

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111219\  
 Data File : BF117762.D  
 Acq On : 12 Nov 2019 16:29  
 Operator : JU  
 Sample : K5789-03  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-17-S

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-BF110619.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 2.216    | 16         | 22       | 33        | rVB   | 1119329     | 1494153    | 26.45%       | 2.707%     |
| 2      | 5.140    | 511        | 519      | 527       | rBV   | 896521      | 1048015    | 18.55%       | 1.898%     |
| 3      | 5.539    | 582        | 587      | 590       | rBV   | 5009927     | 4622345    | 81.82%       | 8.373%     |
| 4      | 6.545    | 752        | 758      | 762       | rBV   | 4719254     | 5060389    | 89.57%       | 9.166%     |
| 5      | 6.692    | 778        | 783      | 786       | rBV   | 5483325     | 5054945    | 89.47%       | 9.157%     |
| 6      | 6.922    | 819        | 822      | 826       | rVB   | 1788437     | 1604065    | 28.39%       | 2.906%     |
| 7      | 7.081    | 845        | 849      | 853       | rBV   | 4475473     | 3896770    | 68.97%       | 7.059%     |
| 8      | 7.486    | 913        | 918      | 921       | rBV   | 3324096     | 3109862    | 55.05%       | 5.633%     |
| 9      | 8.210    | 1037       | 1041     | 1045      | rBV   | 2560674     | 2116934    | 37.47%       | 3.835%     |
| 10     | 9.286    | 1220       | 1224     | 1228      | rBV   | 5835261     | 5649592    | 100.00%      | 10.234%    |
| 11     | 9.404    | 1239       | 1244     | 1255      | rVB   | 379115      | 355435     | 6.29%        | 0.644%     |
| 12     | 9.680    | 1287       | 1291     | 1294      | rBV   | 795847      | 616748     | 10.92%       | 1.117%     |
| 13     | 9.974    | 1337       | 1341     | 1344      | rBV   | 2936428     | 2472892    | 43.77%       | 4.479%     |
| 14     | 10.551   | 1436       | 1439     | 1447      | rBV   | 58596       | 69487      | 1.23%        | 0.126%     |
| 15     | 10.763   | 1470       | 1475     | 1478      | rBV   | 4307993     | 3851291    | 68.17%       | 6.976%     |
| 16     | 11.463   | 1590       | 1594     | 1598      | rBV   | 2923276     | 2544051    | 45.03%       | 4.608%     |
| 17     | 11.527   | 1601       | 1605     | 1608      | rVB2  | 74645       | 80080      | 1.42%        | 0.145%     |
| 18     | 11.692   | 1628       | 1633     | 1640      | rBV5  | 34847       | 78656      | 1.39%        | 0.142%     |
| 19     | 11.986   | 1679       | 1683     | 1688      | rBV   | 622090      | 542394     | 9.60%        | 0.982%     |
| 20     | 12.510   | 1769       | 1772     | 1778      | rBV   | 411103      | 539380     | 9.55%        | 0.977%     |
| 21     | 12.710   | 1803       | 1806     | 1810      | rVB   | 114898      | 98332      | 1.74%        | 0.178%     |
| 22     | 12.774   | 1814       | 1817     | 1822      | rBV   | 156093      | 172948     | 3.06%        | 0.313%     |
| 23     | 13.057   | 1860       | 1865     | 1868      | rBV   | 5928760     | 5285278    | 93.55%       | 9.574%     |
| 24     | 13.957   | 2014       | 2018     | 2026      | rBV2  | 374866      | 704285     | 12.47%       | 1.276%     |
| 25     | 14.110   | 2041       | 2044     | 2048      | rVB   | 1913596     | 1832008    | 32.43%       | 3.319%     |
| 26     | 14.586   | 2122       | 2125     | 2128      | rBV2  | 93535       | 133444     | 2.36%        | 0.242%     |
| 27     | 14.986   | 2190       | 2193     | 2196      | rBV   | 172021      | 175166     | 3.10%        | 0.317%     |
| 28     | 15.621   | 2296       | 2301     | 2305      | rBV   | 1401369     | 1824620    | 32.30%       | 3.305%     |
| 29     | 15.662   | 2305       | 2308     | 2314      | rVB4  | 92207       | 172312     | 3.05%        | 0.312%     |

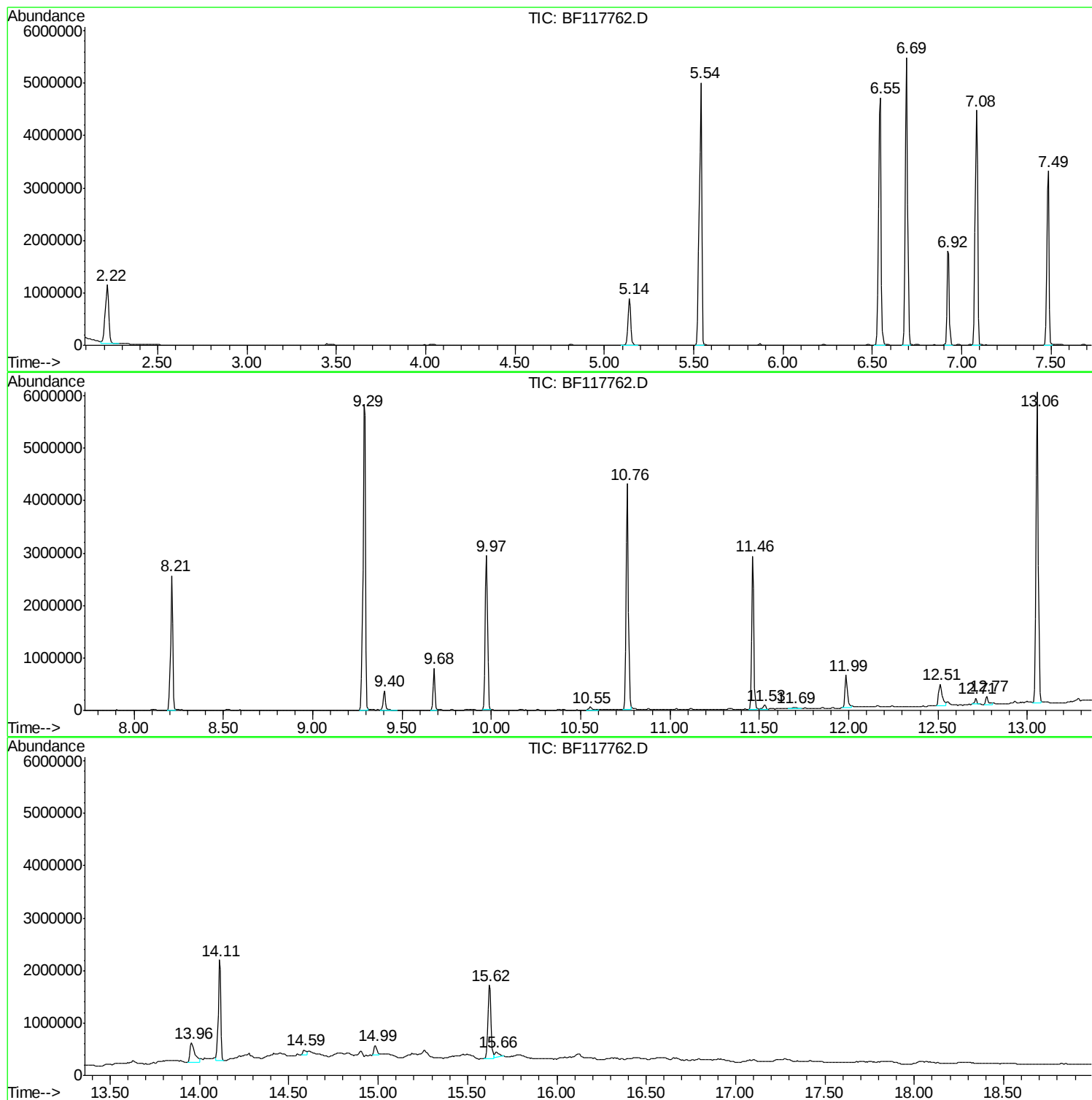
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BNA\_F  
ClientSampled :  
TP-17-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.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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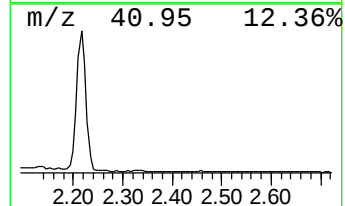
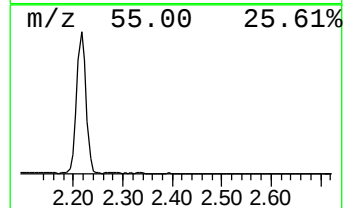
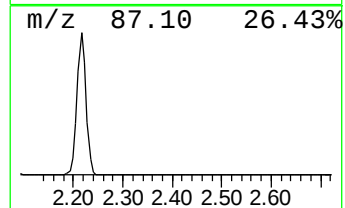
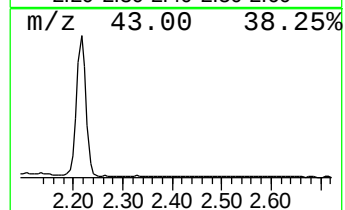
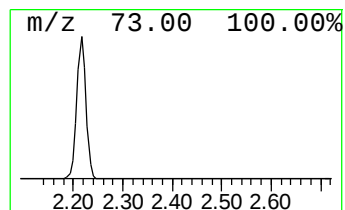
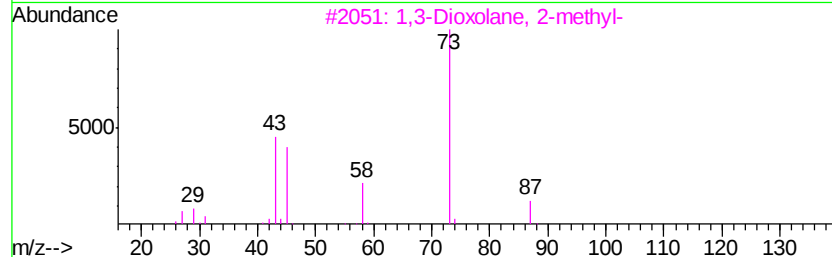
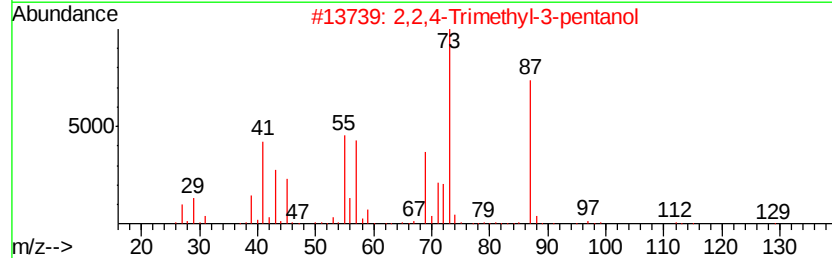
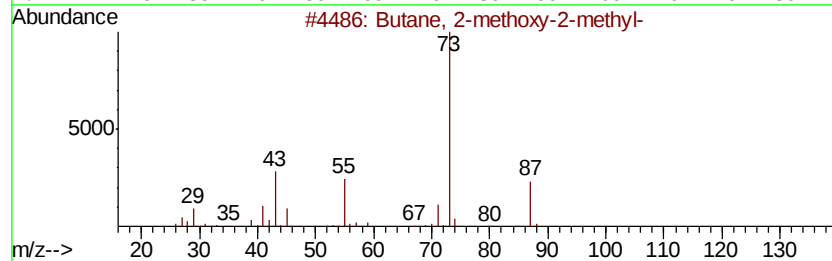
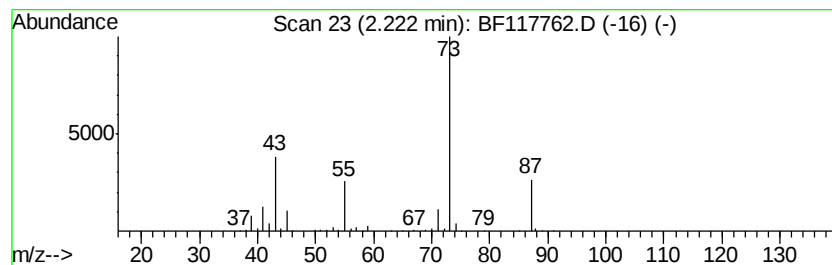
Instrument :  
BNA\_F  
ClientSampleId :  
TP-17-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.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 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T.    | EstConc  | Area                        | Relative to ISTD       | R.T.           |
|---------|----------|-----------------------------|------------------------|----------------|
| 2.22    | 18.63 ng | 1494150                     | 1,4-Dichlorobenzene-d4 | 6.93           |
| Hit# of | 5        | Tentative ID                | MW MolForm             | CAS# Qual      |
| 1       |          | 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 23 |
| 4       |          | Pentane, 3-methoxy-         | 102 C6H14O             | 036839-67-5 12 |
| 5       |          | Silane, tetramethyl-        | 88 C4H12Si             | 000075-76-3 10 |



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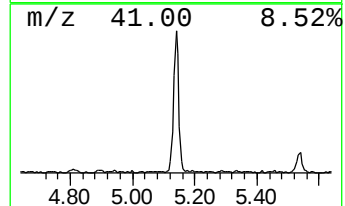
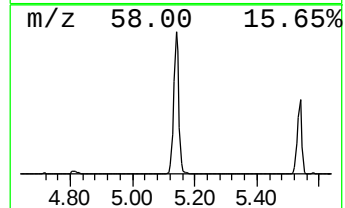
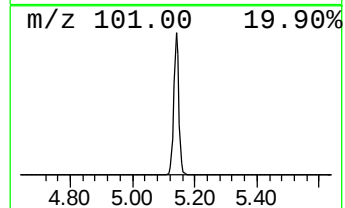
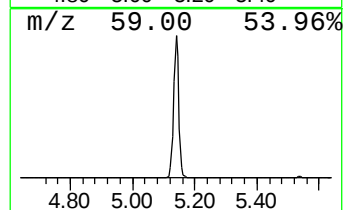
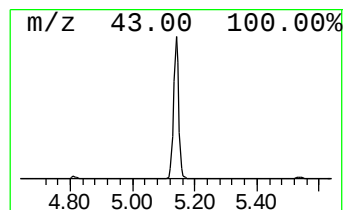
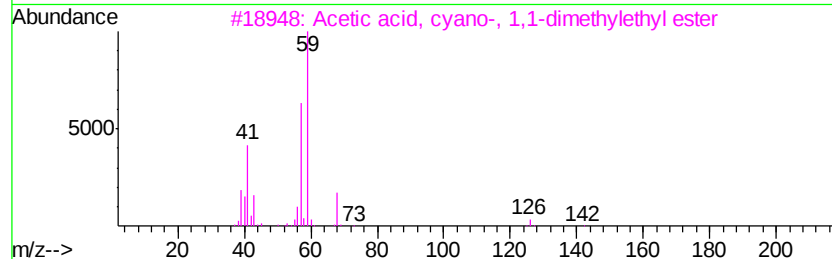
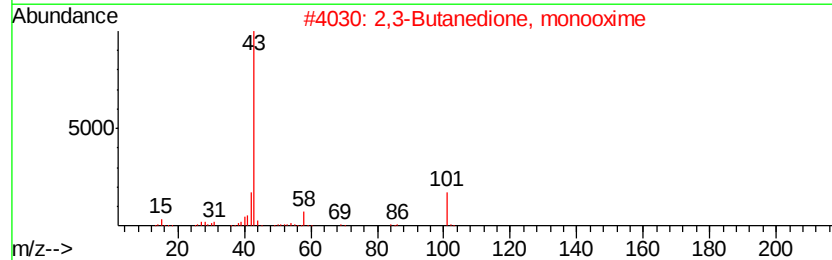
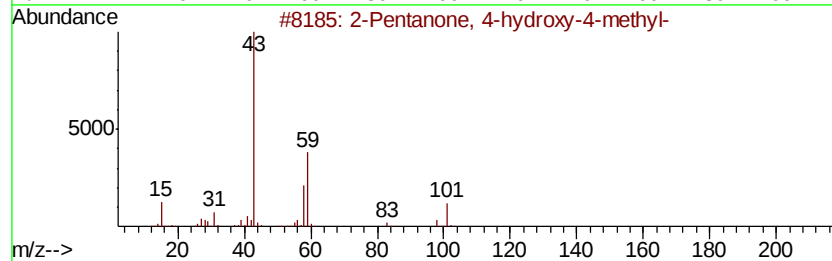
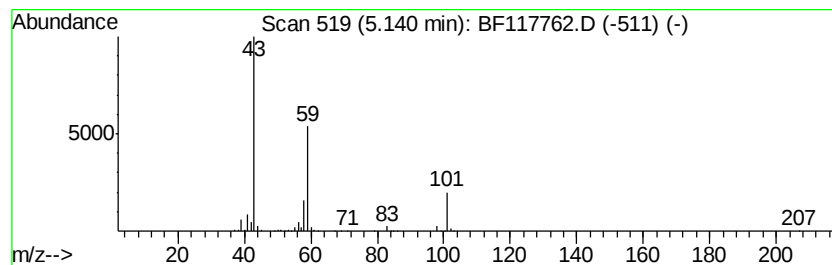
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.14 | 13.07 ng | 1048020 | 1,4-Dichlorobenzene-d4 | 6.93 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 53   |
| 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    |    | 2-Propanone, 1-(1-methylethoxy)-     | 116 | C6H12O2  | 042781-12-4 | 9    |



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Misc :  
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :  
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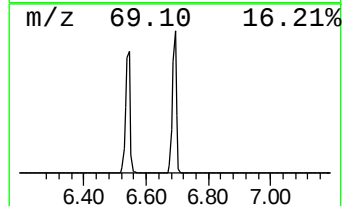
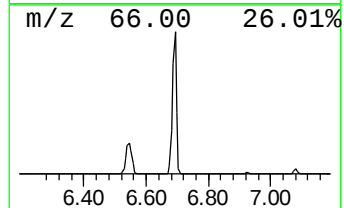
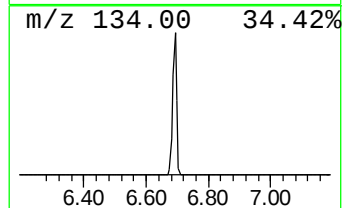
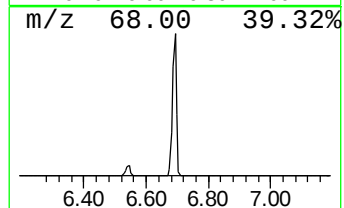
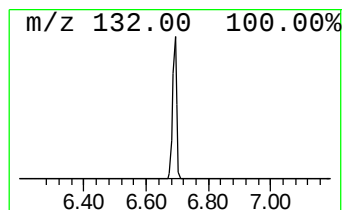
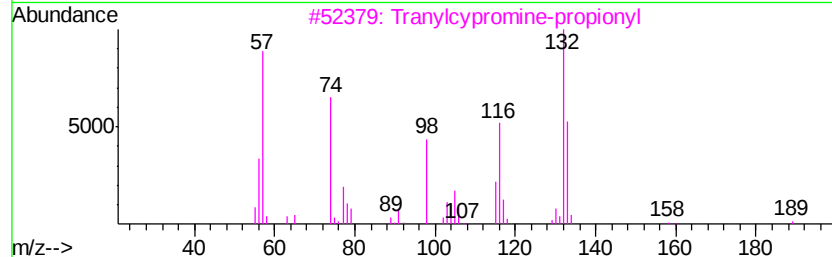
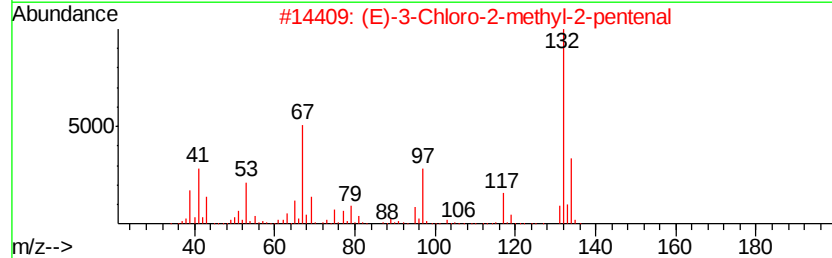
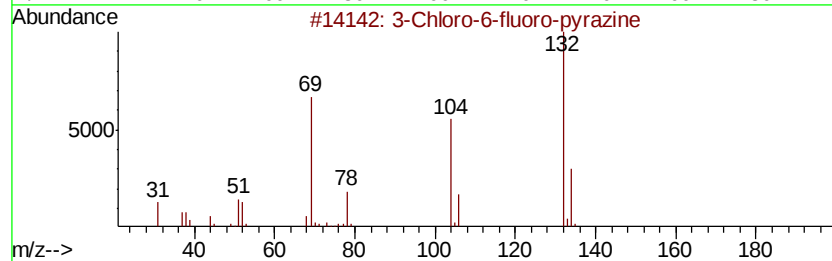
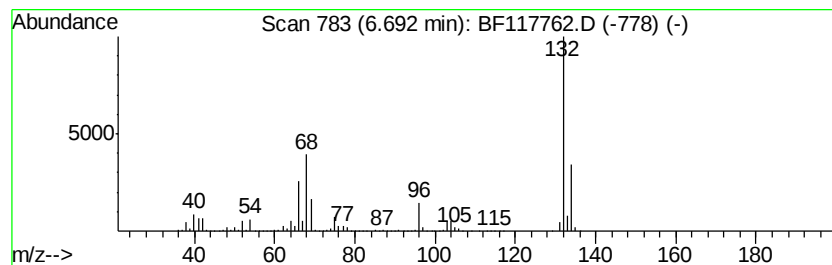
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 unknown6.69 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.69 | 63.03 ng | 5054950 | 1,4-Dichlorobenzene-d4 | 6.93 |

| 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    |



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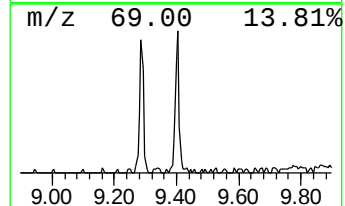
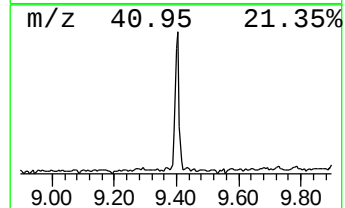
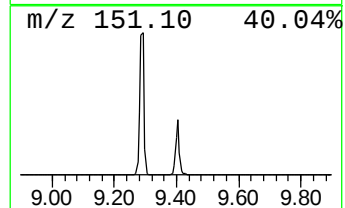
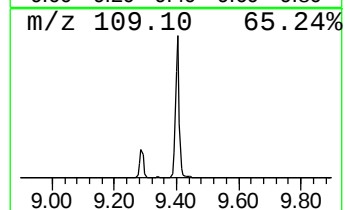
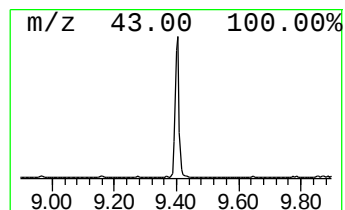
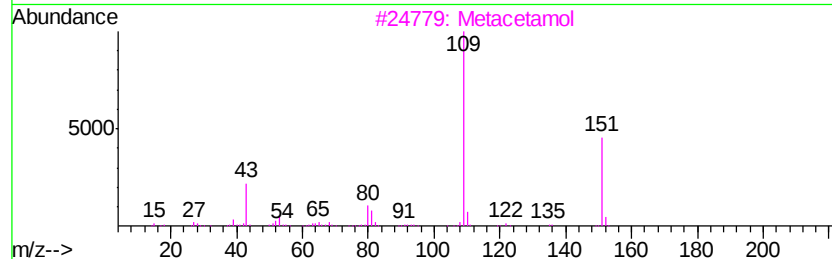
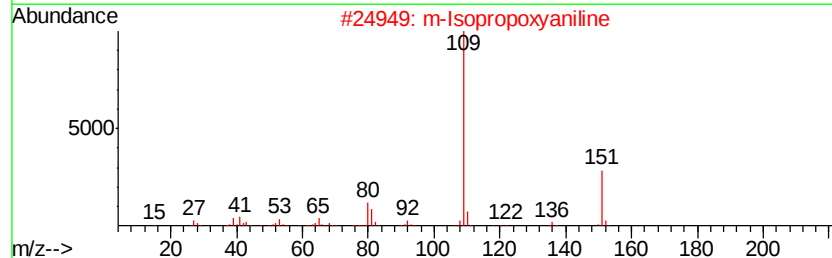
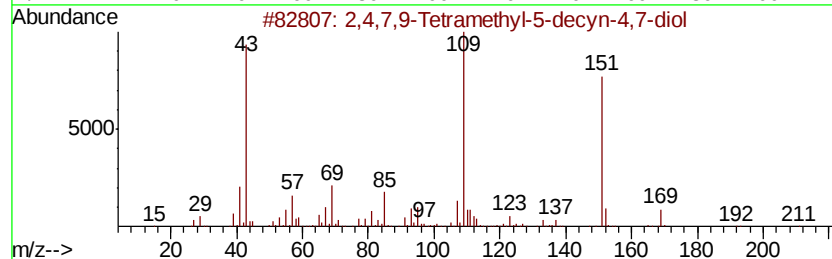
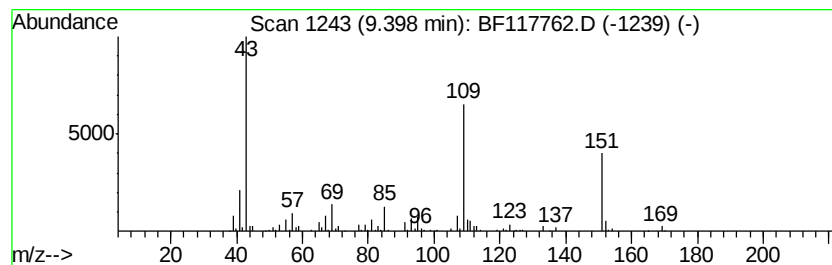
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Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.40 | 2.87 ng | 355435 | 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    |    | m-Isopropoxyvaniline                | 151 | C9H13NO  | 041406-00-2  | 59   |
| 3    |    | Metacetamol                         | 151 | C8H9NO2  | 000621-42-1  | 59   |
| 4    |    | N-Cyclopropyl-p-fluorobenzylamine   | 165 | C10H12FN | 1000250-88-9 | 43   |
| 5    |    | 1-(1,2,3-Trimethyl-cyclopent-2-e... | 152 | C10H16O  | 070987-81-4  | 43   |



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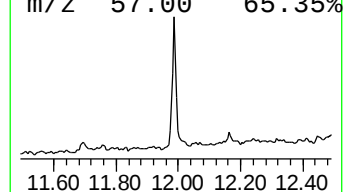
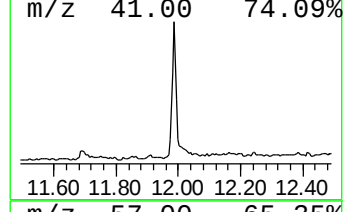
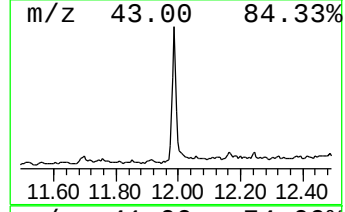
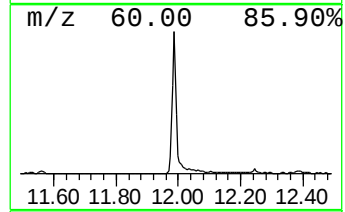
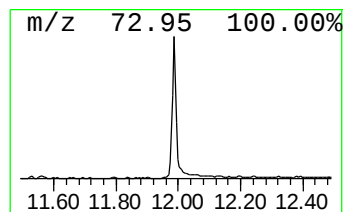
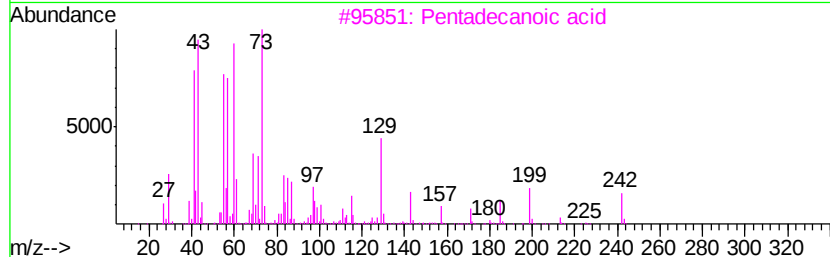
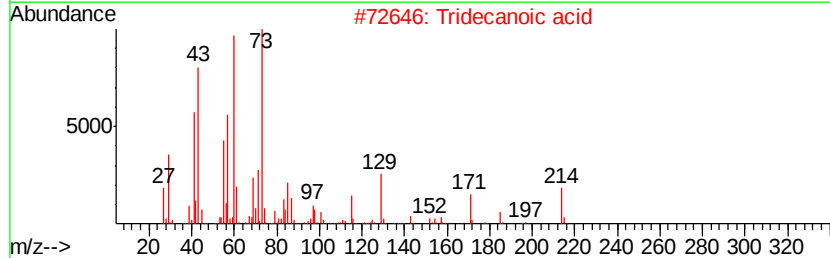
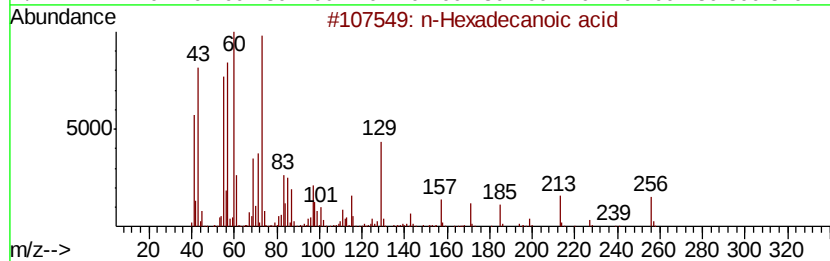
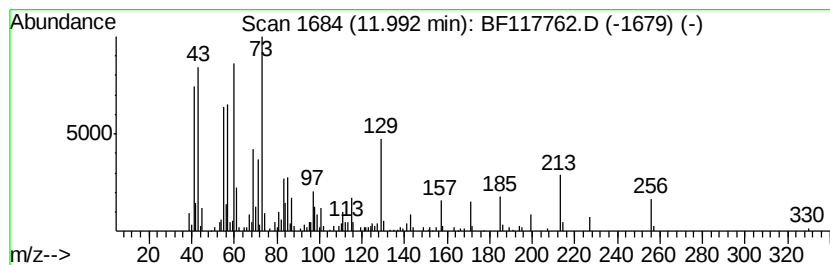
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.99 | 4.26 ng | 542394 | Phenanthrene-d10 | 11.46 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 91   |
| 3    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 91   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 72   |
| 5    |    | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 62   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111219\  
Data File : BF117762.D  
Acq On : 12 Nov 2019 16:29  
Operator : JU  
Sample : K5789-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

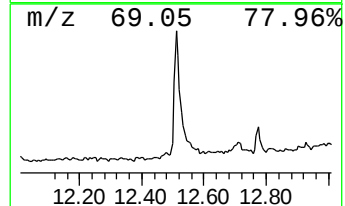
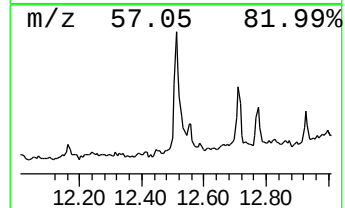
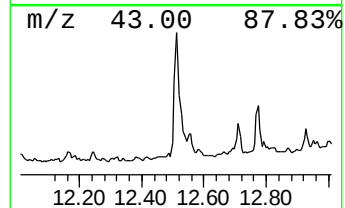
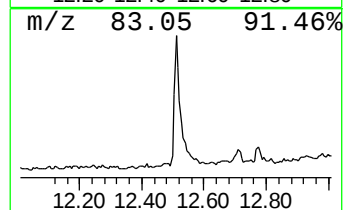
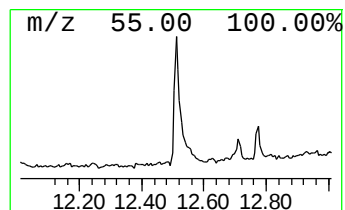
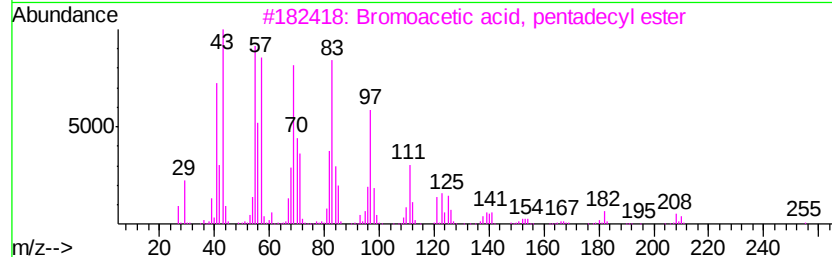
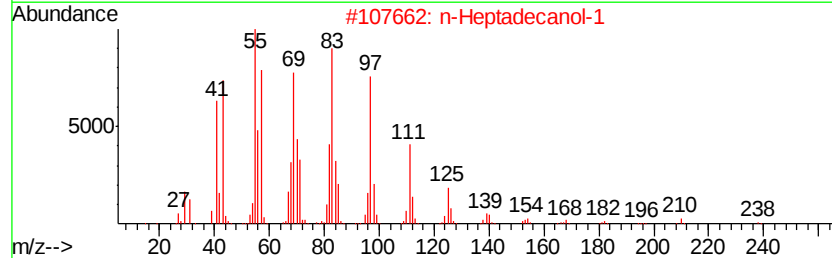
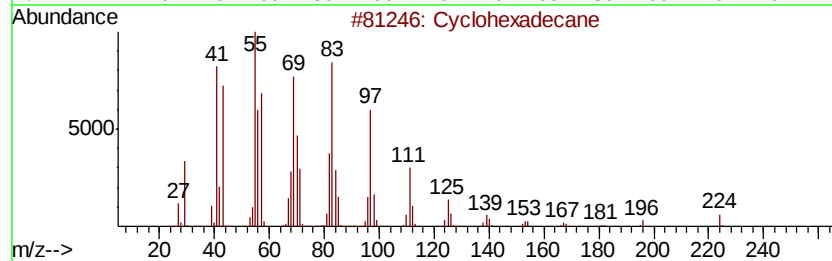
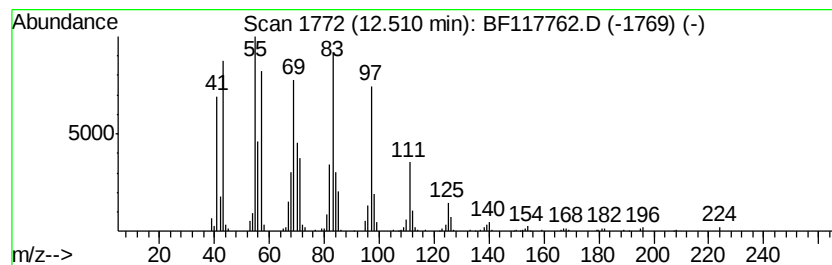
Instrument :  
BNA\_F  
ClientSampled :  
TP-17-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.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 Cyclohexadecane Concentration Rank 7

| R.T.      | EstConc                            | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|--------|------------------|-------------|------|
| 12.51     | 4.24 ng                            | 539380 | Phenanthrene-d10 | 11.46       |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm          | CAS#        | Qual |
| 1         | Cvclohexadecane                    | 224    | C16H32           | 000295-65-8 | 95   |
| 2         | n-Heptadecanol-1                   | 256    | C17H36O          | 001454-85-9 | 95   |
| 3         | Bromoacetic acid, pentadecyl ester | 348    | C17H33BrO2       | 131143-01-6 | 94   |
| 4         | 1-Hexadecanol                      | 242    | C16H34O          | 036653-82-4 | 94   |
| 5         | n-Nonadecanol-1                    | 284    | C19H40O          | 001454-84-8 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111219\  
Data File : BF117762.D  
Acq On : 12 Nov 2019 16:29  
Operator : JU  
Sample : K5789-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-17-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

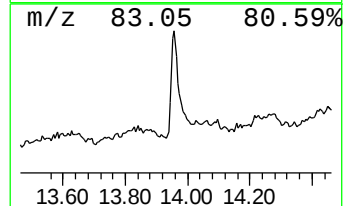
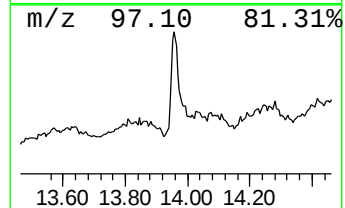
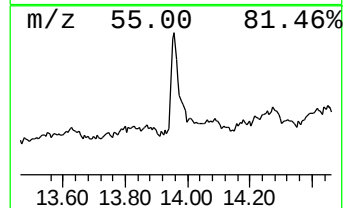
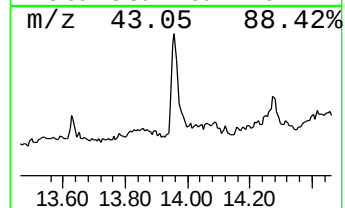
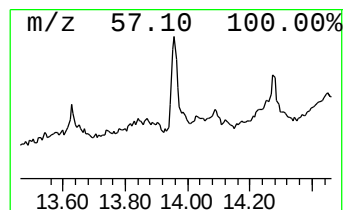
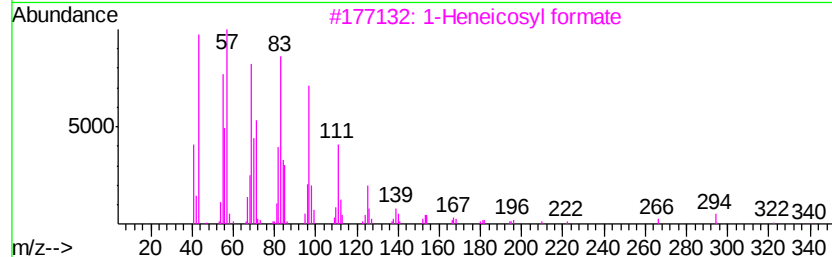
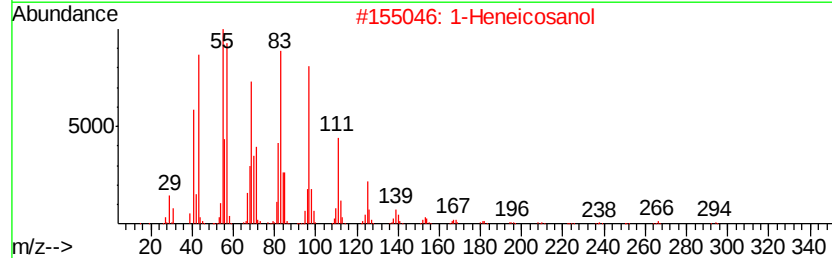
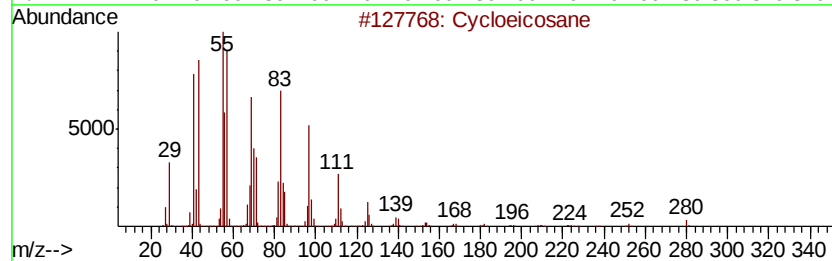
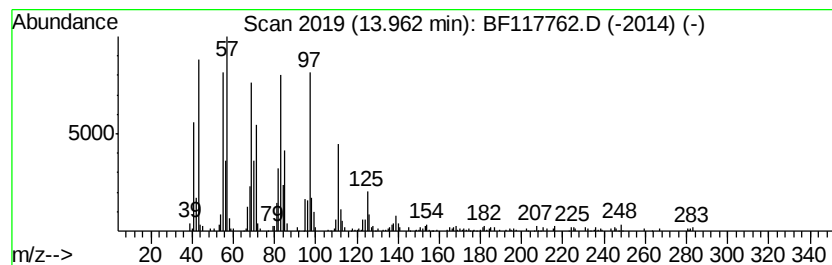
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Cycloeicosane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.96 | 7.69 ng | 704285 | Chrysene-d12     | 14.11 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|---------|---|-------------------------------------|-----|-------------|-------------|------|
| 1       |   | Cycloeicosane                       | 280 | C20H40      | 000296-56-0 | 98   |
| 2       |   | 1-Heneicosanol                      | 312 | C21H44O     | 015594-90-8 | 95   |
| 3       |   | 1-Heneicosyl formate                | 340 | C22H44O2    | 077899-03-7 | 94   |
| 4       |   | n-Tetracosanol-1                    | 354 | C24H50O     | 000506-51-4 | 94   |
| 5       |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF111219\  
Data File : BF117762.D  
Acq On : 12 Nov 2019 16:29  
Operator : JU  
Sample : K5789-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-17-S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110619.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... | 2.22  | 18.6    | ng    | 1494150  | 1                     | 6.93  | 1604070 | 20.0 |
| 2-Pentanone, 4-hy... | 5.14  | 13.1    | ng    | 1048020  | 1                     | 6.93  | 1604070 | 20.0 |
| unknown6.69          | 6.69  | 63.0    | ng    | 5054950  | 1                     | 6.93  | 1604070 | 20.0 |
| 2,4,7,9-Tetrameth... | 9.40  | 2.9     | ng    | 355435   | 3                     | 9.97  | 2472890 | 20.0 |
| n-Hexadecanoic acid  | 11.99 | 4.3     | ng    | 542394   | 4                     | 11.46 | 2544050 | 20.0 |
| Cyclohexadecane      | 12.51 | 4.2     | ng    | 539380   | 4                     | 11.46 | 2544050 | 20.0 |
| Cycloeicosane        | 13.96 | 7.7     | ng    | 704285   | 5                     | 14.11 | 1832010 | 20.0 |