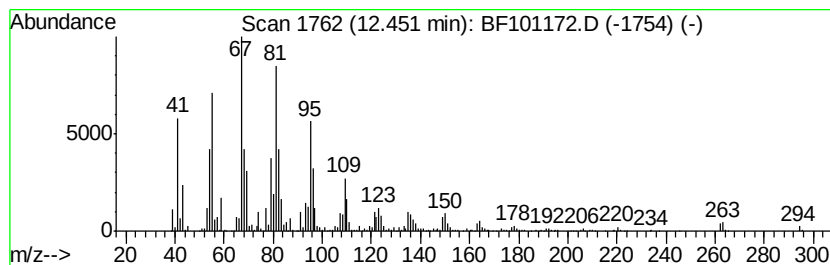
Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.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 9,12-Octadecadienoic acid, ... Concentration Rank 1

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|---------|-------------------------------------|--------------|------------------|-------------|------|------|
| 12.45   | 163.98 ng                           | 2616680      | Phenanthrene-d10 | 11.37       |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1       | 9,12-Octadecadienoic acid, methy... | 294          | C19H34O2         | 002566-97-4 | 99   |      |
| 2       | 10,13-Octadecadienoic acid, meth... | 294          | C19H34O2         | 056554-62-2 | 99   |      |
| 3       | 8,11-Octadecadienoic acid, methy... | 294          | C19H34O2         | 056599-58-7 | 99   |      |
| 4       | 9,15-Octadecadienoic acid, methy... | 294          | C19H34O2         | 017309-05-6 | 99   |      |
| 5       | 9,12-Octadecadienoic acid (Z,Z)-... | 294          | C19H34O2         | 000112-63-0 | 99   |      |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

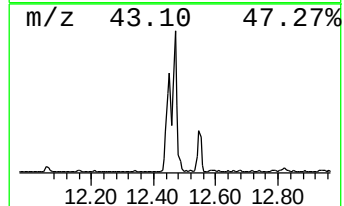
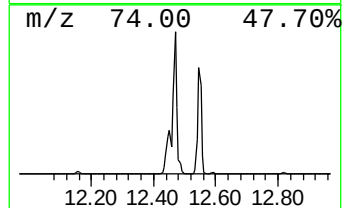
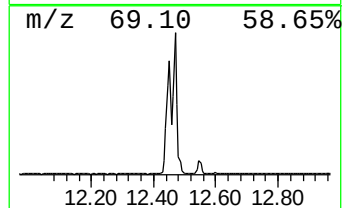
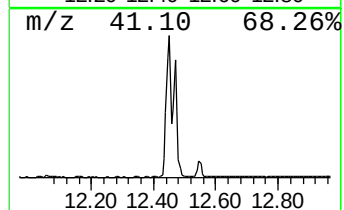
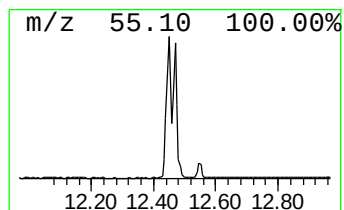
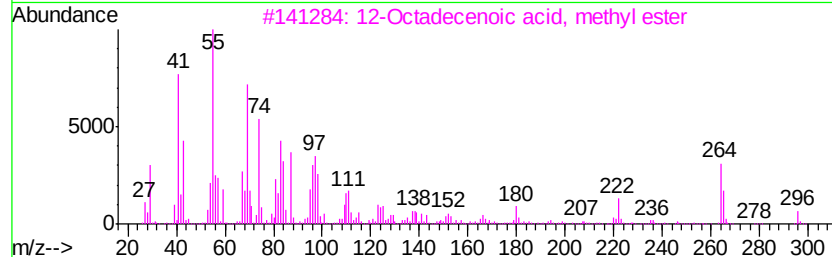
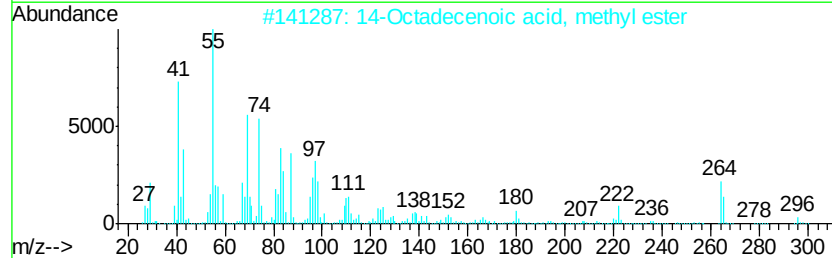
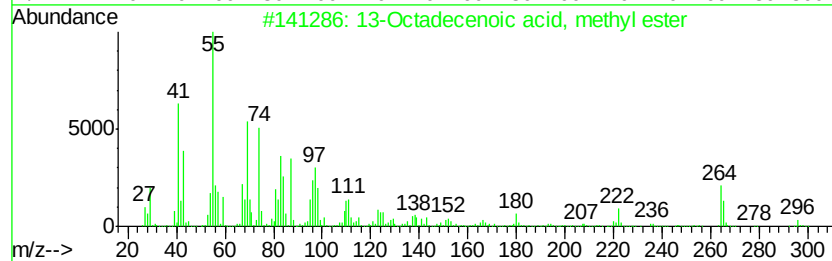
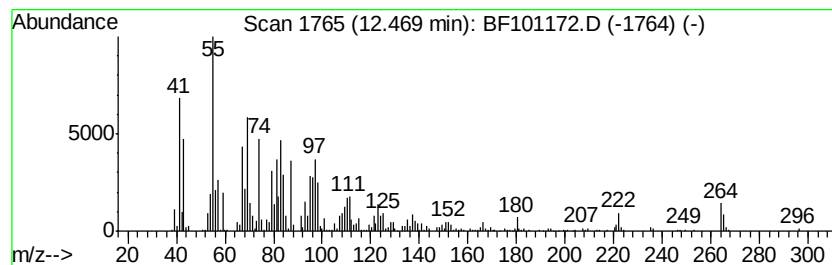
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 13-Octadecenoic acid, methyl... Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.47 | 74.10 ng | 1182460 | Phenanthrene-d10 | 11.37 |

| Hit# | of | Tentative ID                             | MW  | MolForm  | CAS#         | Qual |
|------|----|------------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 13-Octadecenoic acid, methyl ester       | 296 | C19H36O2 | 056554-47-3  | 99   |
| 2    |    | 14-Octadecenoic acid, methyl ester       | 296 | C19H36O2 | 056554-48-4  | 99   |
| 3    |    | 12-Octadecenoic acid, methyl ester       | 296 | C19H36O2 | 056554-46-2  | 96   |
| 4    |    | 15-Octadecenoic acid, methyl ester       | 296 | C19H36O2 | 004764-72-1  | 94   |
| 5    |    | trans-13-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 1000333-61-3 | 94   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

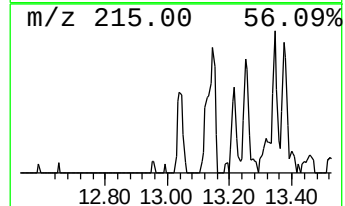
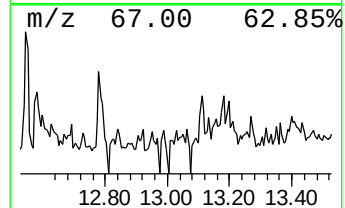
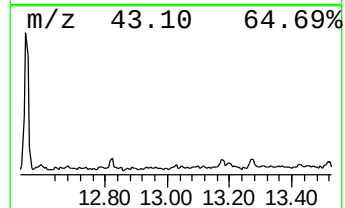
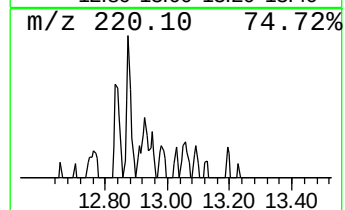
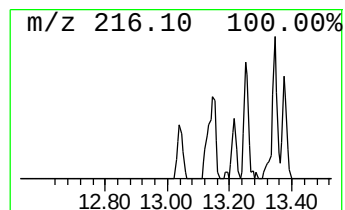
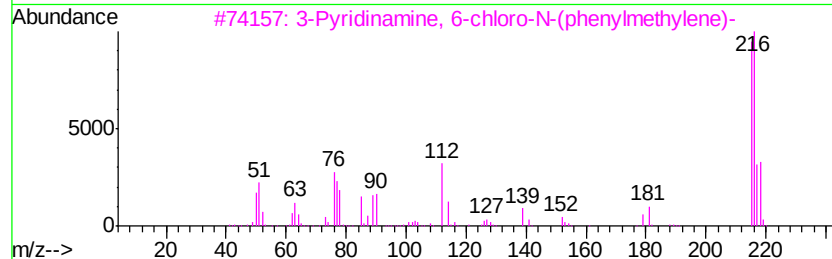
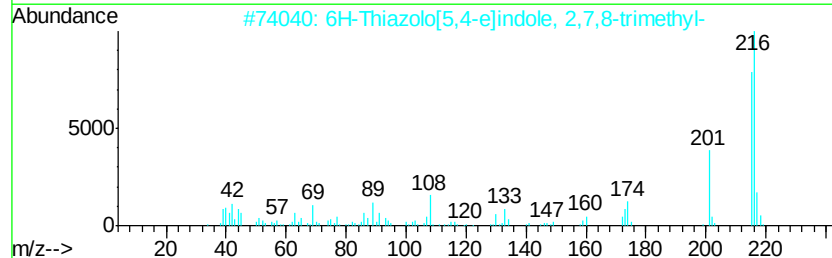
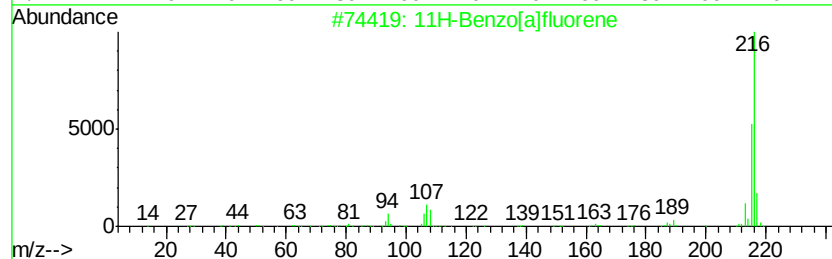
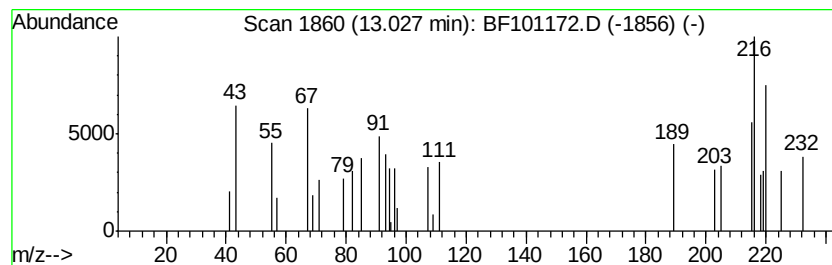
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ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 unknown13.03 Concentration Rank 11

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 13.03     | 2.18 ng                             | 35178 | Chrysene-d12     | 14.02        |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | 11H-Benzo[alfluorene                | 216   | C17H12           | 000238-84-6  | 49   |
| 2         | 6H-Thiazolo[5,4-elindole, 2,7,8-... | 216   | C12H12N2S        | 1000338-32-0 | 47   |
| 3         | 3-Pyridinamine, 6-chloro-N-(phen... | 216   | C12H9ClN2        | 005325-71-3  | 47   |
| 4         | 2H-Benzimidazol-2-one, 6,7-dichl... | 216   | C8H6Cl2N2O       | 077572-79-3  | 32   |
| 5         | 3,4-Dichlorocinnamic acid           | 216   | C9H6Cl2O2        | 001202-39-7  | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

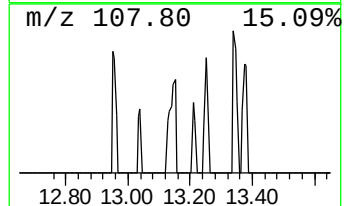
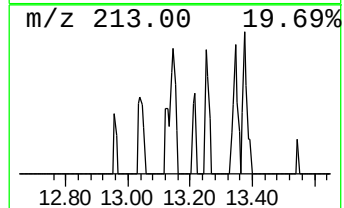
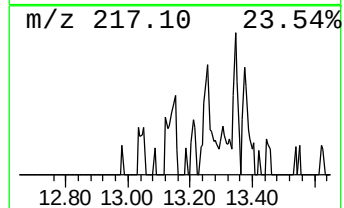
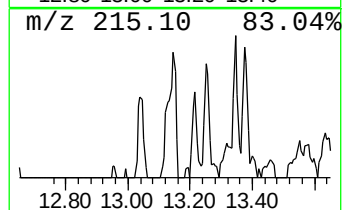
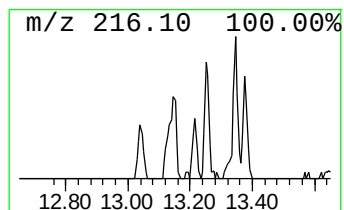
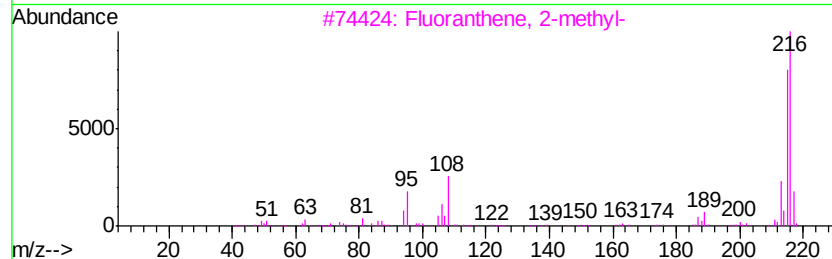
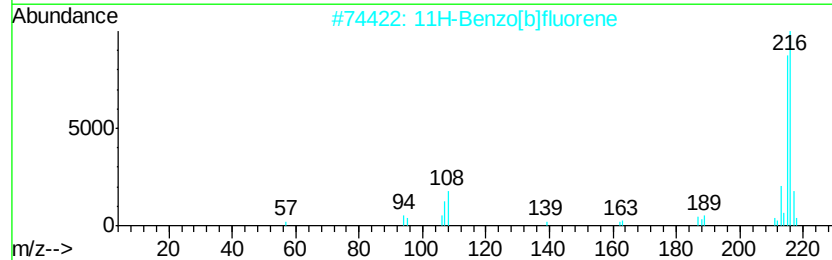
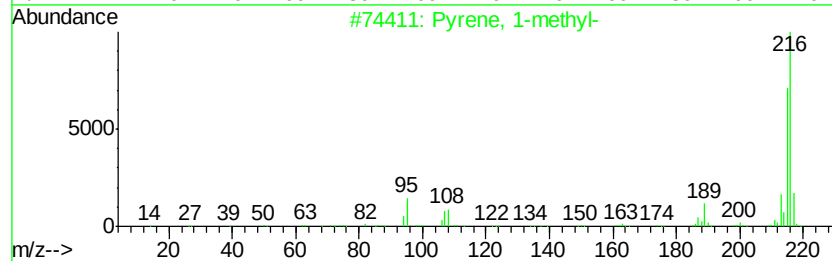
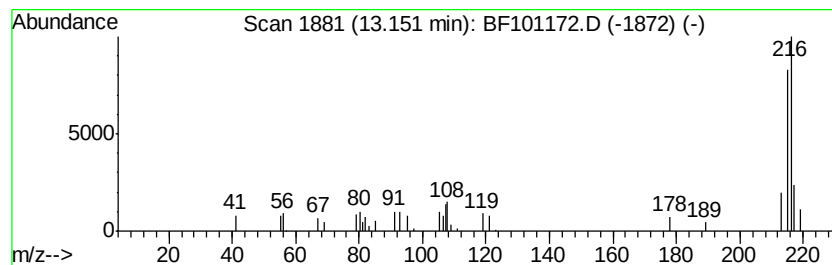
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Pyrene, 1-methyl- Concentration Rank 12

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.15 | 2.06 ng | 33128 | Chrysene-d12     | 14.02 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 74   |
| 2    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 70   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 62   |
| 4    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 58   |
| 5    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 58   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

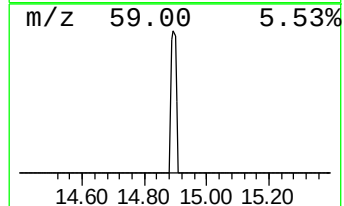
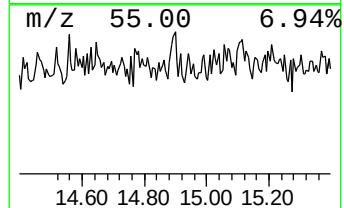
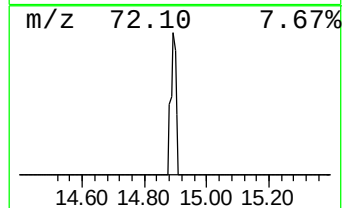
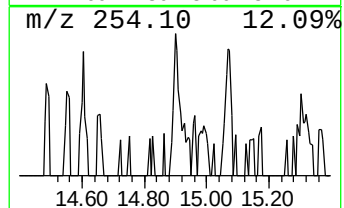
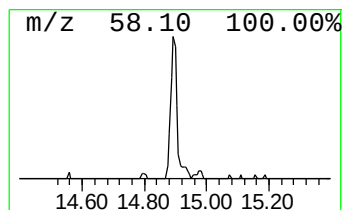
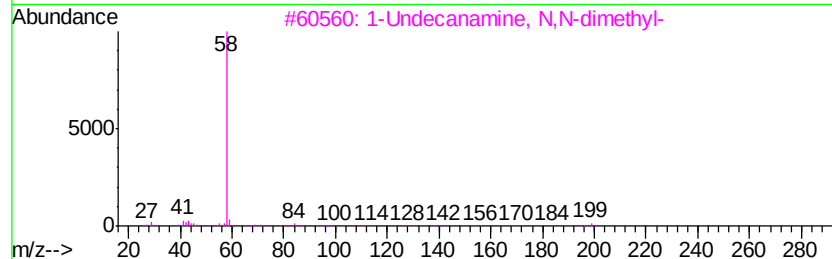
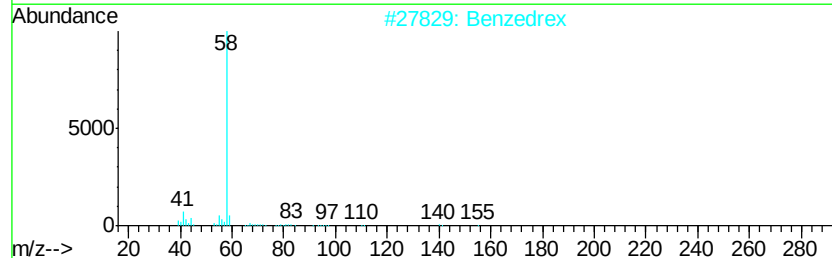
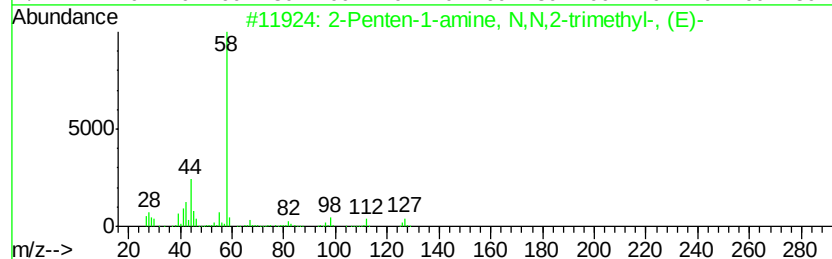
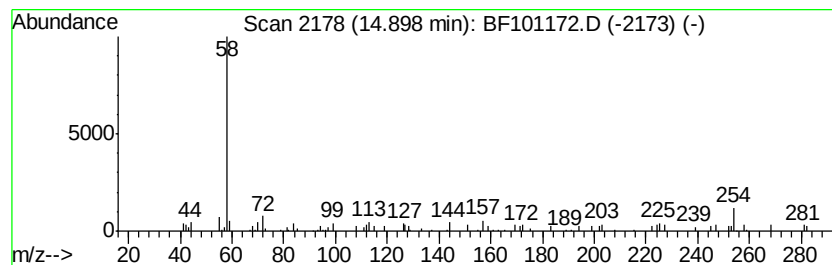
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 2-Penten-1-amine, N,N,2-tri... Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.90 | 2.80 ng | 40241 | Perylene-d12     | 15.50 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Penten-1-amine, N,N,2-trimethy... | 127 | C8H17N  | 055630-70-1 | 58   |
| 2    |    |   | Benzedrex                           | 155 | C10H21N | 000101-40-6 | 53   |
| 3    |    |   | 1-Undecanamine, N,N-dimethyl-       | 199 | C13H29N | 017373-28-3 | 53   |
| 4    |    |   | N,N-Dimethyl-3-butoxypropylamine    | 159 | C9H21NO | 071126-65-3 | 53   |
| 5    |    |   | 1-Hexanamine, 2-ethyl-N,N-dimethyl- | 157 | C10H23N | 028056-87-3 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

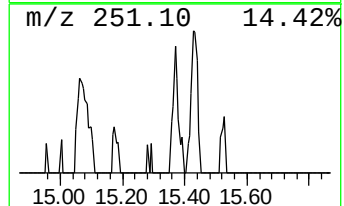
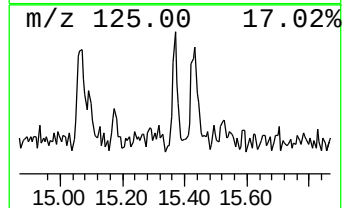
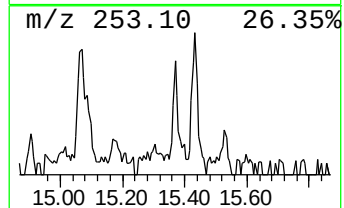
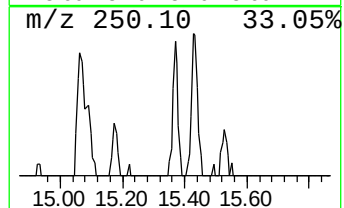
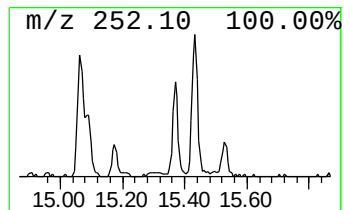
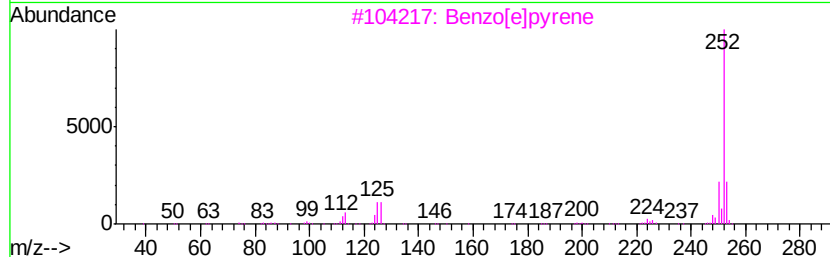
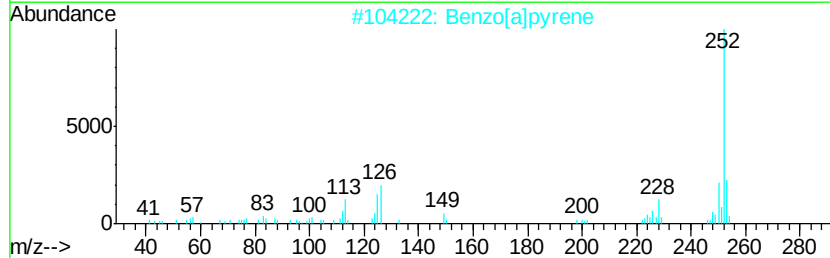
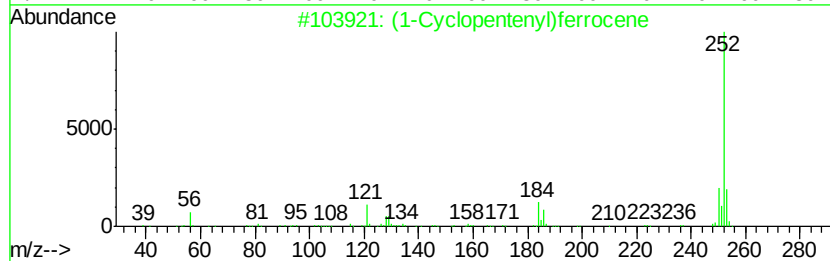
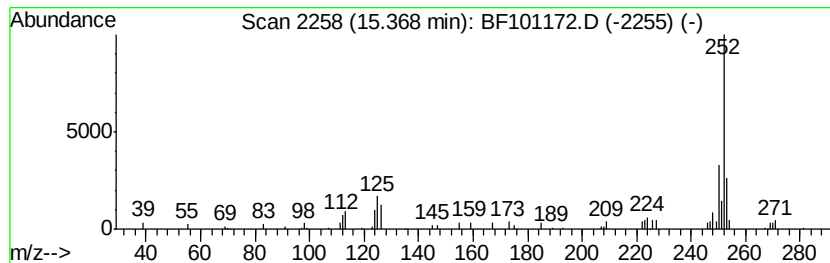
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 (1-Cyclopentenyl)ferrocene Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.37 | 2.45 ng | 35228 | Perylene-d12     | 15.50 |

| Hit# | of | Tentative ID               | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------|-----|----------|-------------|------|
| 1    | 5  | (1-Cyclopentenyl)ferrocene | 252 | C15H16Fe | 012260-67-2 | 86   |
| 2    |    | Benzo[a]pyrene             | 252 | C20H12   | 000050-32-8 | 81   |
| 3    |    | Benzo[e]pyrene             | 252 | C20H12   | 000192-97-2 | 81   |
| 4    |    | Benzo[e]acephenanthrylene  | 252 | C20H12   | 000205-99-2 | 76   |
| 5    |    | Benzo[k]fluoranthene       | 252 | C20H12   | 000207-08-9 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120617\  
Data File : BF101172.D  
Acq On : 6 Dec 2017 17:18  
Operator : SJ/JU  
Sample : I6683-03 10X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EO-01-120517-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF113017.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 |
| unknown6.62          | 6.62  | 7.9     | ng    | 102187   | 1                     | 6.85  | 259818 | 20.0 |
| Hexadecane           | 9.80  | 2.2     | ng    | 38312    | 3                     | 9.89  | 347760 | 20.0 |
| Dodecane             | 10.30 | 2.4     | ng    | 41492    | 3                     | 9.89  | 347760 | 20.0 |
| Octacosane           | 10.77 | 3.0     | ng    | 47239    | 4                     | 11.37 | 319138 | 20.0 |
| Hexadecanoic acid... | 11.76 | 40.1    | ng    | 640143   | 4                     | 11.37 | 319138 | 20.0 |
| 9,12-Octadecadien... | 12.45 | 164.0   | ng    | 2616680  | 4                     | 11.37 | 319138 | 20.0 |
| 13-Octadecenoic a... | 12.47 | 74.1    | ng    | 1182460  | 4                     | 11.37 | 319138 | 20.0 |
| unknown13.03         | 13.03 | 2.2     | ng    | 35178    | 5                     | 14.02 | 322323 | 20.0 |
| Pyrene, 1-methyl-    | 13.15 | 2.1     | ng    | 33128    | 5                     | 14.02 | 322323 | 20.0 |
| 2-Penten-1-amine,... | 14.90 | 2.8     | ng    | 40241    | 6                     | 15.50 | 287422 | 20.0 |
| (1-Cyclopentenyl)... | 15.37 | 2.5     | ng    | 35228    | 6                     | 15.50 | 287422 | 20.0 |