

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF012420\  
 Data File : BF118699.D  
 Acq On : 24 Jan 2020 21:35  
 Operator : CG/JU  
 Sample : L1242-12  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 20200065-AMENDED-COMP-3

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-BF012020.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         | 2.140       | 3             | 9           | 18           | rVB      | 958948         | 1320506       | 11.95%          | 2.704%        |
| 2         | 5.110       | 502           | 514         | 517          | rBV      | 5627194        | 11049306      | 100.00%         | 22.627%       |
| 3         | 5.481       | 572           | 577         | 581          | rVB      | 243068         | 208487        | 1.89%           | 0.427%        |
| 4         | 6.487       | 743           | 748         | 758          | rBV3     | 1626834        | 1829251       | 16.56%          | 3.746%        |
| 5         | 6.863       | 808           | 812         | 816          | rBV      | 1774992        | 1414896       | 12.81%          | 2.897%        |
| 6         | 7.022       | 835           | 839         | 842          | rVB      | 4475211        | 3740247       | 33.85%          | 7.659%        |
| 7         | 7.257       | 873           | 879         | 885          | rVB      | 374181         | 315996        | 2.86%           | 0.647%        |
| 8         | 7.422       | 902           | 907         | 910          | rBV      | 3105281        | 2822803       | 25.55%          | 5.781%        |
| 9         | 7.575       | 929           | 933         | 936          | rBV      | 1206599        | 1003840       | 9.09%           | 2.056%        |
| 10        | 8.145       | 1026          | 1030        | 1033         | rBV      | 2264114        | 1861669       | 16.85%          | 3.812%        |
| 11        | 9.222       | 1208          | 1213        | 1216         | rBV      | 4862190        | 4664126       | 42.21%          | 9.551%        |
| 12        | 9.610       | 1275          | 1279        | 1284         | rVB      | 875739         | 746364        | 6.75%           | 1.528%        |
| 13        | 9.898       | 1324          | 1328        | 1331         | rBV      | 2623054        | 2187951       | 19.80%          | 4.481%        |
| 14        | 10.798      | 1478          | 1481        | 1490         | rVB3     | 248571         | 321948        | 2.91%           | 0.659%        |
| 15        | 10.869      | 1490          | 1493        | 1495         | rBV      | 139850         | 120658        | 1.09%           | 0.247%        |
| 16        | 11.045      | 1520          | 1523        | 1526         | rVV3     | 124219         | 131844        | 1.19%           | 0.270%        |
| 17        | 11.239      | 1552          | 1556        | 1561         | rBV4     | 225839         | 264167        | 2.39%           | 0.541%        |
| 18        | 11.380      | 1576          | 1580        | 1583         | rBV      | 2529268        | 2177250       | 19.70%          | 4.459%        |
| 19        | 11.404      | 1583          | 1584        | 1587         | rVV      | 249321         | 149571        | 1.35%           | 0.306%        |
| 20        | 11.604      | 1615          | 1618        | 1621         | rBV      | 119416         | 120205        | 1.09%           | 0.246%        |
| 21        | 11.904      | 1667          | 1669        | 1674         | rBV3     | 117145         | 131602        | 1.19%           | 0.270%        |
| 22        | 11.998      | 1682          | 1685        | 1693         | rVB      | 111601         | 158466        | 1.43%           | 0.325%        |
| 23        | 12.592      | 1783          | 1786        | 1791         | rBV      | 575637         | 453837        | 4.11%           | 0.929%        |
| 24        | 12.821      | 1819          | 1825        | 1827         | rBV2     | 419839         | 462261        | 4.18%           | 0.947%        |
| 25        | 12.963      | 1845          | 1849        | 1853         | rVB      | 5064596        | 4591126       | 41.55%          | 9.402%        |
| 26        | 13.857      | 1998          | 2001        | 2007         | rBV      | 629481         | 677500        | 6.13%           | 1.387%        |
| 27        | 14.015      | 2022          | 2028        | 2031         | rBV      | 2016298        | 2087249       | 18.89%          | 4.274%        |
| 28        | 14.039      | 2031          | 2032        | 2040         | rVB      | 230255         | 155130        | 1.40%           | 0.318%        |
| 29        | 14.180      | 2054          | 2056        | 2061         | rVB2     | 125779         | 118096        | 1.07%           | 0.242%        |
| 30        | 14.486      | 2105          | 2108        | 2113         | rBV2     | 201746         | 219662        | 1.99%           | 0.450%        |
| 31        | 14.792      | 2157          | 2160        | 2167         | rVB      | 193548         | 227488        | 2.06%           | 0.466%        |
| 32        | 14.868      | 2170          | 2173        | 2177         | rVB      | 193732         | 190743        | 1.73%           | 0.391%        |
| 33        | 15.051      | 2200          | 2204        | 2207         | rBV      | 165183         | 243692        | 2.21%           | 0.499%        |
| 34        | 15.133      | 2214          | 2218        | 2221         | rBV      | 178741         | 216373        | 1.96%           | 0.443%        |

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Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

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-BF012020.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |     |         |         |        |        |
|----|--------|------|------|------|-----|---------|---------|--------|--------|
| 35 | 15.351 | 2253 | 2255 | 2263 | rVB | 104601  | 145399  | 1.32%  | 0.298% |
| 36 | 15.415 | 2263 | 2266 | 2271 | rVB | 128194  | 145801  | 1.32%  | 0.299% |
| 37 | 15.480 | 2272 | 2277 | 2281 | rBV | 1383842 | 1813698 | 16.41% | 3.714% |
| 38 | 15.509 | 2281 | 2282 | 2290 | rVB | 188048  | 207456  | 1.88%  | 0.425% |
| 39 | 15.951 | 2354 | 2357 | 2362 | rVB | 106821  | 134990  | 1.22%  | 0.276% |

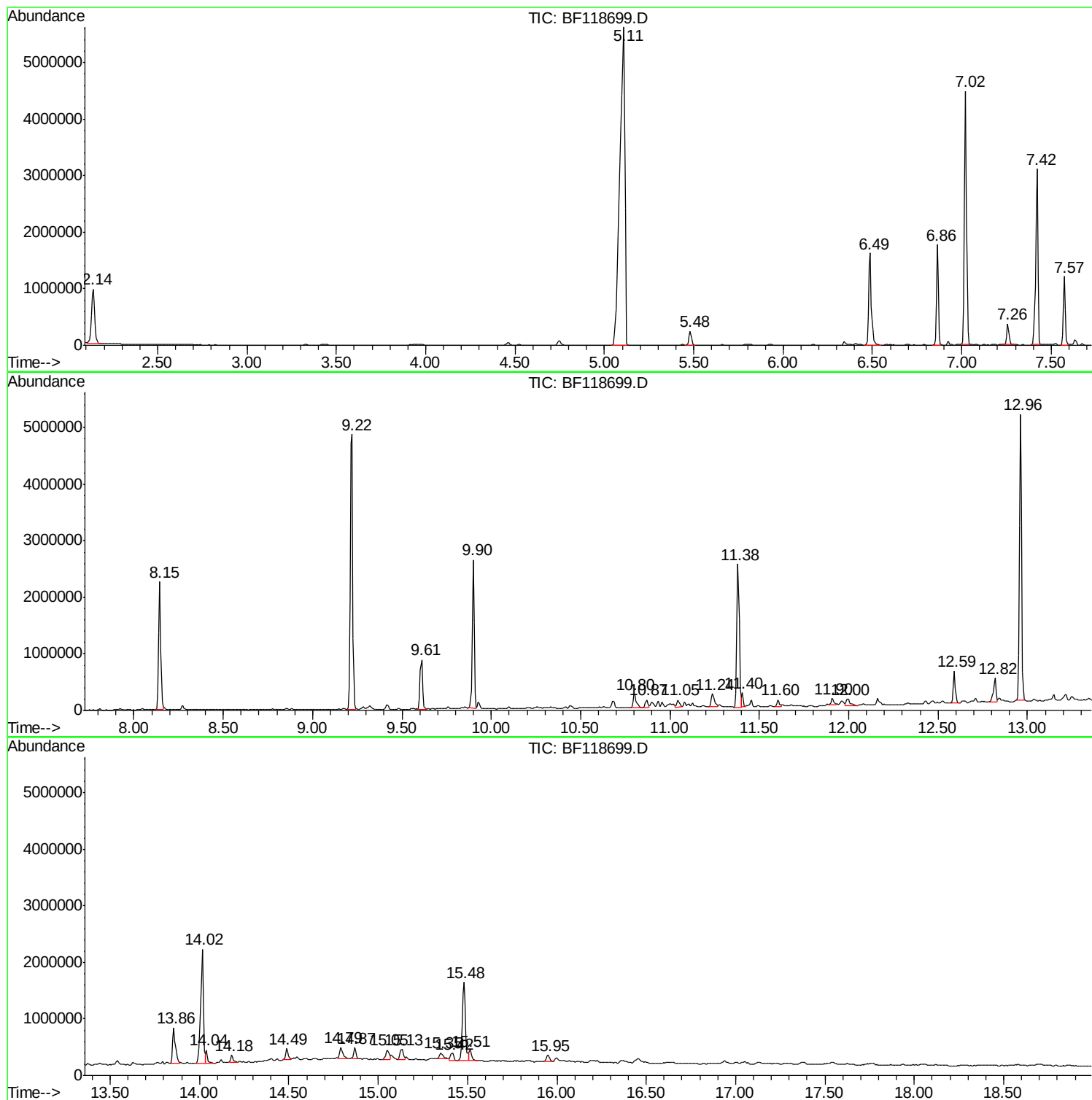
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BNA\_F  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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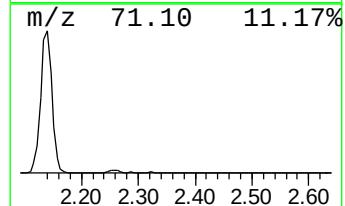
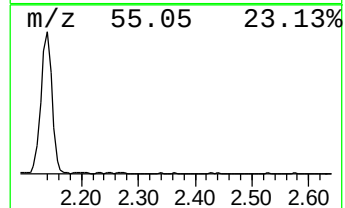
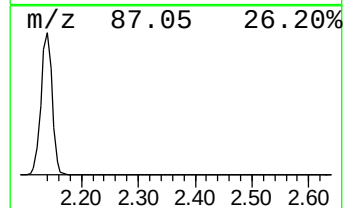
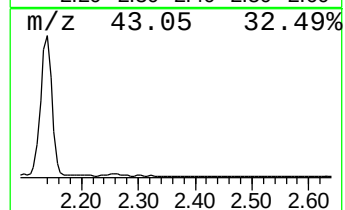
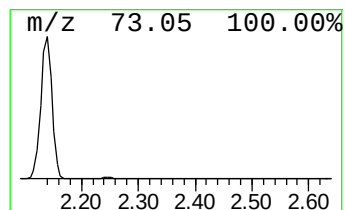
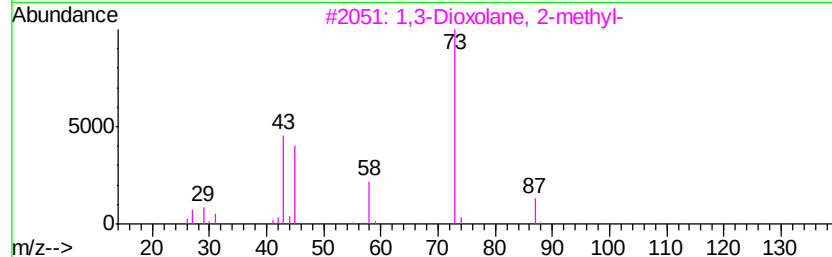
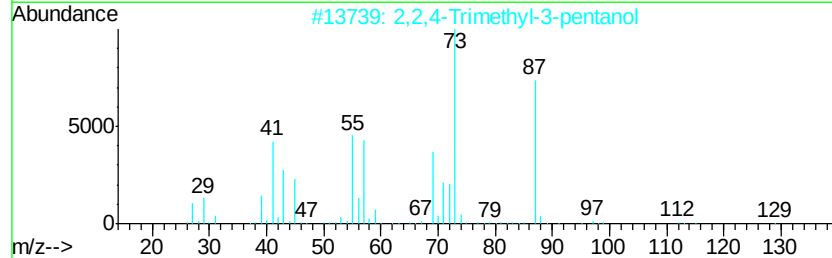
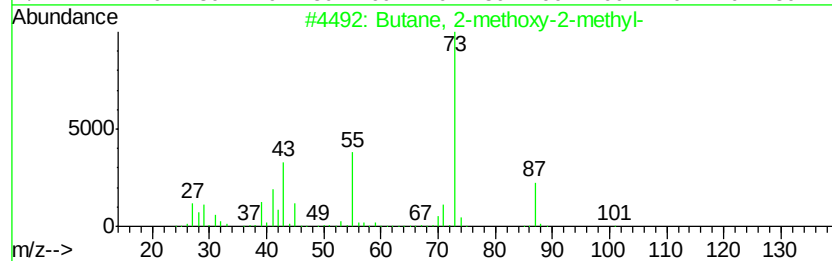
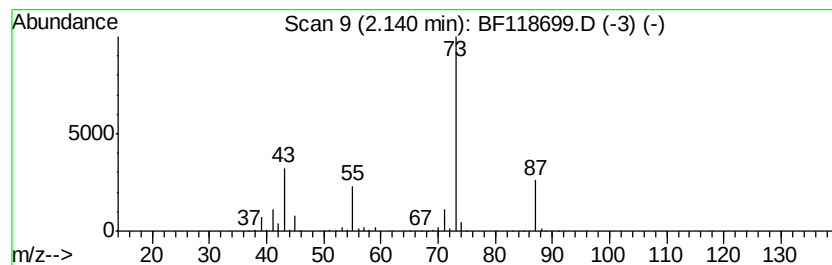
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 2.14 | 18.67 ng | 1320510 | 1,4-Dichlorobenzene-d4 | 6.86 |

| 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 | 39   |
| 3    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 4    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 23   |
| 5    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 17   |



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Instrument :  
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ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

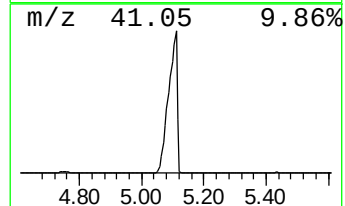
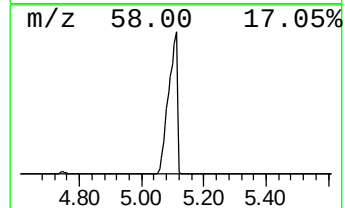
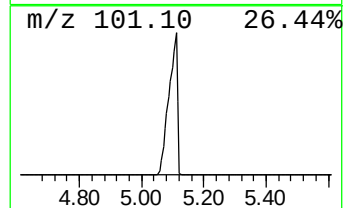
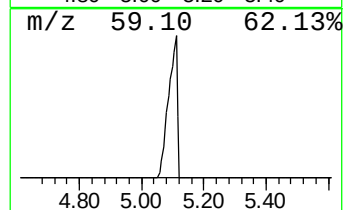
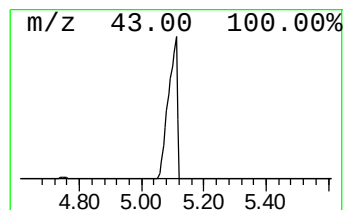
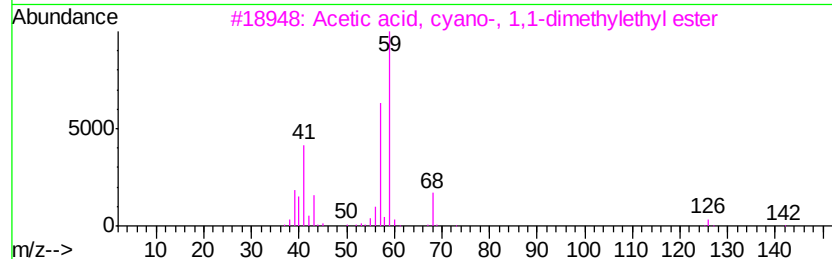
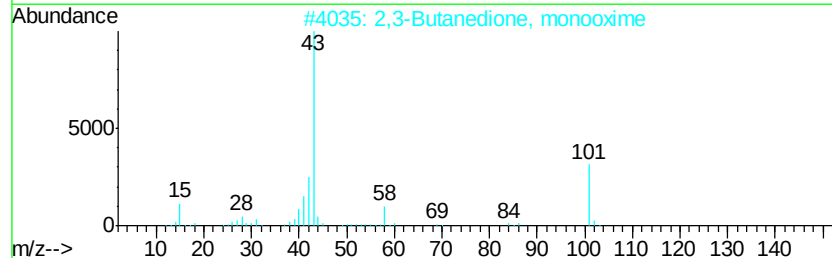
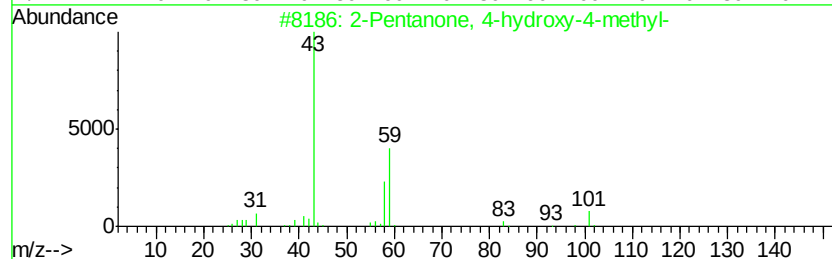
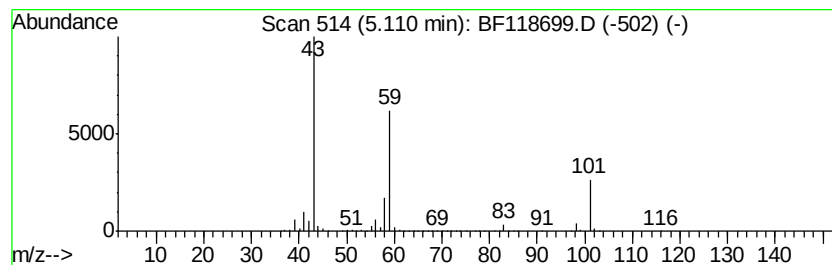
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc   | Area     | Relative to ISTD       | R.T. |
|------|-----------|----------|------------------------|------|
| 5.11 | 156.19 ng | 11049300 | 1,4-Dichlorobenzene-d4 | 6.86 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 17   |
| 3    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 4    |    | Acetone                              | 58  | C3H6O    | 000067-64-1 | 10   |
| 5    |    | 5-Hexen-2-one                        | 98  | C6H10O   | 000109-49-9 | 9    |



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Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

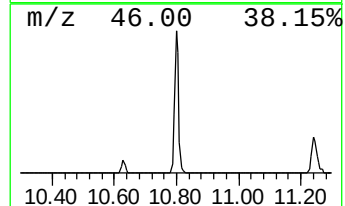
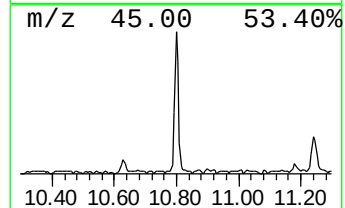
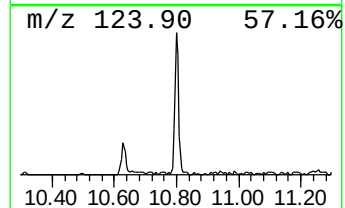
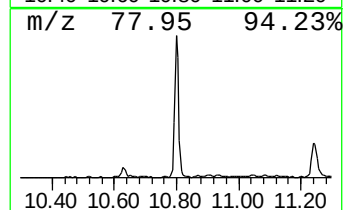
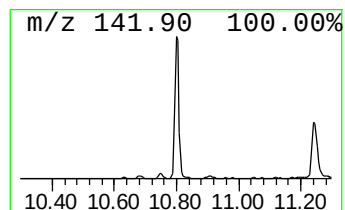
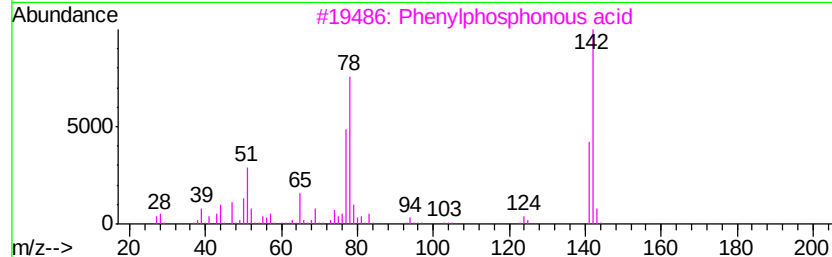
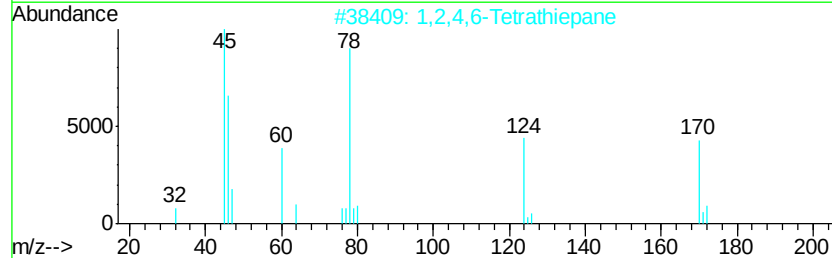
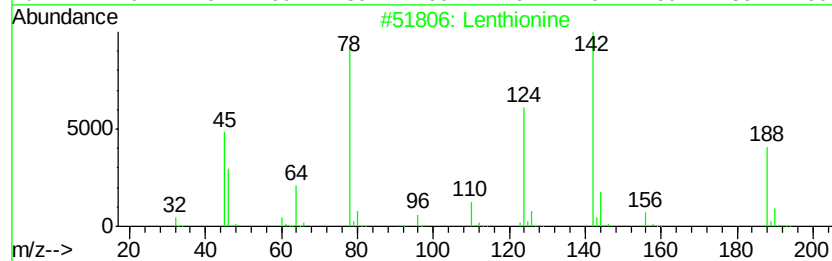
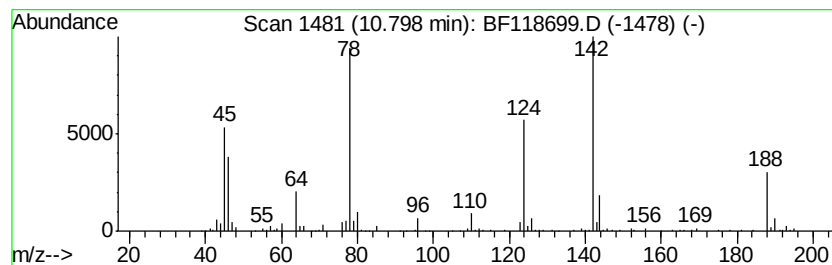
Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Lenthionine Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.80     | 2.96 ng                             | 321948 | Phenanthrene-d10 | 11.38        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Lenthionine                         | 188    | C2H4S5           | 000292-46-6  | 94   |
| 2         | 1,2,4,6-Tetrathiepane               | 170    | C3H6S4           | 000292-45-5  | 45   |
| 3         | Phenvlphosphonous acid              | 142    | C6H7O2P          | 001779-48-2  | 33   |
| 4         | Boronic acid, phenyl-               | 122    | C6H7B02          | 000098-80-6  | 9    |
| 5         | Methylthiophosphonamidic acid, S... | 125    | C2H8NOPS         | 1000306-03-2 | 9    |



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ALS Vial : 19 Sample Multiplier: 1

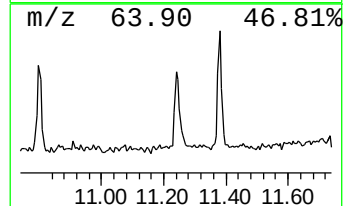
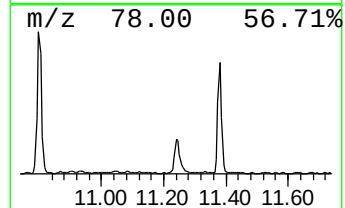
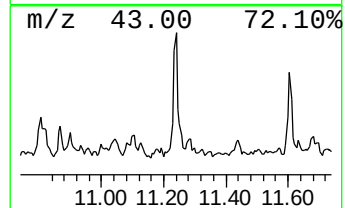
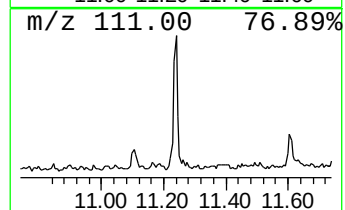
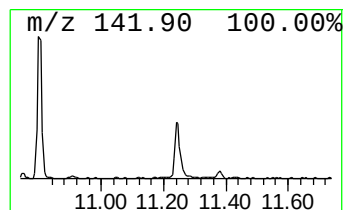
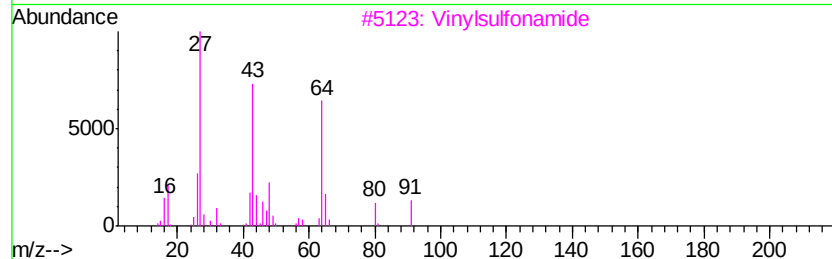
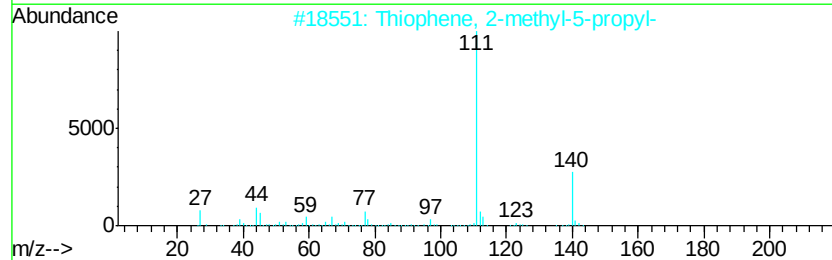
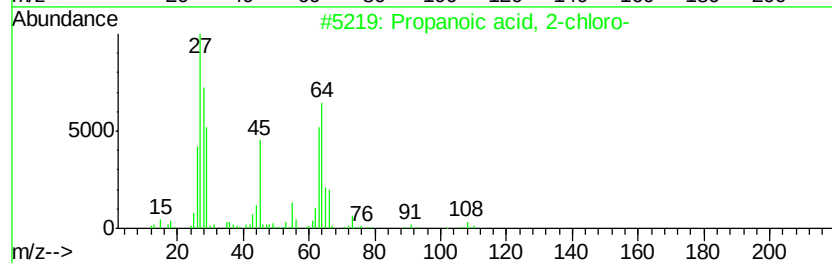
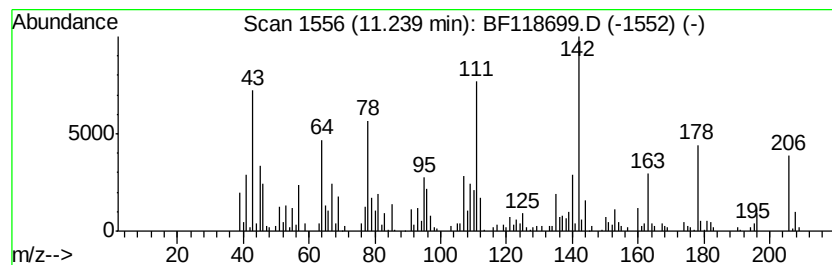
Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 unknown11.24 Concentration Rank 8

| R.T.      | EstConc                        | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------------------------|--------|------------------|-------------|-------|
| 11.24     | 2.43 ng                        | 264167 | Phenanthrene-d10 |             | 11.38 |
| Hit# of 5 | Tentative ID                   | MW     | MolForm          | CAS#        | Qual  |
| 1         | Propanoic acid, 2-chloro-      | 108    | C3H5ClO2         | 000598-78-7 | 14    |
| 2         | Thiophene, 2-methyl-5-propyl-  | 140    | C8H12S           | 033933-73-2 | 10    |
| 3         | Vinylsulfonamide               | 107    | C2H5NO2S         | 002386-58-5 | 10    |
| 4         | (S)-(-)-2-Chloropropionic acid | 108    | C3H5ClO2         | 029617-66-1 | 10    |
| 5         | Tetrazolo[1,5-b]pyridazine     | 121    | C4H3N5           | 000274-89-5 | 10    |



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Misc :  
ALS Vial : 19 Sample Multiplier: 1

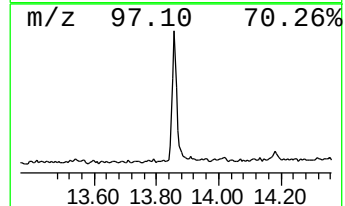
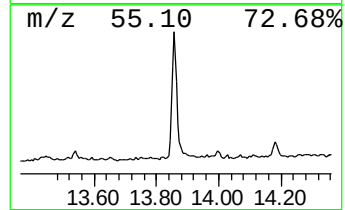
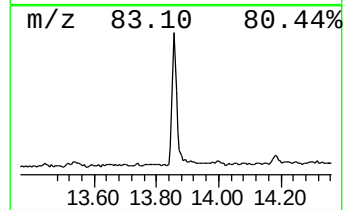
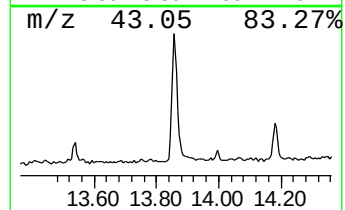
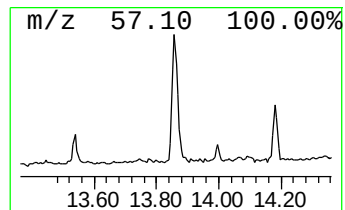
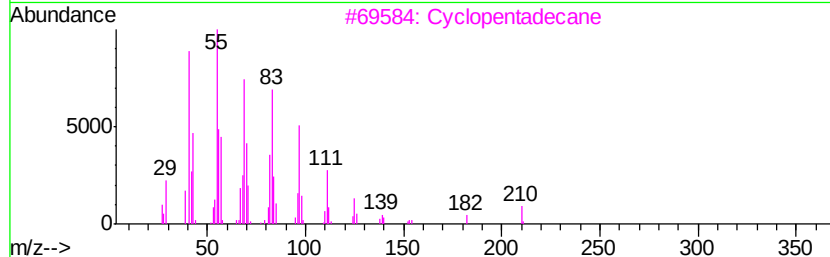
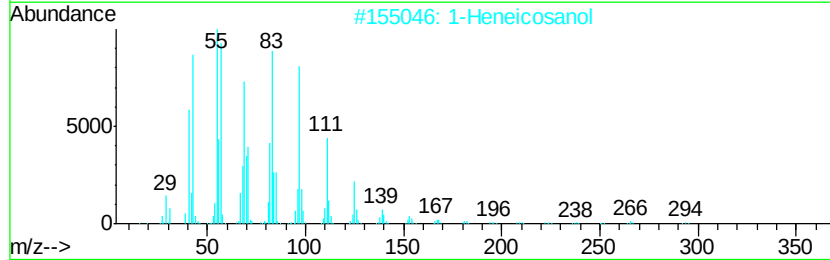
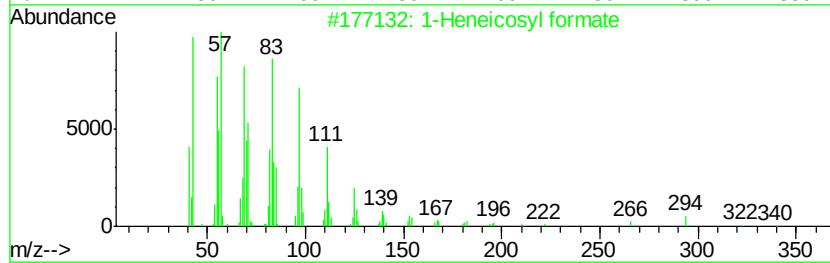
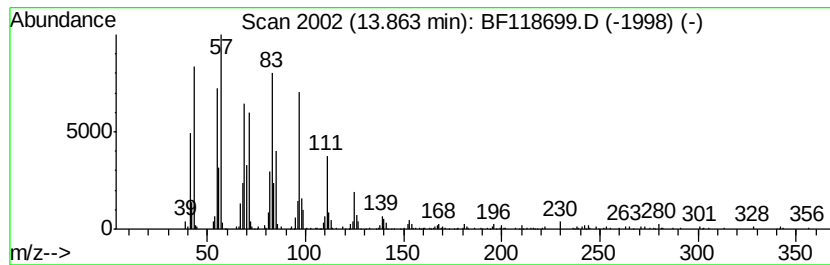
Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.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 1-Heneicosyl formate Concentration Rank 5

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.86     | 6.49 ng                             | 677500 | Chrysene-d12     | 14.02        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Heneicosyl formate                | 340    | C22H44O2         | 077899-03-7  | 98   |
| 2         | 1-Heneicosanol                      | 312    | C21H44O          | 015594-90-8  | 95   |
| 3         | Cyclopentadecane                    | 210    | C15H30           | 000295-48-7  | 94   |
| 4         | Dichloroacetic acid, heptadecyl ... | 366    | C19H36Cl2O2      | 1000282-98-2 | 94   |
| 5         | n-Nonadecanol-1                     | 284    | C19H40O          | 001454-84-8  | 94   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\  
Data File : BF118699.D  
Acq On : 24 Jan 2020 21:35  
Operator : CG/JU  
Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

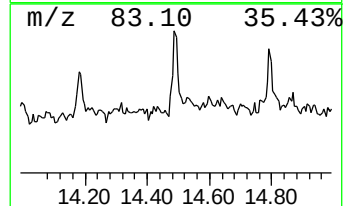
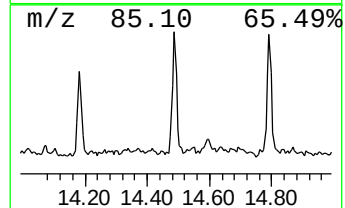
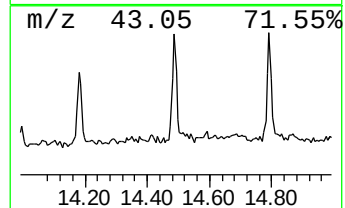
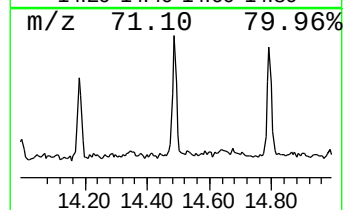
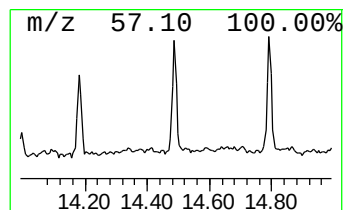
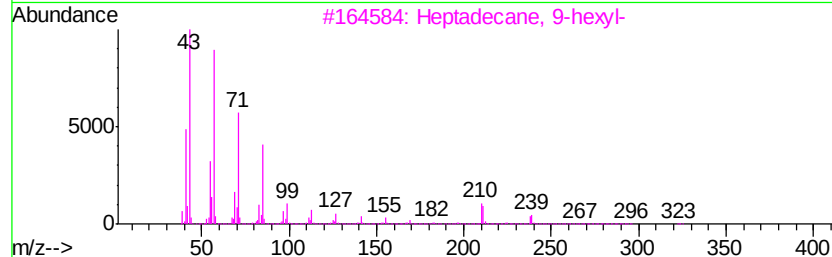
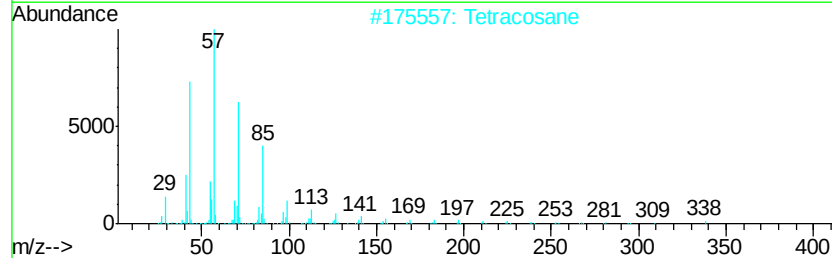
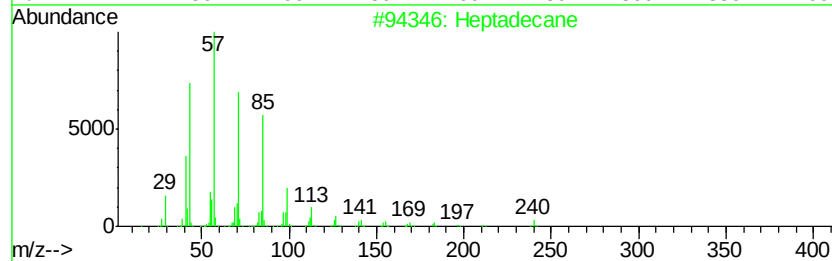
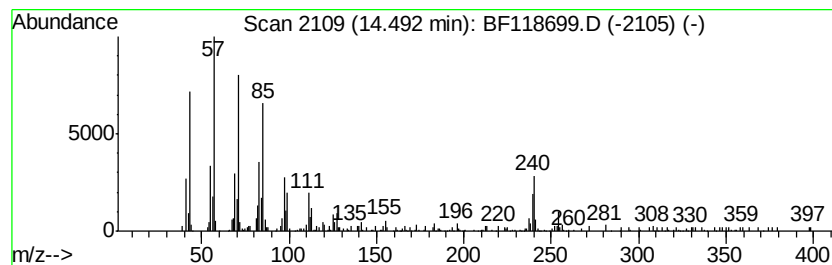
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ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Heptadecane Concentration Rank 10

| R.T.    | EstConc                             | Area          | Relative to ISTD | R.T.      |
|---------|-------------------------------------|---------------|------------------|-----------|
| 14.49   | 2.10 ng                             | 219662        | Chrysene-d12     | 14.02     |
| Hit# of | 5                                   | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | Heptadecane                         | 240 C17H36    | 000629-78-7      | 98        |
| 2       | Tetracosane                         | 338 C24H50    | 000646-31-1      | 92        |
| 3       | Heptadecane, 9-hexyl-               | 324 C23H48    | 055124-79-3      | 90        |
| 4       | Sulfurous acid, dodecyl 2-propyl... | 292 C15H32O3S | 1000309-12-3     | 87        |
| 5       | Sulfurous acid, 2-propyl undecyl... | 278 C14H30O3S | 1000309-12-2     | 87        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\  
Data File : BF118699.D  
Acq On : 24 Jan 2020 21:35  
Operator : CG/JU  
Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

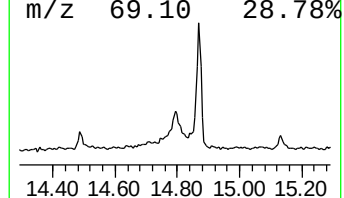
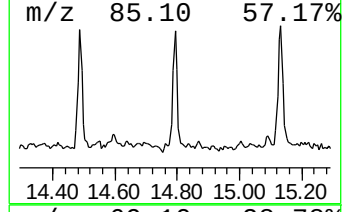
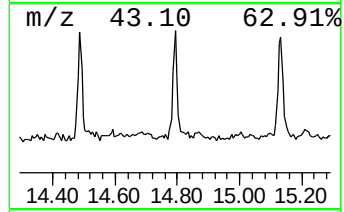
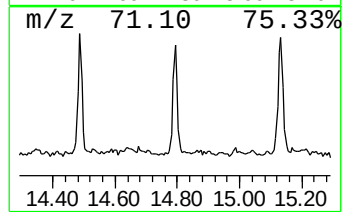
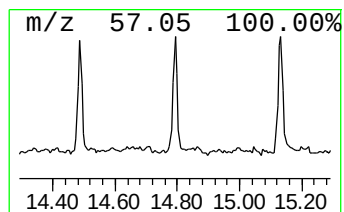
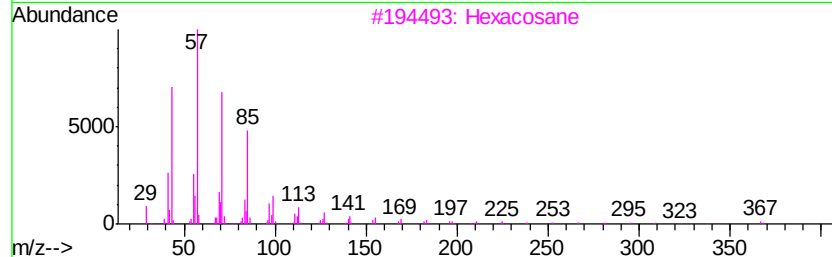
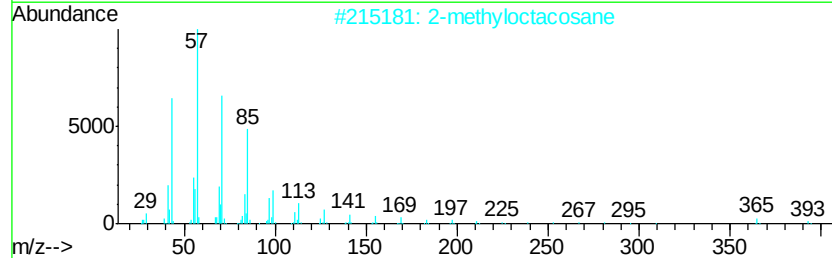
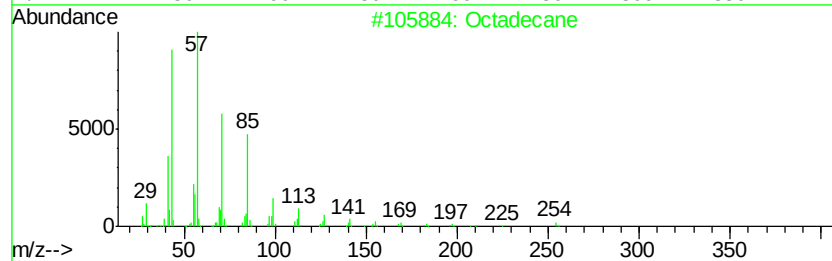
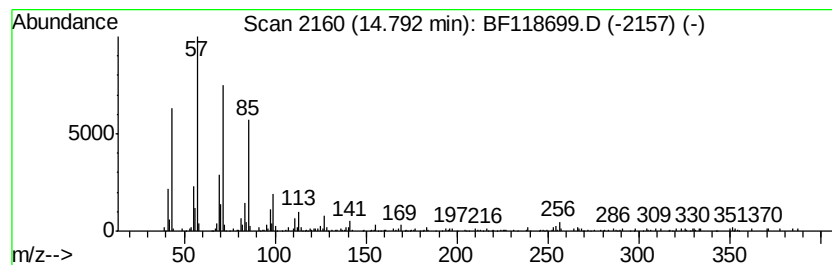
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Octadecane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.79 | 2.51 ng | 227488 | Perylene-d12     | 15.48 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------|-----|---------|--------------|------|
| 1    | 5  | Octadecane         | 254 | C18H38  | 000593-45-3  | 93   |
| 2    |    | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 91   |
| 3    |    | Hexacosane         | 366 | C26H54  | 000630-01-3  | 91   |
| 4    |    | Heneicosane        | 296 | C21H44  | 000629-94-7  | 91   |
| 5    |    | Tetracosane        | 338 | C24H50  | 000646-31-1  | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\  
Data File : BF118699.D  
Acq On : 24 Jan 2020 21:35  
Operator : CG/JU  
Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

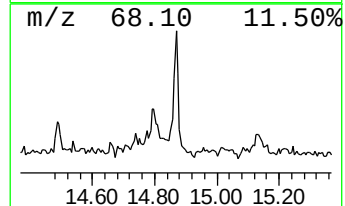
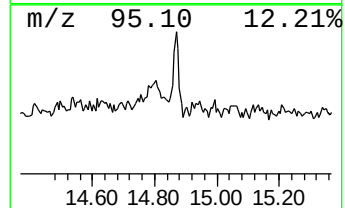
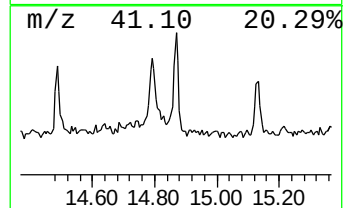
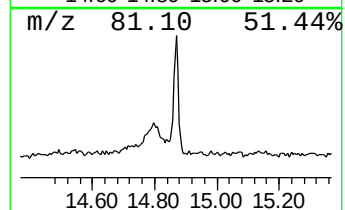
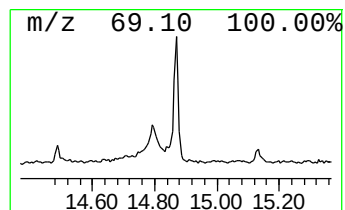
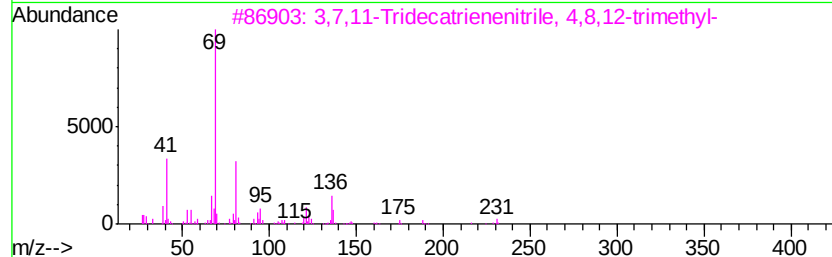
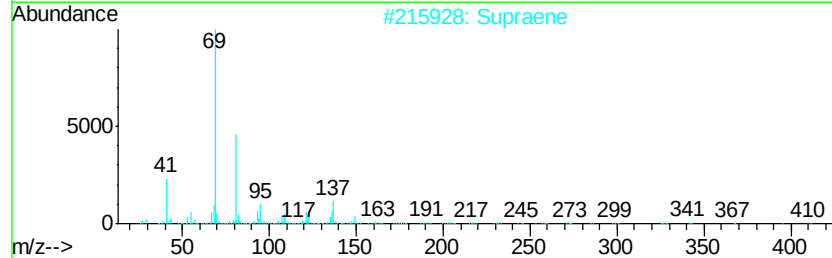
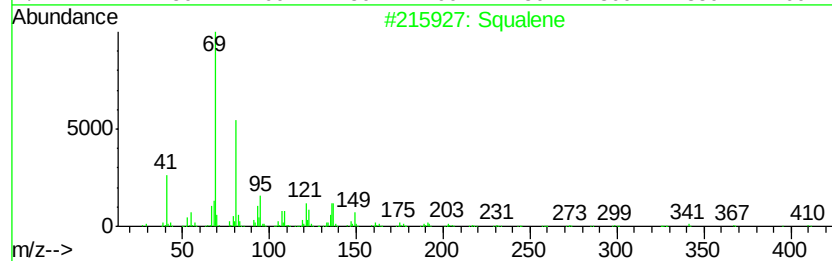
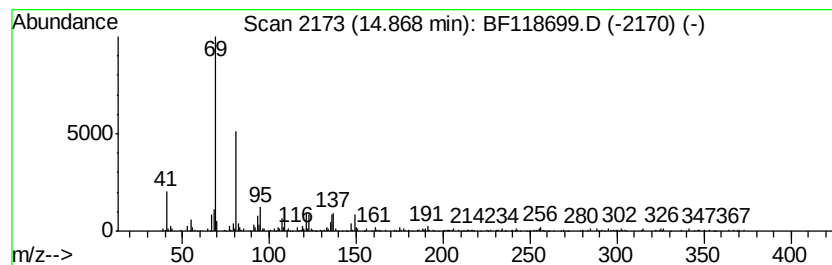
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Squalene Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.87 | 2.10 ng | 190743 | Perylene-d12     | 15.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Squalene                            | 410 | C30H50   | 000111-02-4 | 87   |
| 2    |    | Supraene                            | 410 | C30H50   | 007683-64-9 | 83   |
| 3    |    | 3,7,11-Tridecatrienitrile, 4,8...   | 231 | C16H25N  | 006006-01-5 | 72   |
| 4    |    | 1,6-Octadiene, 3,5-dimethyl-, tr... | 138 | C10H18   | 074630-87-8 | 59   |
| 5    |    | 2,3,5-Trimethyl-2,3,5-hexanetric... | 203 | C12H17N3 | 003974-79-6 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\  
Data File : BF118699.D  
Acq On : 24 Jan 2020 21:35  
Operator : CG/JU  
Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

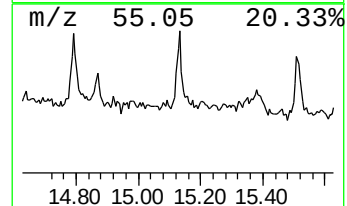
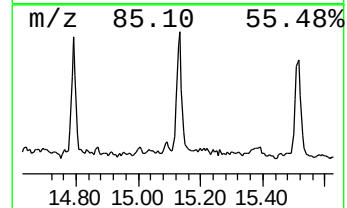
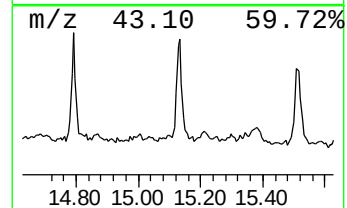
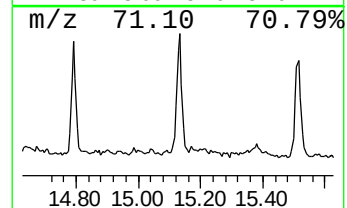
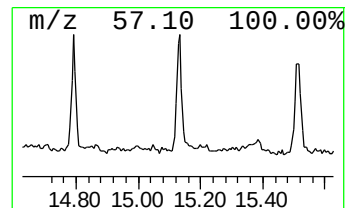
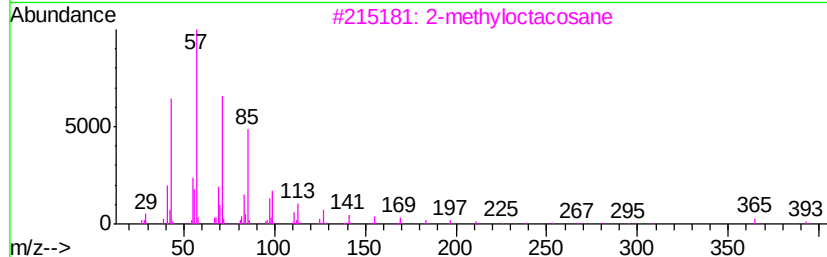
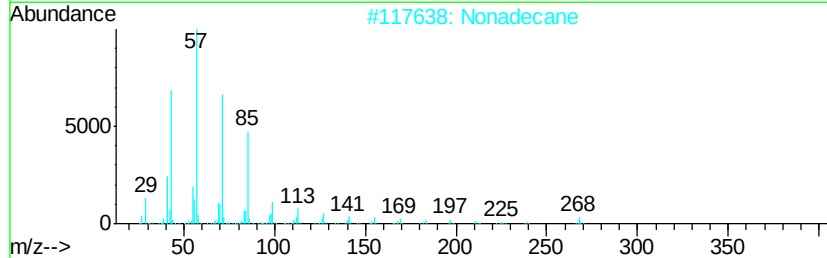
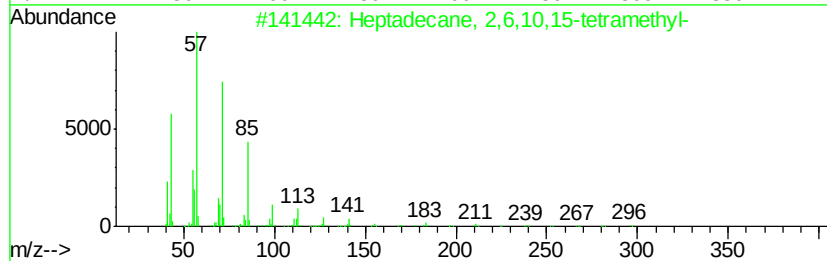
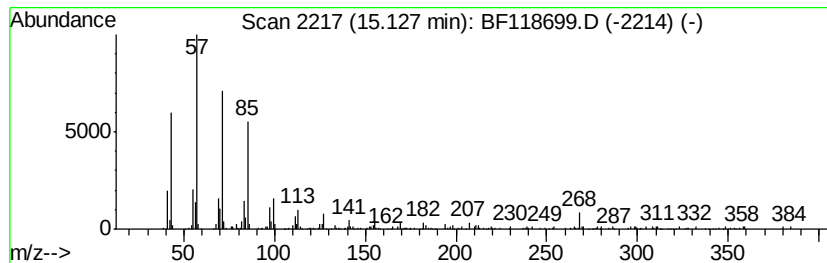
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Heptadecane, 2,6,10,15-tetr... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.13 | 2.39 ng | 216373 | Perylene-d12     | 15.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6  | 93   |
| 2    |    | Nonadecane                          | 268 | C19H40  | 000629-92-5  | 91   |
| 3    |    | 2-methyloctacosane                  | 408 | C29H60  | 1000376-72-8 | 91   |
| 4    |    | Heptadecane, 9-octyl-               | 352 | C25H52  | 007225-64-1  | 91   |
| 5    |    | Docosane, 11-decyl-                 | 451 | C32H66  | 055401-55-3  | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF012420\  
Data File : BF118699.D  
Acq On : 24 Jan 2020 21:35  
Operator : CG/JU  
Sample : L1242-12  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
20200065-AMENDED-COMP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF012020.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... | 2.14  | 18.7    | ng    | 1320510  | 1                     | 6.86  | 1414900 | 20.0 |
| 2-Pentanone, 4-hy... | 5.11  | 156.2   | ng    | 11049300 | 1                     | 6.86  | 1414900 | 20.0 |
| Lenthionine          | 10.80 | 3.0     | ng    | 321948   | 4                     | 11.38 | 2177250 | 20.0 |
| unknown11.24         | 11.24 | 2.4     | ng    | 264167   | 4                     | 11.38 | 2177250 | 20.0 |
| 1-Heneicosyl formate | 13.86 | 6.5     | ng    | 677500   | 5                     | 14.02 | 2087250 | 20.0 |
| Heptadecane          | 14.49 | 2.1     | ng    | 219662   | 5                     | 14.02 | 2087250 | 20.0 |
| Octadecane           | 14.79 | 2.5     | ng    | 227488   | 6                     | 15.48 | 1813700 | 20.0 |
| Squalene             | 14.87 | 2.1     | ng    | 190743   | 6                     | 15.48 | 1813700 | 20.0 |
| Heptadecane, 2,6,... | 15.13 | 2.4     | ng    | 216373   | 6                     | 15.48 | 1813700 | 20.0 |