

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF082820\  
 Data File : BF121461.D  
 Acq On : 28 Aug 2020 15:54  
 Operator : JU/CG  
 Sample : L3824-01 10X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OIL-1

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-BF082720.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           | 32           | rVB4     | 148121         | 366509        | 7.23%           | 1.530%        |
| 2         | 5.463       | 569           | 574         | 578          | rVB      | 86631          | 84540         | 1.67%           | 0.353%        |
| 3         | 6.469       | 741           | 745         | 748          | rBV      | 112742         | 86632         | 1.71%           | 0.362%        |
| 4         | 6.857       | 807           | 811         | 815          | rBV      | 2065020        | 1732403       | 34.16%          | 7.231%        |
| 5         | 8.139       | 1025          | 1029        | 1033         | rBV      | 2442483        | 2308425       | 45.52%          | 9.635%        |
| 6         | 9.216       | 1209          | 1212        | 1215         | rBV      | 196677         | 178437        | 3.52%           | 0.745%        |
| 7         | 9.298       | 1223          | 1226        | 1230         | rBV3     | 229410         | 297648        | 5.87%           | 1.242%        |
| 8         | 9.369       | 1230          | 1238        | 1240         | rBV9     | 117076         | 292300        | 5.76%           | 1.220%        |
| 9         | 9.610       | 1277          | 1279        | 1281         | rBV      | 581766         | 613300        | 12.09%          | 2.560%        |
| 10        | 9.704       | 1293          | 1295        | 1296         | rBV      | 194147         | 151417        | 2.99%           | 0.632%        |
| 11        | 9.775       | 1303          | 1307        | 1309         | rBV2     | 968672         | 1244614       | 24.54%          | 5.195%        |
| 12        | 9.816       | 1311          | 1314        | 1316         | rVB      | 1492396        | 1315551       | 25.94%          | 5.491%        |
| 13        | 9.904       | 1322          | 1329        | 1332         | rBV3     | 4040796        | 5071692       | 100.00%         | 21.168%       |
| 14        | 9.945       | 1334          | 1336        | 1338         | rBV2     | 440966         | 524049        | 10.33%          | 2.187%        |
| 15        | 10.163      | 1371          | 1373        | 1376         | rBV2     | 880291         | 950482        | 18.74%          | 3.967%        |
| 16        | 10.275      | 1390          | 1392        | 1395         | rBV3     | 716304         | 731667        | 14.43%          | 3.054%        |
| 17        | 10.316      | 1396          | 1399        | 1402         | rBV2     | 1891207        | 1878299       | 37.03%          | 7.840%        |
| 18        | 11.392      | 1579          | 1582        | 1585         | rBV      | 1952931        | 1789094       | 35.28%          | 7.467%        |
| 19        | 14.021      | 2026          | 2029        | 2033         | rVB      | 2160922        | 1962602       | 38.70%          | 8.191%        |
| 20        | 14.304      | 2075          | 2077        | 2078         | rVB      | 223826         | 88427         | 1.74%           | 0.369%        |
| 21        | 14.792      | 2157          | 2160        | 2166         | rVB      | 296926         | 323016        | 6.37%           | 1.348%        |
| 22        | 14.886      | 2173          | 2176        | 2180         | rVB      | 283729         | 286966        | 5.66%           | 1.198%        |
| 23        | 15.486      | 2274          | 2278        | 2283         | rVB      | 1353906        | 1681320       | 33.15%          | 7.017%        |

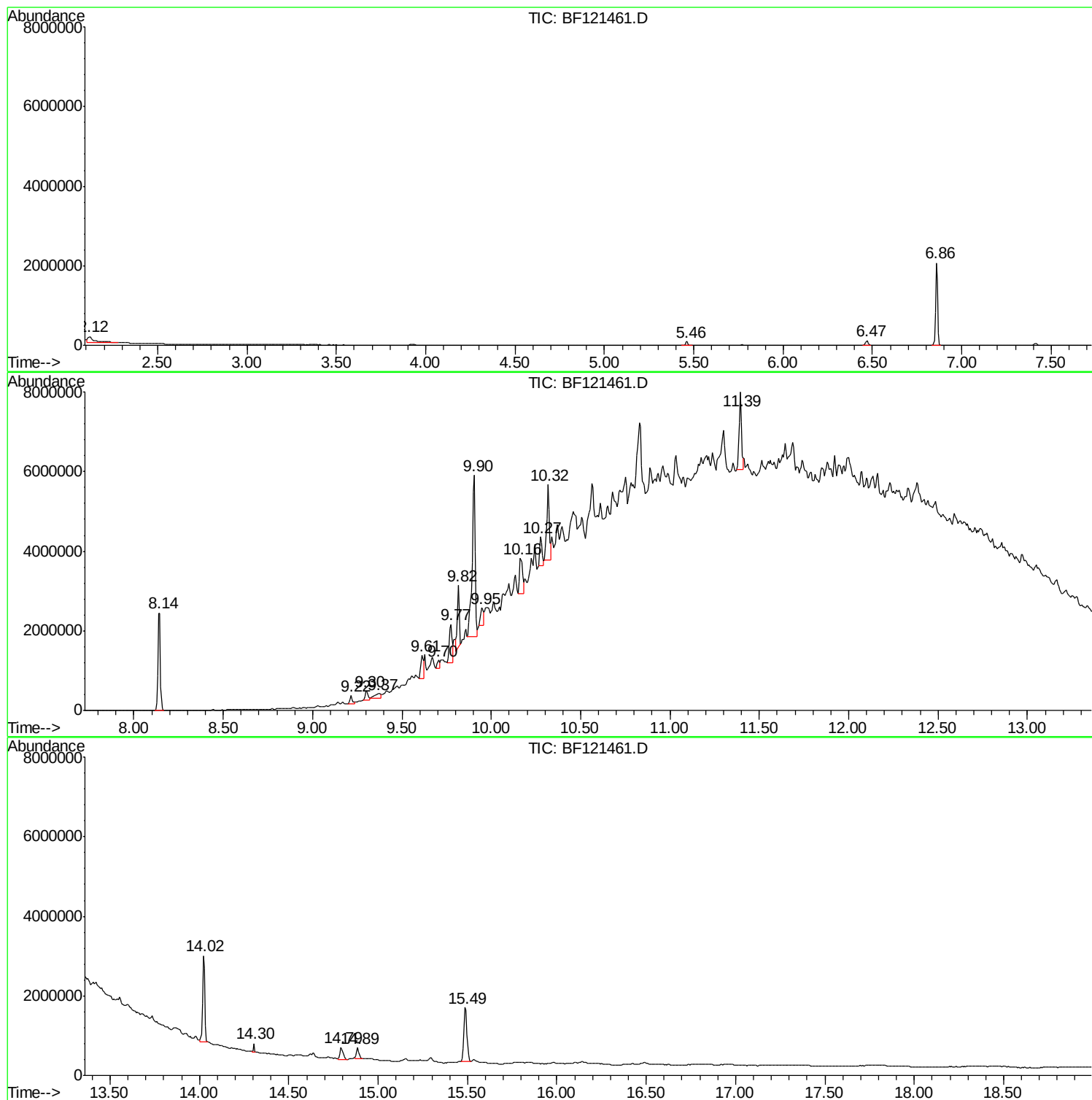
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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.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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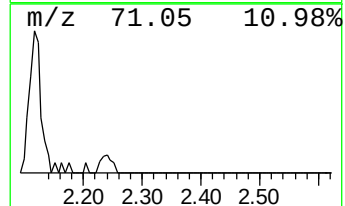
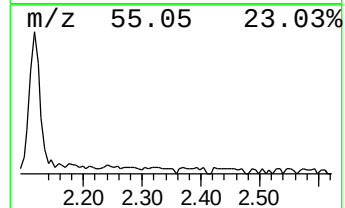
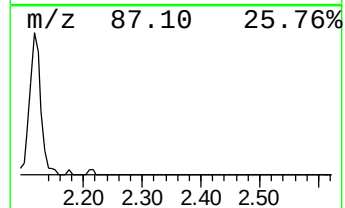
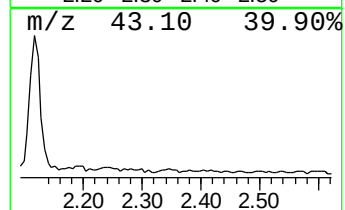
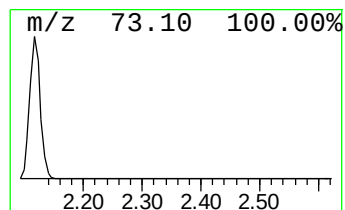
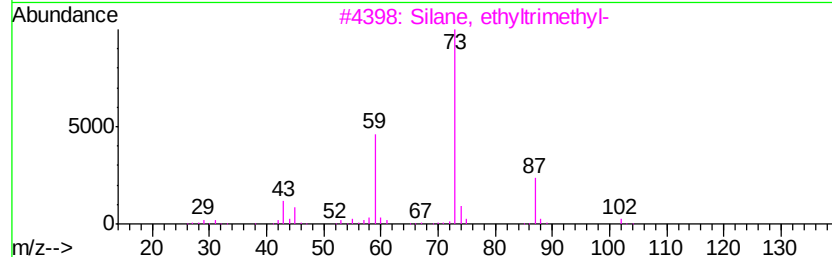
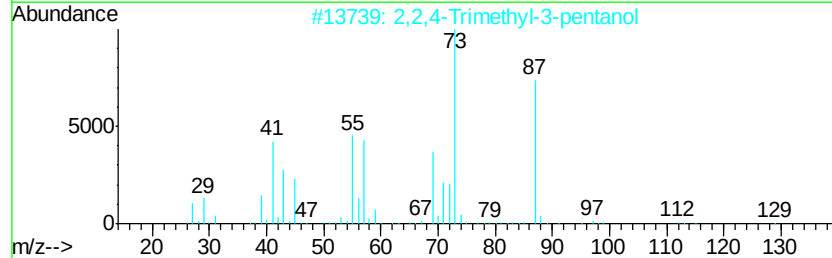
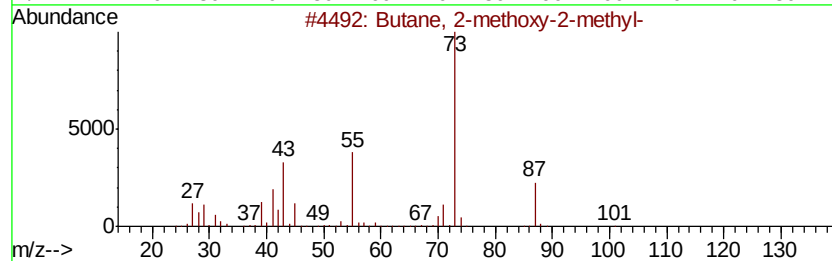
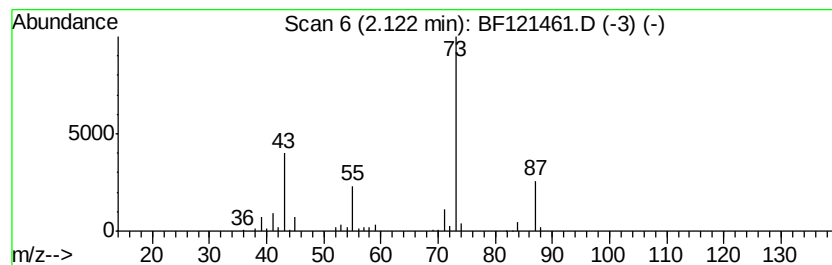
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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 4

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 2.12 | 4.23 ng | 366509 | 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 | 72   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol          | 130 | C8H18O   | 005162-48-1 | 40   |
| 3    |    | Silane, ethyltrimethyl-             | 102 | C5H14Si  | 003439-38-1 | 33   |
| 4    |    | Boronic acid, ethyl-, dimethyl e... | 102 | C4H11BO2 | 007318-82-3 | 10   |
| 5    |    | N-Ethylformamide                    | 73  | C3H7NO   | 000627-45-2 | 9    |



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Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

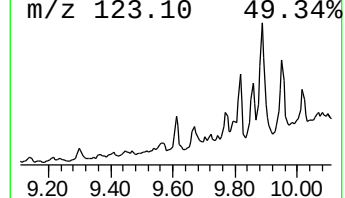
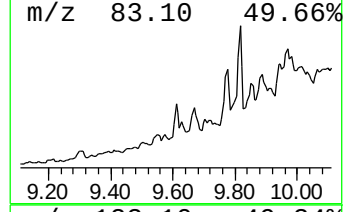
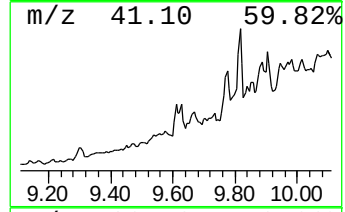
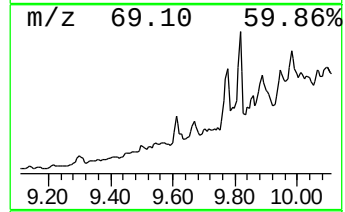
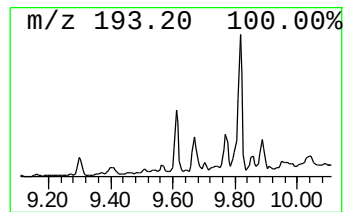
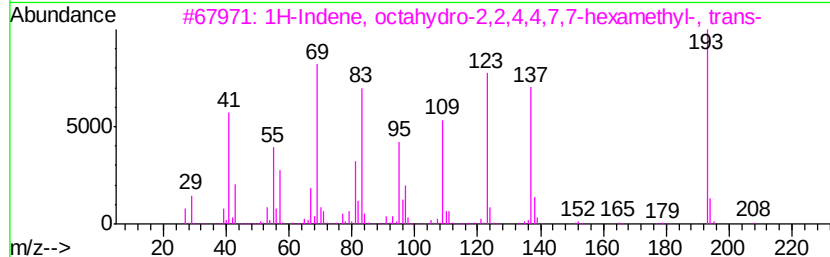
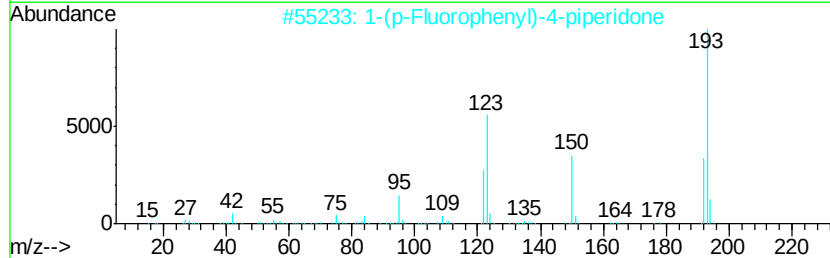
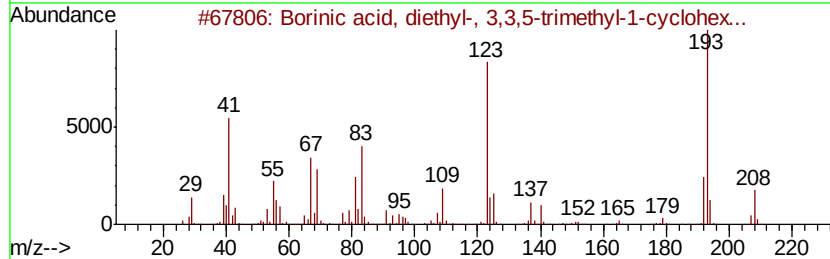
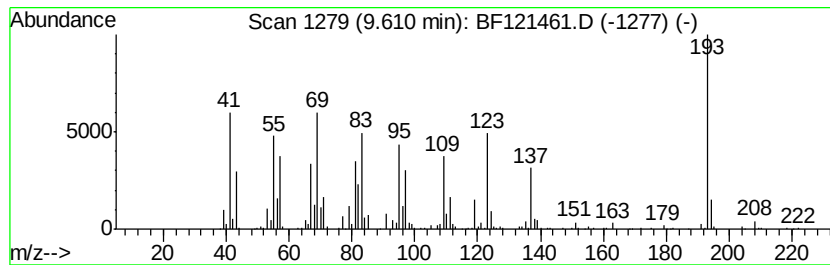
Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.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 unknown9.61 Concentration Rank 9

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.      |              |      |
|---------|---------|-------------------------------------|------------------|-----------|--------------|------|
| 9.61    | 2.42 ng | 613300                              | Acenaphthene-d10 | 9.90      |              |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm   | CAS#         | Qual |
| 1       |         | Borinic acid, diethyl-, 3,3,5-tr... | 208              | C13H25B0  | 057387-76-5  | 47   |
| 2       |         | 1-(p-Fluorophenyl)-4-piperidone     | 193              | C11H12FN0 | 1000238-56-7 | 46   |
| 3       |         | 1H-Indene, octahydro-2,2,4,4,7,7... | 208              | C15H28    | 054832-83-6  | 40   |
| 4       |         | 2-Phenylindolizine                  | 193              | C14H11N   | 025379-20-8  | 25   |
| 5       |         | 1H-Pyrazole, 3,5-bis(1,1-dimethy... | 208              | C13H24N2  | 125281-21-2  | 25   |



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ALS Vial : 11 Sample Multiplier: 1

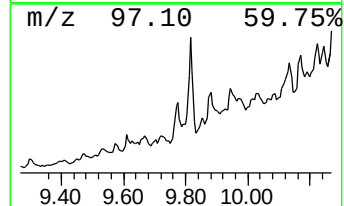
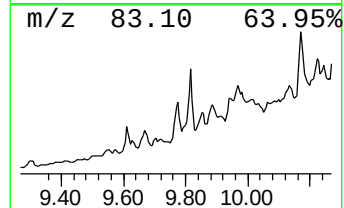
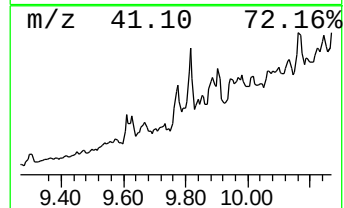
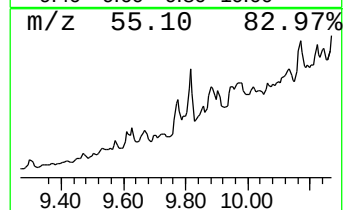
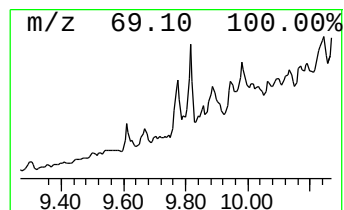
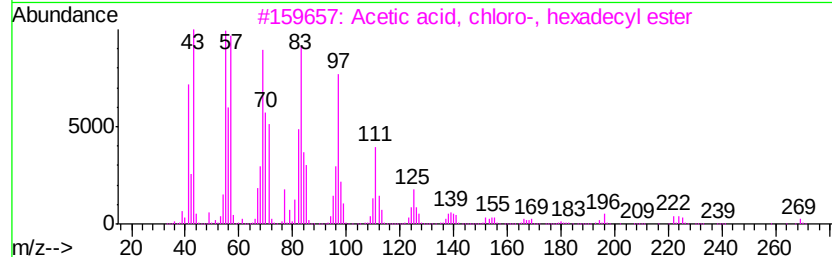
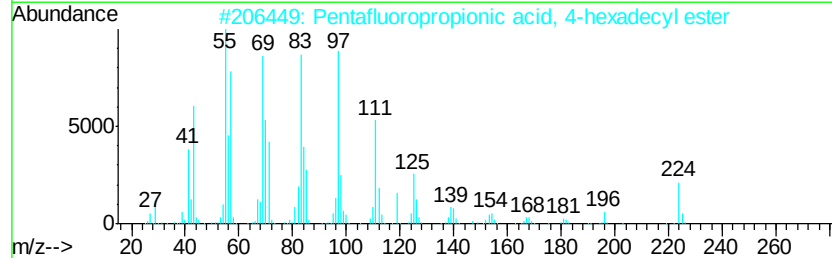
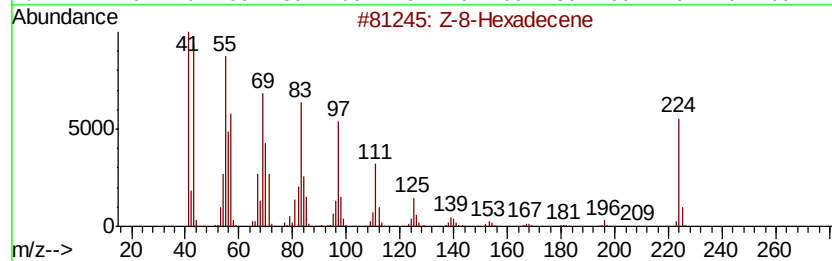
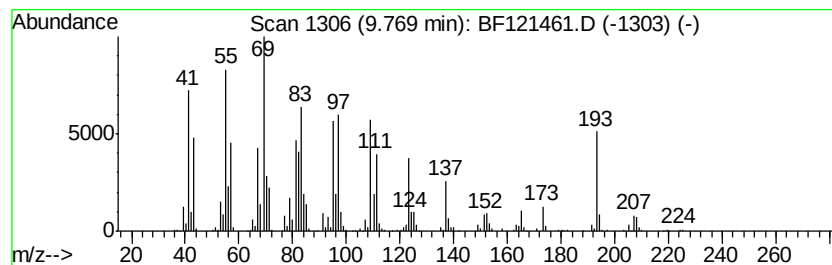
Instrument :  
BNA\_F  
ClientSampled :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.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 unknown9.77 Concentration Rank 3

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.        |              |      |
|---------|---------|-------------------------------------|------------------|-------------|--------------|------|
| 9.77    | 4.91 ng | 1244610                             | Acenaphthene-d10 | 9.90        |              |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm     | CAS#         | Qual |
| 1       |         | Z-8-Hexadecene                      | 224              | C16H32      | 1000130-87-5 | 44   |
| 2       |         | Pentafluoropropionic acid, 4-hex... | 388              | C19H33F5O2  | 1000283-04-1 | 38   |
| 3       |         | Acetic acid, chloro-, hexadecyl ... | 318              | C18H35ClO2  | 052132-58-8  | 38   |
| 4       |         | Carbonic acid, tetradecyl 2,2,2-... | 388              | C17H31Cl3O3 | 1000314-56-0 | 30   |
| 5       |         | n-Nonadecanol-1                     | 284              | C19H40O     | 001454-84-8  | 30   |



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ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

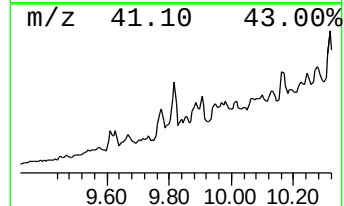
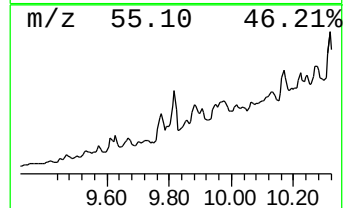
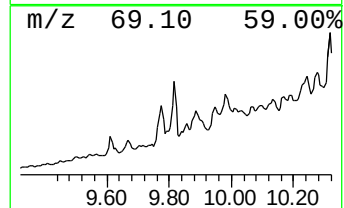
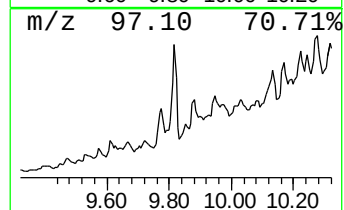
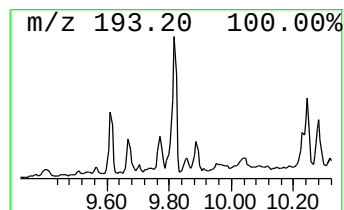
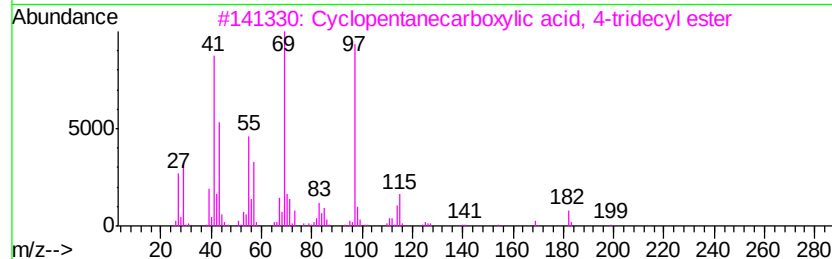
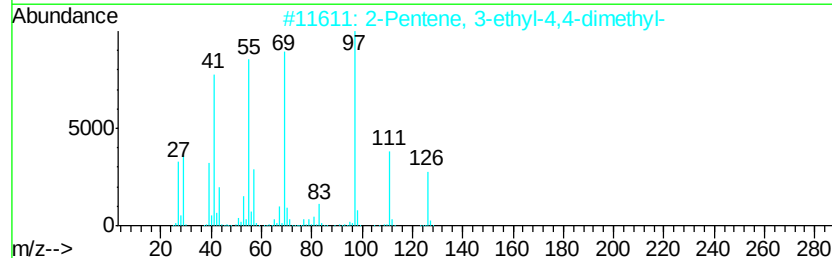
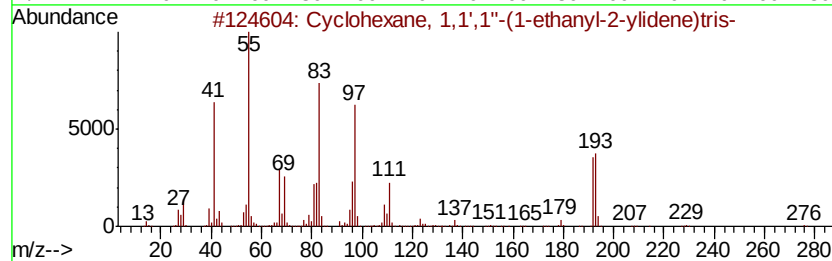
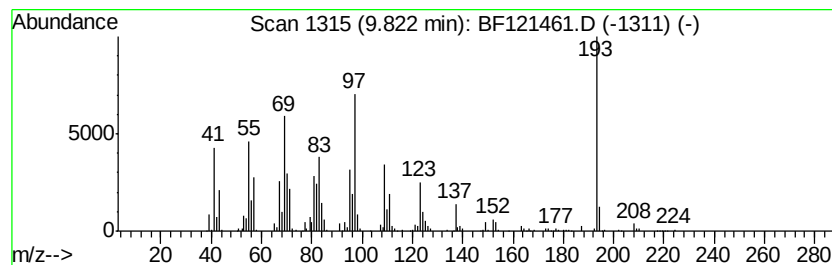
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown9.82 Concentration Rank 2

| R.T. | EstConc | Area    | Relative to ISTD | R.T. |
|------|---------|---------|------------------|------|
| 9.82 | 5.19 ng | 1315550 | Acenaphthene-d10 | 9.90 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | Cyclohexane, 1,1',1''-(1-ethanyl... | 276 | C20H36   | 055682-86-5  | 35   |
| 2       |   | 2-Pentene, 3-ethyl-4,4-dimethyl-    | 126 | C9H18    | 053907-59-8  | 30   |
| 3       |   | Cyclopentanecarboxylic acid, 4-t... | 296 | C19H36O2 | 1000280-58-6 | 27   |
| 4       |   | Cyclopentanecarboxylic acid, 3-p... | 324 | C21H40O2 | 1000280-59-1 | 27   |
| 5       |   | 1,2-Dodecanediol                    | 202 | C12H26O2 | 001119-87-5  | 27   |



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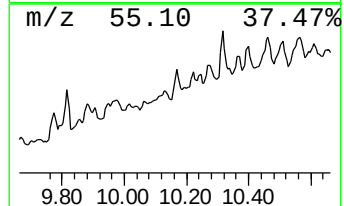
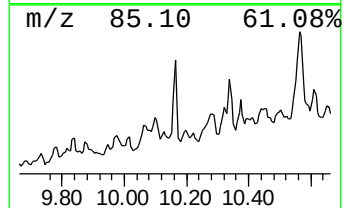
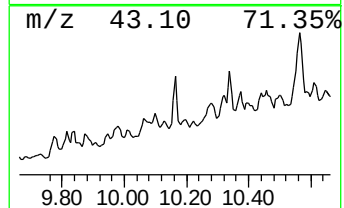
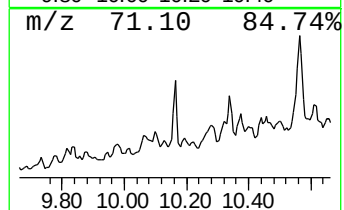
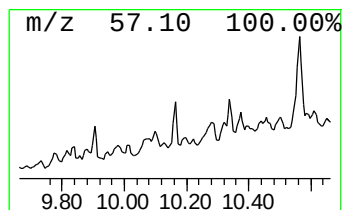
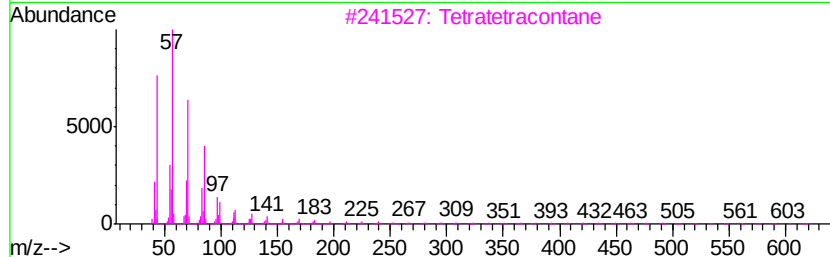
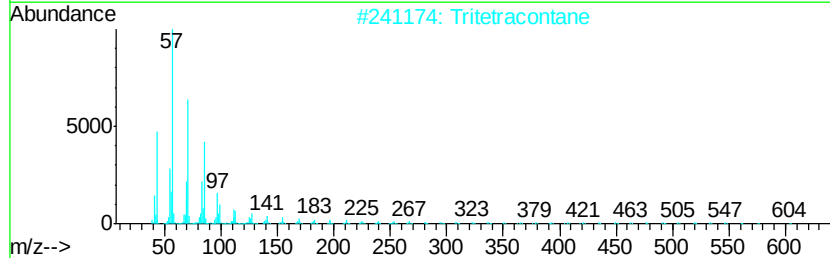
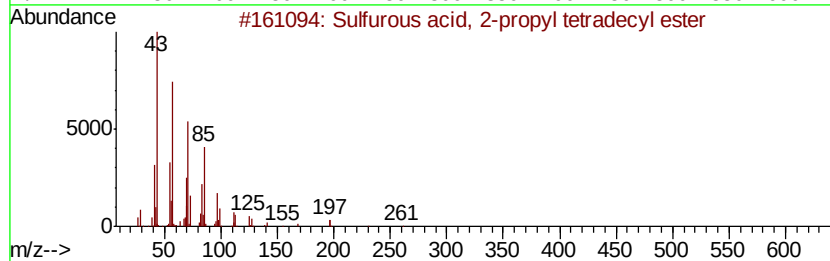
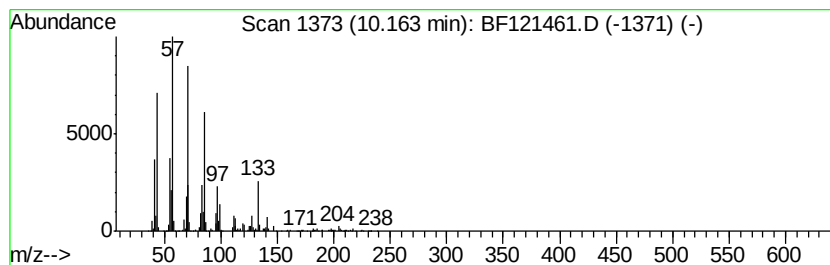
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Sulfurous acid, 2-propyl te... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.16 | 3.75 ng | 950482 | Acenaphthene-d10 | 9.90 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|--------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Sulfurous acid, 2-propyl tetradec... | 320 | C17H36O3S | 1000309-12-5 | 72   |
| 2    |    | Tritetracontane                      | 605 | C43H88    | 007098-21-7  | 72   |
| 3    |    | Tetratetracontane                    | 619 | C44H90    | 007098-22-8  | 58   |
| 4    |    | Tridecane, 7-propyl-                 | 226 | C16H34    | 055045-09-5  | 58   |
| 5    |    | Tridecane                            | 184 | C13H28    | 000629-50-5  | 58   |



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Data File : BF121461.D  
Acq On : 28 Aug 2020 15:54  
Operator : JU/CG  
Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

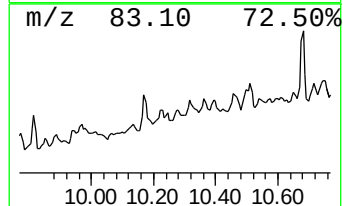
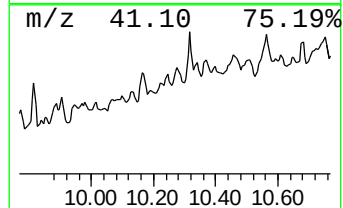
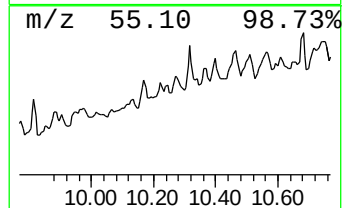
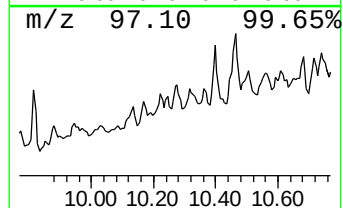
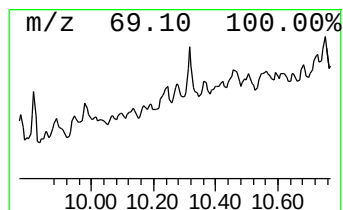
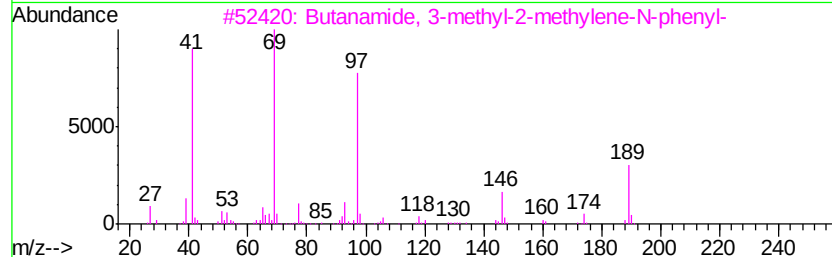
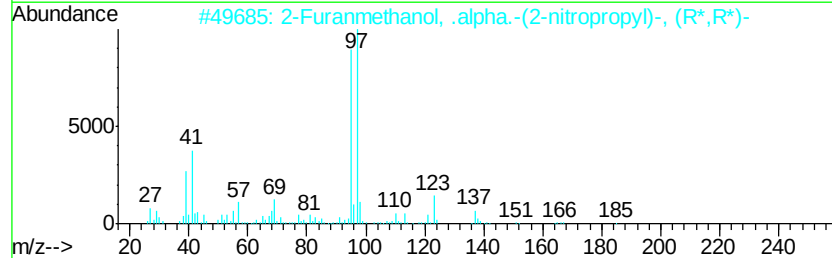
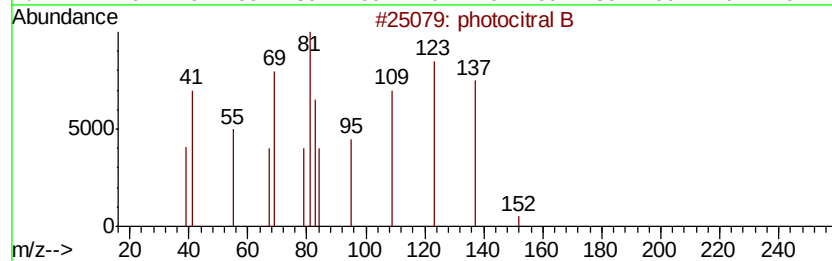
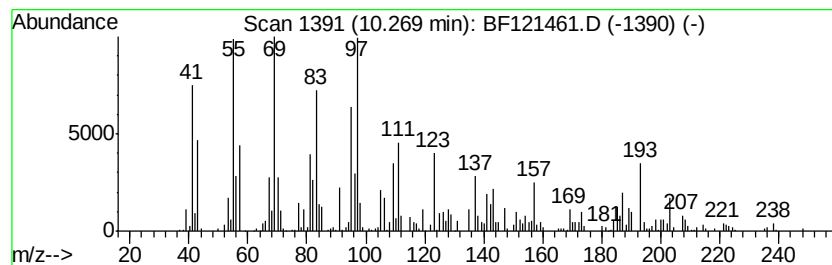
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ClientSampled :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 unknown10.27 Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.27     | 2.89 ng                             | 731667 | Acenaphthene-d10 | 9.90         |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | photocitral B                       | 152    | C10H16O          | 006040-45-5  | 46   |
| 2         | 2-Furanmethanol, .alpha.-(2-nitr... | 185    | C8H11NO4         | 103077-75-4  | 41   |
| 3         | Butanamide, 3-methyl-2-methylene... | 189    | C12H15NO         | 095383-63-4  | 35   |
| 4         | 9-Undecenol, 2,10-dimethyl-         | 198    | C13H26O          | 1000131-86-0 | 30   |
| 5         | Cyclohexane, 1,2,4,5-tetraethyl-... | 196    | C14H28           | 061142-24-3  | 25   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082820\  
Data File : BF121461.D  
Acq On : 28 Aug 2020 15:54  
Operator : JU/CG  
Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

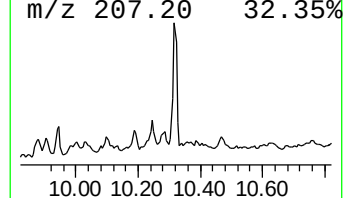
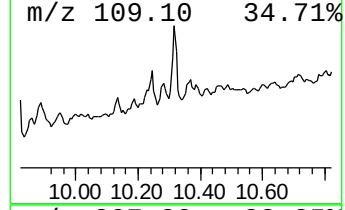
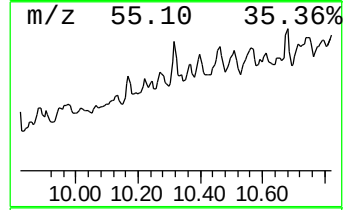
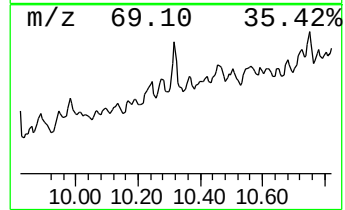
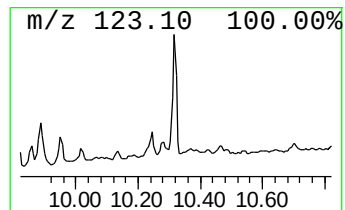
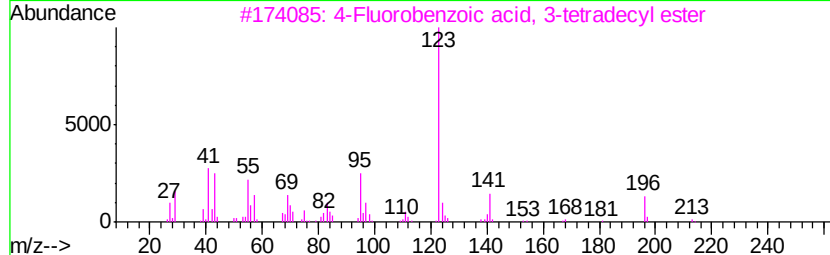
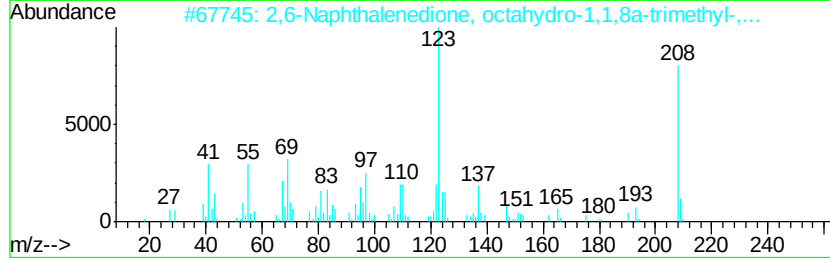
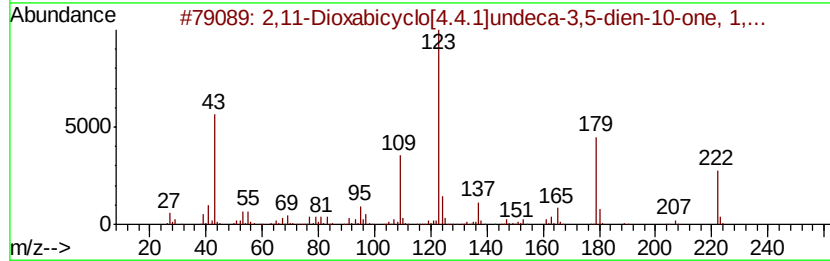
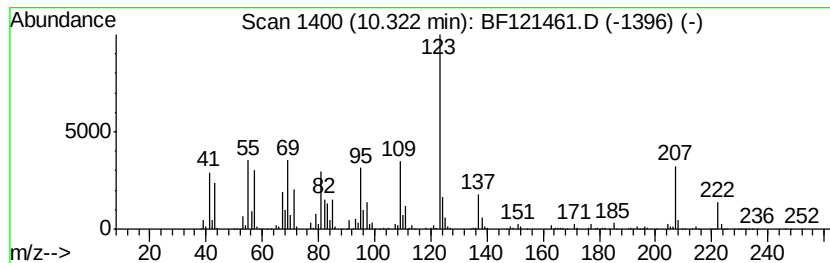
Instrument :  
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ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 1

| R.T.      | EstConc                             | Area    | Relative to ISTD |              | R.T. |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 10.32     | 7.41 ng                             | 1878300 | Acenaphthene-d10 |              | 9.90 |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 2,11-Dioxabicyclo[4.4.1]undeca-3... | 222     | C13H18O3         | 070412-52-1  | 59   |
| 2         | 2,6-Naphthalenedione, octahydro-... | 208     | C13H20O2         | 057289-16-4  | 53   |
| 3         | 4-Fluorobenzoic acid, 3-tetradec... | 336     | C21H33FO2        | 1000280-68-1 | 50   |
| 4         | 4-Fluorobenzoic acid, 5-pentadec... | 350     | C22H35FO2        | 1000280-68-6 | 50   |
| 5         | 3-Fluorobenzoic acid, 4-tridecyl... | 322     | C20H31FO2        | 1000280-60-4 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082820\  
Data File : BF121461.D  
Acq On : 28 Aug 2020 15:54  
Operator : JU/CG  
Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

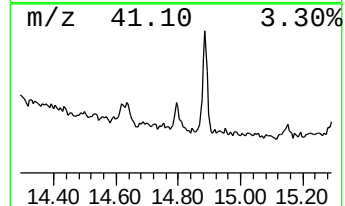
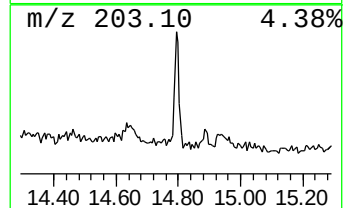
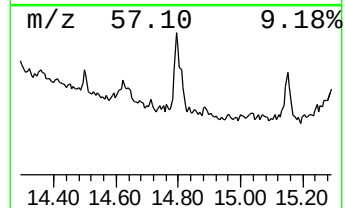
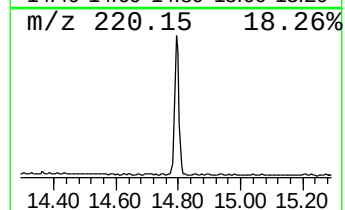
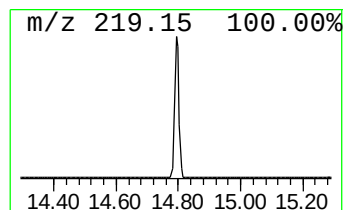
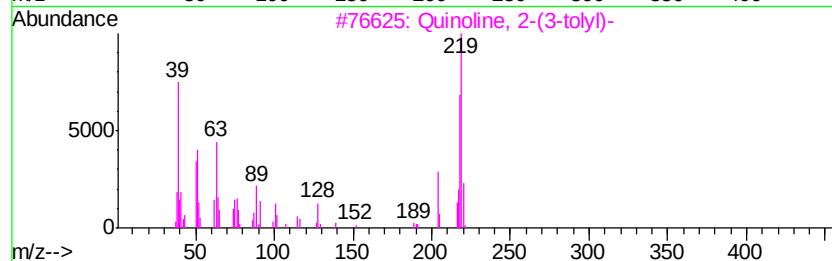
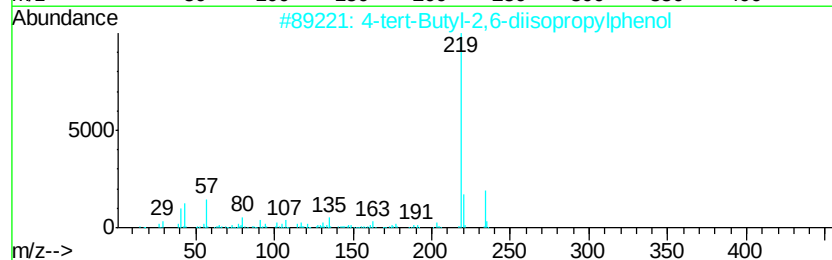
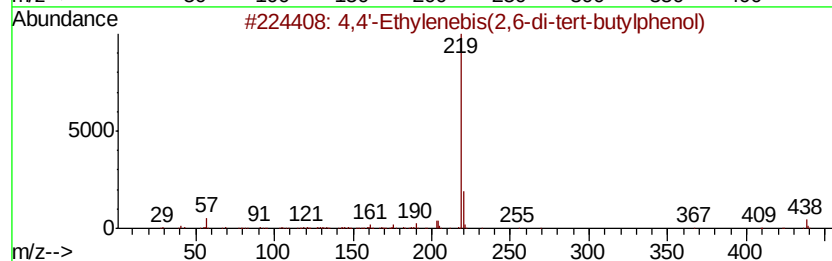
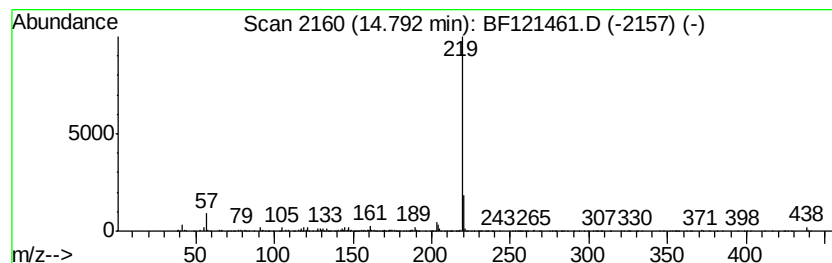
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 4,4'-Ethylenebis(2,6-di-ter... Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.79 | 3.84 ng | 323016 | Perylene-d12     | 15.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 4,4'-Ethylenebis(2,6-di-tert-but... | 438 | C30H46O2   | 001516-94-5  | 90   |
| 2    |    | 4-tert-Butyl-2,6-diisopropylphenol  | 234 | C16H26O    | 057354-65-1  | 72   |
| 3    |    | Quinoline, 2-(3-tolyl)-             | 219 | C16H13N    | 024641-30-3  | 64   |
| 4    |    | Ethyl 5-bromo-2-chlorobenzoate      | 262 | C9H8BrClO2 | 1000343-08-1 | 59   |
| 5    |    | 1-Naphthalenamine, N-phenyl-        | 219 | C16H13N    | 000090-30-2  | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082820\  
Data File : BF121461.D  
Acq On : 28 Aug 2020 15:54  
Operator : JU/CG  
Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

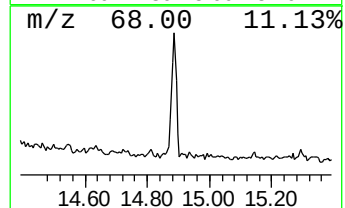
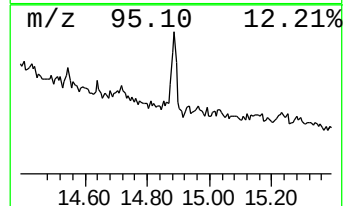
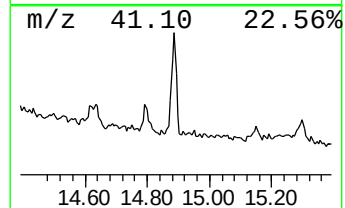
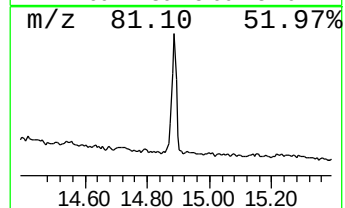
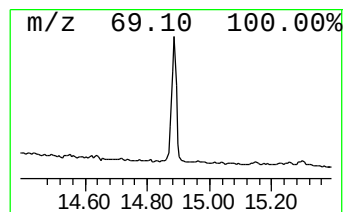
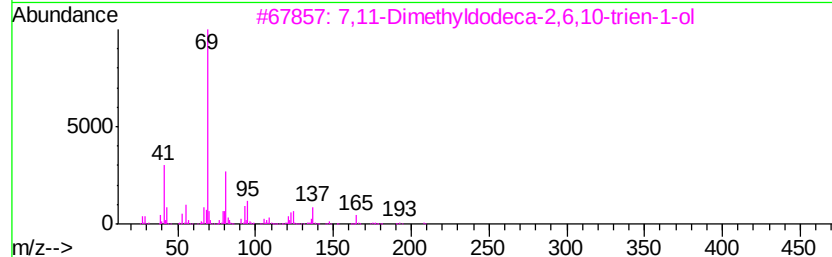
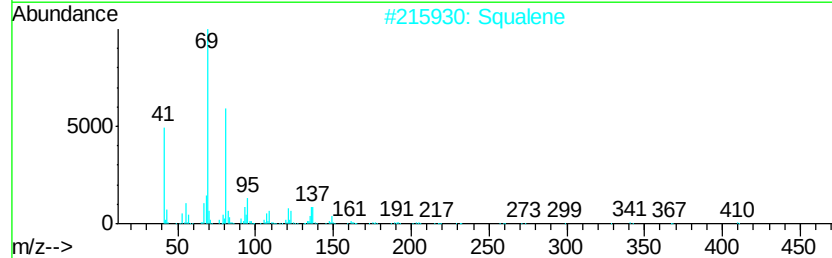
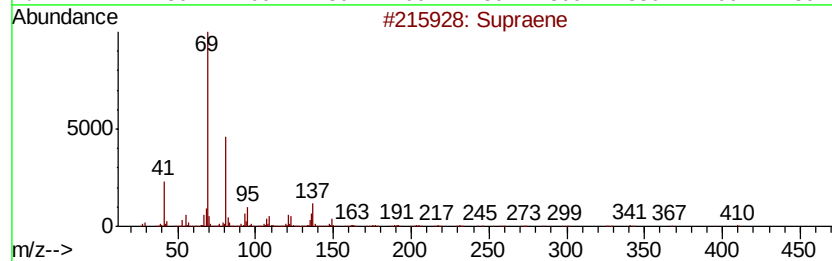
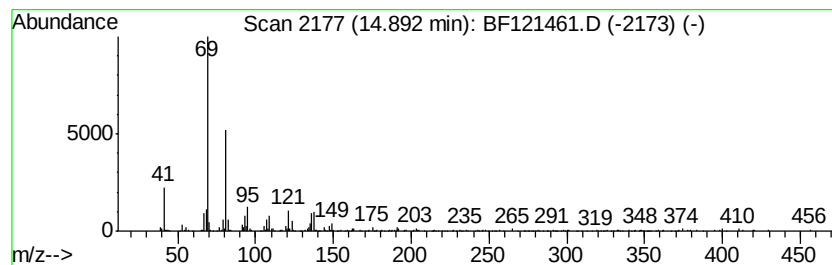
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Supraene Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.89 | 3.41 ng | 286966 | Perylene-d12     | 15.49 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1       | Supraene                            |              | 410 | C30H50  | 007683-64-9 | 87   |
| 2       | Squalene                            |              | 410 | C30H50  | 000111-02-4 | 87   |
| 3       | 7,11-Dimethyldodeca-2,6,10-trien... |              | 208 | C14H24O | 125529-10-4 | 72   |
| 4       | 1,6-Octadiene, 3,5-dimethyl-, tr... |              | 138 | C10H18  | 074630-87-8 | 59   |
| 5       | 1,5-Heptadiene, 3,3,6-trimethyl-    |              | 138 | C10H18  | 035387-63-4 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF082820\  
Data File : BF121461.D  
Acq On : 28 Aug 2020 15:54  
Operator : JU/CG  
Sample : L3824-01 10X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OIL-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF082720.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  | 4.2 ng  |       | 366509   | 1                     | 6.86  | 1732400 | 20.0 |
| unknown9.61          | 9.61  | 2.4 ng  |       | 613300   | 3                     | 9.90  | 5071690 | 20.0 |
| unknown9.77          | 9.77  | 4.9 ng  |       | 1244610  | 3                     | 9.90  | 5071690 | 20.0 |
| unknown9.82          | 9.82  | 5.2 ng  |       | 1315550  | 3                     | 9.90  | 5071690 | 20.0 |
| Sulfurous acid, 2... | 10.16 | 3.8 ng  |       | 950482   | 3                     | 9.90  | 5071690 | 20.0 |
| unknown10.27         | 10.27 | 2.9 ng  |       | 731667   | 3                     | 9.90  | 5071690 | 20.0 |
| 2,11-Dioxabicyclo... | 10.32 | 7.4 ng  |       | 1878300  | 3                     | 9.90  | 5071690 | 20.0 |
| 4,4'-Ethylenebis(... | 14.79 | 3.8 ng  |       | 323016   | 6                     | 15.49 | 1681320 | 20.0 |
| Supraene             | 14.89 | 3.4 ng  |       | 286966   | 6                     | 15.49 | 1681320 | 20.0 |