

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
 Data File : BF117506.D  
 Acq On : 24 Oct 2019 1:37  
 Operator : JU  
 Sample : K5152-10 2X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SU-03-102219

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.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         | 5.351       | 552           | 555         | 574          | rBV      | 248032         | 367530        | 4.96%           | 1.491%        |
| 2         | 6.375       | 726           | 729         | 748          | rBV      | 345042         | 431090        | 5.82%           | 1.749%        |
| 3         | 6.504       | 748           | 751         | 758          | rVV      | 478479         | 445282        | 6.01%           | 1.807%        |
| 4         | 6.728       | 786           | 789         | 803          | rBV      | 226523         | 234938        | 3.17%           | 0.953%        |
| 5         | 6.887       | 812           | 816         | 830          | rBV      | 365103         | 325363        | 4.39%           | 1.320%        |
| 6         | 7.292       | 882           | 885         | 898          | rBV      | 199448         | 252946        | 3.41%           | 1.026%        |
| 7         | 8.010       | 1002          | 1007        | 1014         | rBV      | 357532         | 349546        | 4.72%           | 1.418%        |
| 8         | 9.086       | 1186          | 1190        | 1195         | rBV      | 530880         | 486919        | 6.57%           | 1.976%        |
| 9         | 9.481       | 1253          | 1257        | 1260         | rBV2     | 79068          | 93070         | 1.26%           | 0.378%        |
| 10        | 9.698       | 1291          | 1294        | 1297         | rBV      | 108809         | 80597         | 1.09%           | 0.327%        |
| 11        | 9.763       | 1301          | 1305        | 1309         | rVB      | 473044         | 383777        | 5.18%           | 1.557%        |
| 12        | 10.192      | 1376          | 1378        | 1381         | rVB      | 129995         | 102883        | 1.39%           | 0.417%        |
| 13        | 10.551      | 1437          | 1439        | 1450         | rVB      | 277572         | 284560        | 3.84%           | 1.155%        |
| 14        | 10.669      | 1455          | 1459        | 1467         | rBV      | 147361         | 204221        | 2.75%           | 0.829%        |
| 15        | 11.116      | 1531          | 1535        | 1537         | rBV      | 138046         | 113666        | 1.53%           | 0.461%        |
| 16        | 11.145      | 1537          | 1540        | 1545         | rVB      | 95811          | 98869         | 1.33%           | 0.401%        |
| 17        | 11.245      | 1554          | 1557        | 1560         | rBV      | 367495         | 379487        | 5.12%           | 1.540%        |
| 18        | 11.275      | 1560          | 1562        | 1569         | rVB      | 294466         | 266287        | 3.59%           | 1.080%        |
| 19        | 11.539      | 1604          | 1607        | 1609         | rBV      | 148528         | 112388        | 1.52%           | 0.456%        |
| 20        | 11.645      | 1620          | 1625        | 1628         | rBV      | 2388606        | 1970281       | 26.58%          | 7.994%        |
| 21        | 11.751      | 1640          | 1643        | 1645         | rBV      | 127644         | 112017        | 1.51%           | 0.454%        |
| 22        | 11.780      | 1645          | 1648        | 1653         | rVB      | 222279         | 202706        | 2.73%           | 0.822%        |
| 23        | 11.857      | 1658          | 1661        | 1664         | rVV2     | 125608         | 146089        | 1.97%           | 0.593%        |
| 24        | 11.886      | 1664          | 1666        | 1670         | rVB      | 100552         | 85984         | 1.16%           | 0.349%        |
| 25        | 11.945      | 1672          | 1676        | 1681         | rBV      | 114482         | 112621        | 1.52%           | 0.457%        |
| 26        | 12.045      | 1690          | 1693        | 1697         | rBV      | 78471          | 80713         | 1.09%           | 0.327%        |
| 27        | 12.333      | 1737          | 1742        | 1744         | rVV      | 3543627        | 4527905       | 61.08%          | 18.371%       |
| 28        | 12.369      | 1744          | 1748        | 1752         | rVV      | 5978664        | 7413168       | 100.00%         | 30.078%       |
| 29        | 12.398      | 1752          | 1753        | 1756         | rVV      | 110998         | 83187         | 1.12%           | 0.338%        |
| 30        | 12.433      | 1756          | 1759        | 1762         | rVV      | 942955         | 748570        | 10.10%          | 3.037%        |
| 31        | 12.463      | 1762          | 1764        | 1767         | rVV      | 378420         | 407670        | 5.50%           | 1.654%        |
| 32        | 12.492      | 1767          | 1769        | 1779         | rVB3     | 231776         | 332259        | 4.48%           | 1.348%        |
| 33        | 12.669      | 1796          | 1799        | 1800         | rBV      | 87645          | 83763         | 1.13%           | 0.340%        |
| 34        | 12.692      | 1800          | 1803        | 1811         | rVB      | 408341         | 475708        | 6.42%           | 1.930%        |

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Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 35 | 12.833 | 1823 | 1827 | 1830 | rBV  | 440916 | 392817 | 5.30% | 1.594% |
| 36 | 12.916 | 1837 | 1841 | 1844 | rBV  | 63855  | 108090 | 1.46% | 0.439% |
| 37 | 13.010 | 1855 | 1857 | 1863 | rVV4 | 87868  | 151865 | 2.05% | 0.616% |
| 38 | 13.074 | 1863 | 1868 | 1874 | rVB4 | 167929 | 246786 | 3.33% | 1.001% |
| 39 | 13.157 | 1879 | 1882 | 1885 | rVB  | 145512 | 130737 | 1.76% | 0.530% |
| 40 | 13.251 | 1895 | 1898 | 1906 | rBV2 | 70798  | 104964 | 1.42% | 0.426% |
| 41 | 13.586 | 1950 | 1955 | 1957 | rBV4 | 59300  | 85700  | 1.16% | 0.348% |
| 42 | 13.821 | 1992 | 1995 | 2001 | rVB  | 160136 | 137675 | 1.86% | 0.559% |
| 43 | 13.892 | 2002 | 2007 | 2010 | rBV2 | 370809 | 503745 | 6.80% | 2.044% |
| 44 | 13.916 | 2010 | 2011 | 2016 | rVB  | 159904 | 132496 | 1.79% | 0.538% |
| 45 | 14.351 | 2082 | 2085 | 2089 | rBV2 | 91953  | 125871 | 1.70% | 0.511% |
| 46 | 14.763 | 2151 | 2155 | 2159 | rVB  | 116448 | 155453 | 2.10% | 0.631% |
| 47 | 14.921 | 2179 | 2182 | 2185 | rBV  | 79571  | 100092 | 1.35% | 0.406% |
| 48 | 15.327 | 2247 | 2251 | 2255 | rVB  | 188338 | 236651 | 3.19% | 0.960% |
| 49 | 15.951 | 2355 | 2357 | 2365 | rVB3 | 70055  | 123078 | 1.66% | 0.499% |
| 50 | 16.745 | 2488 | 2492 | 2500 | rVB4 | 58240  | 114641 | 1.55% | 0.465% |

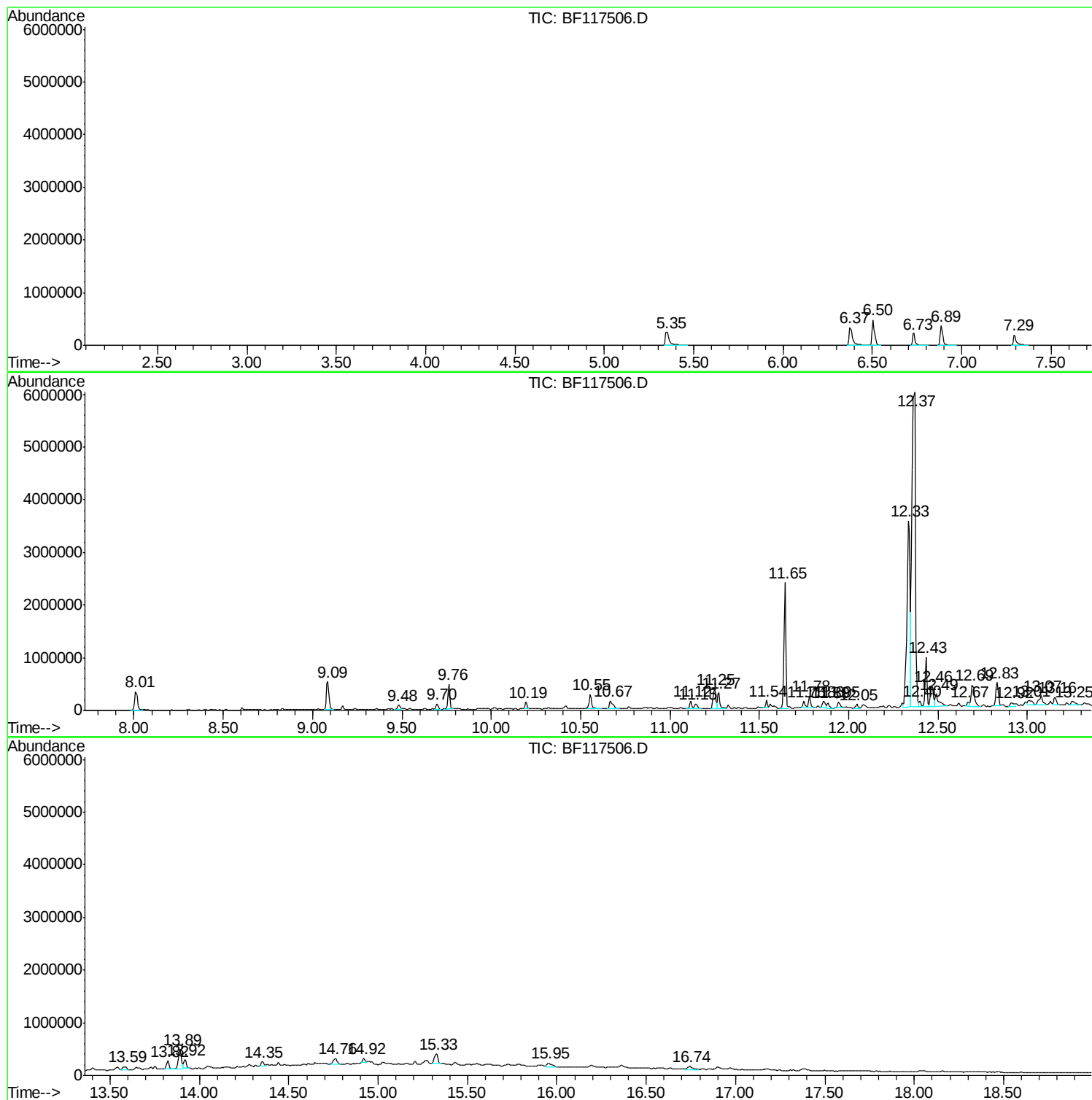
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
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SU-03-102219

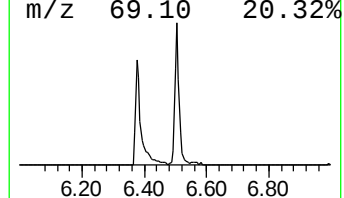
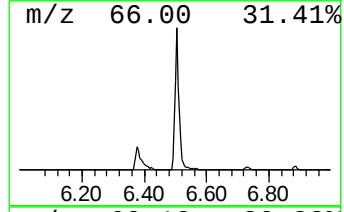
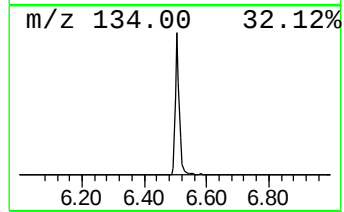
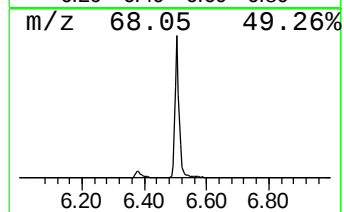
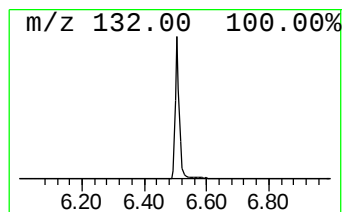
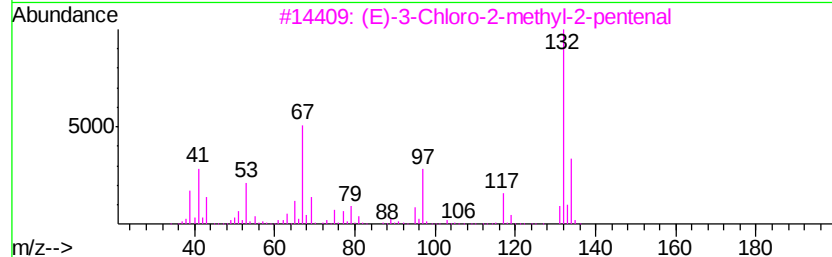
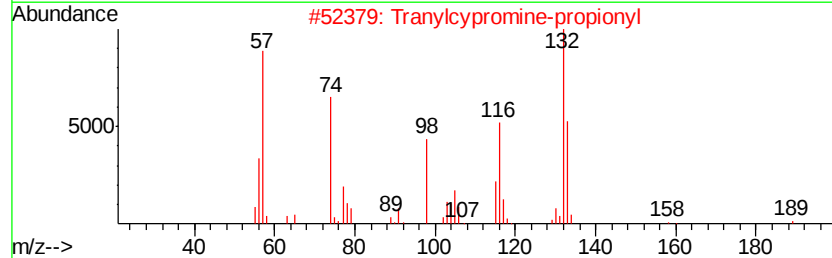
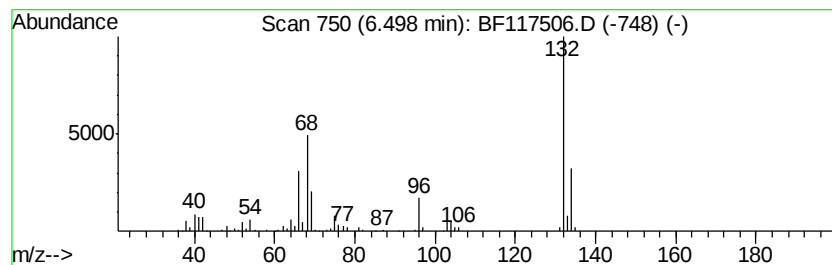
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 unknown6.50 Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.50 | 37.91 ng | 445282 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 2    |    |   | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 3    |    |   | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 4    |    |   | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 10   |
| 5    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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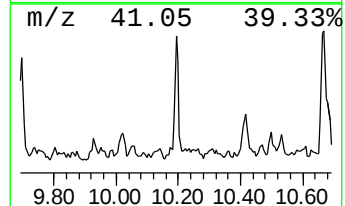
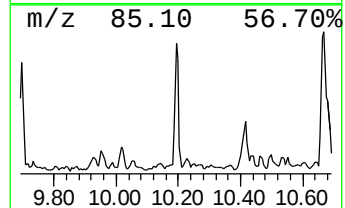
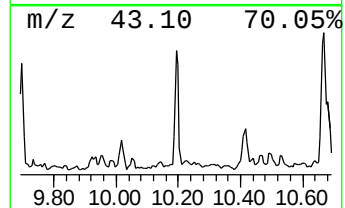
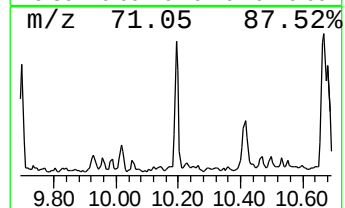
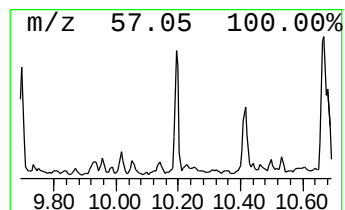
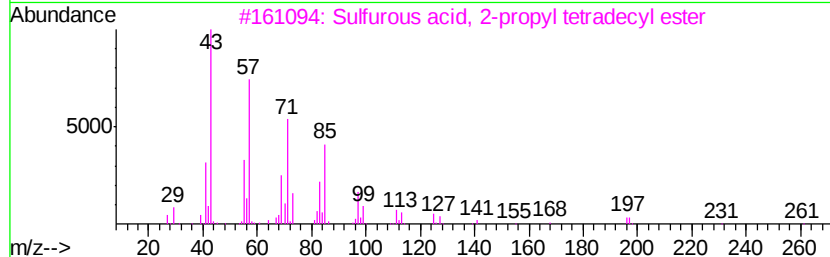
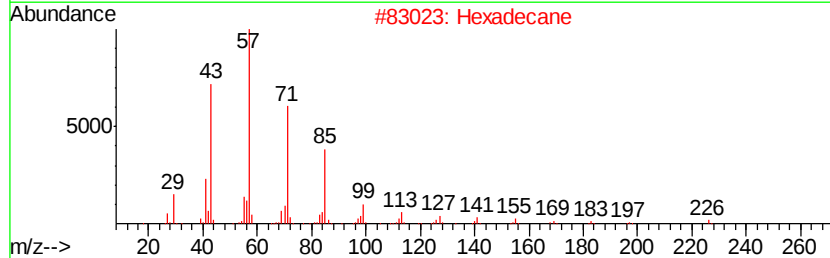
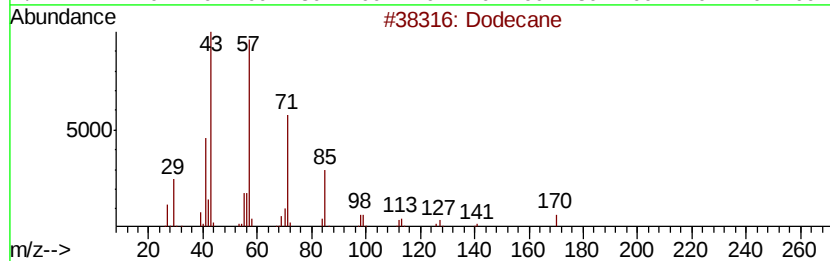
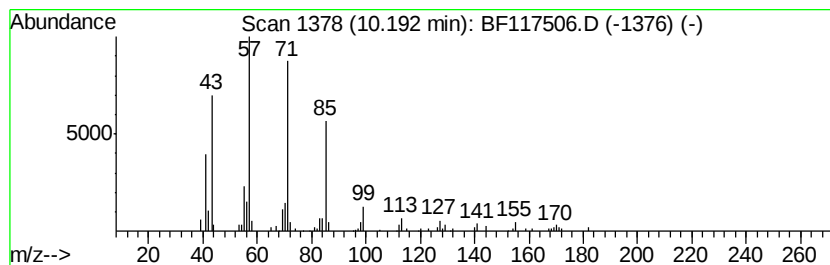
Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.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 Dodecane Concentration Rank 18

| R.T.    | EstConc                                   | Area          | Relative to ISTD | R.T.      |
|---------|-------------------------------------------|---------------|------------------|-----------|
| 10.19   | 5.36 ng                                   | 102883        | Acenaphthene-d10 | 9.76      |
| Hit# of | 5                                         | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | Dodecane                                  | 170 C12H26    | 000112-40-3      | 90        |
| 2       | Hexadecane                                | 226 C16H34    | 000544-76-3      | 87        |
| 3       | Sulfurous acid, 2-propyl tetradecyl ester | 320 C17H36O3S | 1000309-12-5     | 86        |
| 4       | Tetracosane                               | 338 C24H50    | 000646-31-1      | 86        |
| 5       | Dodecane, 1-iodo-                         | 296 C12H25I   | 004292-19-7      | 80        |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

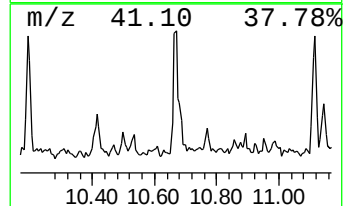
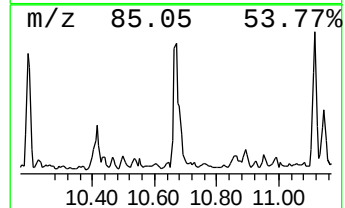
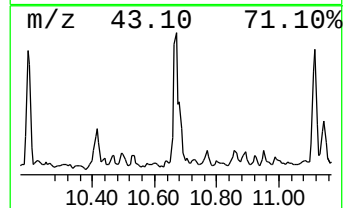
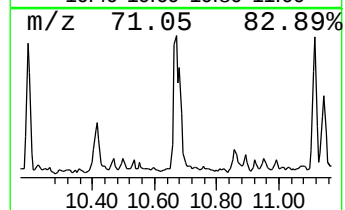
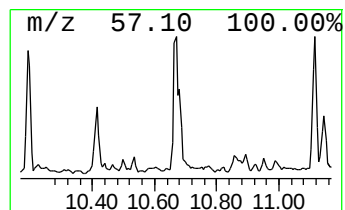
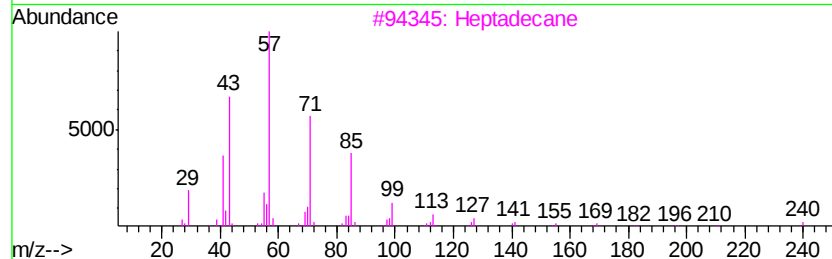
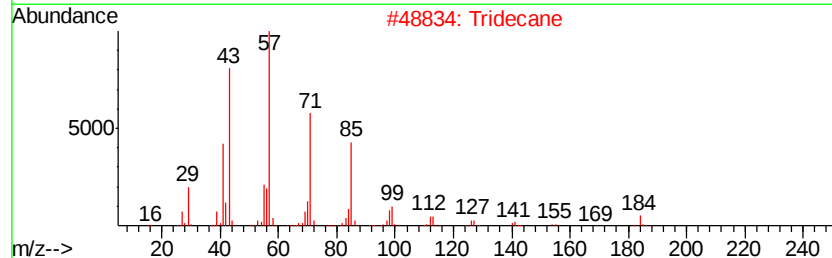
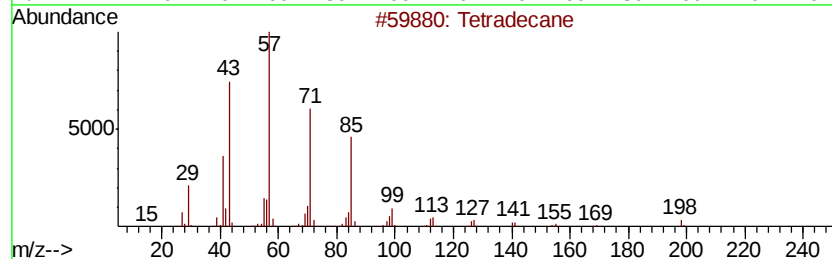
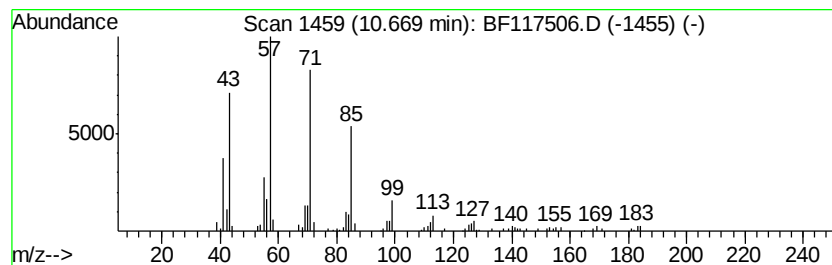
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 4 Tetradecane Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.67 | 10.76 ng | 204221 | Phenanthrene-d10 | 11.25 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|---------|---|-------------------------------------|-----|-----------|--------------|------|
| 1       |   | Tetradecane                         | 198 | C14H30    | 000629-59-4  | 91   |
| 2       |   | Tridecane                           | 184 | C13H28    | 000629-50-5  | 91   |
| 3       |   | Heptadecane                         | 240 | C17H36    | 000629-78-7  | 91   |
| 4       |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 91   |
| 5       |   | Sulfurous acid, 2-propyl tetrade... | 320 | C17H36O3S | 1000309-12-5 | 90   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

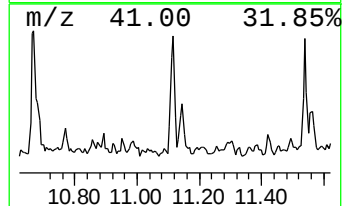
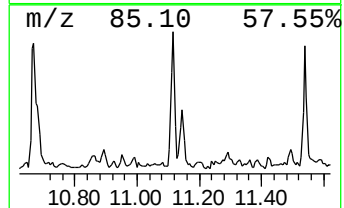
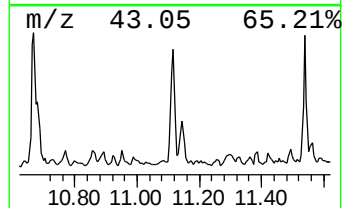
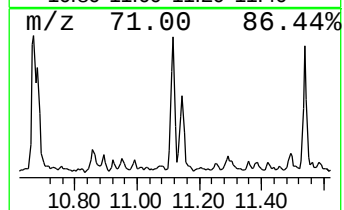
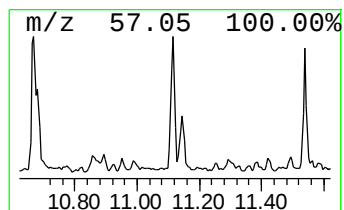
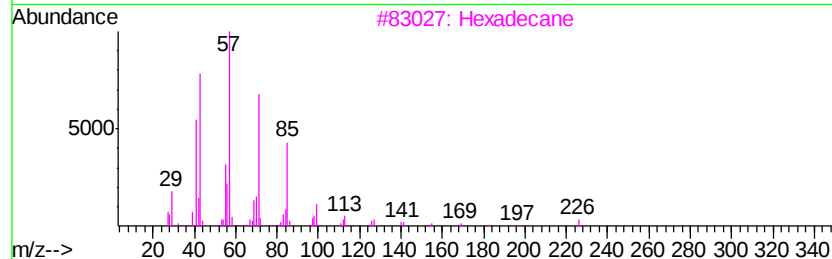
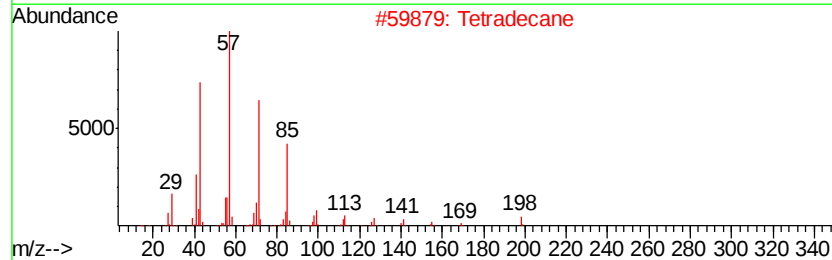
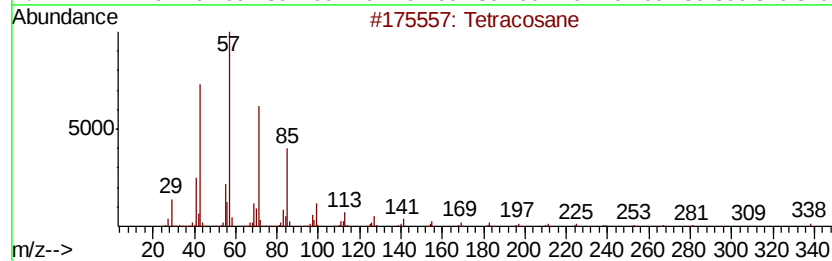
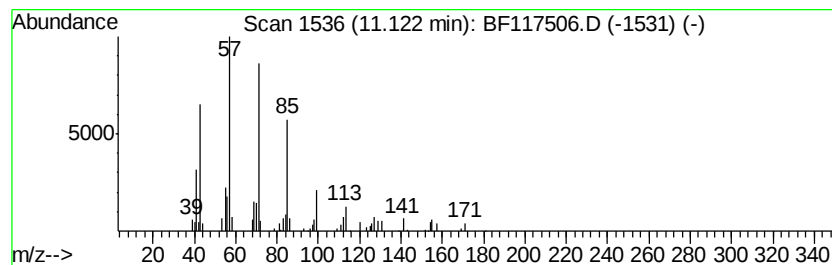
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 5 Tetracosane Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.12 | 5.99 ng | 113666 | Phenanthrene-d10 | 11.25 |

| Hit# of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------------------|-----|---------|-------------|------|
| 1       |   | Tetracosane              | 338 | C24H50  | 000646-31-1 | 90   |
| 2       |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 3       |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 90   |
| 4       |   | Dodecane, 1-iodo-        | 296 | C12H25I | 004292-19-7 | 87   |
| 5       |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

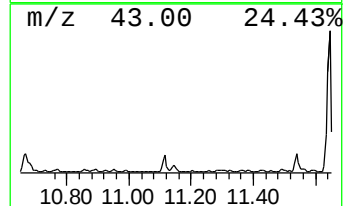
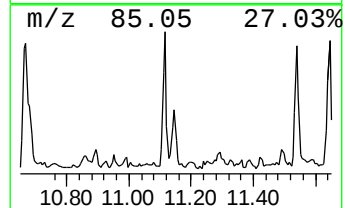
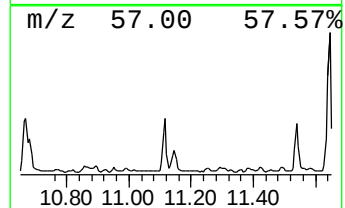
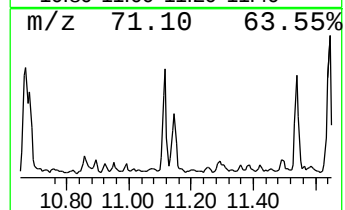
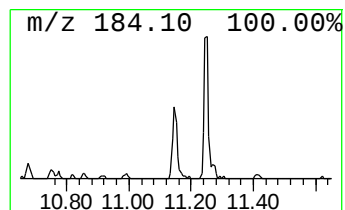
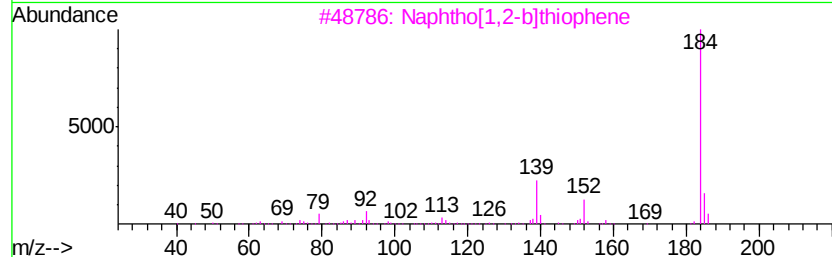
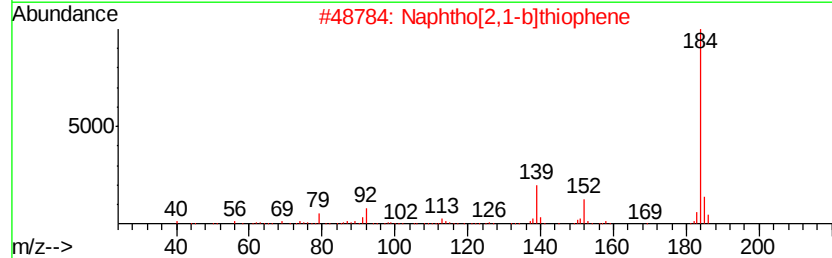
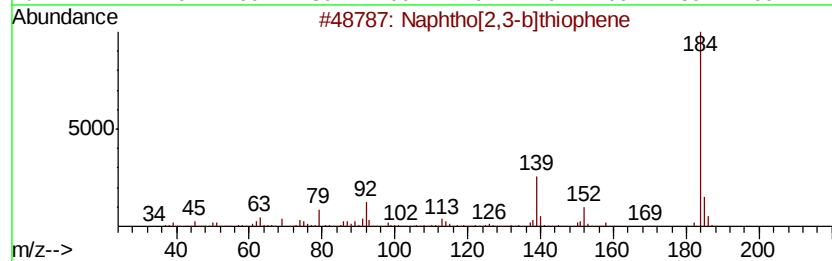
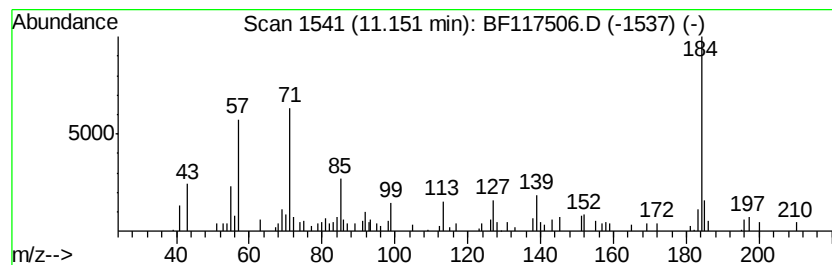
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 unknown11.15 Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.15 | 5.21 ng | 98869 | Phenanthrene-d10 | 11.25 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphtho[2,3-b]thiophene | 184 | C12H8S  | 000268-77-9 | 46   |
| 2    |    | Naphtho[2,1-b]thiophene | 184 | C12H8S  | 000233-02-3 | 46   |
| 3    |    | Naphtho[1,2-b]thiophene | 184 | C12H8S  | 000234-41-3 | 46   |
| 4    |    | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 46   |
| 5    |    | Azuleno(2,1-b)thiophene | 184 | C12H8S  | 000248-13-5 | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

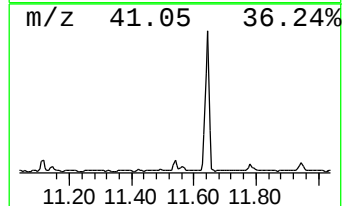
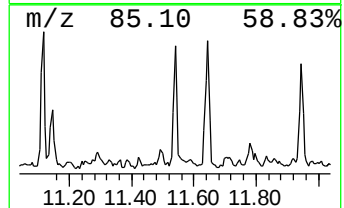
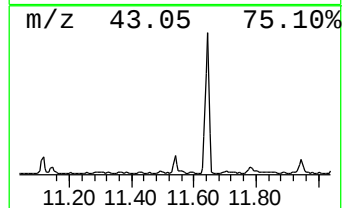
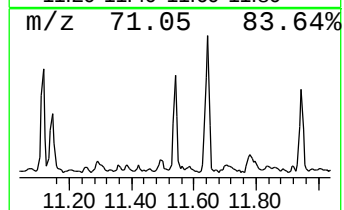
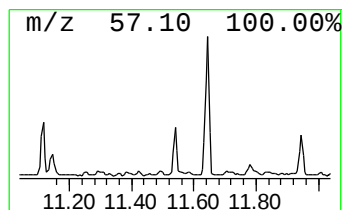
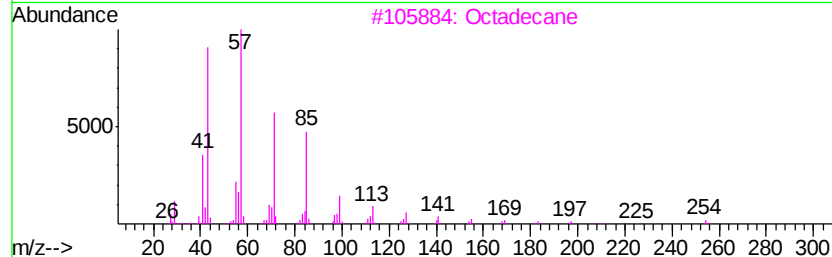
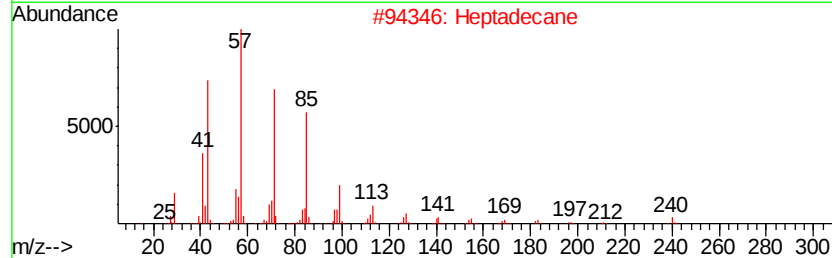
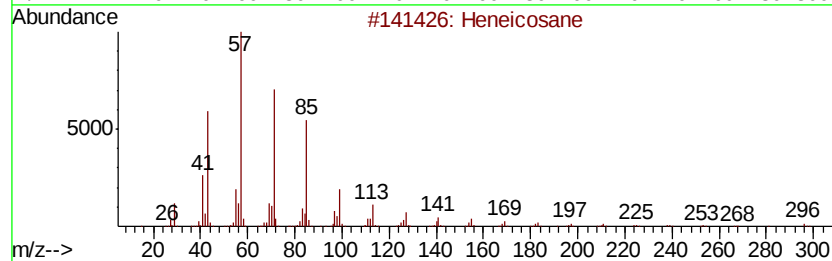
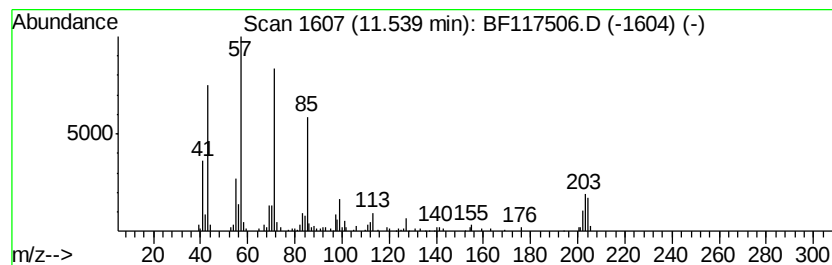
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Heneicosane Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.54 | 5.92 ng | 112388 | Phenanthrene-d10 | 11.25 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Heneicosane  | 296 | C21H44  | 000629-94-7 | 81   |
| 2         | Heptadecane  | 240 | C17H36  | 000629-78-7 | 74   |
| 3         | Octadecane   | 254 | C18H38  | 000593-45-3 | 72   |
| 4         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 72   |
| 5         | Tetracosane  | 338 | C24H50  | 000646-31-1 | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

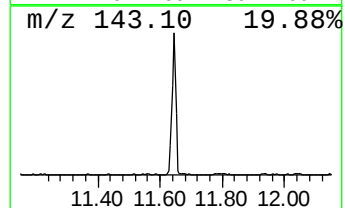
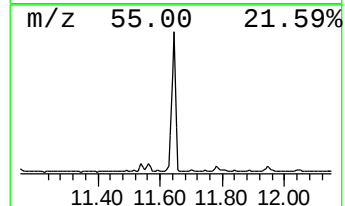
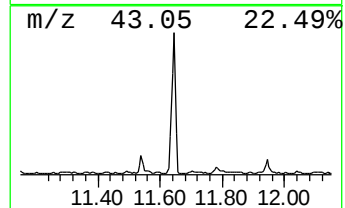
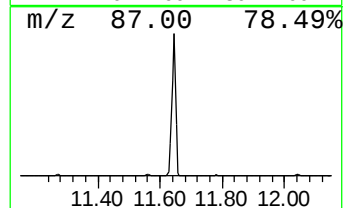
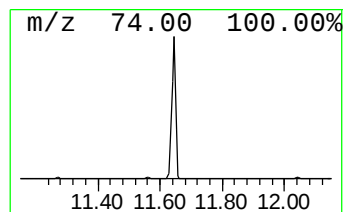
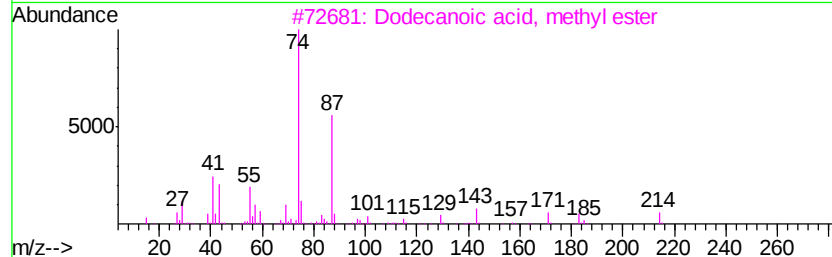
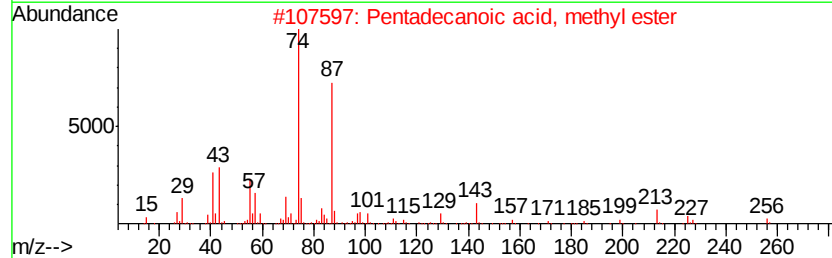
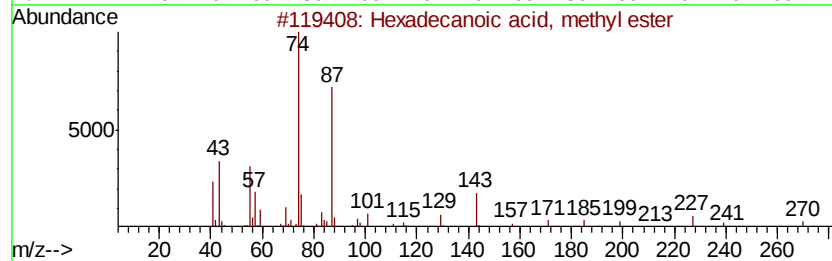
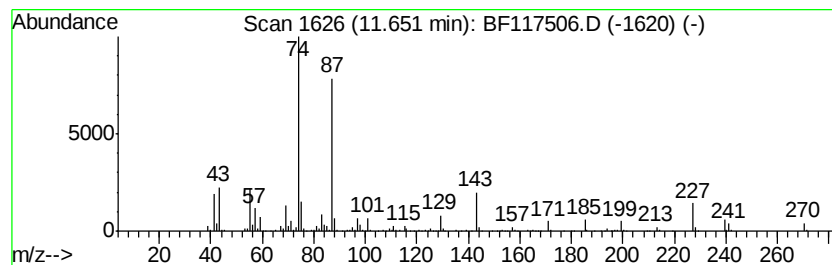
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 11.65 | 103.84 ng | 1970280 | Phenanthrene-d10 | 11.25 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexadecanoic acid, methyl ester  | 270 | C17H34O2 | 000112-39-0 | 97   |
| 2    |    | Pentadecanoic acid, methyl ester | 256 | C16H32O2 | 007132-64-1 | 90   |
| 3    |    | Dodecanoic acid, methyl ester    | 214 | C13H26O2 | 000111-82-0 | 87   |
| 4    |    | Methyl tetradecanoate            | 242 | C15H30O2 | 000124-10-7 | 87   |
| 5    |    | Tridecanoic acid, methyl ester   | 228 | C14H28O2 | 001731-88-0 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

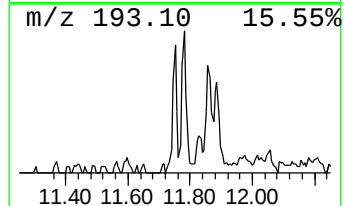
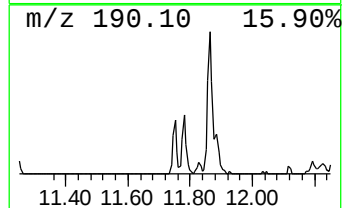
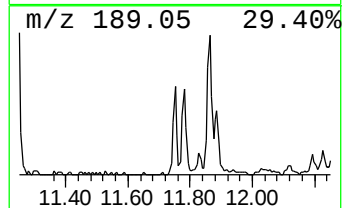
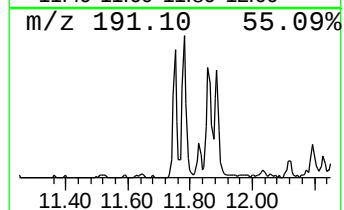
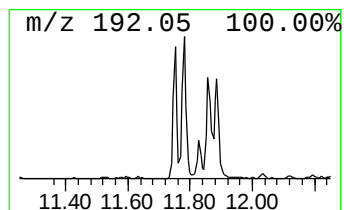
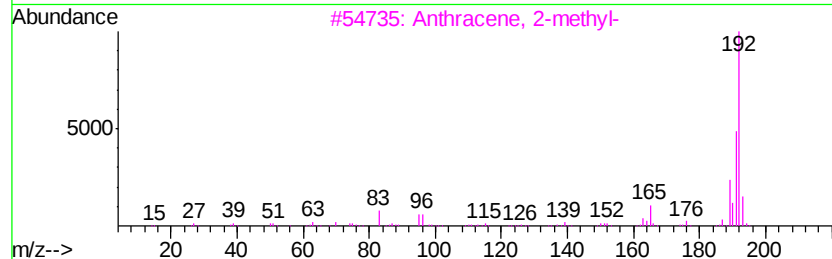
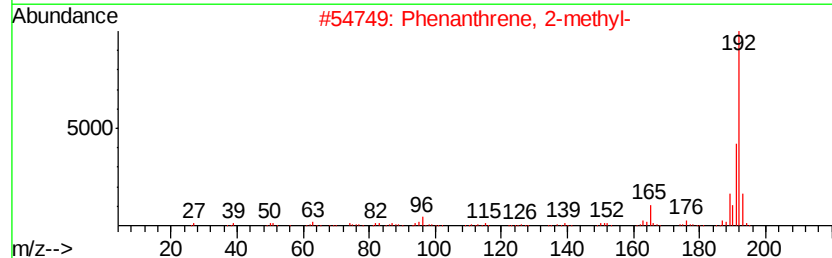
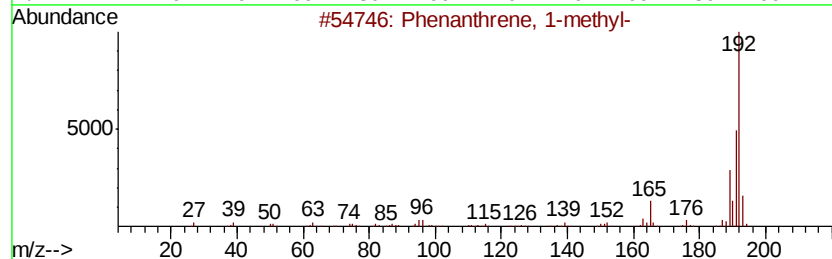
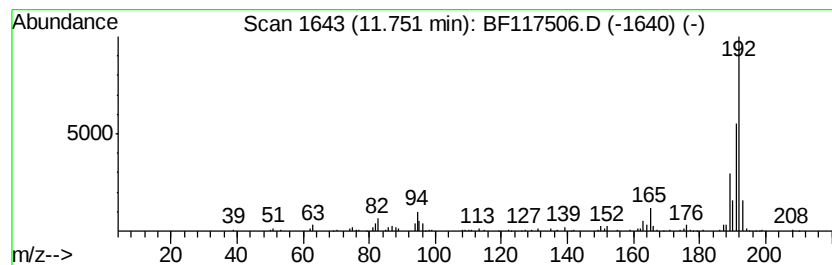
Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 16

| R.T.      | EstConc                               | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------------------|--------|------------------|-------------|------|
| 11.75     | 5.90 ng                               | 112017 | Phenanthrene-d10 | 11.25       |      |
| Hit# of 5 | Tentative ID                          | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 1-methyl-               | 192    | C15H12           | 000832-69-9 | 97   |
| 2         | Phenanthrene, 2-methyl-               | 192    | C15H12           | 002531-84-2 | 96   |
| 3         | Anthracene, 2-methyl-                 | 192    | C15H12           | 000613-12-7 | 95   |
| 4         | Anthracene, 1-methyl-                 | 192    | C15H12           | 000610-48-0 | 94   |
| 5         | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192    | C15H12           | 000949-41-7 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

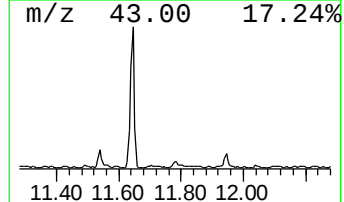
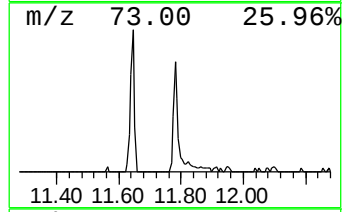
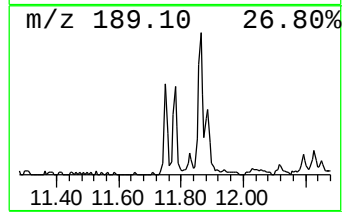
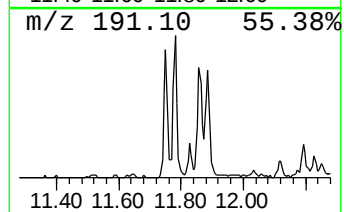
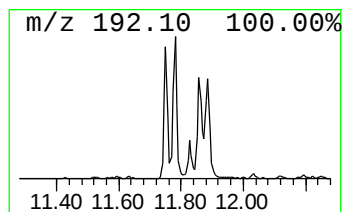
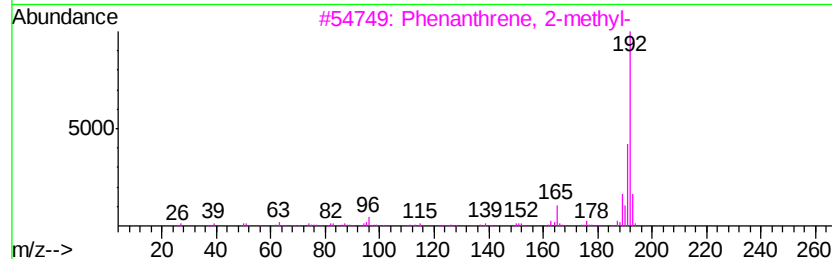
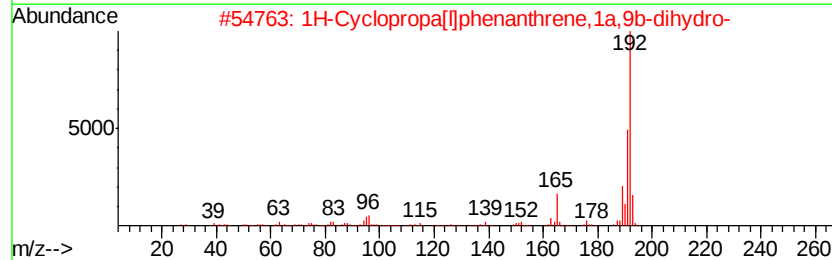
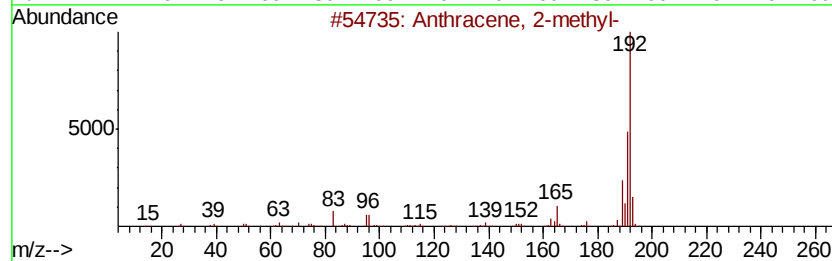
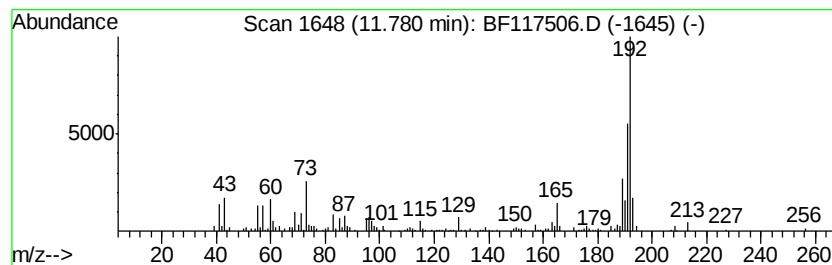
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Anthracene, 2-methyl- Concentration Rank 8

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.78 | 10.68 ng | 202706 | Phenanthrene-d10 | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 96   |
| 2    |    | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 3    |    | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 95   |
| 4    |    | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 95   |
| 5    |    | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

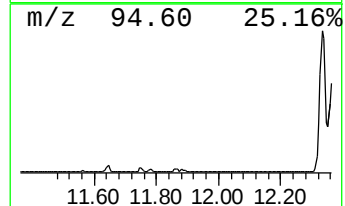
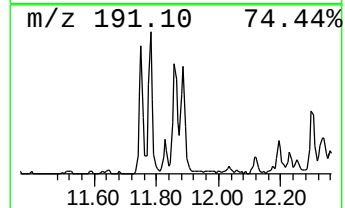
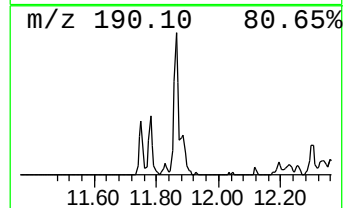
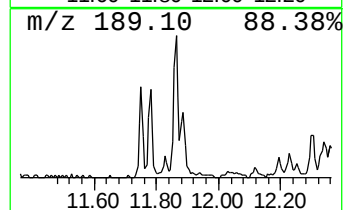
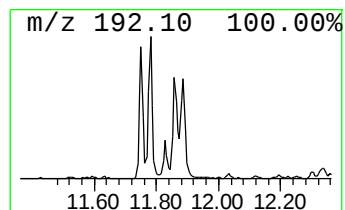
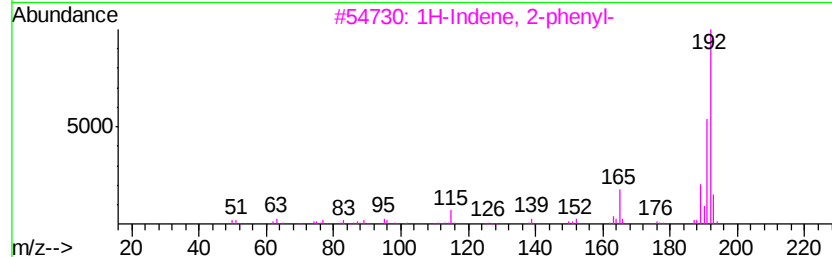
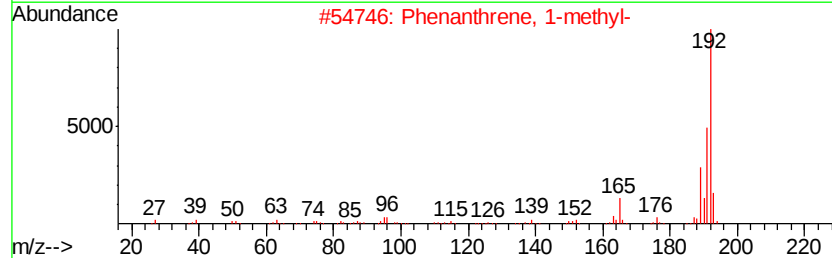
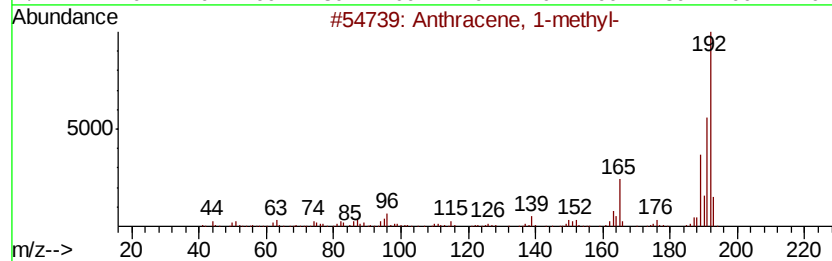
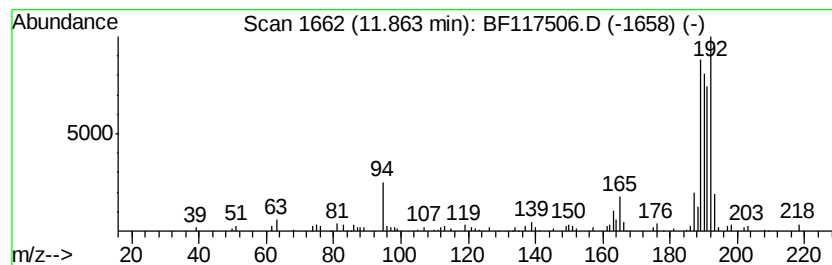
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ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 Anthracene, 1-methyl- Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------------------|--------|------------------|-------------|-------|
| 11.86     | 7.70 ng                             | 146089 | Phenanthrene-d10 |             | 11.25 |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual  |
| 1         | Anthracene, 1-methyl-               | 192    | C15H12           | 000610-48-0 | 91    |
| 2         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 91    |
| 3         | 1H-Indene, 2-phenyl-                | 192    | C15H12           | 004505-48-0 | 70    |
| 4         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 70    |
| 5         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 70    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

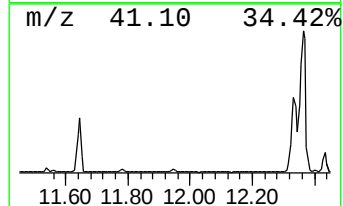
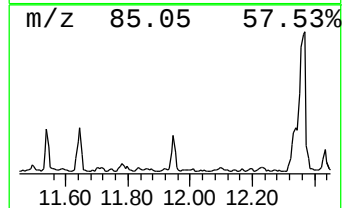
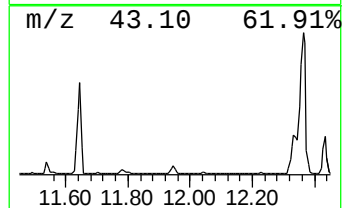
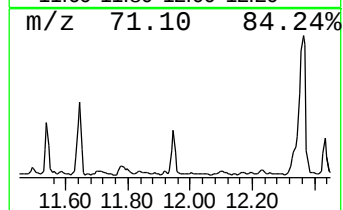
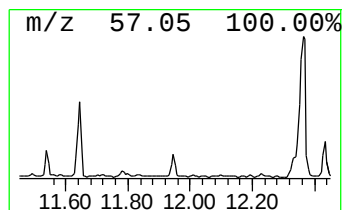
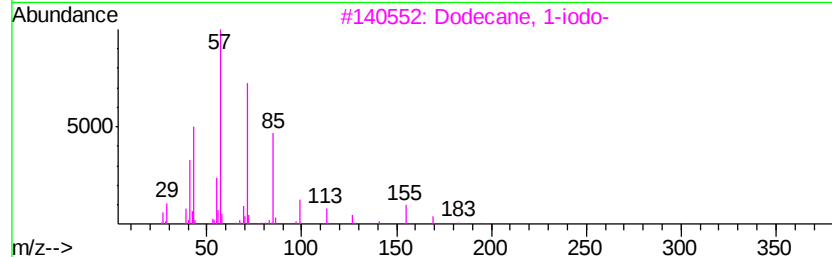
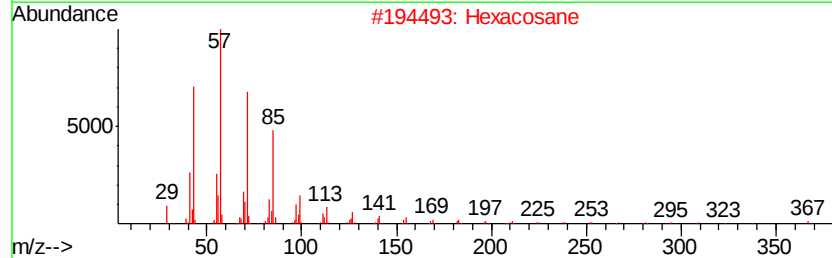
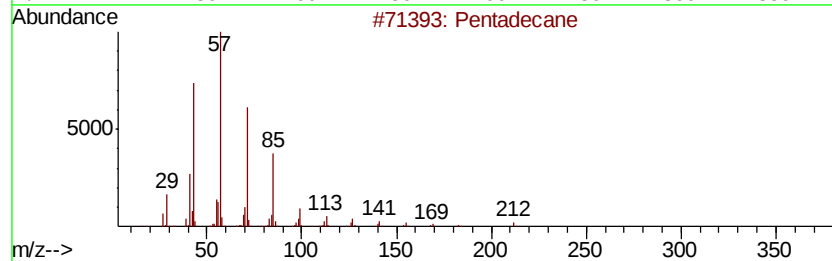
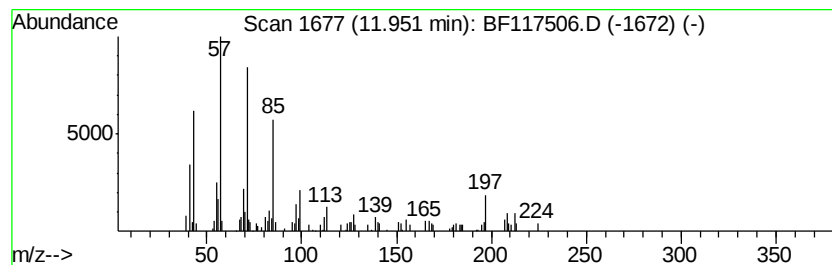
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Pentadecane Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.95 | 5.94 ng | 112621 | Phenanthrene-d10 | 11.25 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane       | 212 | C15H32  | 000629-62-9 | 90   |
| 2    |    | Hexacosane        | 366 | C26H54  | 000630-01-3 | 87   |
| 3    |    | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 80   |
| 4    |    | Heptacosane       | 380 | C27H56  | 000593-49-7 | 74   |
| 5    |    | Hexadecane        | 226 | C16H34  | 000544-76-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

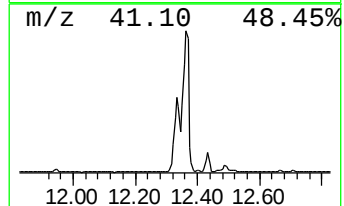
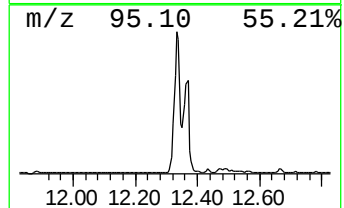
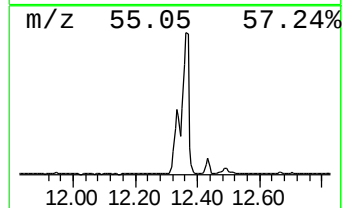
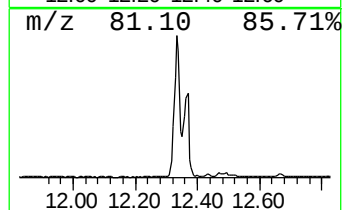
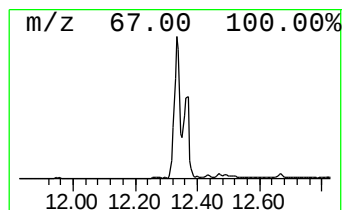
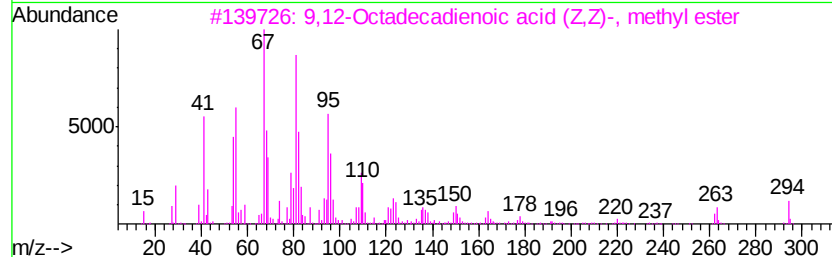
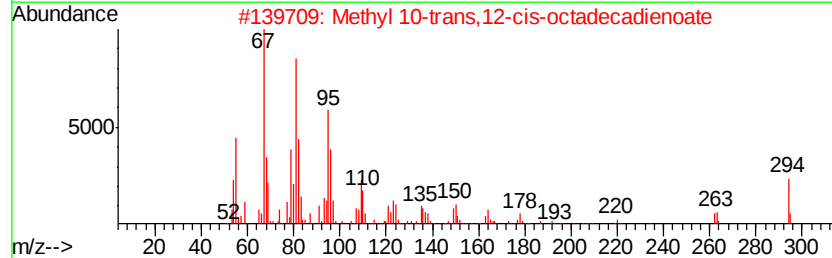
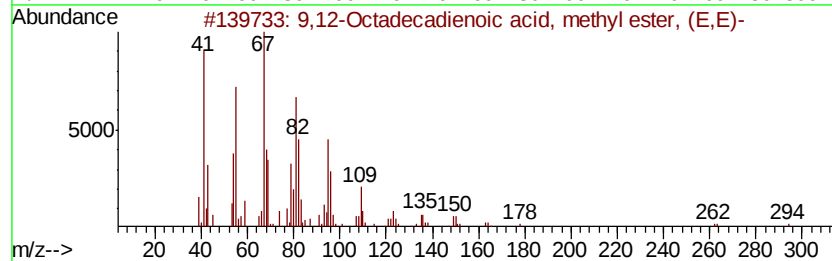
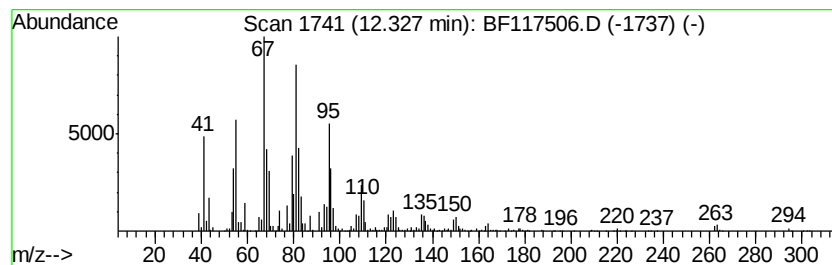
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 9,12-Octadecadienoic acid, ... Concentration Rank 2

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.33 | 238.63 ng | 4527910 | Phenanthrene-d10 | 11.25 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4  | 99   |
| 2    |    |   | Methyl 10-trans,12-cis-octadecad... | 294 | C19H34O2 | 1000336-44-2 | 99   |
| 3    |    |   | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0  | 99   |
| 4    |    |   | 10,13-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-62-2  | 98   |
| 5    |    |   | 11,14-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-61-1  | 98   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

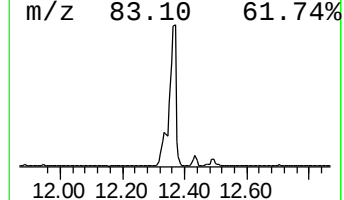
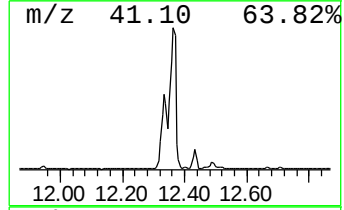
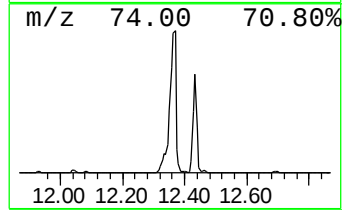
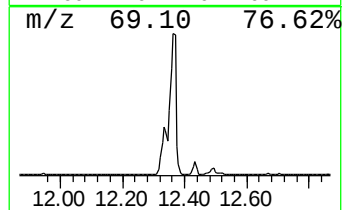
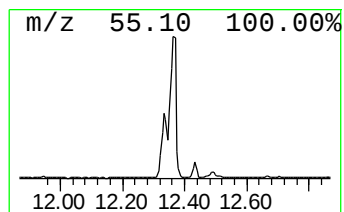
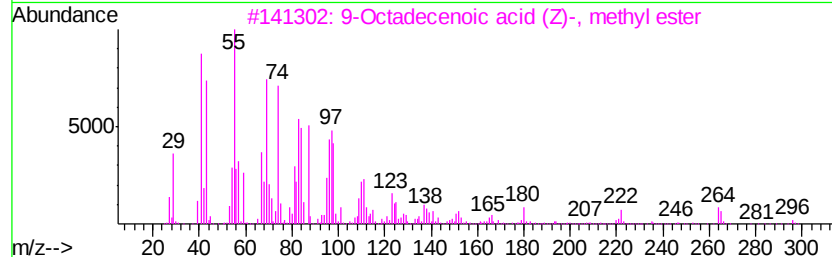
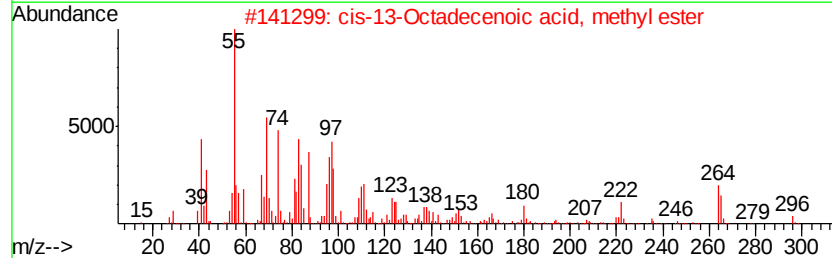
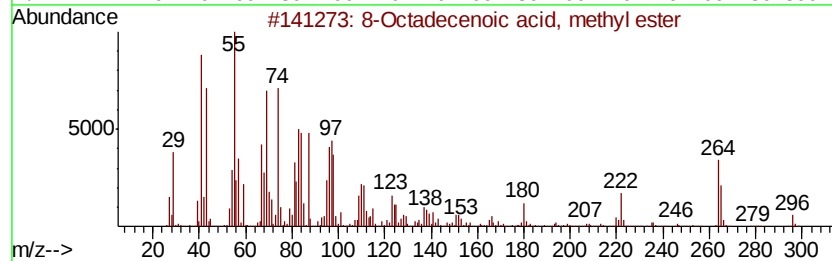
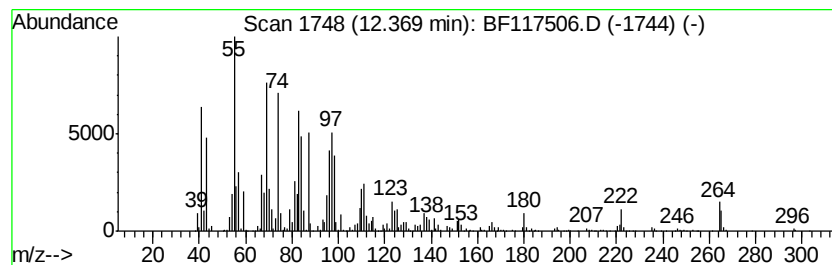
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 8-Octadecenoic acid, methyl... Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 12.37 | 390.69 ng | 7413170 | Phenanthrene-d10 | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 8-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 002345-29-1  | 99   |
| 2    |    | cis-13-Octadecenoic acid, methyl... | 296 | C19H36O2 | 1000333-58-3 | 96   |
| 3    |    | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9  | 91   |
| 4    |    | 9-Octadecenoic acid, methyl este... | 296 | C19H36O2 | 001937-62-8  | 91   |
| 5    |    | trans-13-Octadecenoic acid, meth... | 296 | C19H36O2 | 1000333-61-3 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

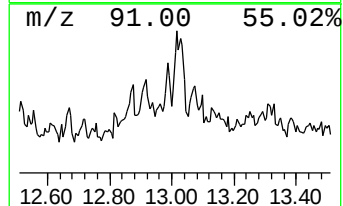
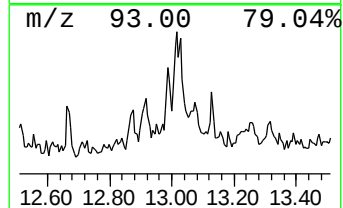
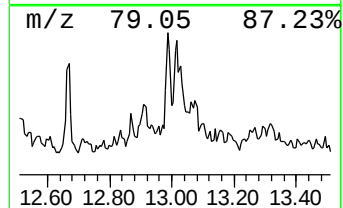
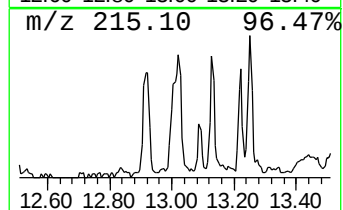
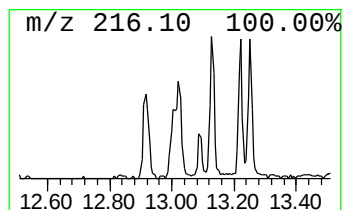
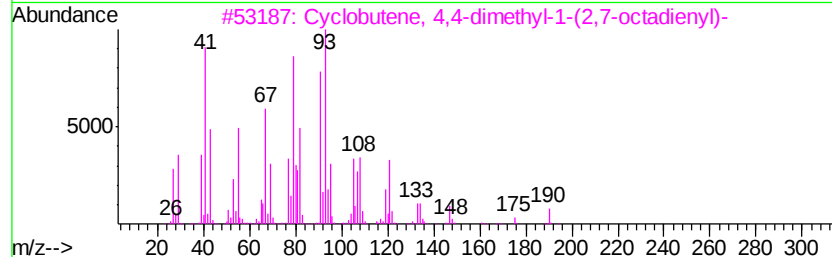
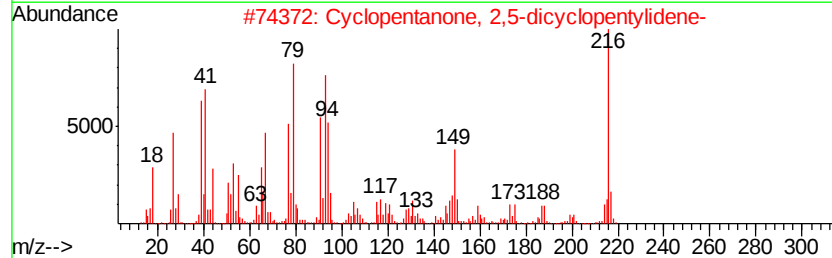
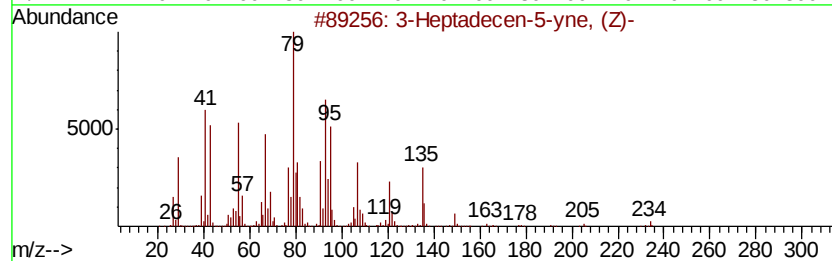
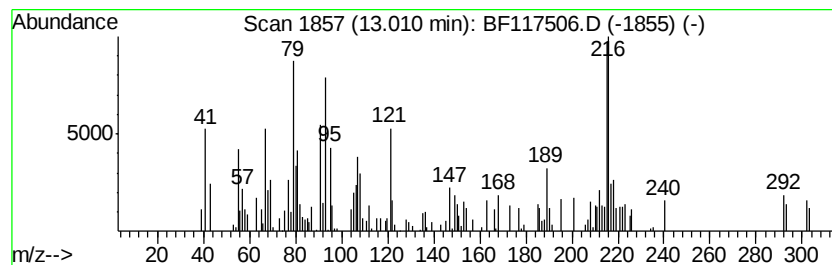
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 unknown13.01 Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.01 | 6.03 ng | 151865 | Chrysene-d12     | 13.89 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 3-Heptadecen-5-yne, (Z)-             | 234 | C17H30  | 074744-55-1 | 43   |
| 2    |    |   | Cyclopentanone, 2,5-dicyclopentyl... | 216 | C15H20O | 005682-82-6 | 43   |
| 3    |    |   | Cyclobutene, 4,4-dimethyl-1-(2,7...  | 190 | C14H22  | 062338-42-5 | 42   |
| 4    |    |   | 1,5-Cycloundecadiene, 9-(1-methy...  | 190 | C14H22  | 062338-55-0 | 35   |
| 5    |    |   | Bicyclo[3.1.1]heptane, 6,6-dimet...  | 136 | C10H16  | 016022-04-1 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

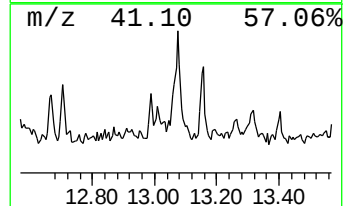
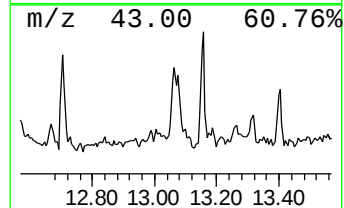
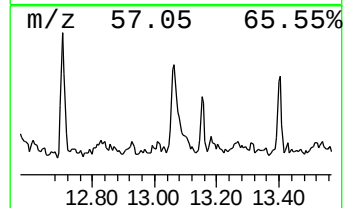
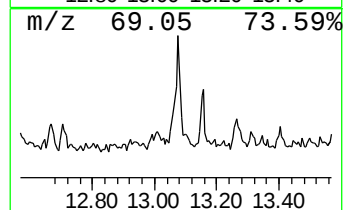
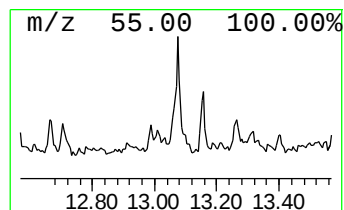
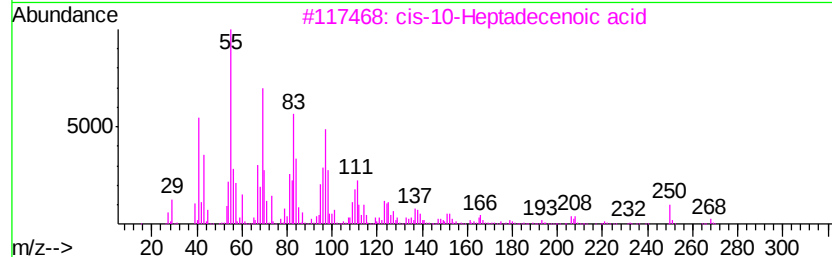
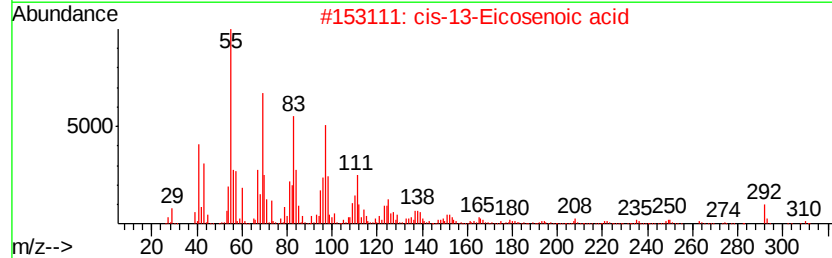
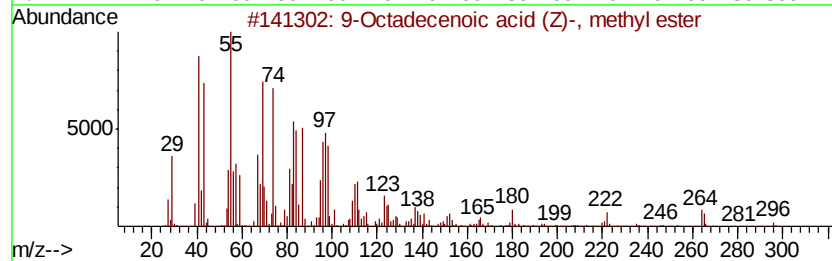
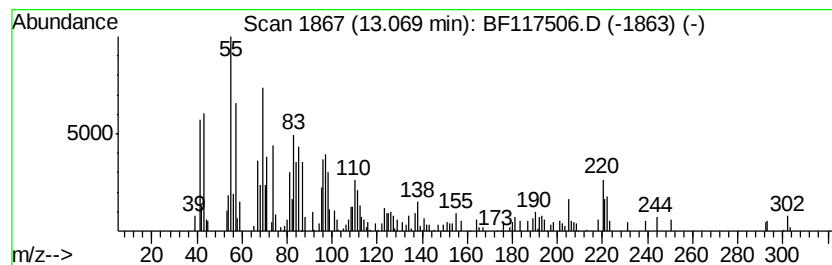
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 9-Octadecenoic acid (Z)-, m... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.07 | 9.80 ng | 246786 | Chrysene-d12     | 13.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9 | 80   |
| 2    |    | cis-13-Eicosenoic acid              | 310 | C20H38O2 | 017735-94-3 | 46   |
| 3    |    | cis-10-Heptadecenoic acid           | 268 | C17H32O2 | 029743-97-3 | 41   |
| 4    |    | cis-11-Eicosenoic acid              | 310 | C20H38O2 | 005561-99-9 | 38   |
| 5    |    | Cyclobutanone, 2-(1,1-dimethylet... | 126 | C8H14O   | 004579-31-1 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

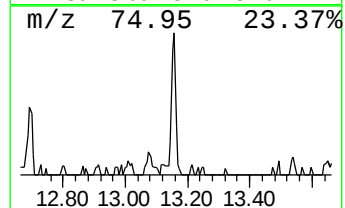
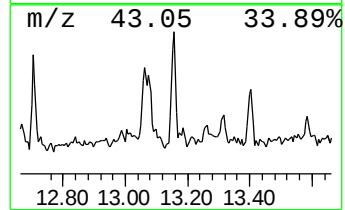
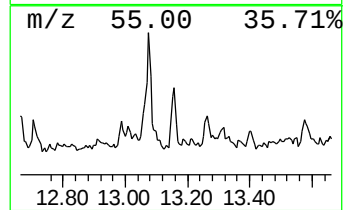
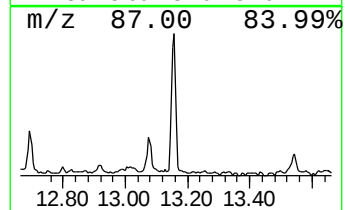
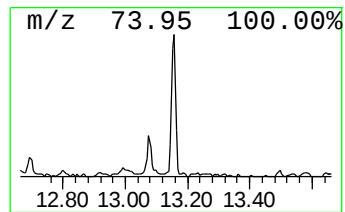
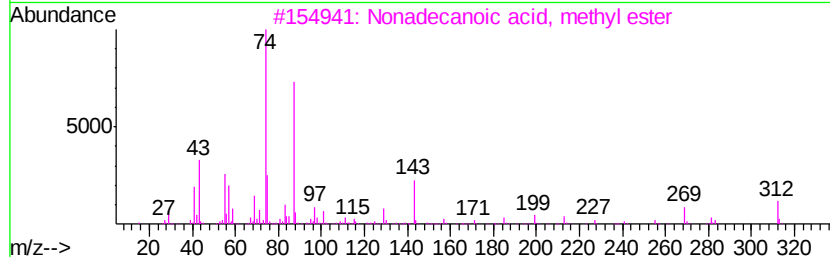
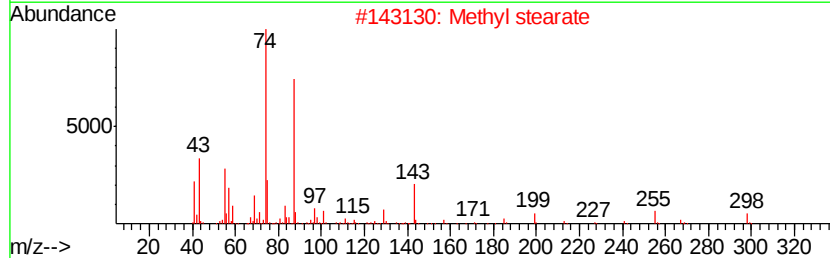
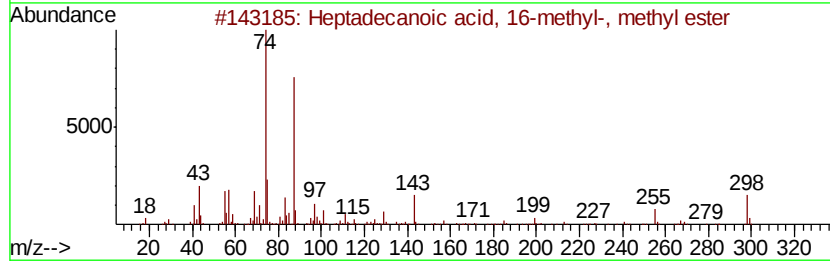
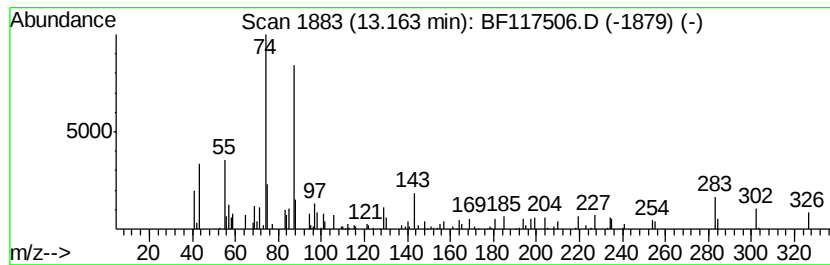
Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Heptadecanoic acid, 16-meth... Concentration Rank 20

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.     |             |      |
|---------|---------|-------------------------------------|------------------|----------|-------------|------|
| 13.16   | 5.19 ng | 130737                              | Chrysene-d12     | 13.89    |             |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1       |         | Heptadecanoic acid, 16-methyl-, ... | 298              | C19H38O2 | 005129-61-3 | 91   |
| 2       |         | Methyl stearate                     | 298              | C19H38O2 | 000112-61-8 | 87   |
| 3       |         | Nonadecanoic acid, methyl ester     | 312              | C20H40O2 | 001731-94-8 | 87   |
| 4       |         | Hexadecanoic acid, methyl ester     | 270              | C17H34O2 | 000112-39-0 | 87   |
| 5       |         | Methyl 9-methyltetradecanoate       | 256              | C16H32O2 | 213617-69-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

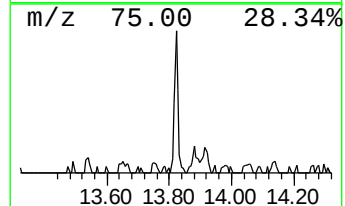
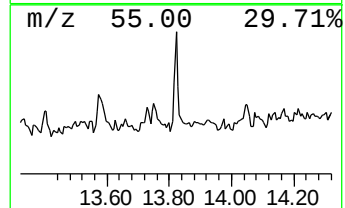
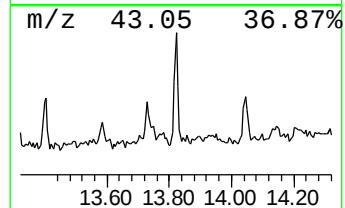
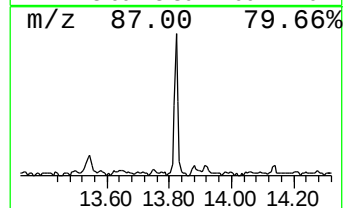
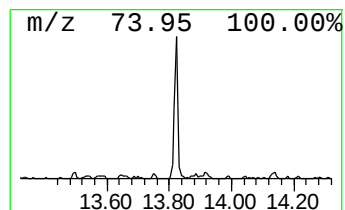
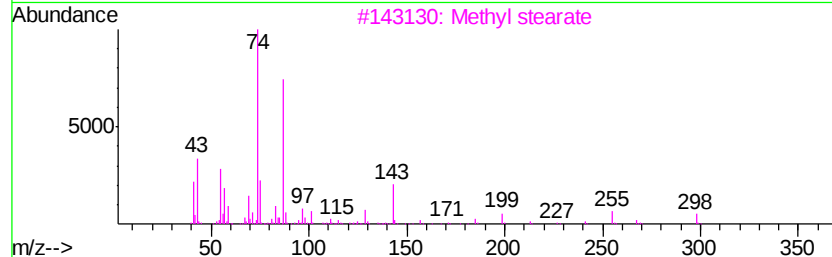
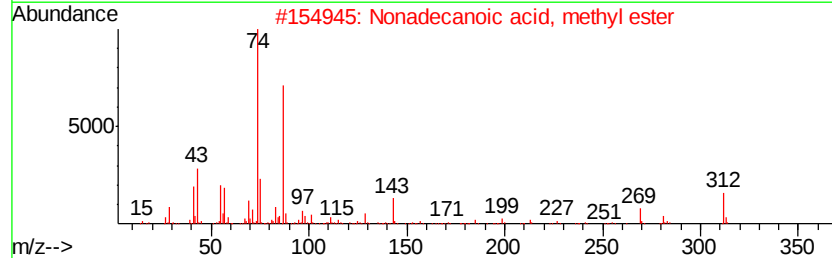
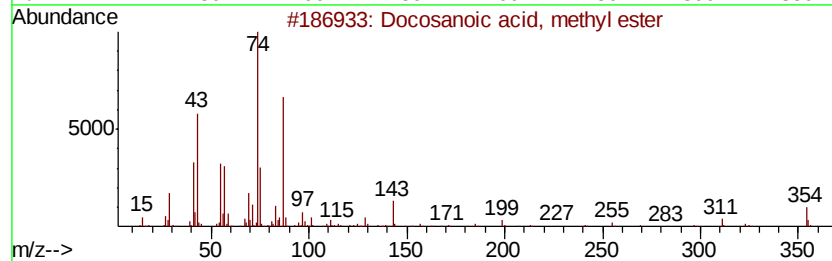
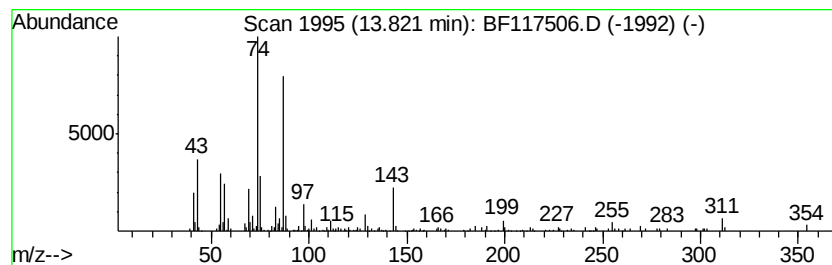
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.82 | 5.47 ng | 137675 | Chrysene-d12     | 13.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Docosanoic acid, methyl ester       | 354 | C23H46O2 | 000929-77-1 | 97   |
| 2    |    | Nonadecanoic acid, methyl ester     | 312 | C20H40O2 | 001731-94-8 | 94   |
| 3    |    | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 93   |
| 4    |    | Octadecanoic acid, 10-methyl-, m... | 312 | C20H40O2 | 002490-19-9 | 91   |
| 5    |    | Hexadecanoic acid, 15-methyl-, m... | 284 | C18H36O2 | 006929-04-0 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

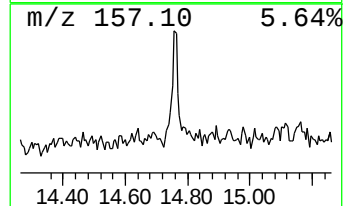
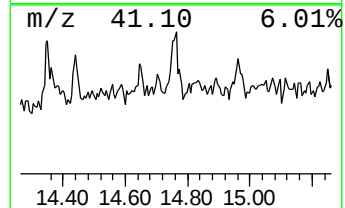
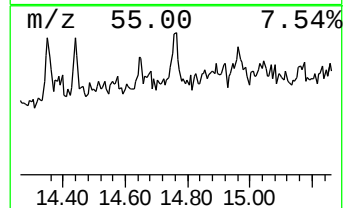
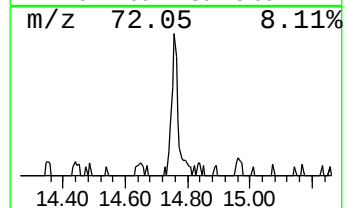
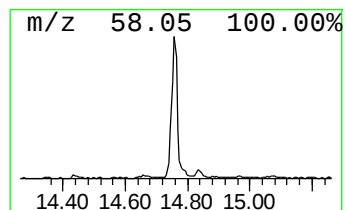
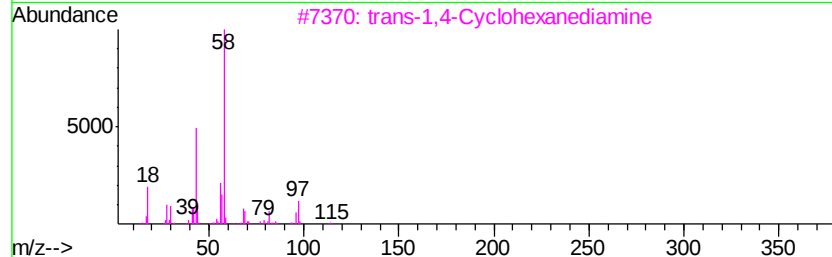
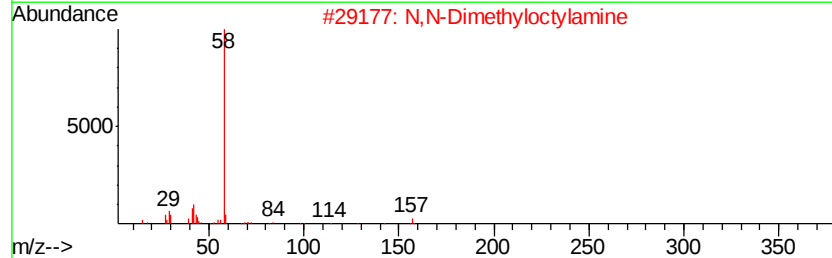
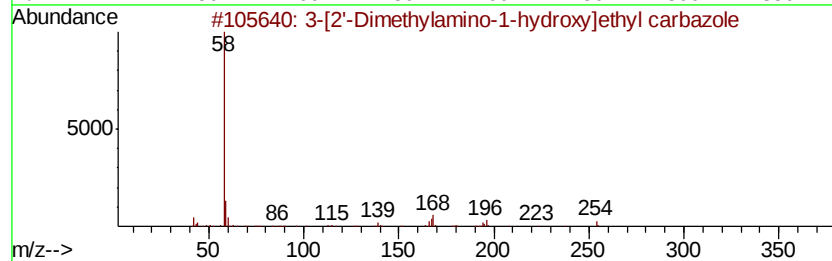
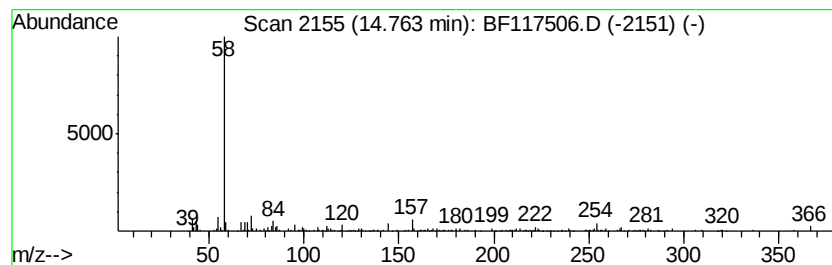
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 3-[2'-Dimethylamino-1-hydro... Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.76 | 13.14 ng | 155453 | Perylene-d12     | 15.33 |

| Hit# | of | Tentative ID                                  | MW  | MolForm     | CAS#         | Qual |
|------|----|-----------------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 3-[2'-Dimethylamino-1-hydroxy]ethyl carbazole | 254 | C16H18N2O   | 1000253-75-3 | 72   |
| 2    |    | N,N-Dimethyloctylamine                        | 157 | C10H23N     | 007378-99-6  | 59   |
| 3    |    | trans-1,4-Cyclohexanediamine                  | 114 | C6H14N2     | 003114-70-3  | 59   |
| 4    |    | Meclofenoxate                                 | 257 | C12H16ClN03 | 000051-68-3  | 59   |
| 5    |    | 4-Pentenylamine, N,N-dimethyl-                | 113 | C7H15N      | 001001-91-8  | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102319\  
Data File : BF117506.D  
Acq On : 24 Oct 2019 1:37  
Operator : JU  
Sample : K5152-10 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

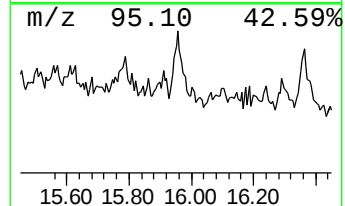
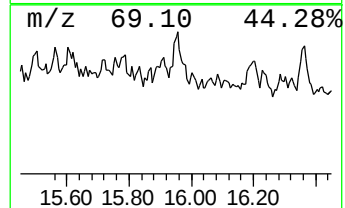
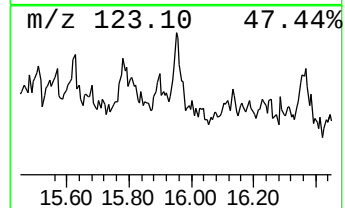
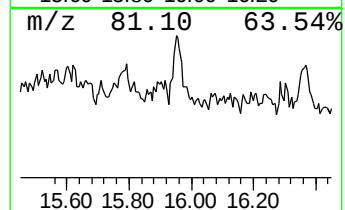
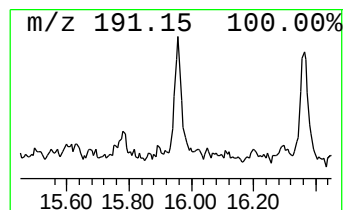
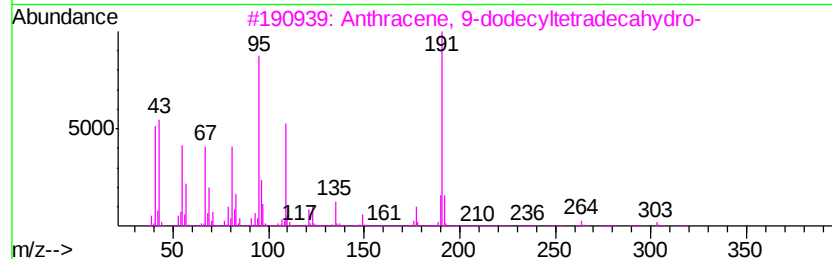
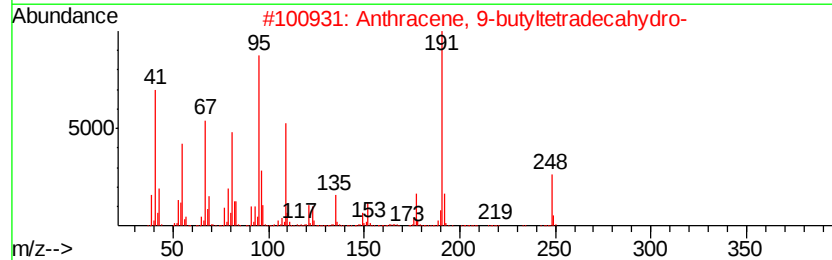
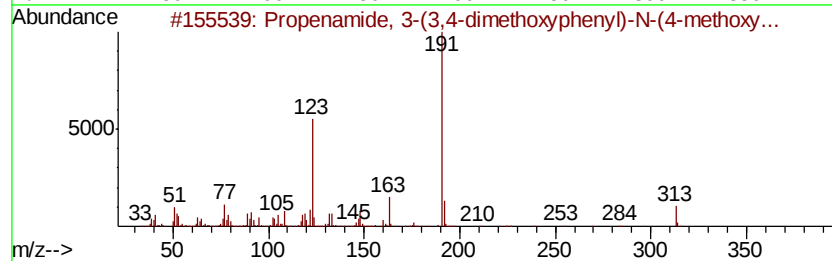
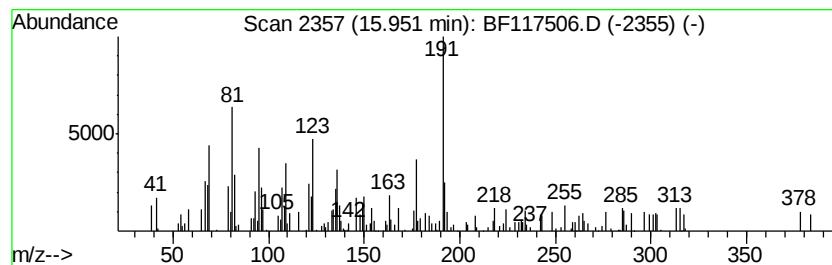
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 unknown15.95 Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.95 | 10.40 ng | 123078 | Perylene-d12     | 15.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Propenamide, 3-(3,4-dimethoxyphe... | 313 | C18H19N04 | 339165-72-9 | 46   |
| 2    |    | Anthracene, 9-butyltetradecahydro-  | 248 | C18H32    | 055133-89-6 | 43   |
| 3    |    | Anthracene, 9-dodecyltetradecahv... | 360 | C26H48    | 055401-75-7 | 43   |
| 4    |    | 2H-1,2,3-Triazole-4-carboxaldehv... | 191 | C9H6FN3O  | 051306-43-5 | 38   |
| 5    |    | Anthracene, 9-butyl-                | 234 | C18H18    | 001498-69-7 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF102319\  
 Data File : BF117506.D  
 Acq On : 24 Oct 2019 1:37  
 Operator : JU  
 Sample : K5152-10 2X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SU-03-102219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.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 |
| unknown6.50          | 6.50  | 37.9    | ng    | 445282   | 1                     | 6.73  | 234938 | 20.0 |
| Dodecane             | 10.19 | 5.4     | ng    | 102883   | 3                     | 9.76  | 383777 | 20.0 |
| Tetradecane          | 10.67 | 10.8    | ng    | 204221   | 4                     | 11.25 | 379487 | 20.0 |
| Tetracosane          | 11.12 | 6.0     | ng    | 113666   | 4                     | 11.25 | 379487 | 20.0 |
| unknown11.15         | 11.15 | 5.2     | ng    | 98869    | 4                     | 11.25 | 379487 | 20.0 |
| Heneicosane          | 11.54 | 5.9     | ng    | 112388   | 4                     | 11.25 | 379487 | 20.0 |
| Hexadecanoic acid... | 11.65 | 103.8   | ng    | 1970280  | 4                     | 11.25 | 379487 | 20.0 |
| Phenanthrene, 1-m... | 11.75 | 5.9     | ng    | 112017   | 4                     | 11.25 | 379487 | 20.0 |
| Anthracene, 2-met... | 11.78 | 10.7    | ng    | 202706   | 4                     | 11.25 | 379487 | 20.0 |
| Anthracene, 1-met... | 11.86 | 7.7     | ng    | 146089   | 4                     | 11.25 | 379487 | 20.0 |
| Pentadecane          | 11.95 | 5.9     | ng    | 112621   | 4                     | 11.25 | 379487 | 20.0 |
| 9,12-Octadecadien... | 12.33 | 238.6   | ng    | 4527910  | 4                     | 11.25 | 379487 | 20.0 |
| 8-Octadecenoic ac... | 12.37 | 390.7   | ng    | 7413170  | 4                     | 11.25 | 379487 | 20.0 |
| unknown13.01         | 13.01 | 6.0     | ng    | 151865   | 5                     | 13.89 | 503745 | 20.0 |
| 9-Octadecenoic ac... | 13.07 | 9.8     | ng    | 246786   | 5                     | 13.89 | 503745 | 20.0 |
| Heptadecanoic aci... | 13.16 | 5.2     | ng    | 130737   | 5                     | 13.89 | 503745 | 20.0 |
| Docosanoic acid, ... | 13.82 | 5.5     | ng    | 137675   | 5                     | 13.89 | 503745 | 20.0 |
| 3-[2'-Dimethylami... | 14.76 | 13.1    | ng    | 155453   | 6                     | 15.33 | 236651 | 20.0 |
| unknown15.95         | 15.95 | 10.4    | ng    | 123078   | 6                     | 15.33 | 236651 | 20.0 |