

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092419\  
 Data File : BF117220.D  
 Acq On : 24 Sep 2019 17:20  
 Operator : JU  
 Sample : PB123268BL  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB123268BL

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-BF091319.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         | 2.116       | 3             | 5           | 14           | rVB      | 290092         | 293173        | 20.98%          | 2.878%        |
| 2         | 5.446       | 567           | 571         | 588          | rBV      | 971479         | 921856        | 65.96%          | 9.048%        |
| 3         | 6.457       | 738           | 743         | 746          | rBV      | 1002867        | 1000848       | 71.61%          | 9.824%        |
| 4         | 6.587       | 761           | 765         | 768          | rBV      | 1170368        | 1071669       | 76.68%          | 10.519%       |
| 5         | 6.810       | 800           | 803         | 812          | rBV      | 224174         | 204174        | 14.61%          | 2.004%        |
| 6         | 6.969       | 826           | 830         | 839          | rBV      | 923756         | 801497        | 57.35%          | 7.867%        |
| 7         | 7.375       | 895           | 899         | 916          | rBV      | 589324         | 639894        | 45.78%          | 6.281%        |
| 8         | 8.092       | 1017          | 1021        | 1028         | rBV      | 241990         | 247108        | 17.68%          | 2.425%        |
| 9         | 9.169       | 1199          | 1204        | 1207         | rBV      | 1390243        | 1259192       | 90.10%          | 12.360%       |
| 10        | 9.563       | 1267          | 1271        | 1283         | rBB      | 78230          | 76267         | 5.46%           | 0.749%        |
| 11        | 9.845       | 1315          | 1319        | 1330         | rBB      | 337371         | 301219        | 21.55%          | 2.957%        |
| 12        | 10.639      | 1450          | 1454        | 1470         | rBV      | 839621         | 807656        | 57.79%          | 7.928%        |
| 13        | 11.333      | 1568          | 1572        | 1584         | rBV      | 289300         | 286238        | 20.48%          | 2.810%        |
| 14        | 11.857      | 1659          | 1661        | 1665         | rBV      | 35653          | 33449         | 2.39%           | 0.328%        |
| 15        | 12.416      | 1749          | 1756        | 1771         | rBV4     | 16322          | 53689         | 3.84%           | 0.527%        |
| 16        | 12.922      | 1836          | 1842        | 1845         | rBV      | 1364834        | 1397607       | 100.00%         | 13.718%       |
| 17        | 13.804      | 1988          | 1992        | 1994         | rBV2     | 11979          | 15625         | 1.12%           | 0.153%        |
| 18        | 13.833      | 1994          | 1997        | 2014         | rVV2     | 65953          | 147071        | 10.52%          | 1.444%        |
| 19        | 13.974      | 2017          | 2021        | 2033         | rVB      | 209365         | 211374        | 15.12%          | 2.075%        |
| 20        | 14.427      | 2094          | 2098        | 2101         | rBV2     | 20613          | 18648         | 1.33%           | 0.183%        |
| 21        | 14.727      | 2145          | 2149        | 2154         | rBV      | 36958          | 34886         | 2.50%           | 0.342%        |
| 22        | 15.057      | 2196          | 2205        | 2211         | rBV      | 45067          | 51720         | 3.70%           | 0.508%        |
| 23        | 15.427      | 2262          | 2268        | 2280         | rBV      | 145977         | 215805        | 15.44%          | 2.118%        |
| 24        | 15.845      | 2330          | 2339        | 2347         | rBV      | 36398          | 58360         | 4.18%           | 0.573%        |
| 25        | 16.339      | 2414          | 2423        | 2427         | rBV2     | 21841          | 38941         | 2.79%           | 0.382%        |

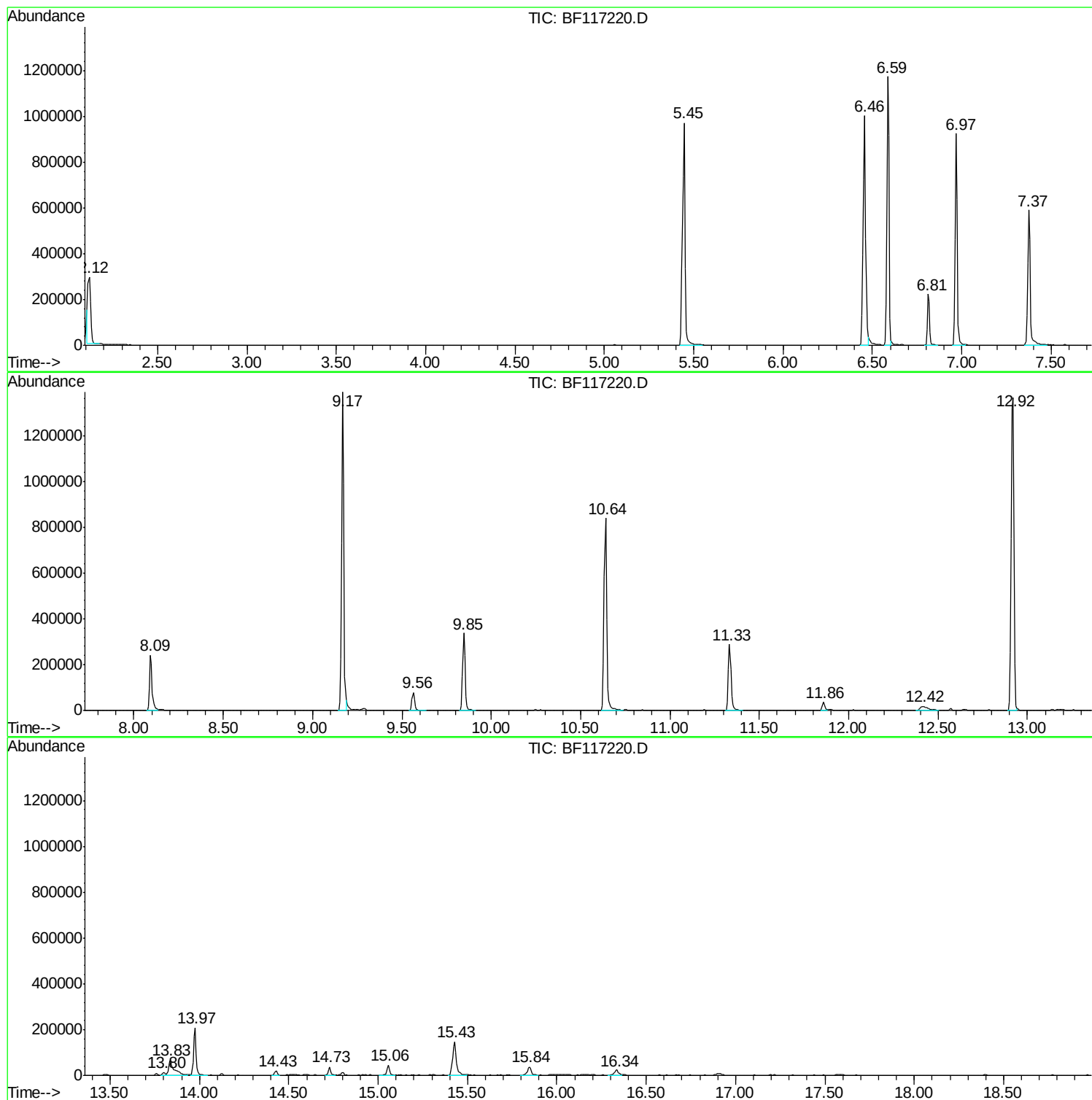
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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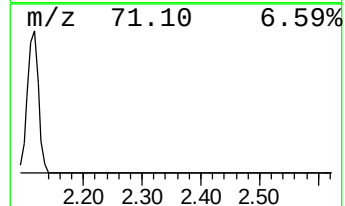
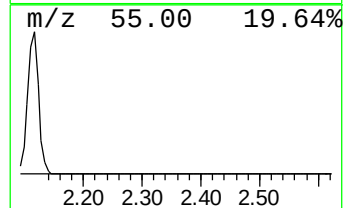
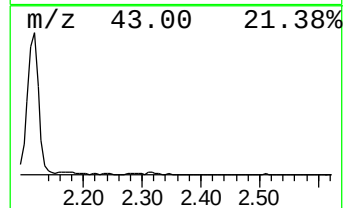
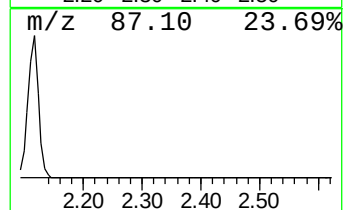
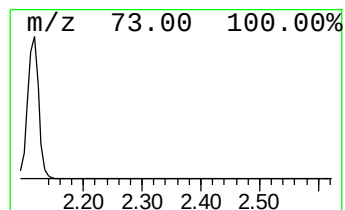
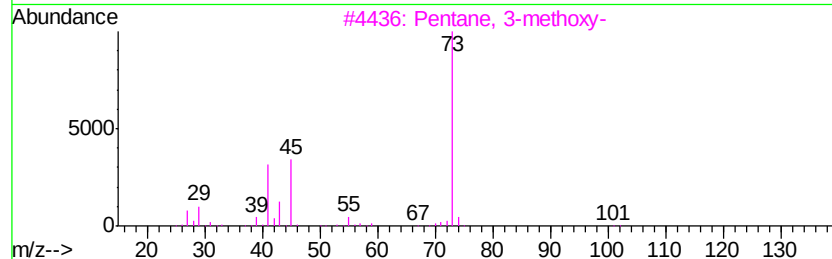
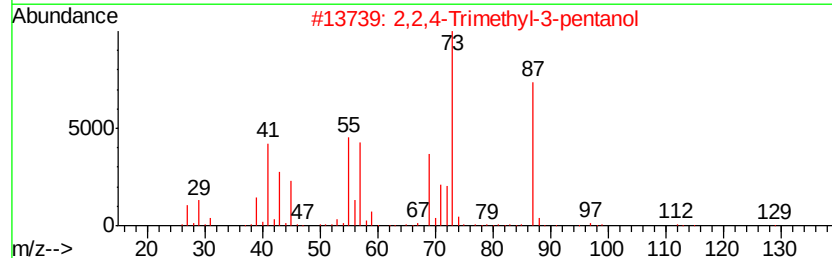
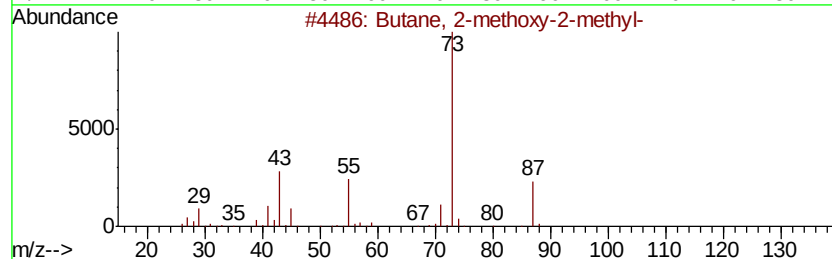
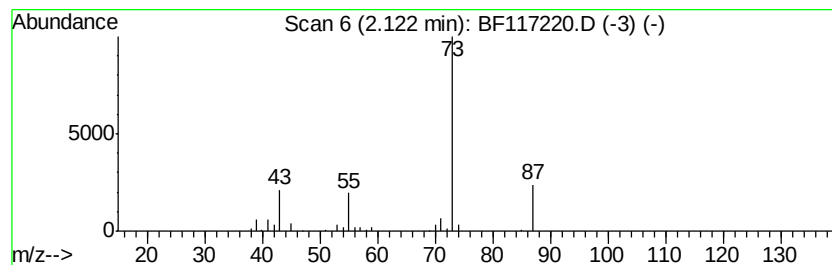
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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. |
|------|----------|--------|------------------------|------|
| 2.12 | 28.72 ng | 293173 | 1,4-Dichlorobenzene-d4 | 6.81 |

| 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    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



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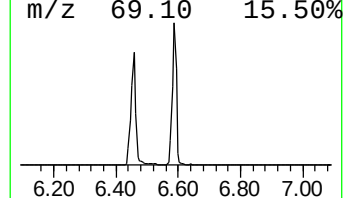
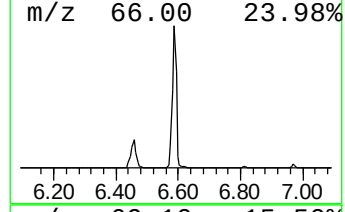
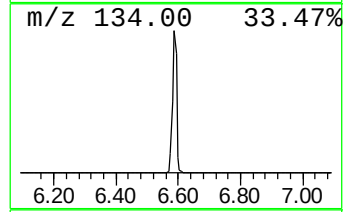
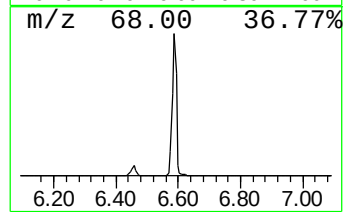
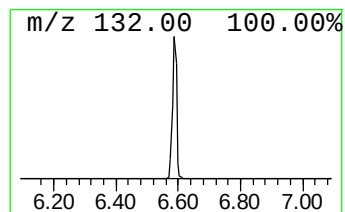
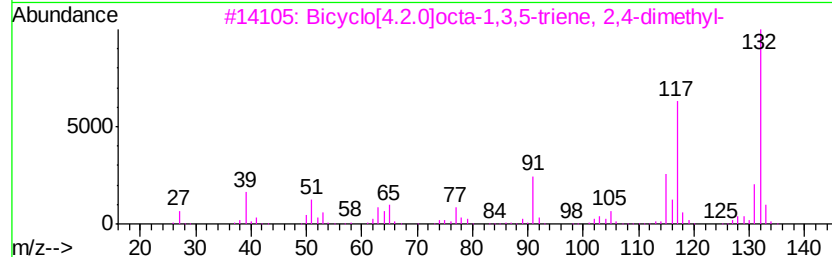
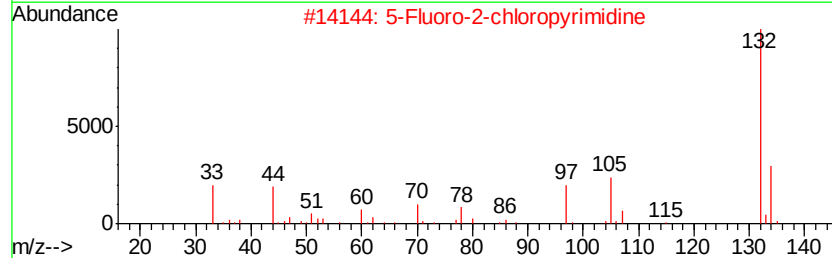
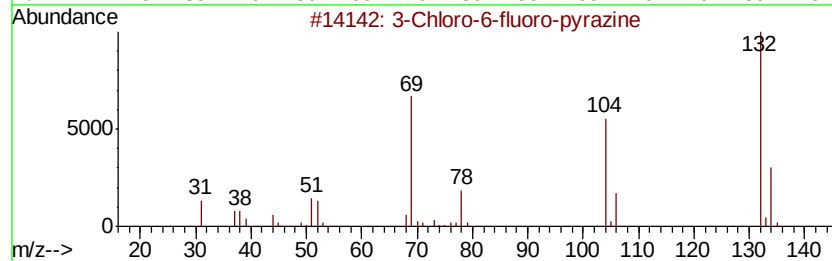
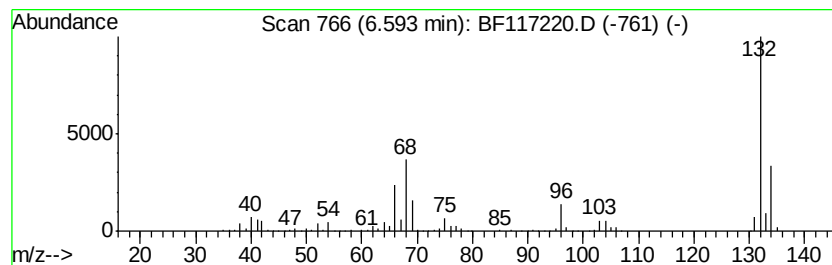
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.59 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 6.59 | 104.98 ng | 1071670 | 1,4-Dichlorobenzene-d4 | 6.81 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 2    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 12   |
| 3    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 4    |    | 1,3,4-Thiadiazole-2(3H)-thione, ... | 132 | C3H4N2S2  | 029490-19-5  | 9    |
| 5    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 9    |



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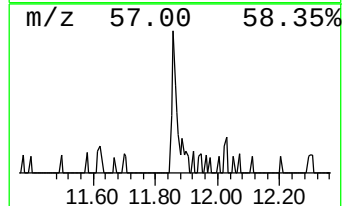
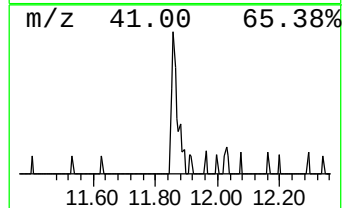
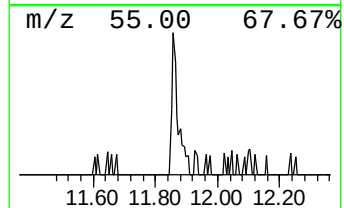
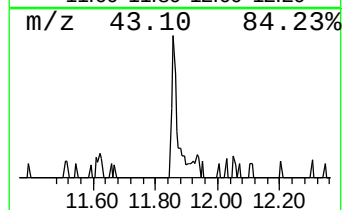
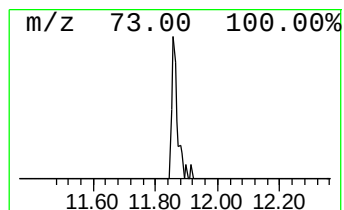
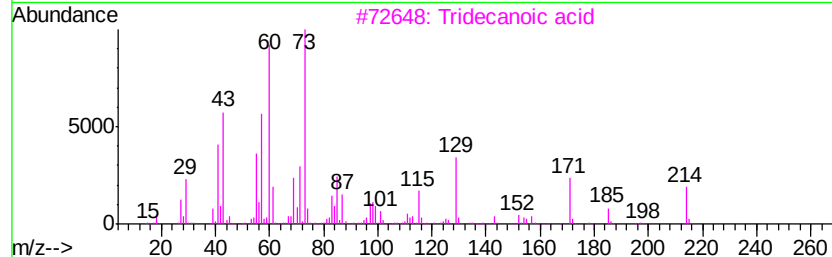
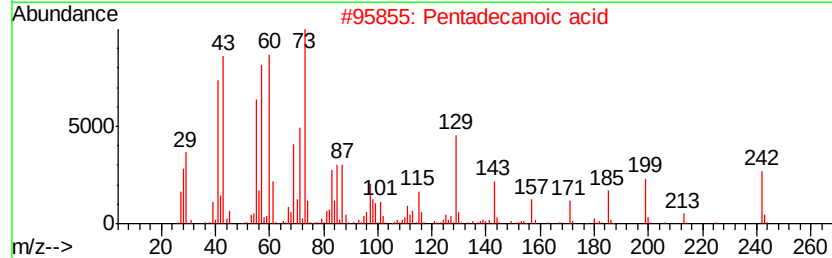
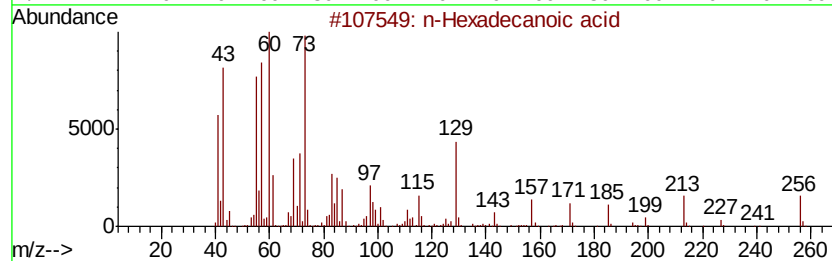
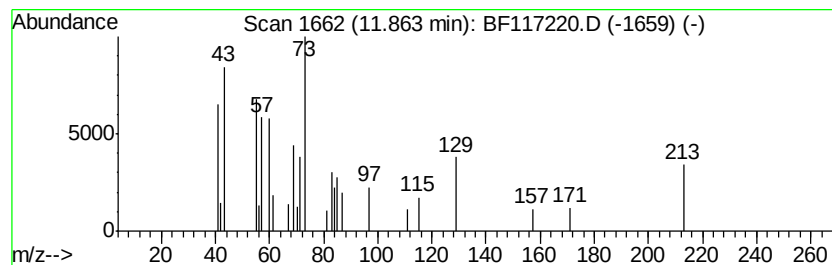
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.86 | 2.34 ng | 33449 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 86   |
| 2    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 72   |
| 3    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 53   |
| 4    |    | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 53   |
| 5    |    | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 53   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
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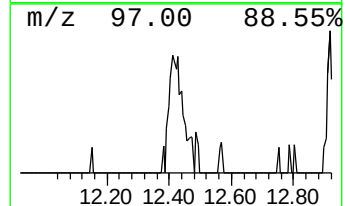
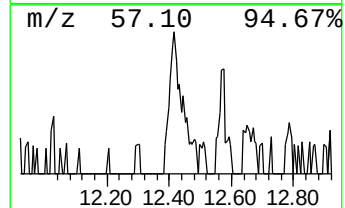
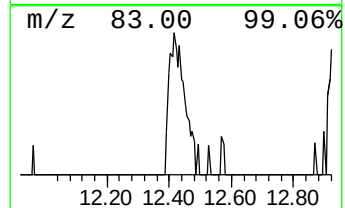
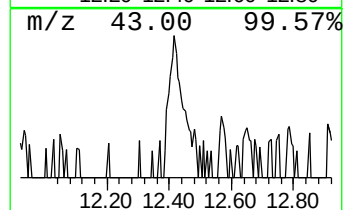
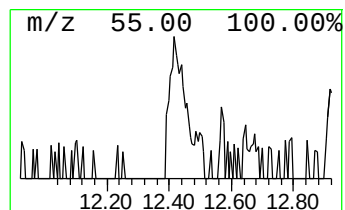
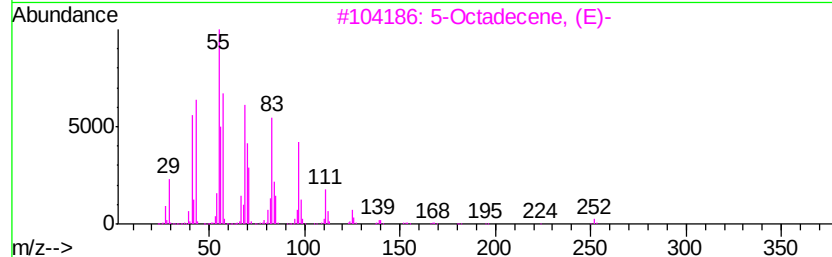
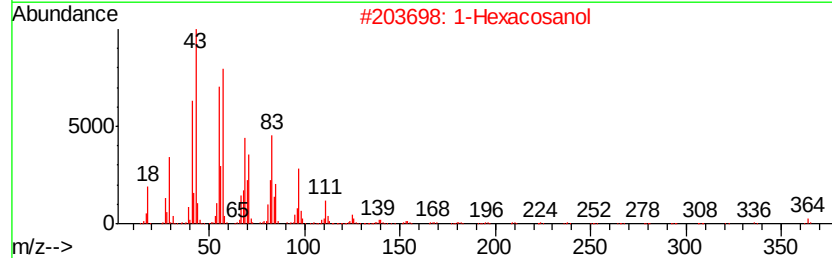
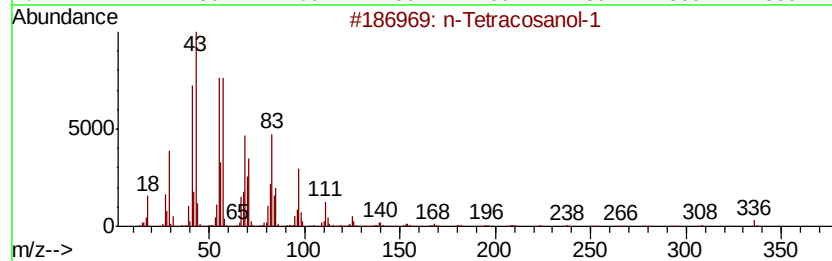
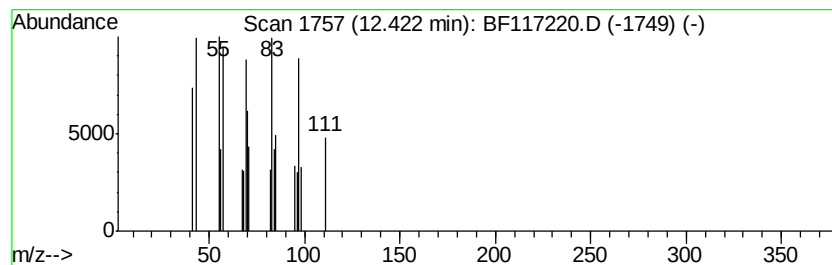
TIC Library : C:\DATABASE\NIST11.L  
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Peak Number 5 n-Tetracosanol-1 Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.42 | 3.75 ng | 53689 | Phenanthrene-d10 | 11.33 |

| Hit# of | 5 | Tentative ID                  | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------|-----|----------|-------------|------|
| 1       |   | n-Tetracosanol-1              | 354 | C24H50O  | 000506-51-4 | 91   |
| 2       |   | 1-Hexacosanol                 | 382 | C26H54O  | 000506-52-5 | 78   |
| 3       |   | 5-Octadecene, (E)-            | 252 | C18H36   | 007206-21-5 | 78   |
| 4       |   | Isobutyl tetradecyl carbonate | 314 | C19H38O3 | 959275-58-2 | 53   |
| 5       |   | 1-Tridecene                   | 182 | C13H26   | 002437-56-1 | 50   |



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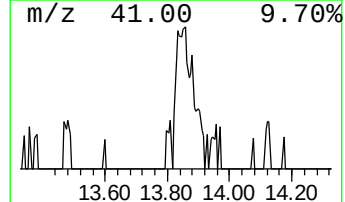
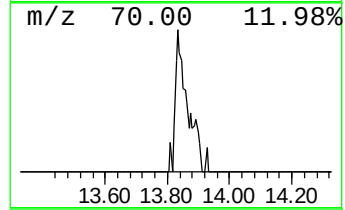
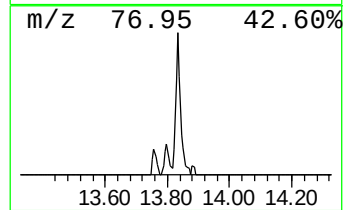
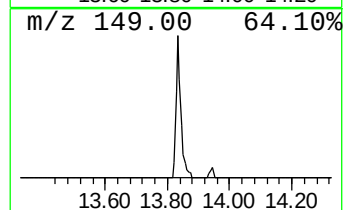
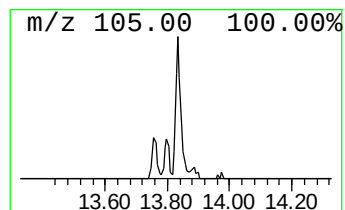
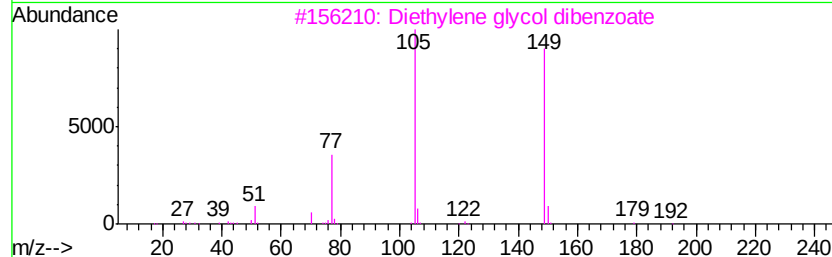
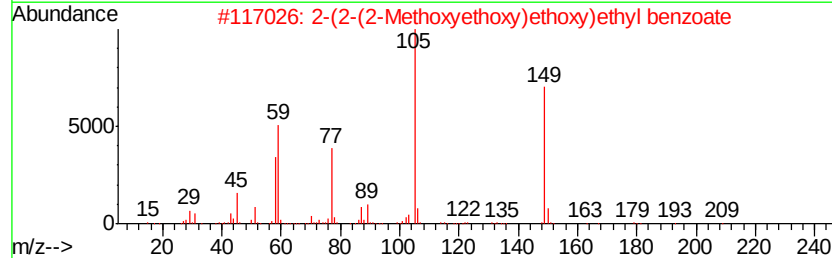
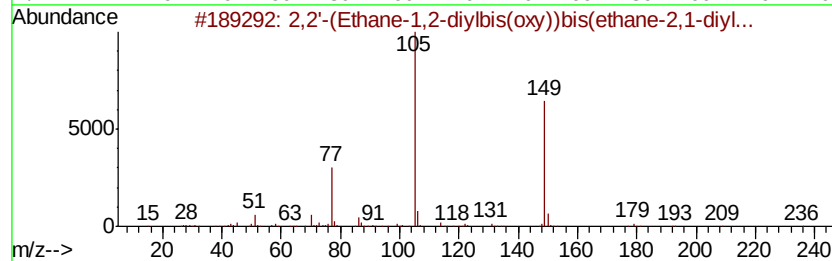
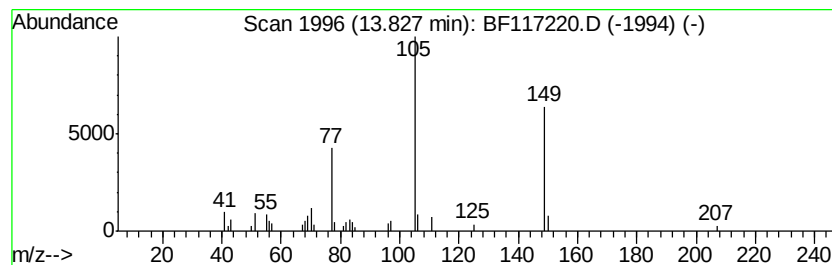
TIC Library : C:\DATABASE\NIST11.L  
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Peak Number 6 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.83 | 13.92 ng | 147071 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 2,2'-(Ethane-1,2-diylbis(oxv))bi... | 358 | C20H22O6   | 1000366-99-4 | 72   |
| 2    |    | 2-(2-(2-Methoxyethoxy)ethoxy)eth... | 268 | C14H20O5   | 1000366-99-1 | 64   |
| 3    |    | Diethylene glycol dibenzoate        | 314 | C18H18O5   | 000120-55-8  | 59   |
| 4    |    | 1,3-Dioxolane, 2-(methoxymethyl)... | 194 | C11H14O3   | 1000156-71-2 | 53   |
| 5    |    | Benzoic acid, 2-(2-chlorophenoxy... | 276 | C15H13ClO3 | 1000368-25-9 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092419\  
Data File : BF117220.D  
Acq On : 24 Sep 2019 17:20  
Operator : JU  
Sample : PB123268BL  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB123268BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

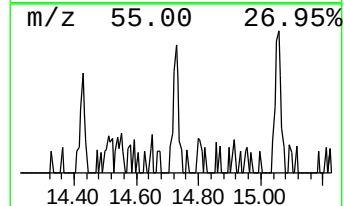
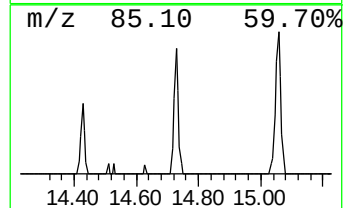
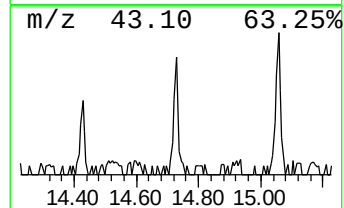
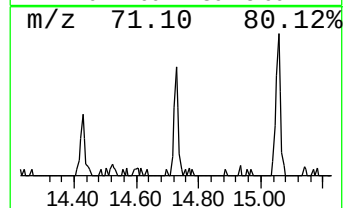
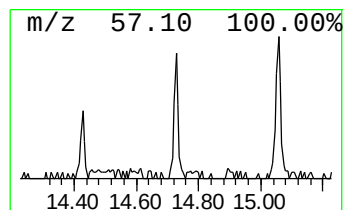
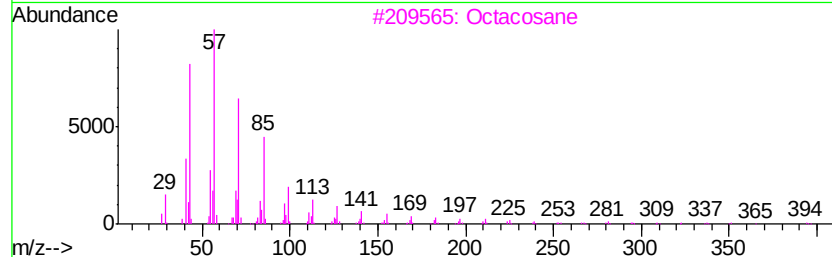
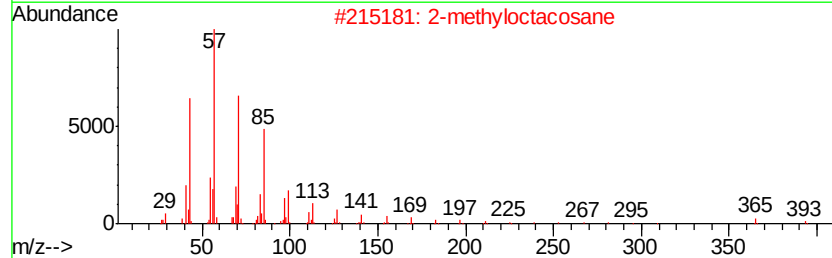
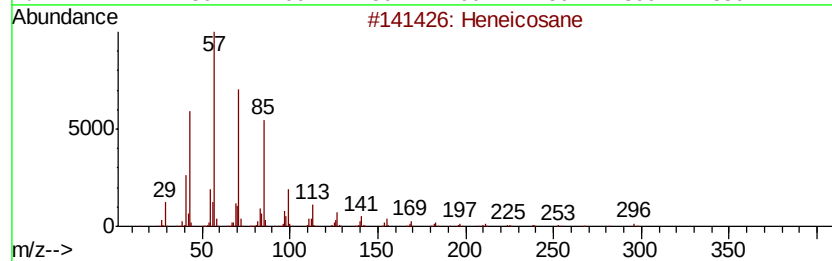
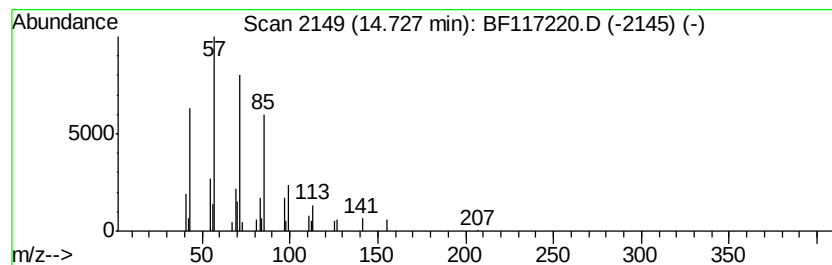
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Heneicosane Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.73 | 3.23 ng | 34886 | Perylene-d12     | 15.43 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|--------------------|--------------|-----|---------|--------------|------|
| 1       | Heneicosane        |              | 296 | C21H44  | 000629-94-7  | 91   |
| 2       | 2-methyloctacosane |              | 408 | C29H60  | 1000376-72-8 | 90   |
| 3       | Octacosane         |              | 394 | C28H58  | 000630-02-4  | 90   |
| 4       | Hexacosane         |              | 366 | C26H54  | 000630-01-3  | 87   |
| 5       | Heptacosane        |              | 380 | C27H56  | 000593-49-7  | 86   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092419\  
Data File : BF117220.D  
Acq On : 24 Sep 2019 17:20  
Operator : JU  
Sample : PB123268BL  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

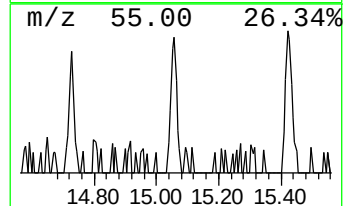
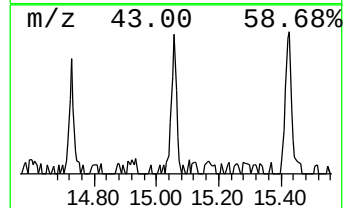
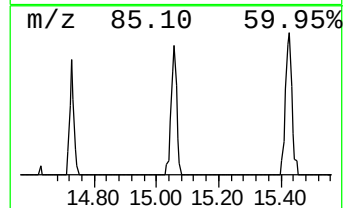
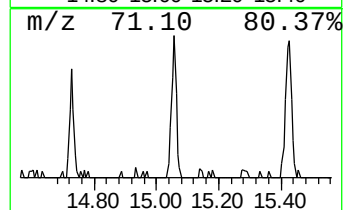
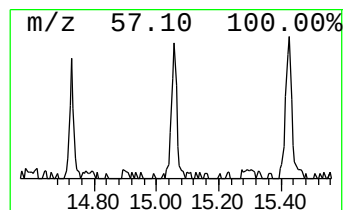
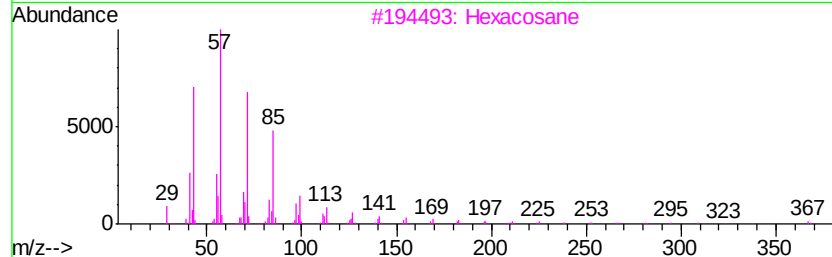
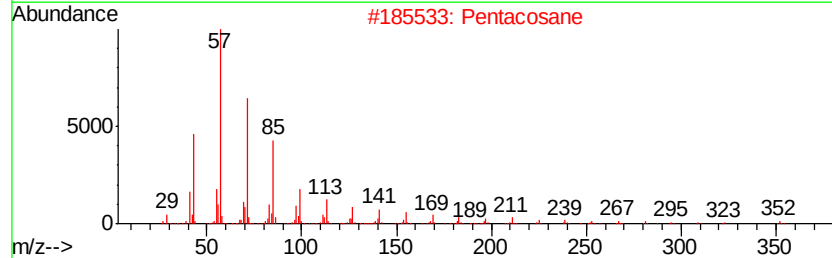
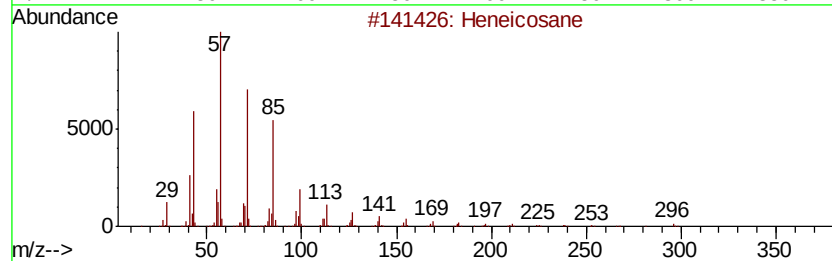
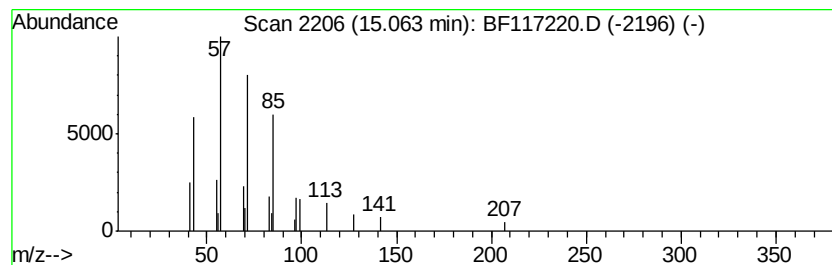
Instrument :  
BNA\_F  
ClientSampleId :  
PB123268BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.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 Pentacosane Concentration Rank 6

| R.T.    | EstConc            | Area         | Relative to ISTD | R.T.      |
|---------|--------------------|--------------|------------------|-----------|
| 15.06   | 4.79 ng            | 51720        | Perylene-d12     | 15.43     |
| Hit# of | 5                  | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Heneicosane        | 296 C21H44   | 000629-94-7      | 91        |
| 2       | Pentacosane        | 352 C25H52   | 000629-99-2      | 90        |
| 3       | Hexacosane         | 366 C26H54   | 000630-01-3      | 87        |
| 4       | Heptacosane        | 380 C27H56   | 000593-49-7      | 87        |
| 5       | 2-methyloctacosane | 408 C29H60   | 1000376-72-8     | 86        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092419\  
Data File : BF117220.D  
Acq On : 24 Sep 2019 17:20  
Operator : JU  
Sample : PB123268BL  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB123268BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

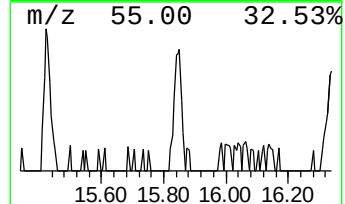
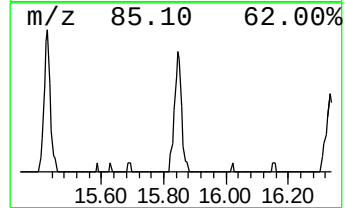
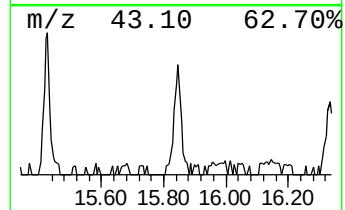
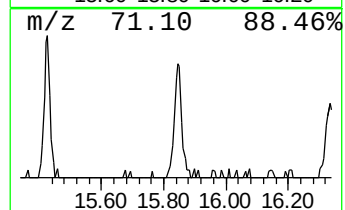
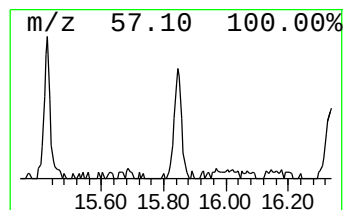
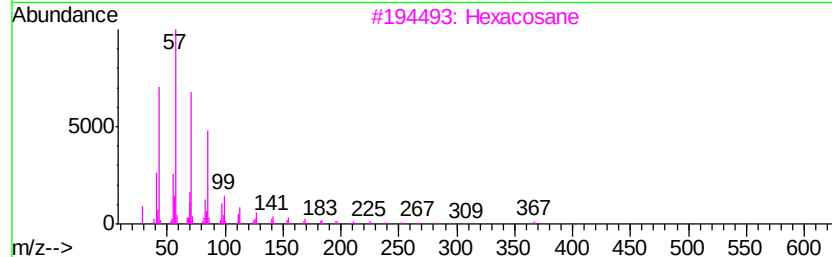
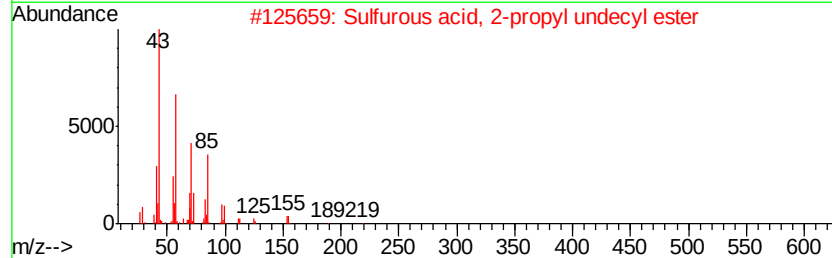
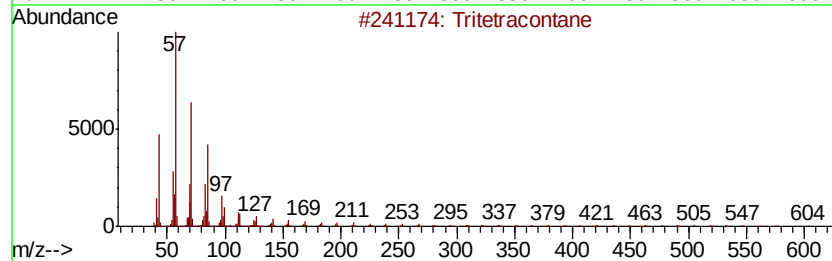
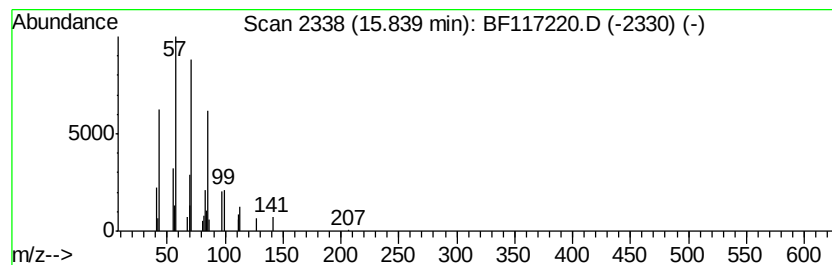
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Tritetracontane Concentration Rank 5

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.84 | 5.41 ng | 58360 | Perylene-d12     | 15.43 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Tritetracontane                     | 605 | C43H88    | 007098-21-7  | 90   |
| 2    |    | Sulfurous acid, 2-propyl undecyl... | 278 | C14H30O3S | 1000309-12-2 | 86   |
| 3    |    | Hexacosane                          | 366 | C26H54    | 000630-01-3  | 86   |
| 4    |    | Hentriacontane                      | 437 | C31H64    | 000630-04-6  | 86   |
| 5    |    | Pentatriacontane                    | 493 | C35H72    | 000630-07-9  | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092419\  
Data File : BF117220.D  
Acq On : 24 Sep 2019 17:20  
Operator : JU  
Sample : PB123268BL  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB123268BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

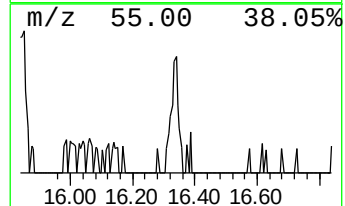
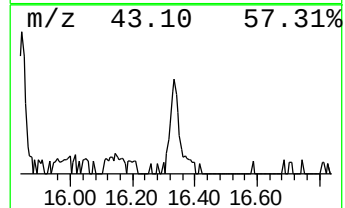
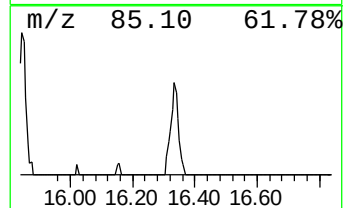
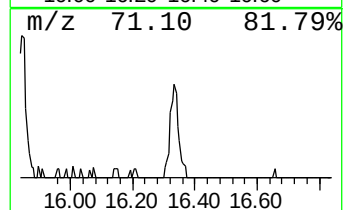
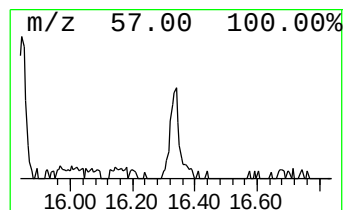
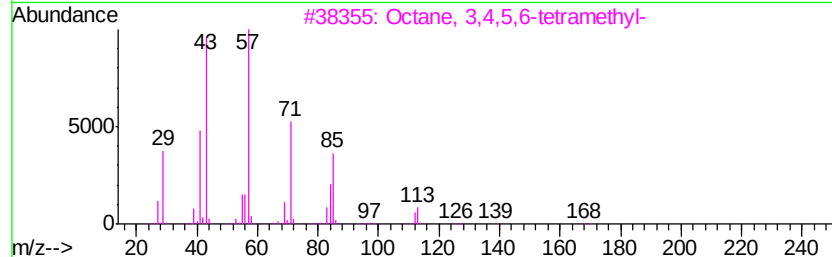
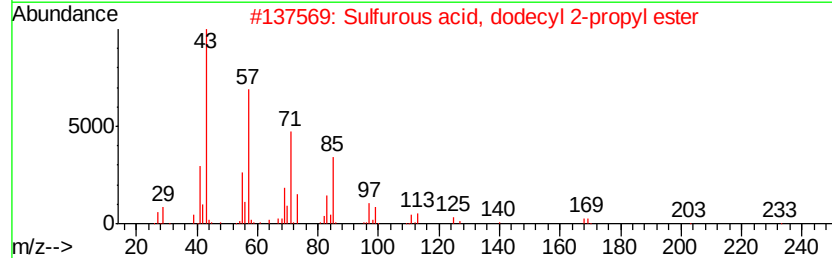
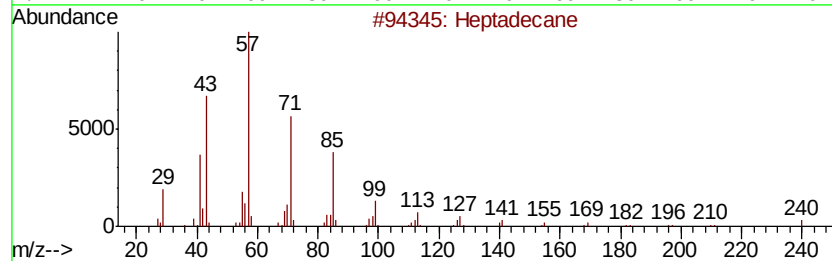
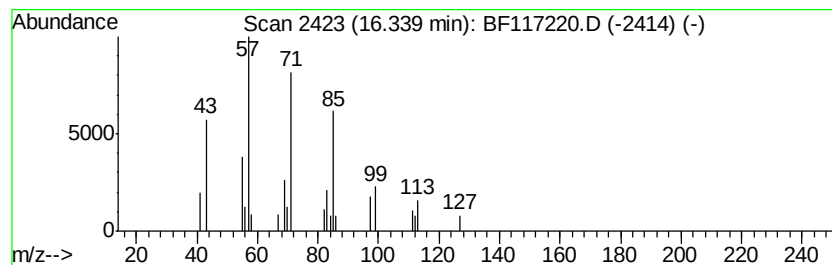
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Heptadecane Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.34 | 3.61 ng | 38941 | Perylene-d12     | 15.43 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Heptadecane                         | 240 | C17H36    | 000629-78-7  | 72   |
| 2    |    | Sulfurous acid, dodecyl 2-propyl... | 292 | C15H32O3S | 1000309-12-3 | 72   |
| 3    |    | Octane, 3,4,5,6-tetramethyl-        | 170 | C12H26    | 062185-21-1  | 59   |
| 4    |    | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 53   |
| 5    |    | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF092419\  
Data File : BF117220.D  
Acq On : 24 Sep 2019 17:20  
Operator : JU  
Sample : PB123268BL  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB123268BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF091319.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... | 2.12  | 28.7    | ng    | 293173   | 1                     | 6.81  | 204174 | 20.0 |
| unknown6.59          | 6.59  | 105.0   | ng    | 1071670  | 1                     | 6.81  | 204174 | 20.0 |
| n-Hexadecanoic acid  | 11.86 | 2.3     | ng    | 33449    | 4                     | 11.33 | 286238 | 20.0 |
| n-Tetracosanol-1     | 12.42 | 3.8     | ng    | 53689    | 4                     | 11.33 | 286238 | 20.0 |
| 2,2'-(Ethane-1,2-... | 13.83 | 13.9    | ng    | 147071   | 5                     | 13.97 | 211374 | 20.0 |
| Heneicosane          | 14.73 | 3.2     | ng    | 34886    | 6                     | 15.43 | 215805 | 20.0 |
| Pentacosane          | 15.06 | 4.8     | ng    | 51720    | 6                     | 15.43 | 215805 | 20.0 |
| Tritetracontane      | 15.84 | 5.4     | ng    | 58360    | 6                     | 15.43 | 215805 | 20.0 |
| Heptadecane          | 16.34 | 3.6     | ng    | 38941    | 6                     | 15.43 | 215805 | 20.0 |