

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
 Data File : BF117812.D  
 Acq On : 15 Nov 2019 19:30  
 Operator : JU  
 Sample : K5861-14  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EP-11

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-BF111319.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         | 2.181       | 10            | 16          | 24           | rVB      | 816986         | 1139592       | 16.33%          | 1.854%        |
| 2         | 3.416       | 223           | 226         | 239          | rBV      | 30365          | 71640         | 1.03%           | 0.117%        |
| 3         | 5.128       | 509           | 517         | 525          | rBV      | 882086         | 1156781       | 16.58%          | 1.882%        |
| 4         | 5.528       | 580           | 585         | 589          | rBV      | 4802372        | 5196711       | 74.49%          | 8.455%        |
| 5         | 6.539       | 751           | 757         | 760          | rBV      | 5295949        | 5481198       | 78.57%          | 8.917%        |
| 6         | 6.686       | 777           | 782         | 785          | rBV      | 5837713        | 5552328       | 79.59%          | 9.033%        |
| 7         | 6.916       | 818           | 821         | 825          | rVB      | 1406937        | 1207634       | 17.31%          | 1.965%        |
| 8         | 7.075       | 844           | 848         | 852          | rBV      | 5060515        | 4340663       | 62.22%          | 7.062%        |
| 9         | 7.481       | 912           | 917         | 920          | rBV      | 3657960        | 3463531       | 49.65%          | 5.635%        |
| 10        | 8.204       | 1036          | 1040        | 1044         | rBV      | 1915275        | 1559818       | 22.36%          | 2.538%        |
| 11        | 9.286       | 1219          | 1224        | 1227         | rBV      | 7034500        | 6361026       | 91.18%          | 10.349%       |
| 12        | 9.398       | 1238          | 1243        | 1251         | rVB      | 243322         | 227268        | 3.26%           | 0.370%        |
| 13        | 9.674       | 1286          | 1290        | 1294         | rBV      | 842106         | 664972        | 9.53%           | 1.082%        |
| 14        | 9.969       | 1336          | 1340        | 1343         | rBV      | 2233253        | 1817442       | 26.05%          | 2.957%        |
| 15        | 10.551      | 1435          | 1439        | 1448         | rBV      | 84296          | 132314        | 1.90%           | 0.215%        |
| 16        | 10.757      | 1469          | 1474        | 1477         | rBV      | 4888292        | 4412748       | 63.25%          | 7.179%        |
| 17        | 11.457      | 1589          | 1593        | 1597         | rBV      | 2328036        | 1963878       | 28.15%          | 3.195%        |
| 18        | 11.986      | 1679          | 1683        | 1698         | rBV2     | 843047         | 975681        | 13.99%          | 1.587%        |
| 19        | 12.504      | 1768          | 1771        | 1787         | rBV      | 716086         | 1383433       | 19.83%          | 2.251%        |
| 20        | 12.704      | 1797          | 1805        | 1808         | rBV      | 221450         | 259856        | 3.72%           | 0.423%        |
| 21        | 12.727      | 1808          | 1809        | 1813         | rVB      | 105166         | 86269         | 1.24%           | 0.140%        |
| 22        | 12.768      | 1813          | 1816        | 1828         | rBV      | 252384         | 314028        | 4.50%           | 0.511%        |
| 23        | 12.868      | 1828          | 1833        | 1836         | rVB2     | 130822         | 114165        | 1.64%           | 0.186%        |
| 24        | 12.927      | 1839          | 1843        | 1852         | rVV2     | 200030         | 284928        | 4.08%           | 0.464%        |
| 25        | 13.051      | 1859          | 1864        | 1867         | rBV      | 7368935        | 6976425       | 100.00%         | 11.350%       |
| 26        | 13.280      | 1897          | 1903        | 1915         | rBV      | 231553         | 513536        | 7.36%           | 0.835%        |
| 27        | 13.904      | 2004          | 2009        | 2012         | rBV      | 115916         | 138256        | 1.98%           | 0.225%        |
| 28        | 13.951      | 2012          | 2017        | 2037         | rVV3     | 526178         | 1460169       | 20.93%          | 2.376%        |
| 29        | 14.104      | 2037          | 2043        | 2047         | rVB      | 1899483        | 1803012       | 25.84%          | 2.933%        |
| 30        | 14.574      | 2119          | 2123        | 2126         | rBV      | 94450          | 96995         | 1.39%           | 0.158%        |
| 31        | 14.615      | 2126          | 2130        | 2139         | rVB10    | 35298          | 99188         | 1.42%           | 0.161%        |
| 32        | 14.892      | 2174          | 2177        | 2182         | rVB      | 85280          | 94135         | 1.35%           | 0.153%        |
| 33        | 15.251      | 2234          | 2238        | 2245         | rVB      | 109153         | 126167        | 1.81%           | 0.205%        |
| 34        | 15.615      | 2294          | 2300        | 2304         | rBV      | 1357206        | 1765112       | 25.30%          | 2.872%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

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-BF111319.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |     |       |        |       |        |
|----|--------|------|------|------|-----|-------|--------|-------|--------|
| 35 | 15.651 | 2304 | 2306 | 2311 | rVB | 92828 | 105646 | 1.51% | 0.172% |
| 36 | 16.109 | 2380 | 2384 | 2389 | rBV | 76947 | 119269 | 1.71% | 0.194% |

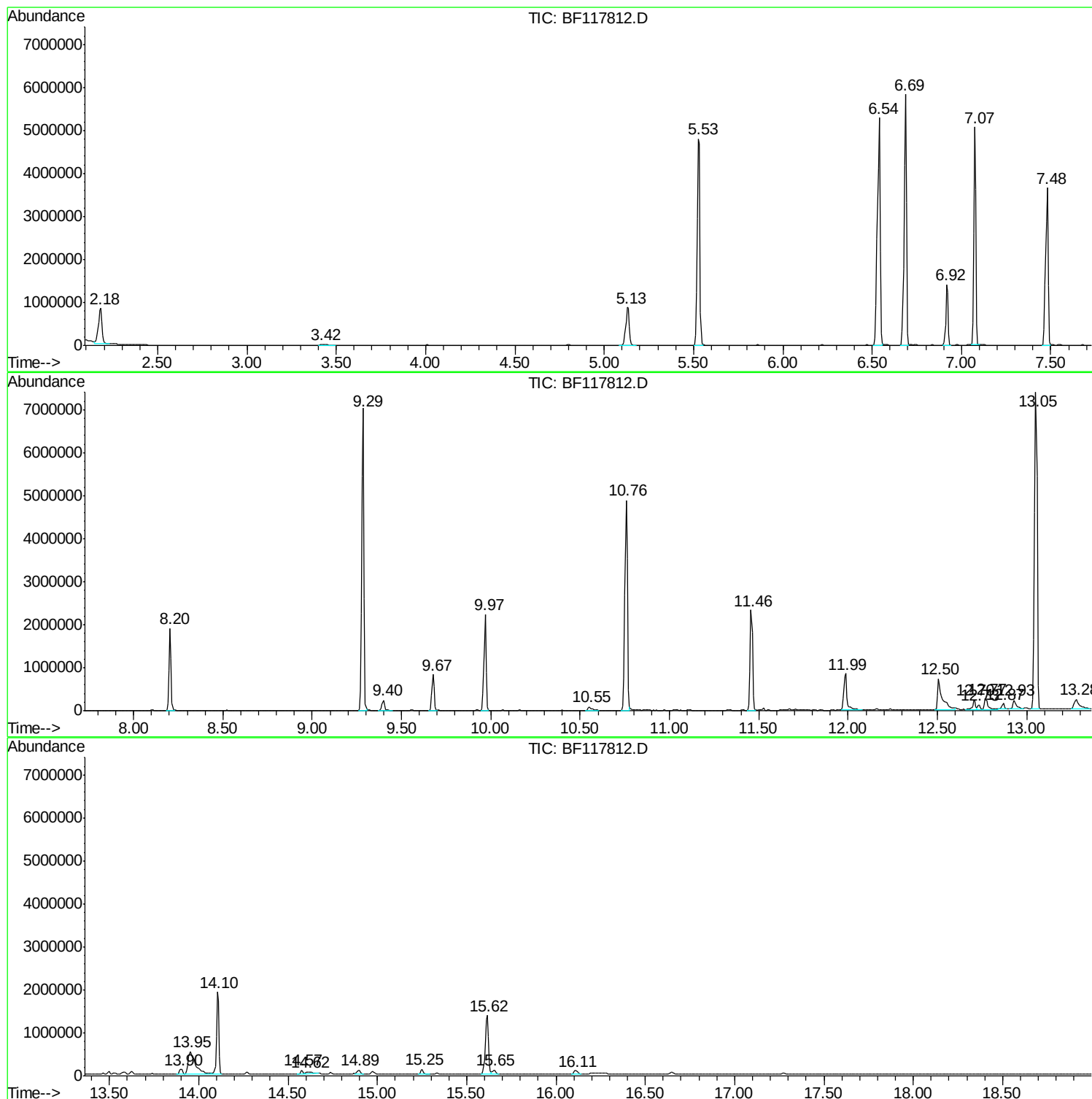
Sum of corrected areas: 61465814

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.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\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

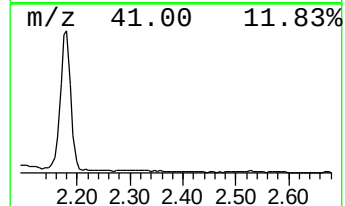
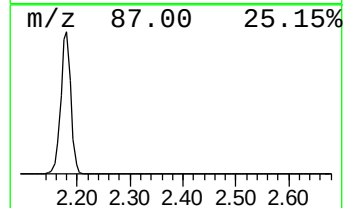
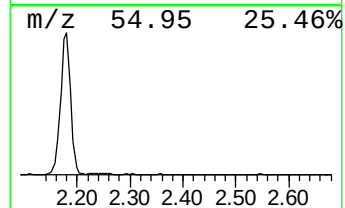
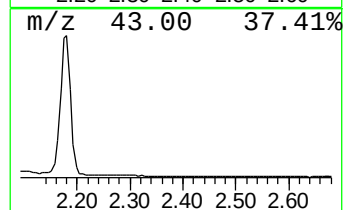
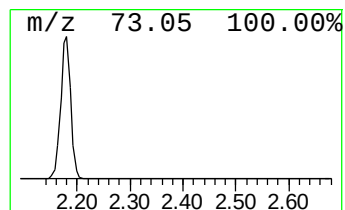
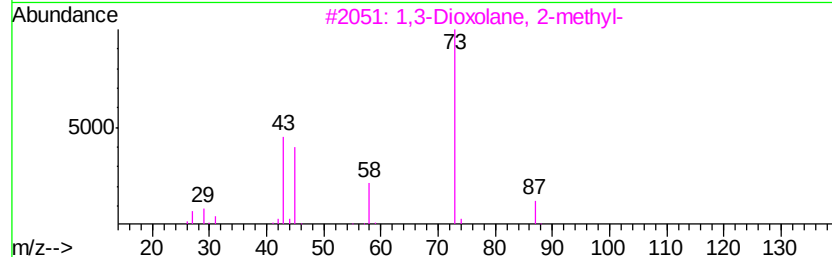
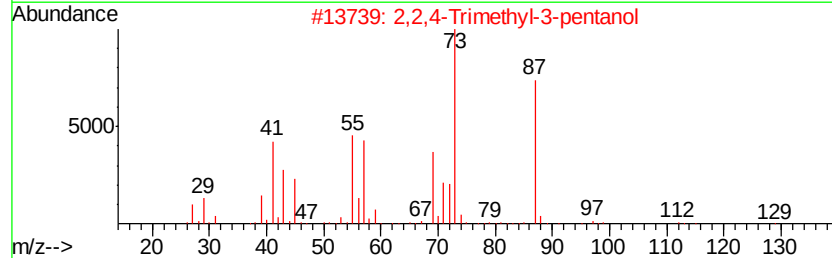
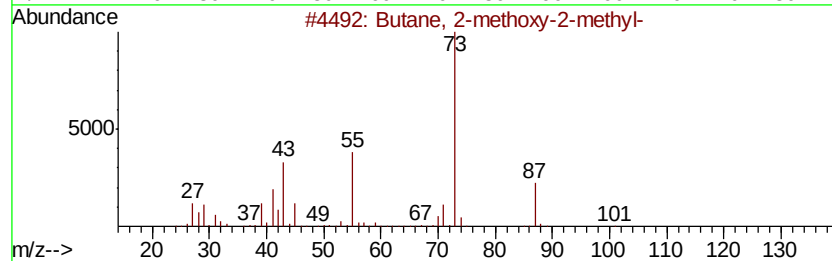
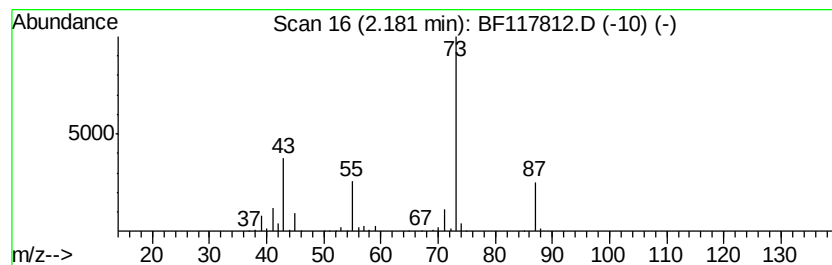
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 2.18 | 18.87 ng | 1139590 | 1,4-Dichlorobenzene-d4 | 6.92 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 4    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 5    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

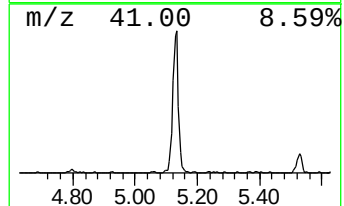
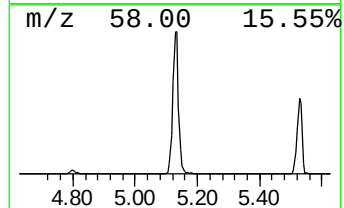
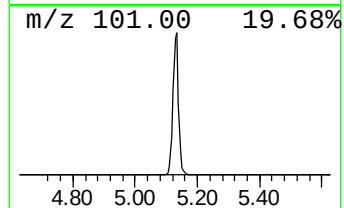
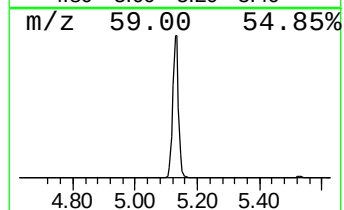
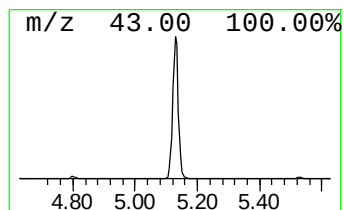
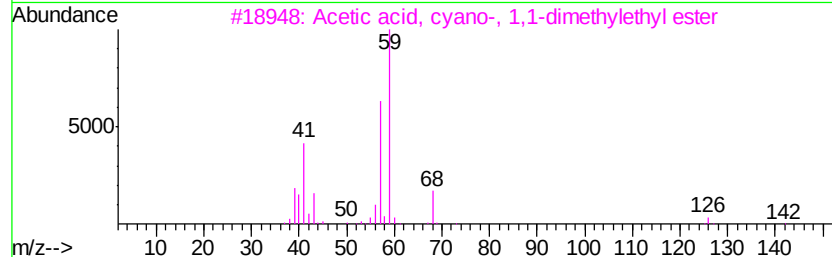
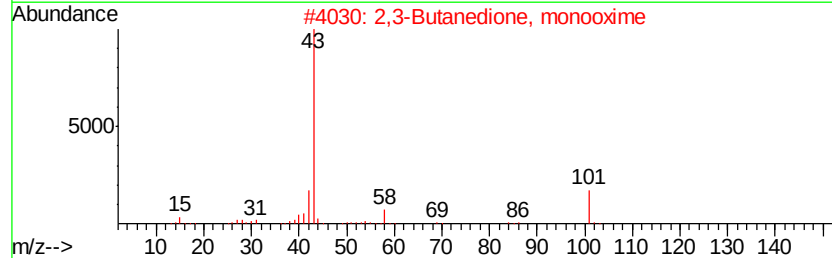
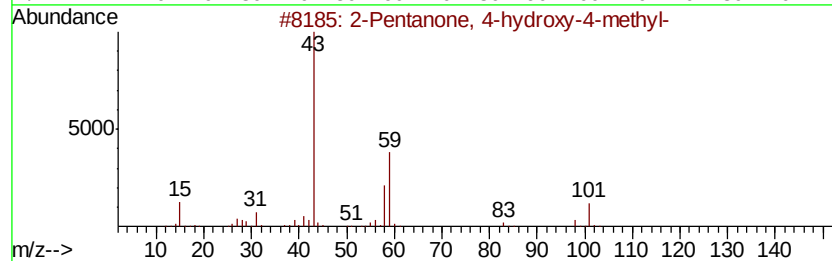
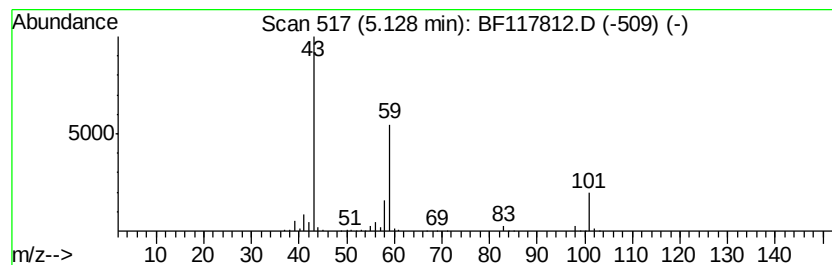
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown5.13 Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.13 | 19.16 ng | 1156780 | 1,4-Dichlorobenzene-d4 | 6.92 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 47   |
| 2    |    |   | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 25   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 4    |    |   | Morpholine, 4-methyl-                | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    |   | Butane, 1-ethoxy-                    | 102 | C6H14O   | 000628-81-9 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

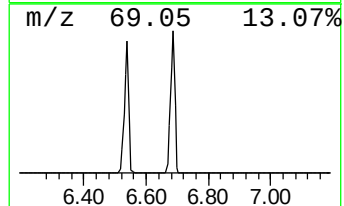
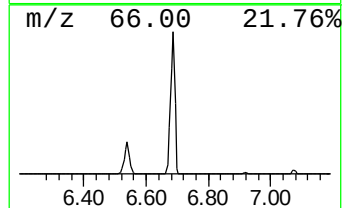
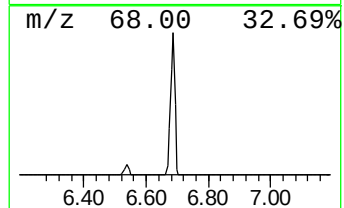
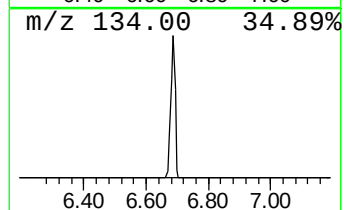
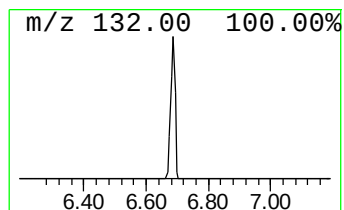
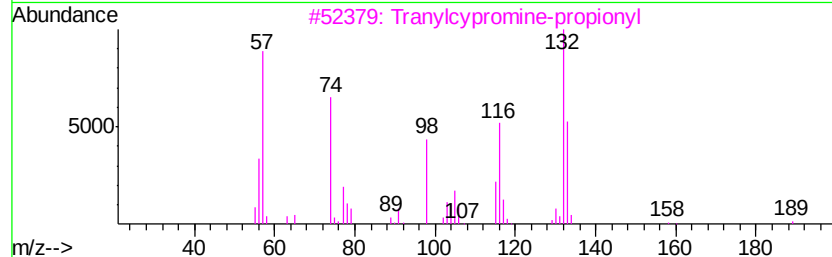
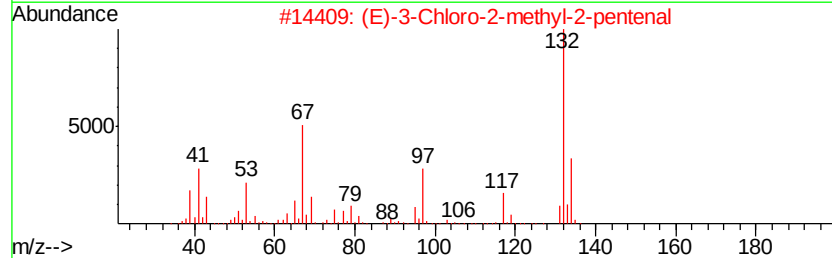
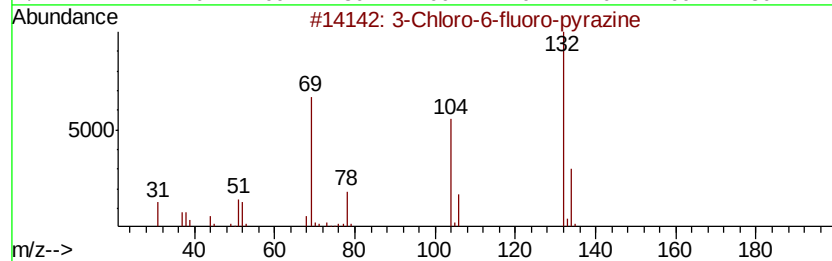
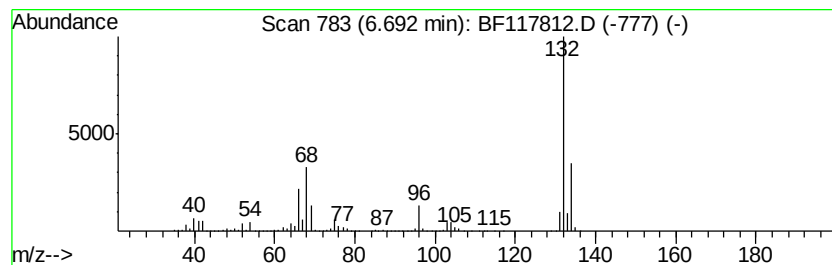
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown6.69 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.69 | 91.95 ng | 5552330 | 1,4-Dichlorobenzene-d4 | 6.92 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | Benzene, 1-ethenyl-3,5-dimethyl-    | 132 | C10H12    | 005379-20-4  | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

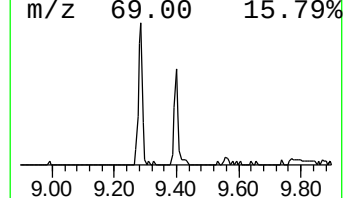
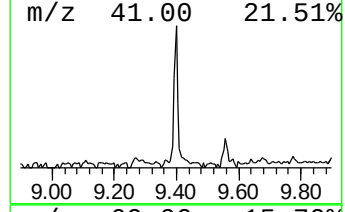
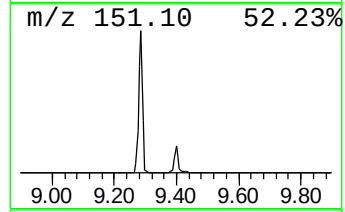
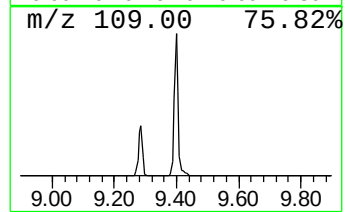
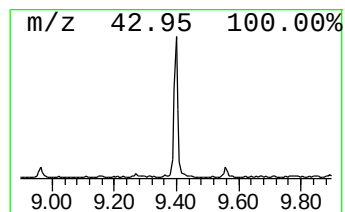
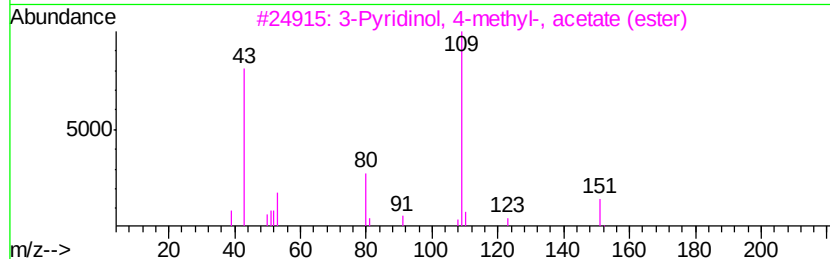
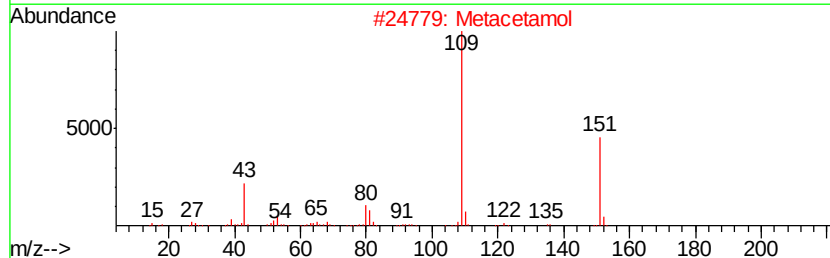
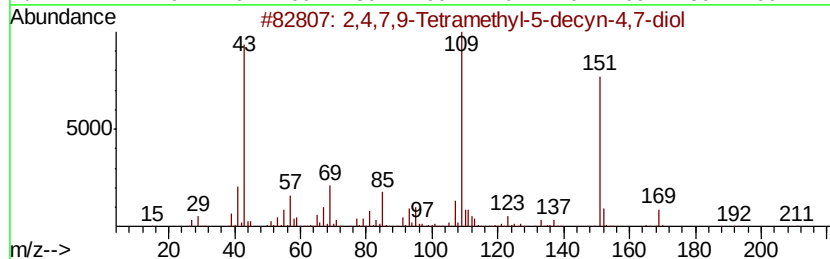
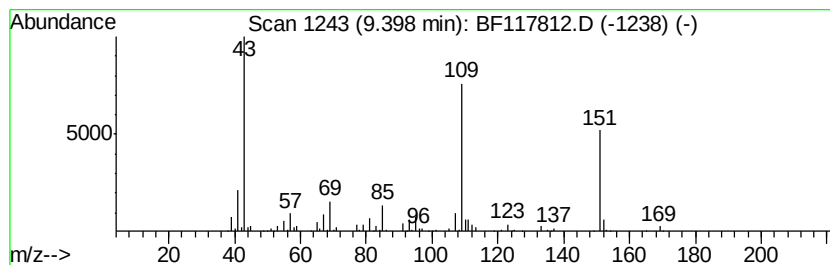
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 12

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.40 | 2.50 ng | 227268 | Acenaphthene-d10 | 9.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226 | C14H26O2 | 000126-86-3 | 91   |
| 2    |    | Metacetamol                         | 151 | C8H9NO2  | 000621-42-1 | 59   |
| 3    |    | 3-Pyridinol, 4-methyl-, acetate ... | 151 | C8H9NO2  | 001006-96-8 | 59   |
| 4    |    | m-Isopropoxyaniline                 | 151 | C9H13NO  | 041406-00-2 | 45   |
| 5    |    | Cyclohexanol, 3,3,5-trimethyl-      | 142 | C9H18O   | 000116-02-9 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
 Data File : BF117812.D  
 Acq On : 15 Nov 2019 19:30  
 Operator : JU  
 Sample : K5861-14  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

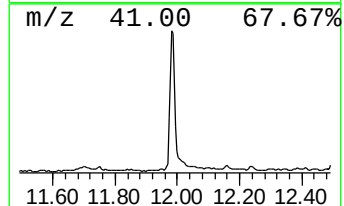
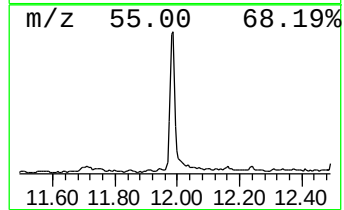
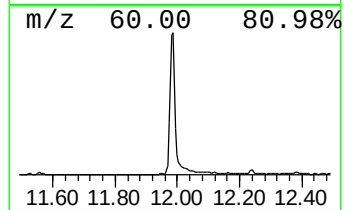
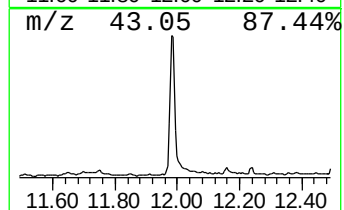
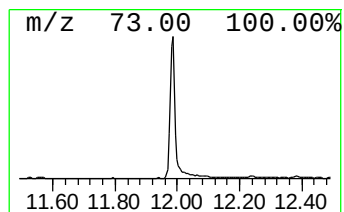
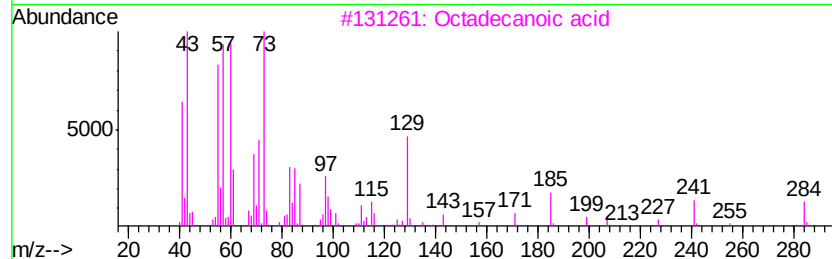
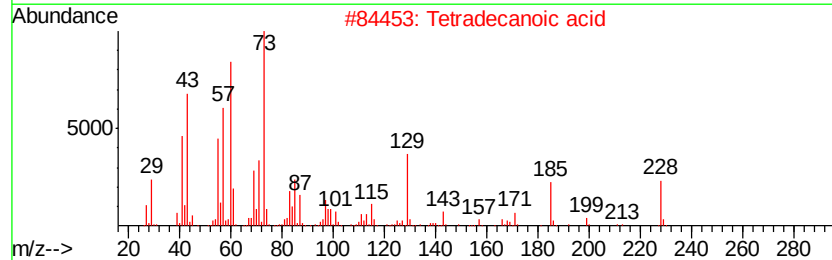
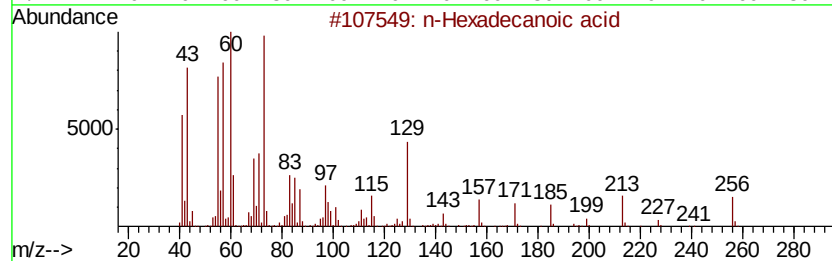
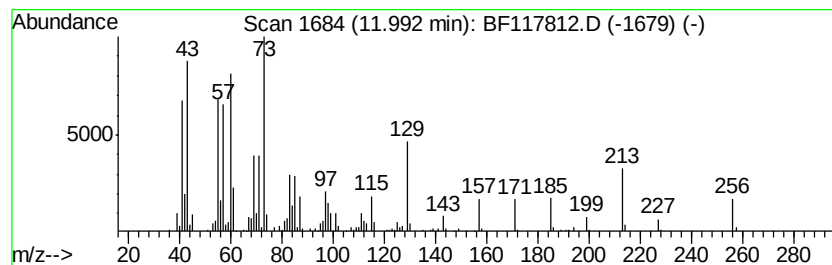
Instrument :  
 BNA\_F  
 ClientSampleId :  
 EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
 Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

| R.T.    | EstConc | Area                | Relative to ISTD |          | R.T.        |      |
|---------|---------|---------------------|------------------|----------|-------------|------|
| 11.99   | 9.94 ng | 975681              | Phenanthrene-d10 |          | 11.46       |      |
| Hit# of | 5       | Tentative ID        | MW               | MolForm  | CAS#        | Qual |
| 1       |         | n-Hexadecanoic acid | 256              | C16H32O2 | 000057-10-3 | 99   |
| 2       |         | Tetradecanoic acid  | 228              | C14H28O2 | 000544-63-8 | 94   |
| 3       |         | Octadecanoic acid   | 284              | C18H36O2 | 000057-11-4 | 93   |
| 4       |         | Pentadecanoic acid  | 242              | C15H30O2 | 001002-84-2 | 91   |
| 5       |         | Tridecanoic acid    | 214              | C13H26O2 | 000638-53-9 | 90   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

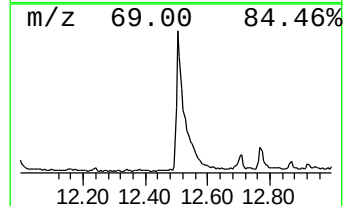
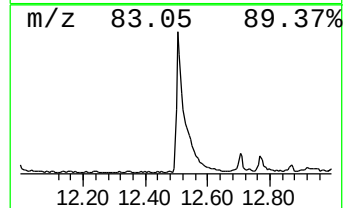
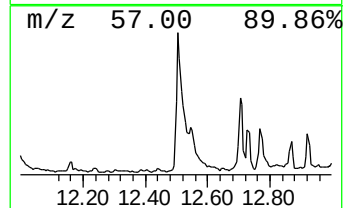
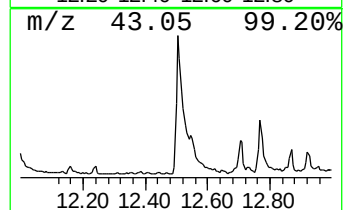
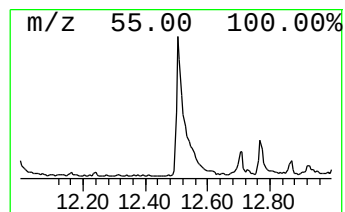
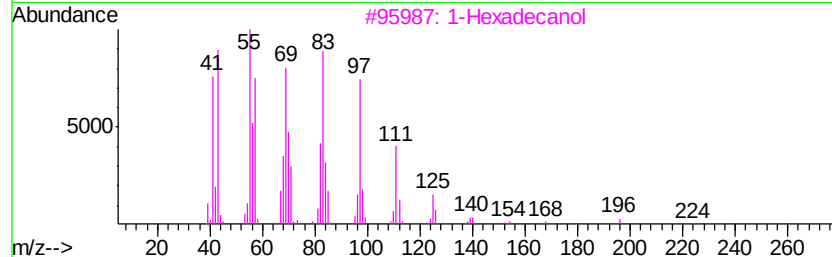
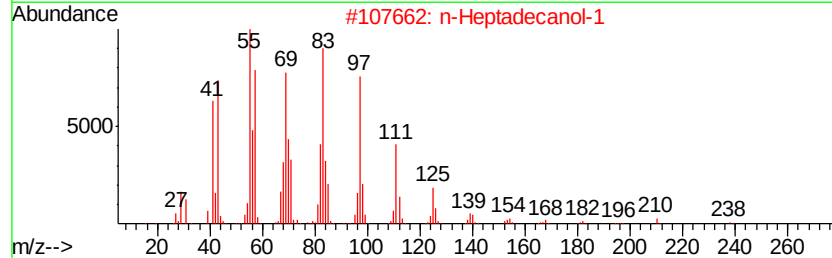
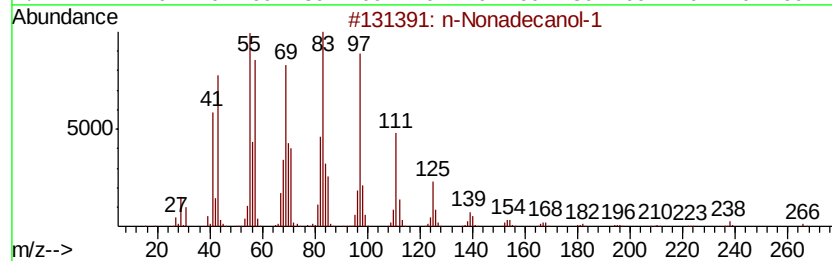
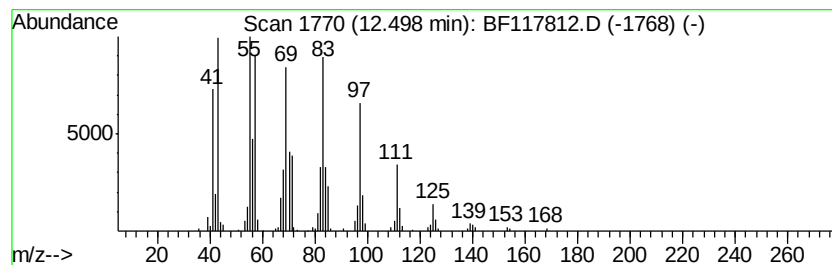
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 n-Nonadecanol-1 Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD |  | R.T.  |
|-------|----------|---------|------------------|--|-------|
| 12.50 | 14.09 ng | 1383430 | Phenanthrene-d10 |  | 11.46 |

| Hit# | of                                 | 5 | Tentative ID | MW  | MolForm    | CAS#        | Qual |
|------|------------------------------------|---|--------------|-----|------------|-------------|------|
| 1    | n-Nonadecanol-1                    |   |              | 284 | C19H40O    | 001454-84-8 | 95   |
| 2    | n-Heptadecanol-1                   |   |              | 256 | C17H36O    | 001454-85-9 | 95   |
| 3    | 1-Hexadecanol                      |   |              | 242 | C16H34O    | 036653-82-4 | 94   |
| 4    | Bromoacetic acid, pentadecyl ester |   |              | 348 | C17H33BrO2 | 131143-01-6 | 94   |
| 5    | n-Pentadecanol                     |   |              | 228 | C15H32O    | 000629-76-5 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

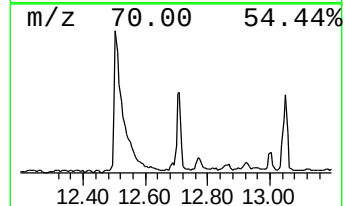
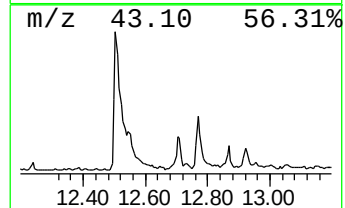
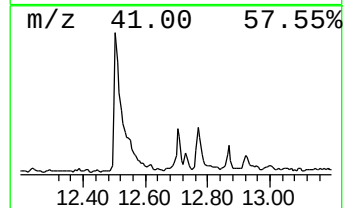
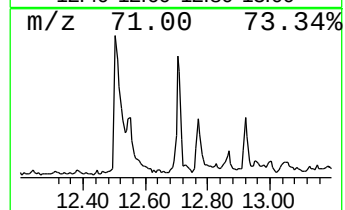
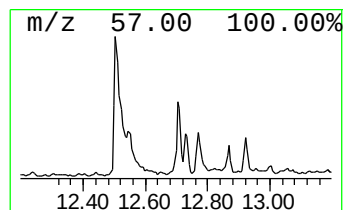
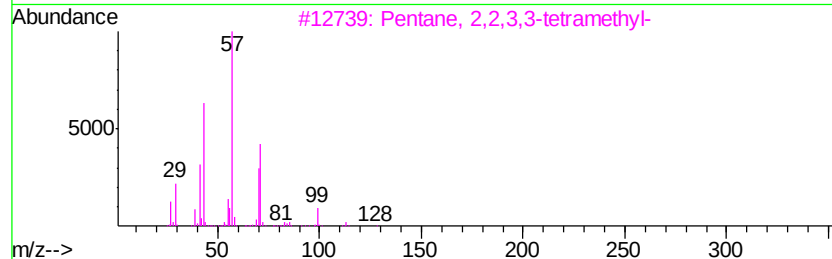
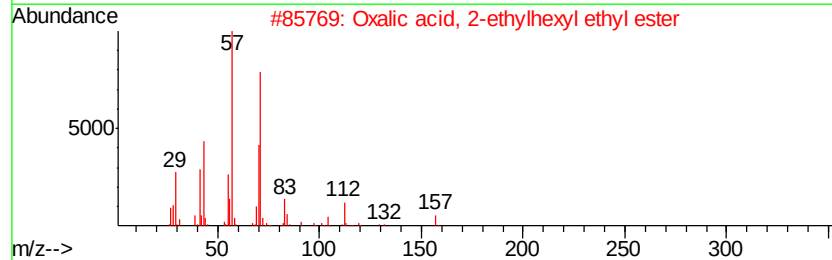
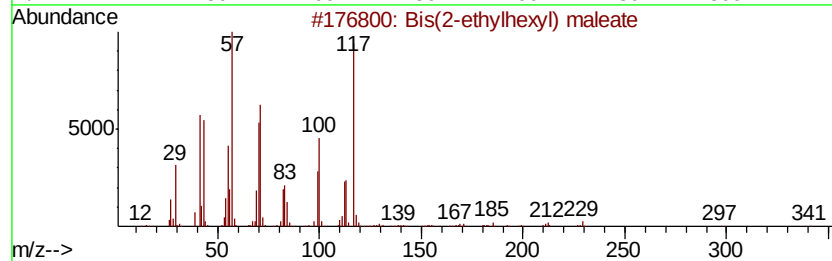
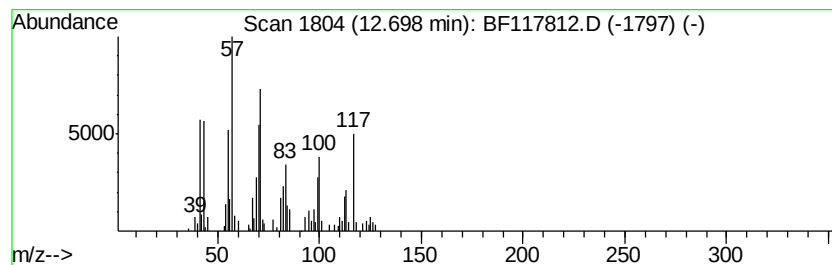
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Bis(2-ethylhexyl) maleate Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.70 | 2.65 ng | 259856 | Phenanthrene-d10 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Bis(2-ethylhexyl) maleate           | 340 | C20H36O4 | 000142-16-5  | 91   |
| 2    |    | Oxalic acid, 2-ethylhexyl ethyl ... | 230 | C12H22O4 | 1000309-38-4 | 43   |
| 3    |    | Pentane, 2,2,3,3-tetramethyl-       | 128 | C9H20    | 007154-79-2  | 43   |
| 4    |    | Oxalic acid, 2-ethylhexyl octyl ... | 314 | C18H34O4 | 1000309-39-1 | 38   |
| 5    |    | 3-Hexanone                          | 100 | C6H12O   | 000589-38-8  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

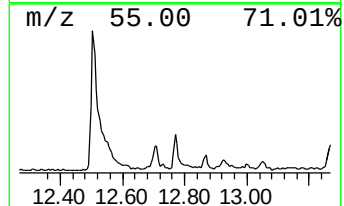
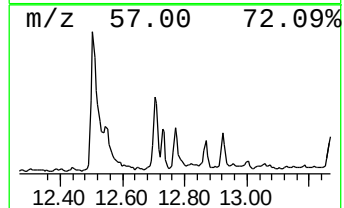
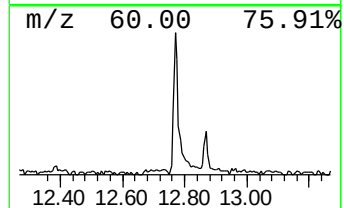
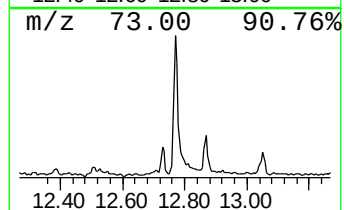
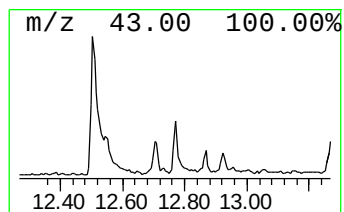
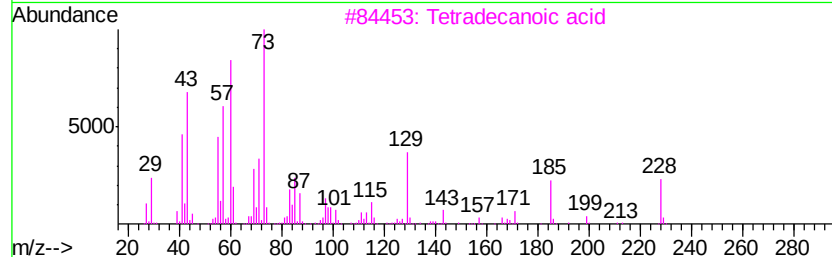
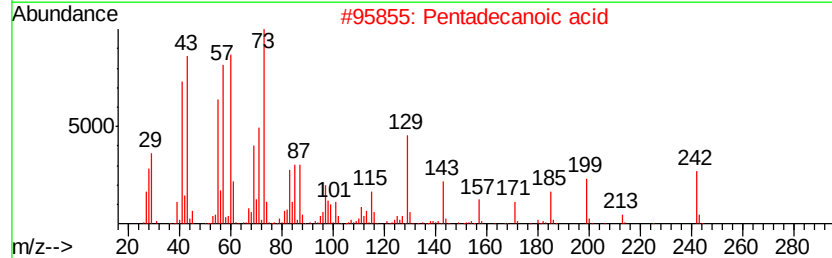
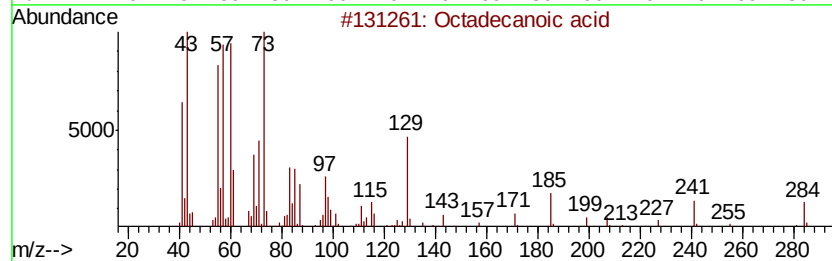
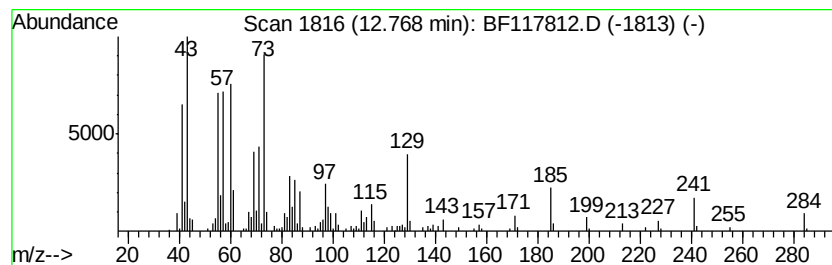
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Octadecanoic acid Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.77 | 3.20 ng | 314028 | Phenanthrene-d10 | 11.46 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|-----------|-------------------------------------|-----|----------|-------------|------|
| 1         | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2         | Pentadecanoic acid                  | 242 | C15H30O2 | 001002-84-2 | 89   |
| 3         | Tetradecanoic acid                  | 228 | C14H28O2 | 000544-63-8 | 74   |
| 4         | Octadecanoic acid, 2-(2-hydroxye... | 372 | C22H44O4 | 000106-11-6 | 72   |
| 5         | n-Decanoic acid                     | 172 | C10H20O2 | 000334-48-5 | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

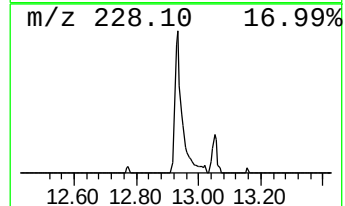
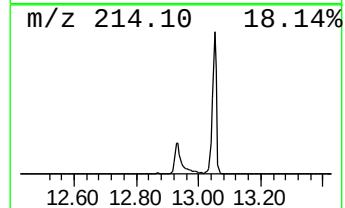
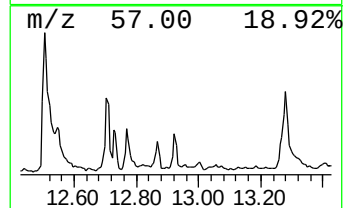
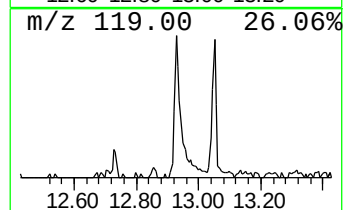
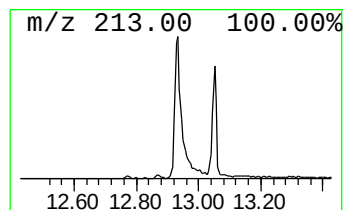
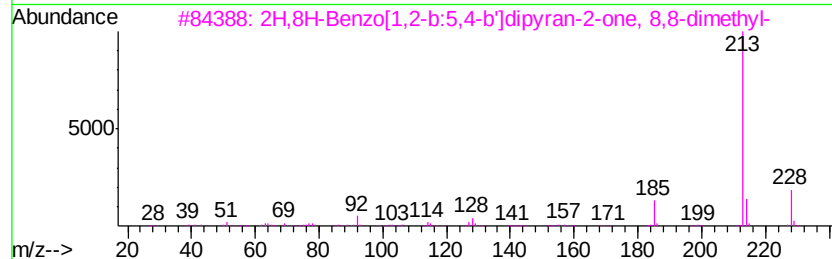
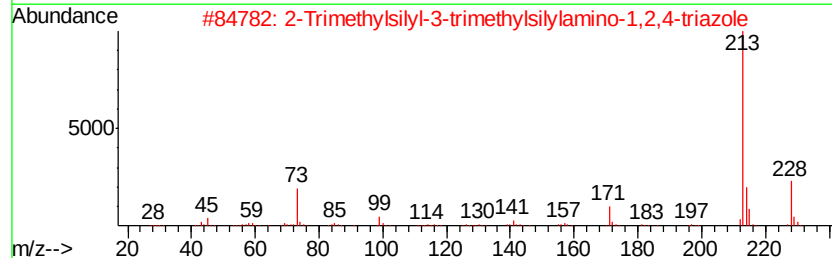
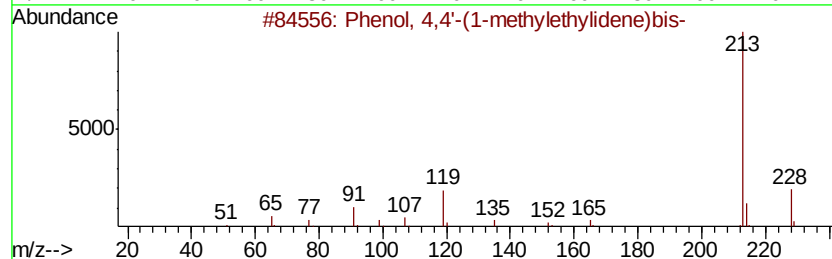
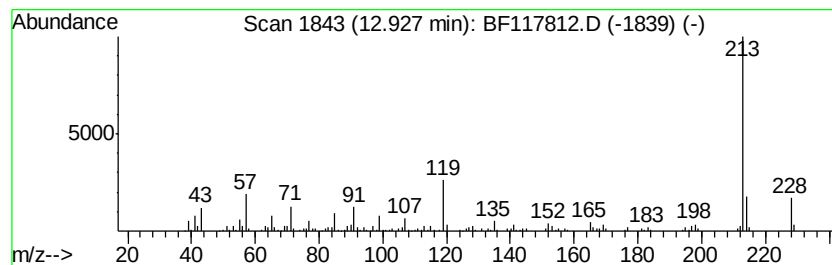
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.93 | 3.16 ng | 284928 | Chrysene-d12     | 14.10 |

| Hit# | of | 5 | Tentative ID                                     | MW  | MolForm    | CAS#         | Qual |
|------|----|---|--------------------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Phenol, 4,4'-(1-methylethylidene)bis-            | 228 | C15H16O2   | 000080-05-7  | 96   |
| 2    |    |   | 2-Trimethylsilyl-3-trimethylsilyl-1,2,4-triazole | 228 | C8H20N4Si2 | 1000373-05-5 | 64   |
| 3    |    |   | 2H,8H-Benzo[1,2-b:5,4-b']dipyrans                | 228 | C14H12O3   | 000553-19-5  | 58   |
| 4    |    |   | Carbanilic acid, p-phenyl-                       | 213 | C13H11NO2  | 004474-53-7  | 58   |
| 5    |    |   | 2H,8H-Benzo[1,2-b:3,4-b']dipyrans                | 228 | C14H12O3   | 000523-59-1  | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

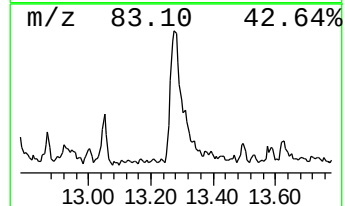
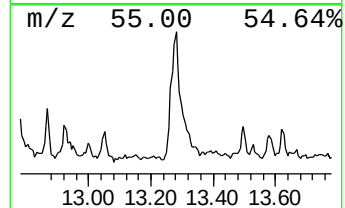
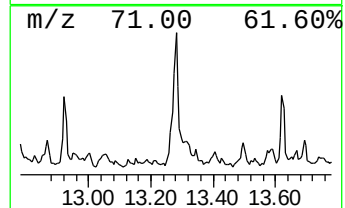
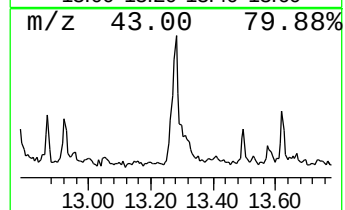
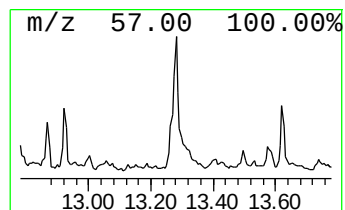
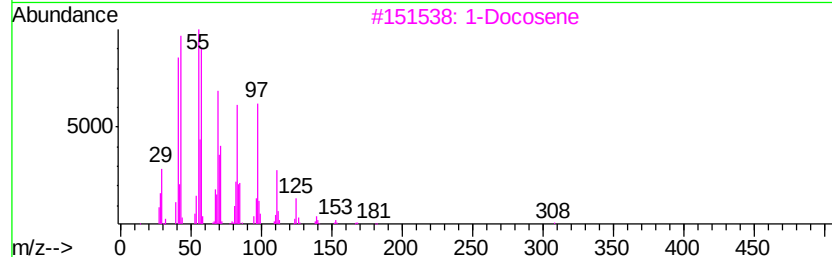
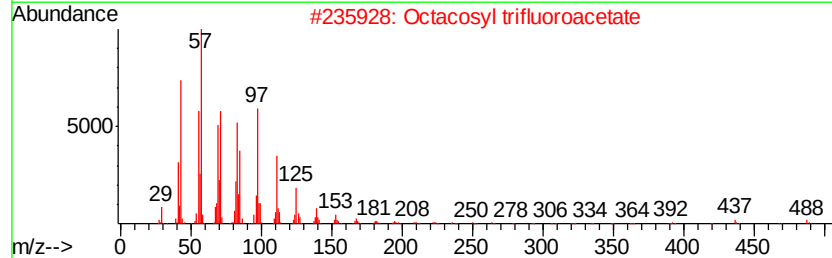
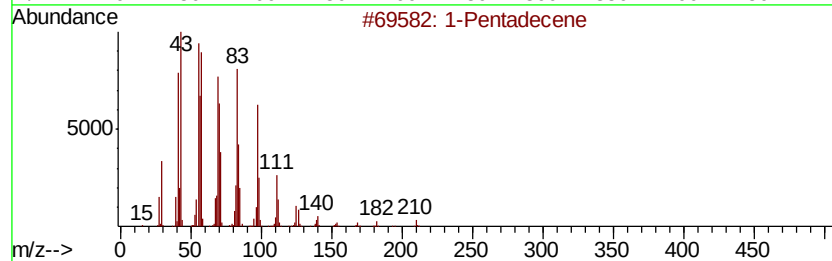
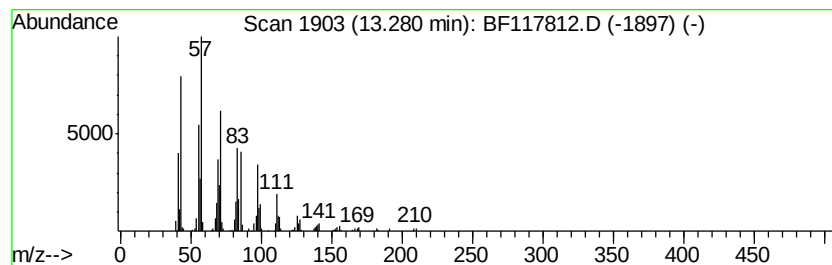
Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 1-Pentadecene Concentration Rank 8

| R.T.    | EstConc                             | Area            | Relative to ISTD | R.T.      |
|---------|-------------------------------------|-----------------|------------------|-----------|
| 13.28   | 5.70 ng                             | 513536          | Chrysene-d12     | 14.10     |
| Hit# of | 5                                   | Tentative ID    | MW MolForm       | CAS# Qual |
| 1       | 1-Pentadecene                       | 210 C15H30      | 013360-61-7      | 92        |
| 2       | Octacosyl trifluoroacetate          | 506 C30H57F3O2  | 1000351-74-9     | 83        |
| 3       | 1-Docosene                          | 308 C22H44      | 001599-67-3      | 70        |
| 4       | Heptadecyl trifluoroacetate         | 352 C19H35F3O2  | 1000351-87-0     | 64        |
| 5       | Trichloroacetic acid, pentadecyl... | 372 C17H31Cl3O2 | 074339-53-0      | 64        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

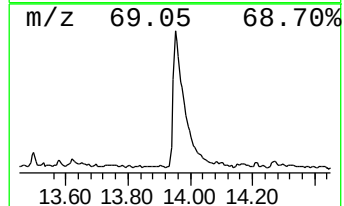
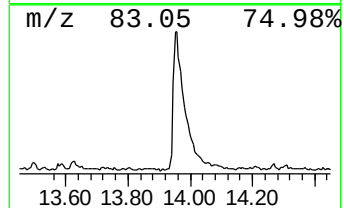
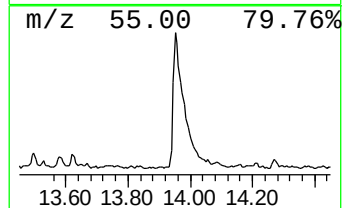
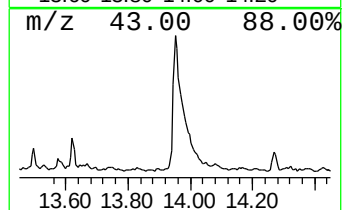
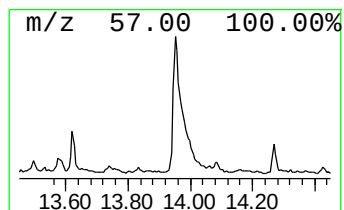
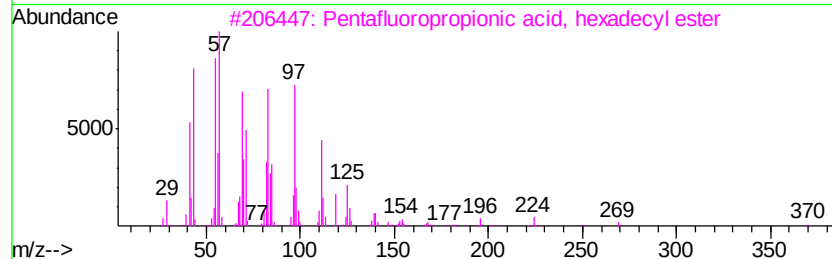
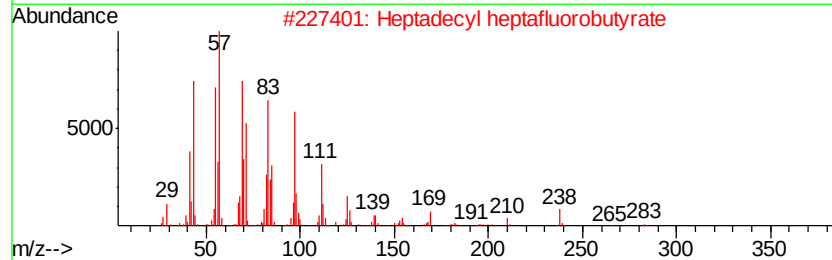
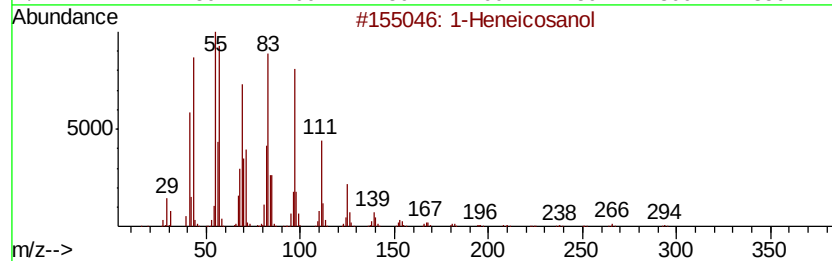
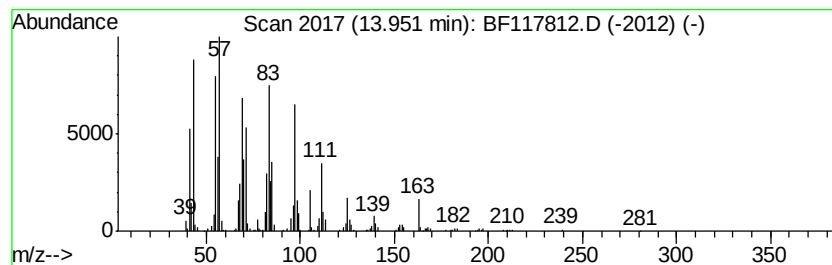
Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 1-Heneicosanol Concentration Rank 5

| R.T.      | EstConc                                    | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------------------|---------|------------------|-------------|------|
| 13.95     | 16.20 ng                                   | 1460170 | Chrysene-d12     | 14.10       |      |
| Hit# of 5 | Tentative ID                               | MW      | MolForm          | CAS#        | Qual |
| 1         | 1-Heneicosanol                             | 312     | C21H44O          | 015594-90-8 | 93   |
| 2         | Heptadecyl heptafluorobutyrate             | 452     | C21H35F7O2       | 959085-66-6 | 91   |
| 3         | Pentafluoropropionic acid, hexadecyl ester | 388     | C19H33F5O2       | 006222-07-7 | 91   |
| 4         | Trichloroacetic acid, hexadecyl ester      | 386     | C18H33Cl3O2      | 074339-54-1 | 91   |
| 5         | Heptafluorobutyric acid, pentadecyl ester  | 424     | C19H31F7O2       | 959261-23-5 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF111519\  
Data File : BF117812.D  
Acq On : 15 Nov 2019 19:30  
Operator : JU  
Sample : K5861-14  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF111319.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 |
| Butane, 2-methoxy... | 2.18  | 18.9    | ng    | 1139590  | 1                     | 6.92  | 1207630 | 20.0 |
| unknown5.13          | 5.13  | 19.2    | ng    | 1156780  | 1                     | 6.92  | 1207630 | 20.0 |
| unknown6.69          | 6.69  | 92.0    | ng    | 5552330  | 1                     | 6.92  | 1207630 | 20.0 |
| 2,4,7,9-Tetrameth... | 9.40  | 2.5     | ng    | 227268   | 3                     | 9.97  | 1817440 | 20.0 |
| n-Hexadecanoic acid  | 11.99 | 9.9     | ng    | 975681   | 4                     | 11.46 | 1963880 | 20.0 |
| n-Nonadecanol-1      | 12.50 | 14.1    | ng    | 1383430  | 4                     | 11.46 | 1963880 | 20.0 |
| Bis(2-ethylhexyl)... | 12.70 | 2.6     | ng    | 259856   | 4                     | 11.46 | 1963880 | 20.0 |
| Octadecanoic acid    | 12.77 | 3.2     | ng    | 314028   | 4                     | 11.46 | 1963880 | 20.0 |
| Phenol, 4,4'-(1-m... | 12.93 | 3.2     | ng    | 284928   | 5                     | 14.10 | 1803010 | 20.0 |
| 1-Pentadecene        | 13.28 | 5.7     | ng    | 513536   | 5                     | 14.10 | 1803010 | 20.0 |
| 1-Heneicosanol       | 13.95 | 16.2    | ng    | 1460170  | 5                     | 14.10 | 1803010 | 20.0 |