

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
 Data File : BF106744.D  
 Acq On : 23 Jun 2018 7:46  
 Operator : JU/SJ  
 Sample : J3596-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-01

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-BF061918.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         | 5.201       | 481           | 487         | 490          | rBV      | 503716         | 603258        | 60.45%          | 3.835%        |
| 2         | 5.572       | 546           | 550         | 552          | rBV      | 128061         | 111937        | 11.22%          | 0.712%        |
| 3         | 5.601       | 552           | 555         | 565          | rVB      | 546031         | 483469        | 48.45%          | 3.073%        |
| 4         | 6.119       | 636           | 643         | 649          | rBV      | 99413          | 128653        | 12.89%          | 0.818%        |
| 5         | 6.319       | 673           | 677         | 680          | rBV      | 241801         | 185010        | 18.54%          | 1.176%        |
| 6         | 6.607       | 722           | 726         | 732          | rBV      | 270459         | 297278        | 29.79%          | 1.890%        |
| 7         | 6.737       | 744           | 748         | 752          | rBV      | 902570         | 845840        | 84.76%          | 5.377%        |
| 8         | 6.801       | 757           | 759         | 761          | rVB      | 28896          | 21167         | 2.12%           | 0.135%        |
| 9         | 6.837       | 761           | 765         | 767          | rBV3     | 13500          | 14641         | 1.47%           | 0.093%        |
| 10        | 6.960       | 782           | 786         | 789          | rBV      | 326165         | 266381        | 26.69%          | 1.693%        |
| 11        | 7.048       | 799           | 801         | 804          | rVB      | 22546          | 15933         | 1.60%           | 0.101%        |
| 12        | 7.119       | 806           | 813         | 816          | rBV      | 898689         | 838896        | 84.07%          | 5.332%        |
| 13        | 7.178       | 819           | 823         | 829          | rVV      | 110751         | 115782        | 11.60%          | 0.736%        |
| 14        | 7.243       | 829           | 834         | 838          | rVV7     | 13618          | 24621         | 2.47%           | 0.157%        |
| 15        | 7.307       | 843           | 845         | 850          | rVB3     | 30838          | 30213         | 3.03%           | 0.192%        |
| 16        | 7.395       | 855           | 860         | 862          | rBV      | 65503          | 64227         | 6.44%           | 0.408%        |
| 17        | 7.448       | 866           | 869         | 871          | rBV      | 26114          | 24616         | 2.47%           | 0.156%        |
| 18        | 7.478       | 871           | 874         | 877          | rBV3     | 34676          | 47028         | 4.71%           | 0.299%        |
| 19        | 7.525       | 877           | 882         | 884          | rVV      | 689978         | 707419        | 70.89%          | 4.497%        |
| 20        | 7.548       | 884           | 886         | 888          | rVV      | 223650         | 179702        | 18.01%          | 1.142%        |
| 21        | 7.584       | 888           | 892         | 894          | rVB2     | 56527          | 60068         | 6.02%           | 0.382%        |
| 22        | 7.637       | 899           | 901         | 904          | rVB2     | 44956          | 35401         | 3.55%           | 0.225%        |
| 23        | 7.713       | 911           | 914         | 916          | rVB2     | 29054          | 24778         | 2.48%           | 0.157%        |
| 24        | 7.737       | 916           | 918         | 920          | rVB2     | 42490          | 31092         | 3.12%           | 0.198%        |
| 25        | 7.778       | 920           | 925         | 929          | rBV2     | 79009          | 104640        | 10.49%          | 0.665%        |
| 26        | 7.819       | 929           | 932         | 935          | rBV4     | 43124          | 66349         | 6.65%           | 0.422%        |
| 27        | 7.848       | 935           | 937         | 939          | rVB      | 21246          | 15246         | 1.53%           | 0.097%        |
| 28        | 7.913       | 943           | 948         | 951          | rBV3     | 92515          | 143186        | 14.35%          | 0.910%        |
| 29        | 7.948       | 951           | 954         | 956          | rVB2     | 61773          | 57042         | 5.72%           | 0.363%        |
| 30        | 7.990       | 958           | 961         | 965          | rBV3     | 65632          | 89198         | 8.94%           | 0.567%        |
| 31        | 8.066       | 971           | 974         | 976          | rVB3     | 97250          | 87793         | 8.80%           | 0.558%        |
| 32        | 8.125       | 976           | 984         | 986          | rBV2     | 140506         | 168629        | 16.90%          | 1.072%        |
| 33        | 8.154       | 986           | 989         | 991          | rVB2     | 74743          | 60661         | 6.08%           | 0.386%        |
| 34        | 8.184       | 991           | 994         | 996          | rVB2     | 112102         | 112733        | 11.30%          | 0.717%        |

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 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-01

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-BF061918.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |        |        |        |
|----|--------|------|------|------|------|---------|--------|--------|--------|
| 35 | 8.207  | 996  | 998  | 1001 | rVB3 | 65156   | 52067  | 5.22%  | 0.331% |
| 36 | 8.242  | 1001 | 1004 | 1007 | rBV  | 390997  | 341704 | 34.24% | 2.172% |
| 37 | 8.284  | 1007 | 1011 | 1015 | rVB3 | 126777  | 166449 | 16.68% | 1.058% |
| 38 | 8.319  | 1015 | 1017 | 1019 | rBV3 | 16210   | 16461  | 1.65%  | 0.105% |
| 39 | 8.366  | 1021 | 1025 | 1027 | rBV3 | 117414  | 184437 | 18.48% | 1.172% |
| 40 | 8.437  | 1032 | 1037 | 1039 | rBV3 | 66841   | 92032  | 9.22%  | 0.585% |
| 41 | 8.472  | 1039 | 1043 | 1049 | rVV7 | 119746  | 235662 | 23.62% | 1.498% |
| 42 | 8.519  | 1049 | 1051 | 1052 | rBV2 | 37365   | 22566  | 2.26%  | 0.143% |
| 43 | 8.566  | 1056 | 1059 | 1062 | rVB4 | 75715   | 92912  | 9.31%  | 0.591% |
| 44 | 8.625  | 1067 | 1069 | 1072 | rBV2 | 127143  | 99910  | 10.01% | 0.635% |
| 45 | 8.689  | 1072 | 1080 | 1083 | rBV4 | 212437  | 377727 | 37.85% | 2.401% |
| 46 | 8.725  | 1083 | 1086 | 1090 | rVB5 | 96639   | 141376 | 14.17% | 0.899% |
| 47 | 8.772  | 1091 | 1094 | 1098 | rVB3 | 134096  | 227592 | 22.81% | 1.447% |
| 48 | 8.825  | 1098 | 1103 | 1108 | rBV7 | 102069  | 161680 | 16.20% | 1.028% |
| 49 | 8.866  | 1108 | 1110 | 1112 | rBV2 | 90515   | 76614  | 7.68%  | 0.487% |
| 50 | 8.925  | 1115 | 1120 | 1122 | rBV3 | 115763  | 205088 | 20.55% | 1.304% |
| 51 | 9.013  | 1131 | 1135 | 1137 | rBV4 | 161668  | 209192 | 20.96% | 1.330% |
| 52 | 9.048  | 1139 | 1141 | 1146 | rVB4 | 237504  | 311004 | 31.17% | 1.977% |
| 53 | 9.137  | 1154 | 1156 | 1159 | rBV2 | 126909  | 109582 | 10.98% | 0.697% |
| 54 | 9.207  | 1166 | 1168 | 1170 | rBV3 | 51741   | 43034  | 4.31%  | 0.274% |
| 55 | 9.260  | 1170 | 1177 | 1180 | rBV5 | 173492  | 360275 | 36.10% | 2.290% |
| 56 | 9.289  | 1180 | 1182 | 1184 | rVV  | 90488   | 74082  | 7.42%  | 0.471% |
| 57 | 9.319  | 1184 | 1187 | 1190 | rBV  | 1065457 | 970669 | 97.27% | 6.170% |
| 58 | 9.354  | 1190 | 1193 | 1194 | rBV2 | 126840  | 98590  | 9.88%  | 0.627% |
| 59 | 9.454  | 1208 | 1210 | 1212 | rVB  | 110074  | 77071  | 7.72%  | 0.490% |
| 60 | 9.495  | 1215 | 1217 | 1219 | rVB3 | 192873  | 158361 | 15.87% | 1.007% |
| 61 | 9.525  | 1219 | 1222 | 1227 | rBV4 | 239810  | 307408 | 30.81% | 1.954% |
| 62 | 9.736  | 1255 | 1258 | 1262 | rVB2 | 204256  | 237491 | 23.80% | 1.510% |
| 63 | 9.819  | 1268 | 1272 | 1275 | rBV4 | 131407  | 237354 | 23.79% | 1.509% |
| 64 | 9.978  | 1296 | 1299 | 1301 | rBV2 | 170083  | 208520 | 20.90% | 1.325% |
| 65 | 10.007 | 1301 | 1304 | 1307 | rVB  | 311570  | 287838 | 28.84% | 1.830% |
| 66 | 10.736 | 1425 | 1428 | 1431 | rBV3 | 116905  | 148434 | 14.87% | 0.944% |
| 67 | 10.807 | 1435 | 1440 | 1443 | rBV  | 578254  | 618951 | 62.03% | 3.934% |
| 68 | 10.995 | 1470 | 1472 | 1478 | rBV4 | 100732  | 129712 | 13.00% | 0.825% |
| 69 | 11.154 | 1497 | 1499 | 1507 | rVB6 | 92507   | 145275 | 14.56% | 0.923% |
| 70 | 11.501 | 1555 | 1558 | 1561 | rVB  | 320389  | 261274 | 26.18% | 1.661% |
| 71 | 12.025 | 1645 | 1647 | 1655 | rVB2 | 79867   | 93107  | 9.33%  | 0.592% |

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Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

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-BF061918.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |        |         |        |
|----|--------|------|------|------|------|---------|--------|---------|--------|
| 72 | 13.083 | 1823 | 1827 | 1830 | rBV  | 1111425 | 997897 | 100.00% | 6.343% |
| 73 | 13.960 | 1974 | 1976 | 1981 | rVB  | 59268   | 55715  | 5.58%   | 0.354% |
| 74 | 14.148 | 2005 | 2008 | 2012 | rBV  | 318184  | 314170 | 31.48%  | 1.997% |
| 75 | 14.283 | 2029 | 2031 | 2034 | rVB  | 40145   | 32135  | 3.22%   | 0.204% |
| 76 | 14.601 | 2082 | 2085 | 2090 | rVB  | 64721   | 65765  | 6.59%   | 0.418% |
| 77 | 14.918 | 2136 | 2139 | 2146 | rVB  | 89548   | 98313  | 9.85%   | 0.625% |
| 78 | 15.277 | 2197 | 2200 | 2205 | rVB  | 96489   | 110092 | 11.03%  | 0.700% |
| 79 | 15.689 | 2265 | 2270 | 2277 | rVB2 | 194165  | 335291 | 33.60%  | 2.131% |
| 80 | 16.142 | 2343 | 2347 | 2352 | rVB  | 49113   | 78378  | 7.85%   | 0.498% |

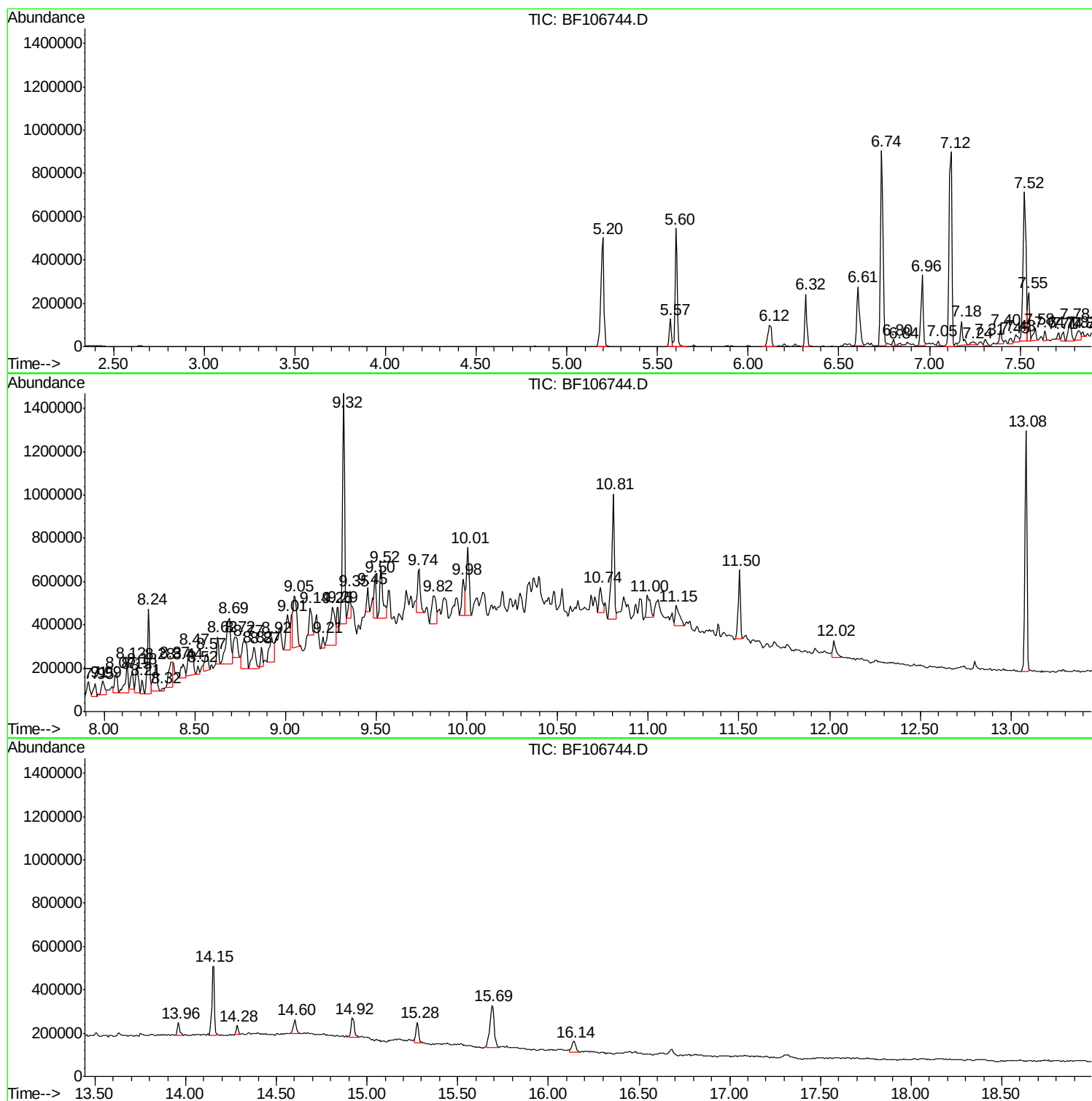
Sum of corrected areas: 15732144

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Data File : BF106744.D  
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Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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 :  
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ClientSampleId :  
MW-01

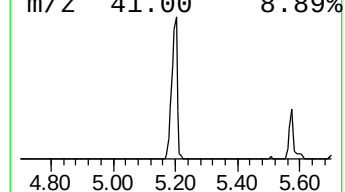
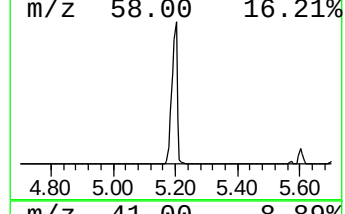
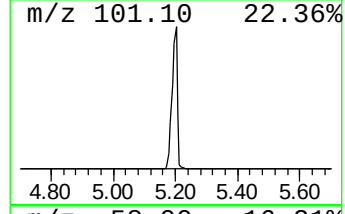
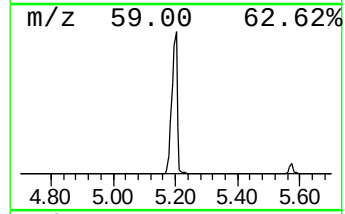
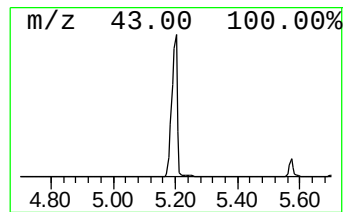
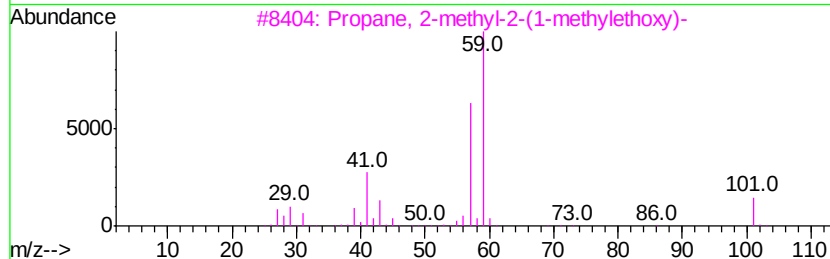
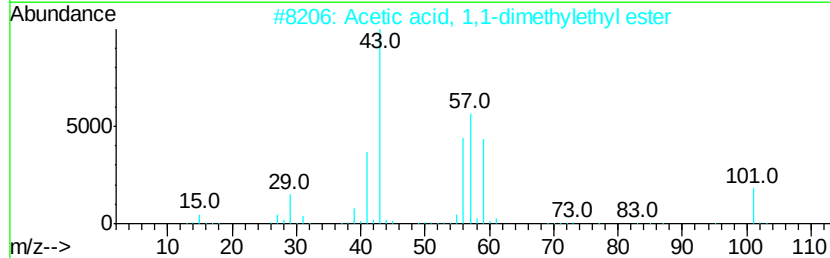
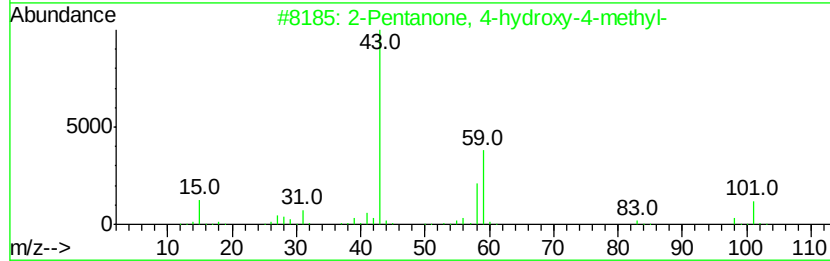
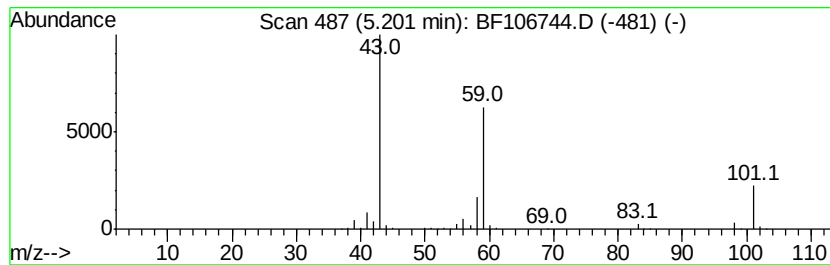
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.20 | 45.29 ng | 603258 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 72   |
| 2    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    |   | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 28   |
| 4    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 5    |    |   | Butane, 1-ethoxy-                   | 102 | C6H14O   | 000628-81-9 | 9    |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

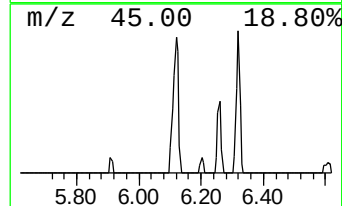
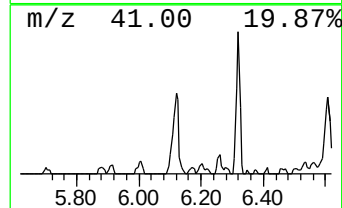
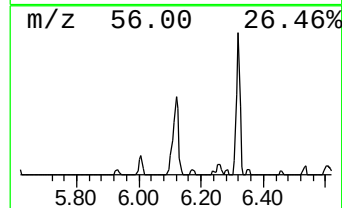
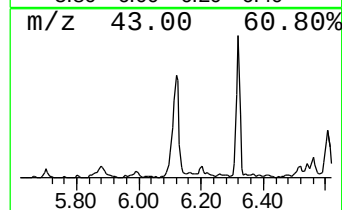
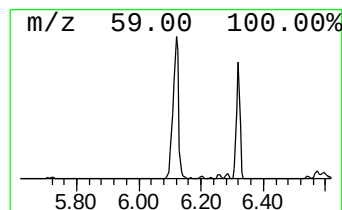
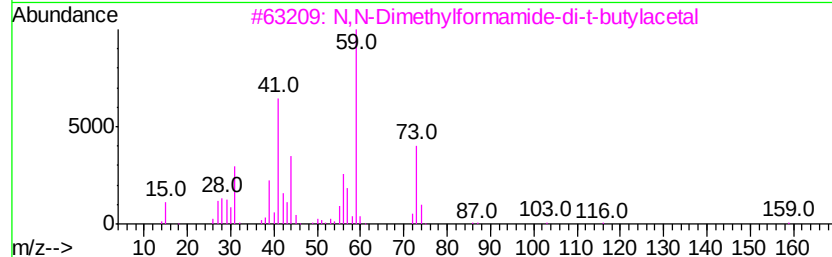
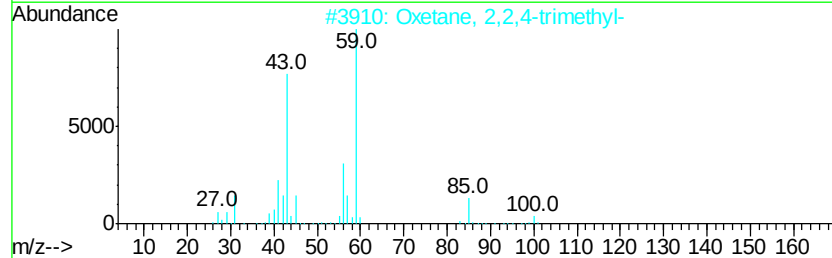
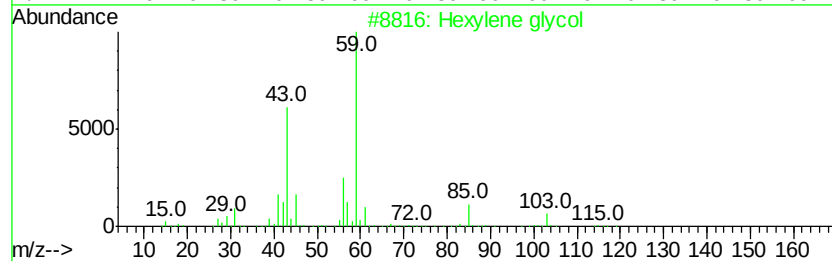
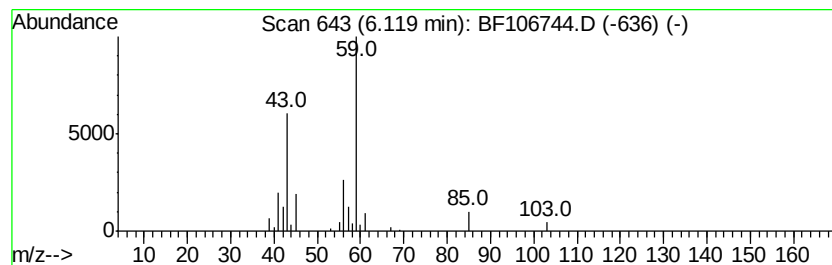
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Hexylene glycol Concentration Rank 20

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.12 | 9.66 ng | 128653 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Hexylene glycol                     | 118 | C6H14O2   | 000107-41-5  | 83   |
| 2    |    | Oxetane, 2,2,4-trimethyl-           | 100 | C6H12O    | 023120-44-7  | 78   |
| 3    |    | N,N-Dimethylformamide-di-t-butyl... | 203 | C11H25NO2 | 036805-97-7  | 50   |
| 4    |    | dl-Erythro-0-methylthreonine        | 133 | C5H11NO3  | 1000214-70-7 | 45   |
| 5    |    | 2-Propanol, 1-methoxy-2-methyl-     | 104 | C5H12O2   | 003587-64-2  | 40   |



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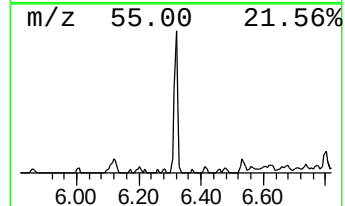
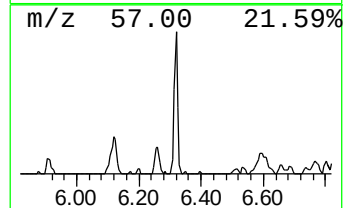
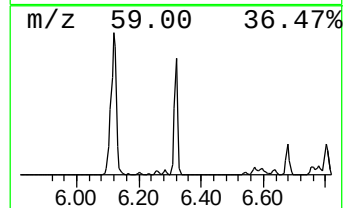
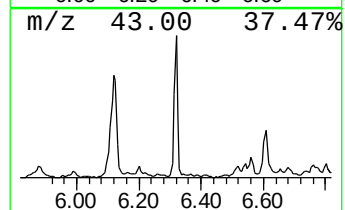
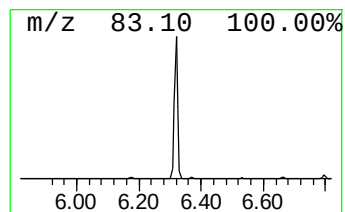
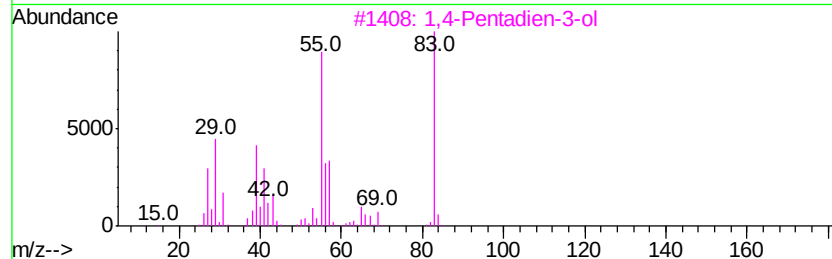
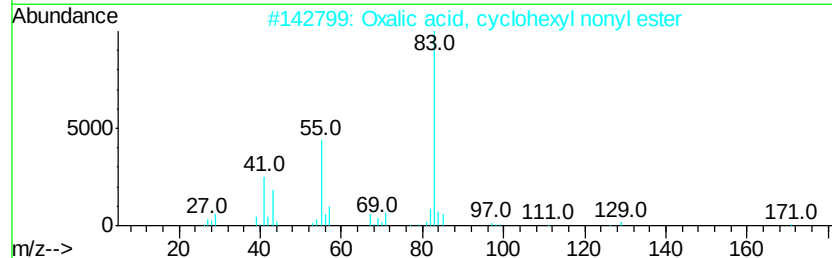
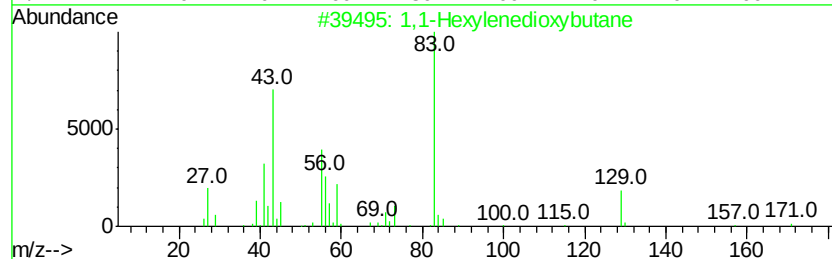
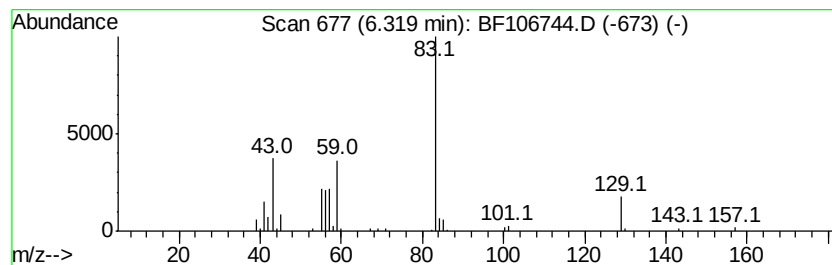
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 1,1-Hexylenedioxybutane Concentration Rank 9

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.32 | 13.89 ng | 185010 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1,1-Hexylenedioxybutane             | 172 | C10H20O2 | 1000249-77-4 | 72   |
| 2    |    | Oxalic acid, cyclohexyl nonyl ester | 298 | C17H30O4 | 1000309-31-1 | 40   |
| 3    |    | 1,4-Pentadien-3-ol                  | 84  | C5H8O    | 000922-65-6  | 9    |
| 4    |    | .alpha.-Amino-2,5-dihydro-5-meth... | 157 | C7H11NO3 | 018455-25-9  | 9    |
| 5    |    | 1H-Imidazol-2-amine                 | 83  | C3H5N3   | 007720-39-0  | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

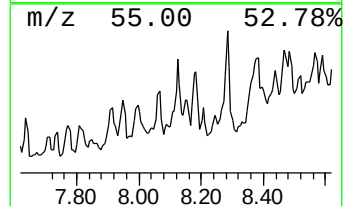
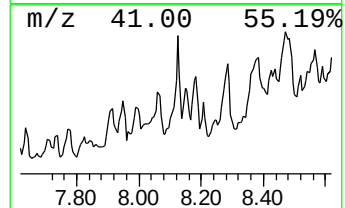
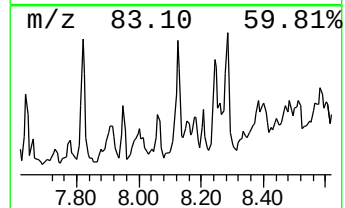
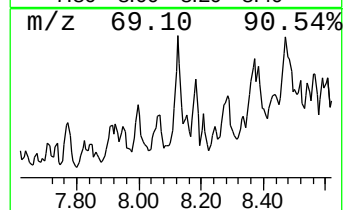
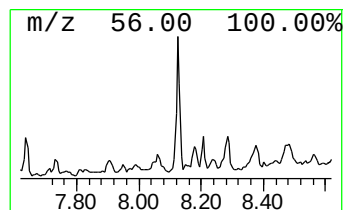
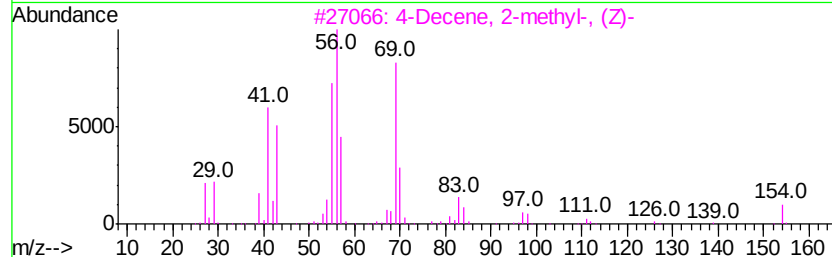
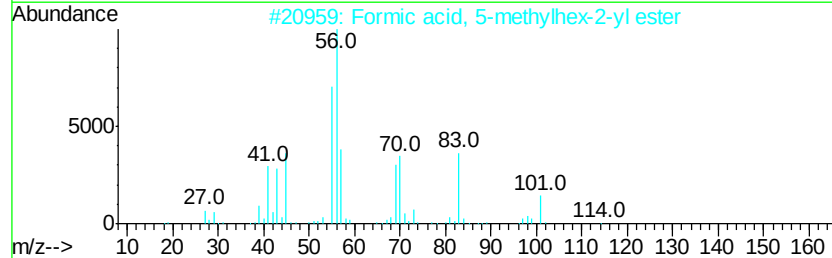
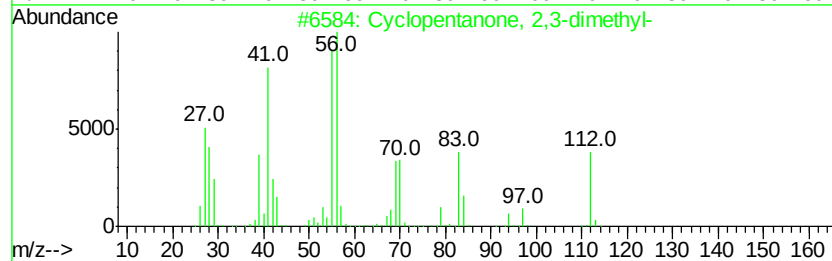
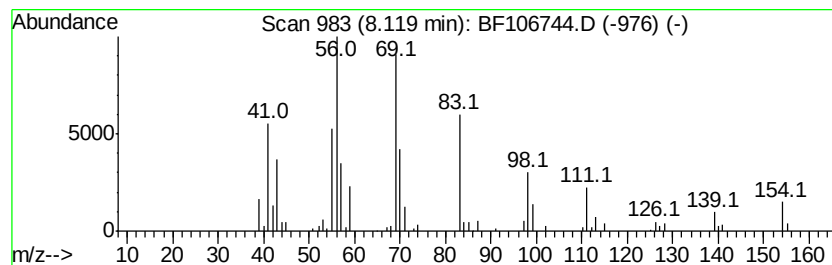
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown8.12 Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 8.12 | 9.87 ng | 168629 | Naphthalene-d8   | 8.24 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|-----------|-------------------------------------|-----|---------|--------------|------|
| 1         | Cyclopentanone, 2,3-dimethyl-       | 112 | C7H12O  | 1000151-06-7 | 42   |
| 2         | Formic acid, 5-methylhex-2-yl ester | 144 | C8H16O2 | 1000368-88-7 | 38   |
| 3         | 4-Decene, 2-methyl-, (Z)-           | 154 | C11H22  | 055499-07-5  | 38   |
| 4         | 2-Heptene                           | 98  | C7H14   | 000592-77-8  | 38   |
| 5         | Cyclopentane, 1,1-dimethyl-         | 98  | C7H14   | 001638-26-2  | 38   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

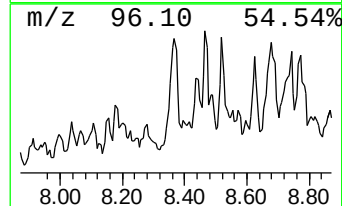
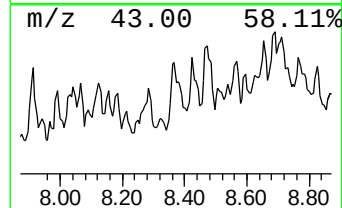
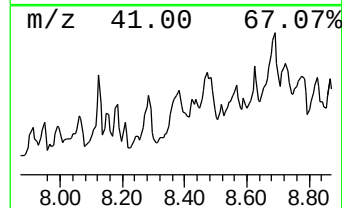
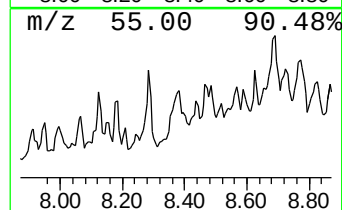
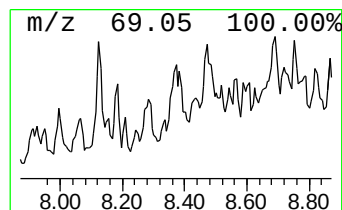
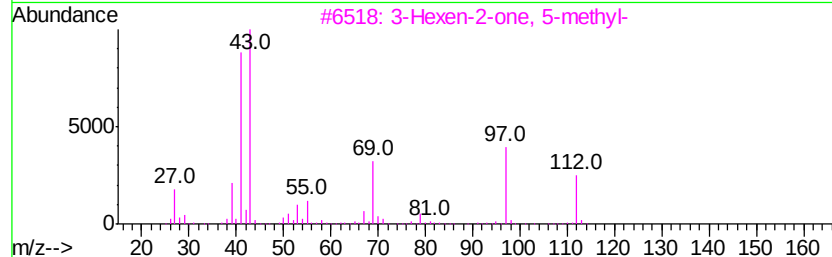
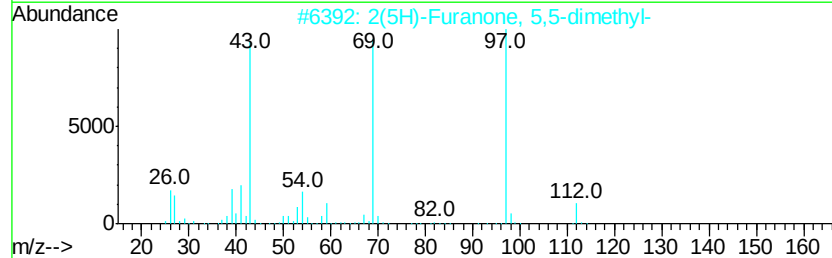
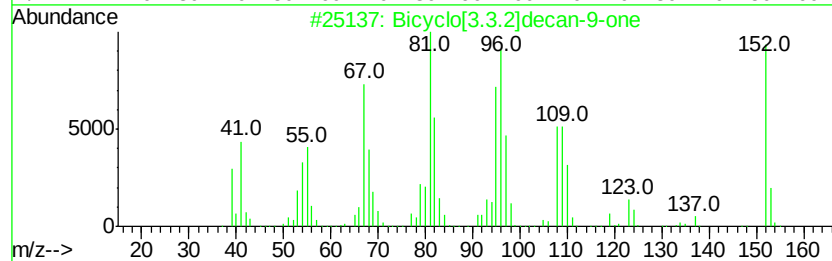
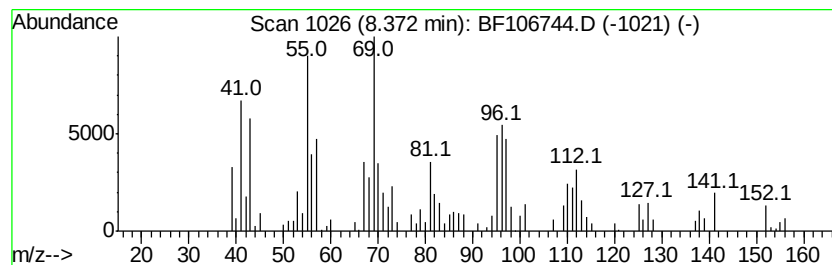
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ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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 Bicyclo[3.3.2]decan-9-one Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 8.37      | 10.80 ng                            | 184437 | Napthalene-d8    | 8.24        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Bicyclo[3.3.2]decan-9-one           | 152    | C10H16O          | 028054-91-3 | 51   |
| 2         | 2(5H)-Furanone, 5,5-dimethyl-       | 112    | C6H8O2           | 020019-64-1 | 38   |
| 3         | 3-Hexen-2-one, 5-methyl-            | 112    | C7H12O           | 005166-53-0 | 38   |
| 4         | Cyclohexanone, 4-methyl-, 0-meth... | 141    | C8H15NO          | 039477-43-5 | 30   |
| 5         | Cyclopentane, (2-methylpropyl)-     | 126    | C9H18            | 003788-32-7 | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

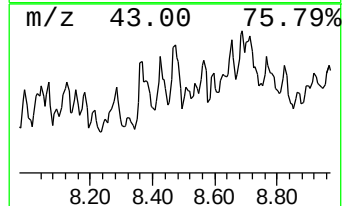
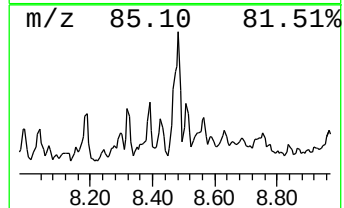
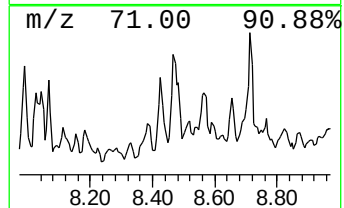
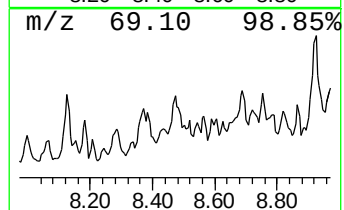
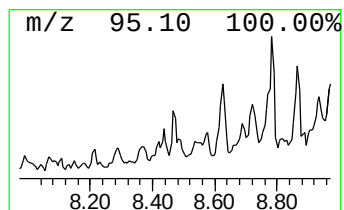
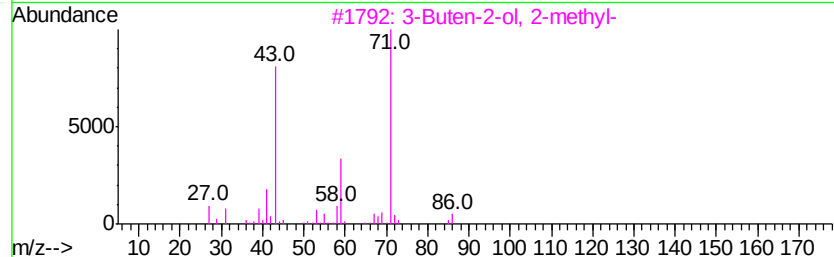
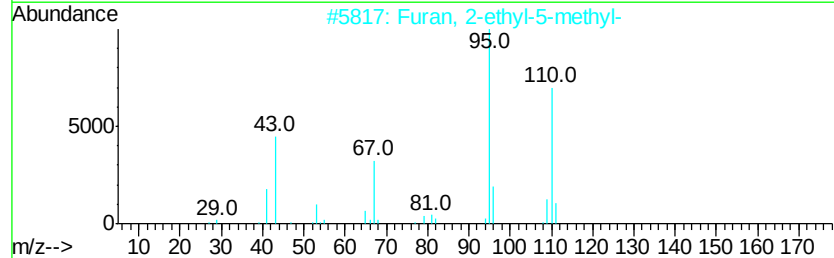
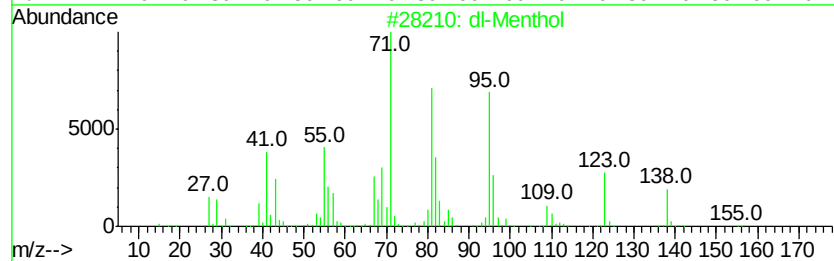
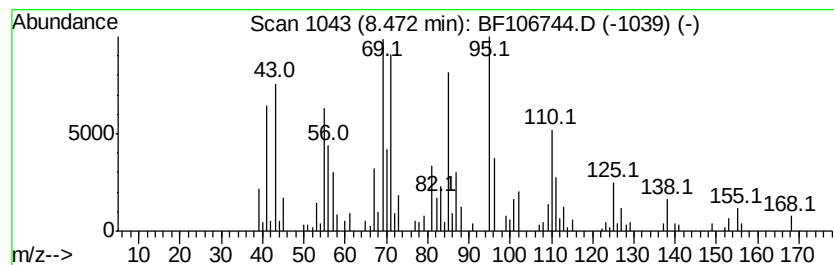
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown8.47 Concentration Rank 10

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.47 | 13.79 ng | 235662 | Naphthalene-d8   | 8.24 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | dl-Menthol                 | 156 | C10H20O | 000089-78-1 | 27   |
| 2    |    | Furan, 2-ethyl-5-methyl-   | 110 | C7H10O  | 001703-52-2 | 18   |
| 3    |    | 3-Buten-2-ol, 2-methyl-    | 86  | C5H10O  | 000115-18-4 | 11   |
| 4    |    | 4-Hexen-3-ol               | 100 | C6H12O  | 004798-58-7 | 11   |
| 5    |    | 1,3-Dimethyl-1-cyclohexene | 110 | C8H14   | 002808-76-6 | 11   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

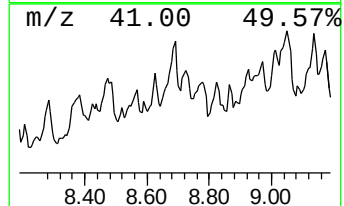
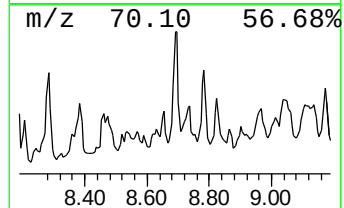
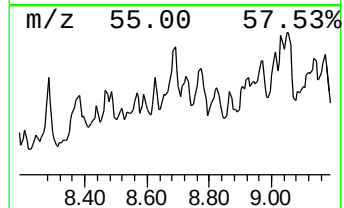
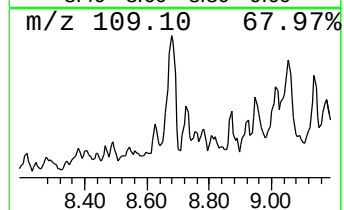
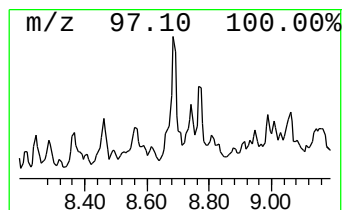
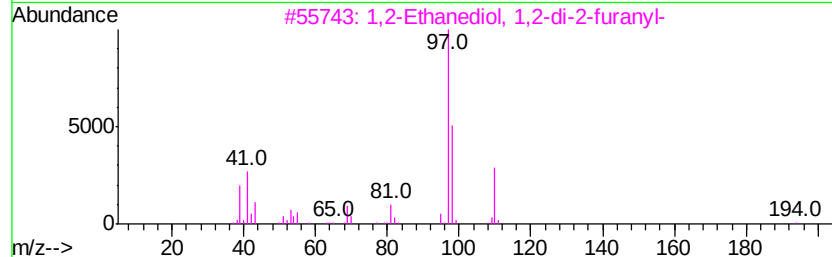
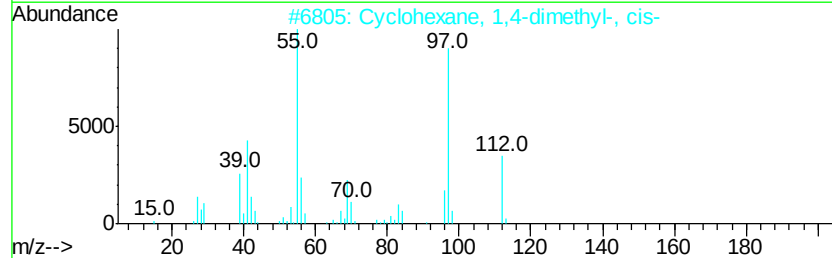
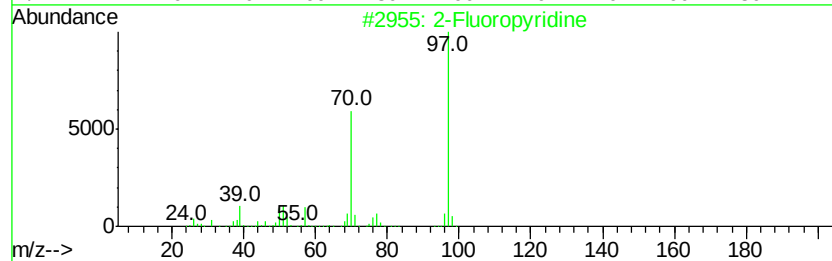
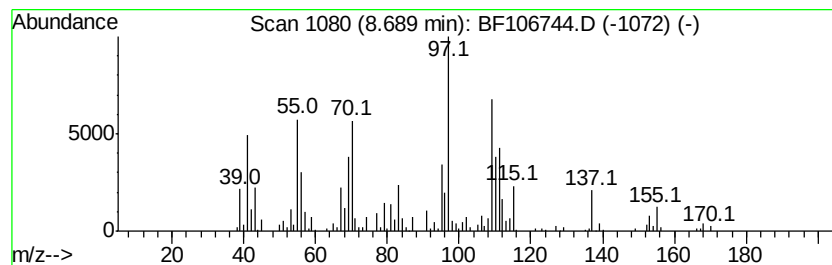
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown8.69 Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.69 | 22.11 ng | 377727 | Naphthalene-d8   | 8.24 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Fluoropyridine                   | 97  | C5H4FN   | 000372-48-5 | 30   |
| 2    |    | Cyclohexane, 1,4-dimethyl-, cis-   | 112 | C8H16    | 000624-29-3 | 22   |
| 3    |    | 1,2-Ethanediol, 1,2-di-2-furanyl-  | 194 | C10H10O4 | 004464-77-1 | 22   |
| 4    |    | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16    | 002207-03-6 | 18   |
| 5    |    | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16    | 002207-04-7 | 18   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

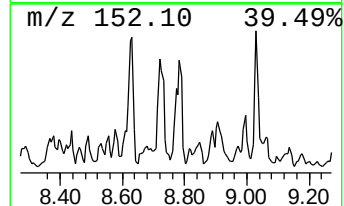
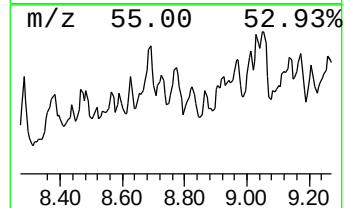
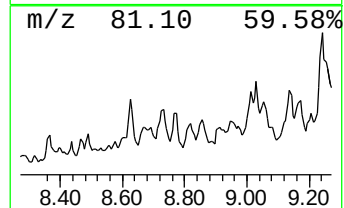
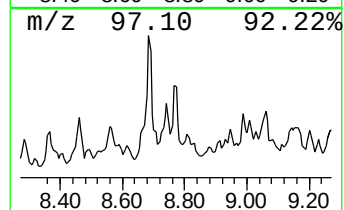
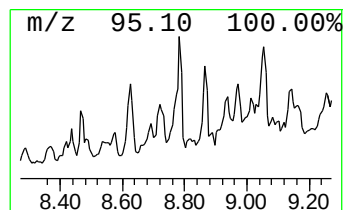
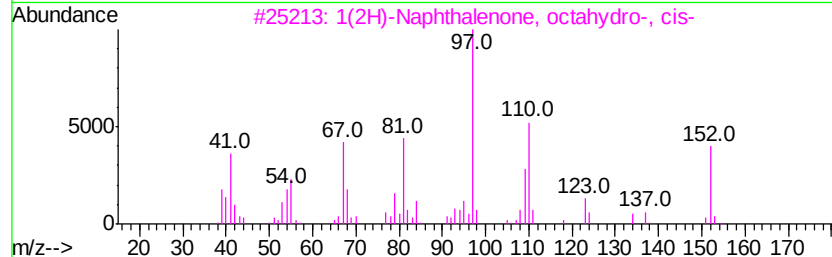
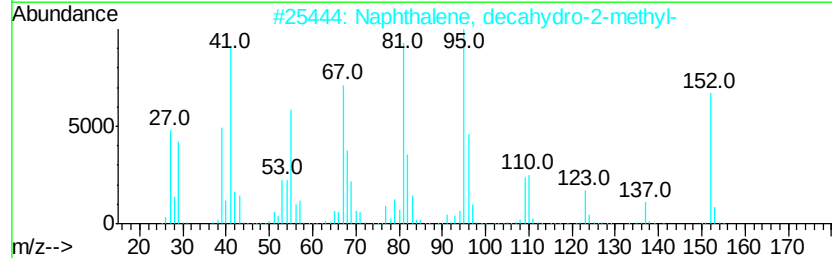
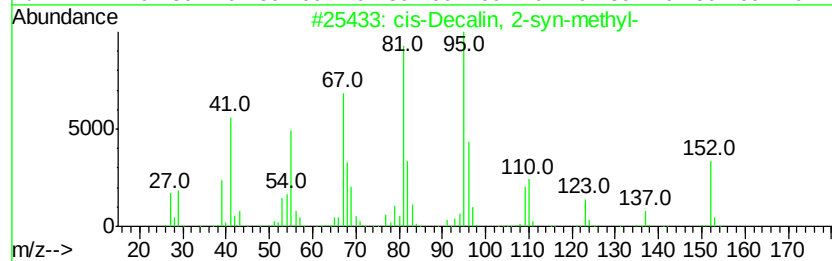
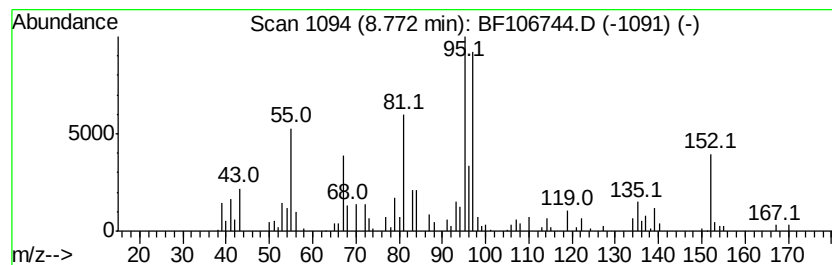
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown8.77 Concentration Rank 11

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.77 | 13.32 ng | 227592 | Naphthalene-d8   | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 49   |
| 2    |    | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 46   |
| 3    |    | 1(2H)-Naphthalenone, octahydro-,... | 152 | C10H16O | 1000141-71-1 | 46   |
| 4    |    | 2-Oxatricyclo[3.3.1.1(3,7)]decan... | 152 | C10H16O | 006508-22-1  | 41   |
| 5    |    | Camphor                             | 152 | C10H16O | 000076-22-2  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

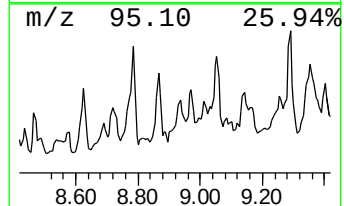
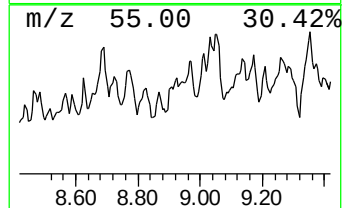
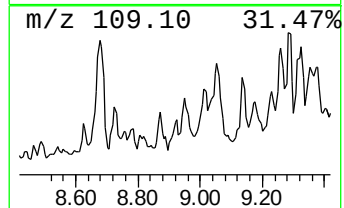
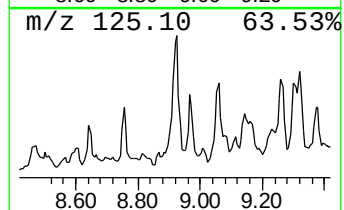
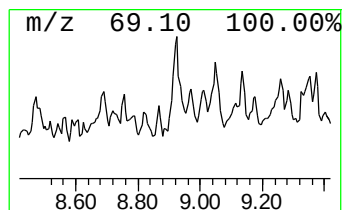
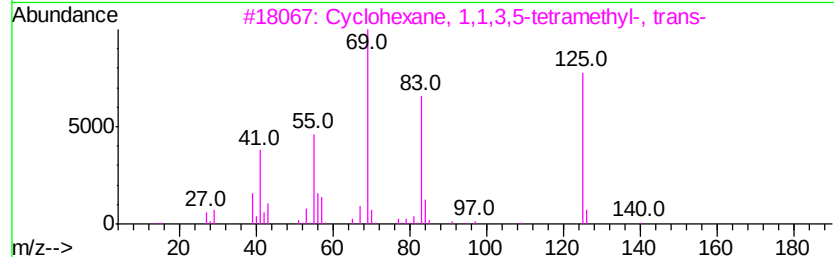
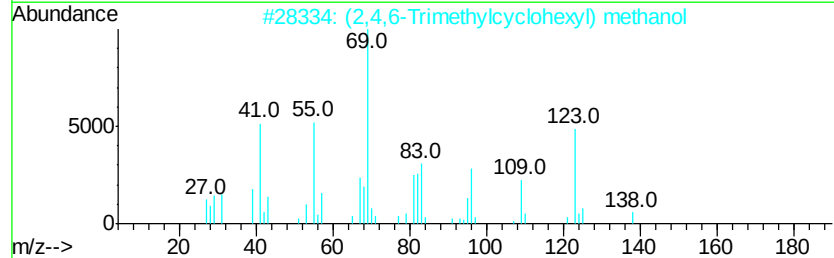
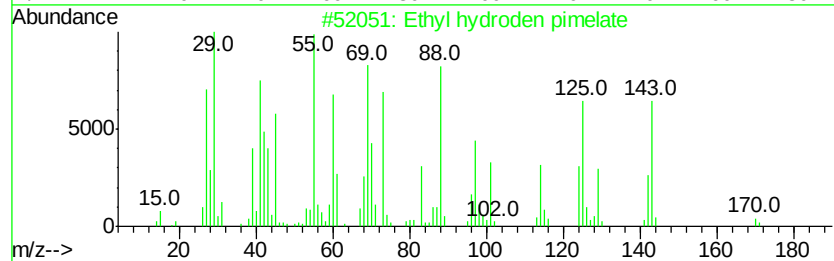
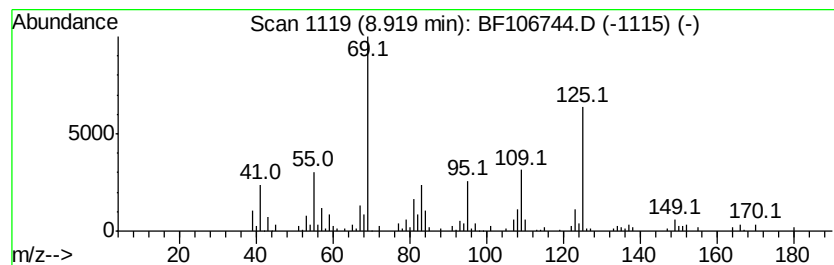
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown8.92 Concentration Rank 13

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.92 | 12.00 ng | 205088 | Napthalene-d8    | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Ethyl hydroden pimelate             | 188 | C9H16O4 | 1000342-35-7 | 38   |
| 2    |    | (2,4,6-Trimethylcyclohexyl) meth... | 156 | C10H20O | 013702-56-2  | 37   |
| 3    |    | Cyclohexane, 1,1,3,5-tetramethyl... | 140 | C10H20  | 050876-31-8  | 37   |
| 4    |    | 1-Isopropyl-1,4,5-trimethylcyclo... | 168 | C12H24  | 219783-06-9  | 37   |
| 5    |    | Methanone, dicyclopropyl-           | 110 | C7H10O  | 001121-37-5  | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

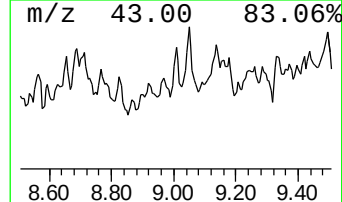
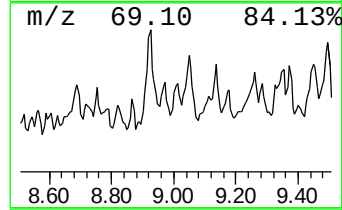
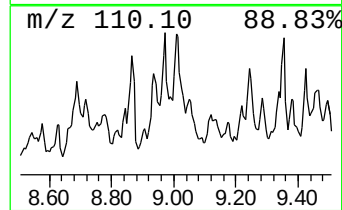
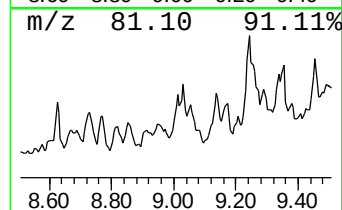
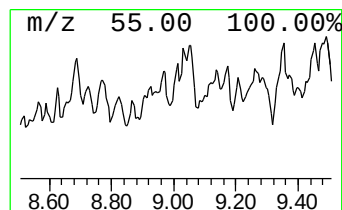
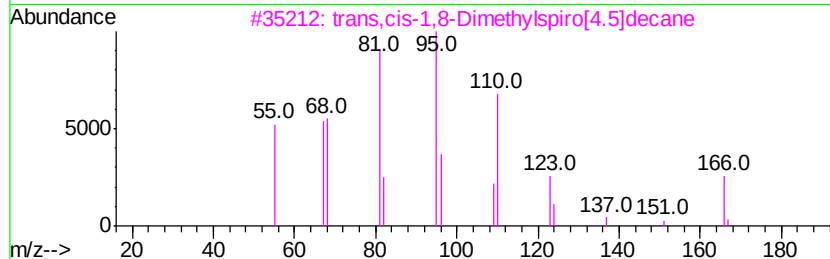
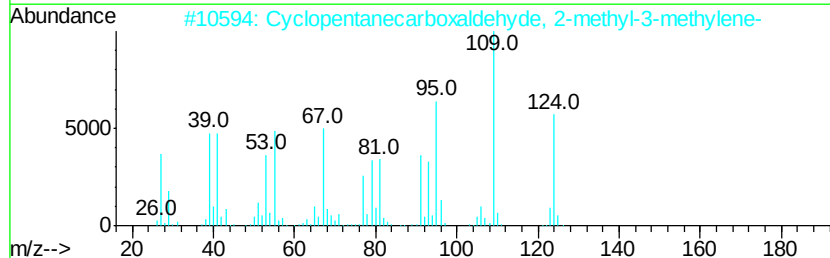
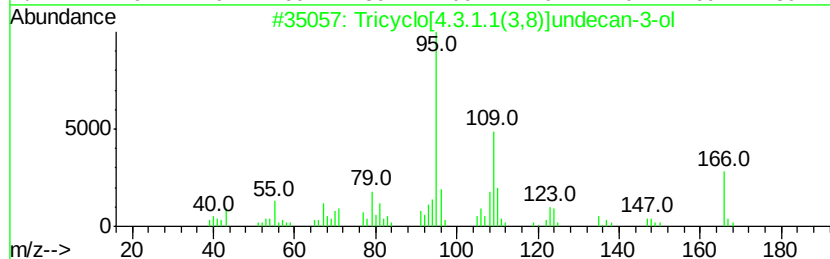
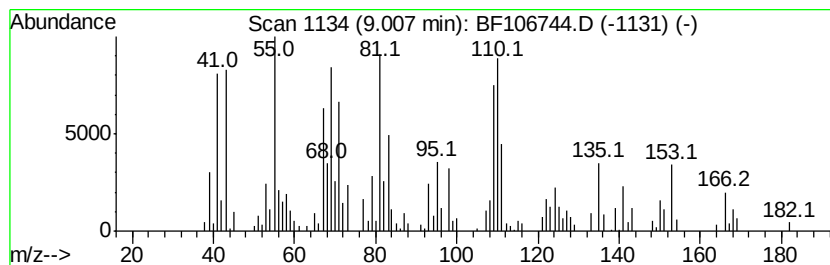
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Tricyclo[4.3.1.1(3,8)]undec... Concentration Rank 12

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.01 | 12.24 ng | 209192 | Napthalene-d8    | 8.24 |

| Hit# | of | Tentative ID                                      | MW  | MolForm  | CAS#         | Qual |
|------|----|---------------------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Tricyclo[4.3.1.1(3,8)]undecan-3-ol                | 166 | C11H18O  | 014504-80-4  | 64   |
| 2    |    | Cyclopentanecarboxaldehyde, 2-methyl-3-methylene- | 124 | C8H12O   | 1000154-24-0 | 30   |
| 3    |    | trans,cis-1,8-Dimethylspiro[4.5]decane            | 166 | C12H22   | 1000111-72-9 | 30   |
| 4    |    | Cyclohexene, 3,5,5-trimethyl-                     | 124 | C9H16    | 000933-12-0  | 25   |
| 5    |    | Iridomyrmecin                                     | 168 | C10H16O2 | 000485-43-8  | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

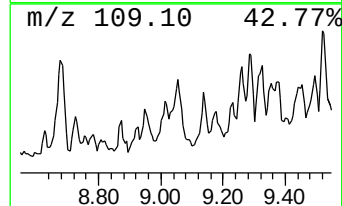
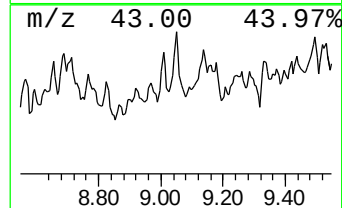
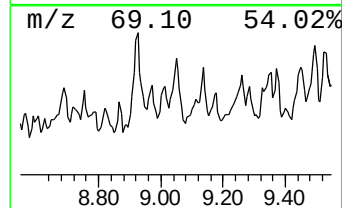
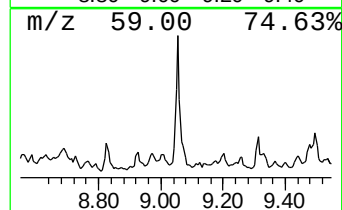
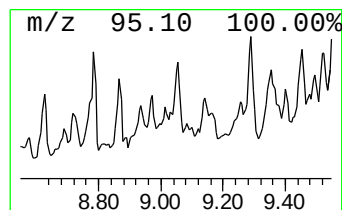
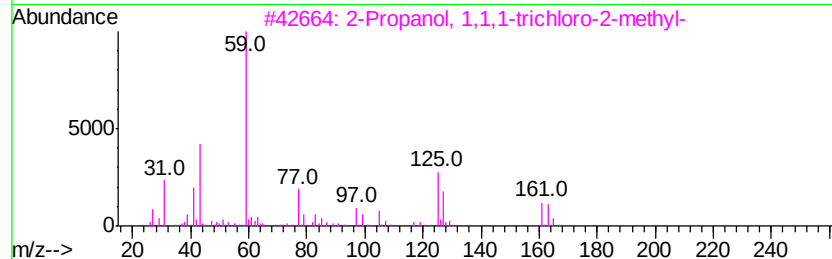
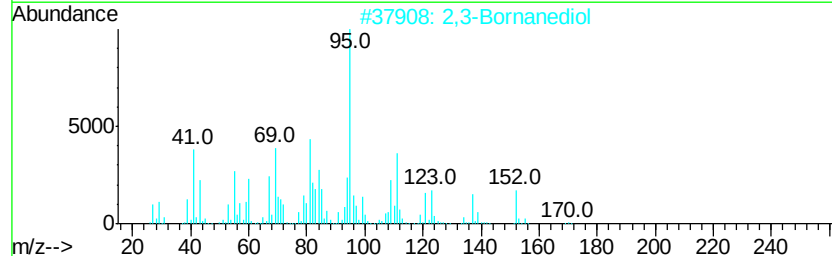
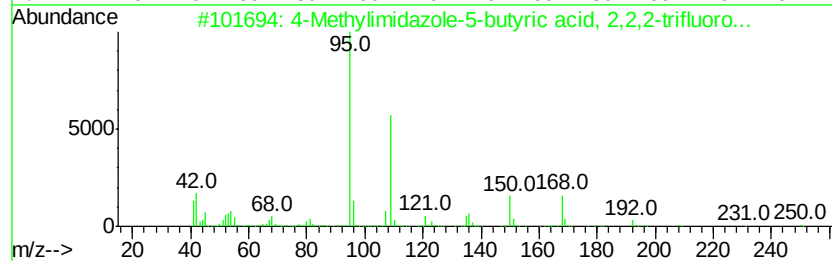
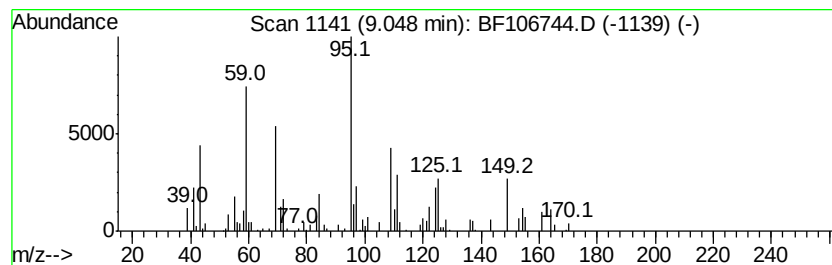
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown9.05 Concentration Rank 7

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.05 | 18.20 ng | 311004 | Naphthalene-d8   | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 4-Methylimidazole-5-butyric acid... | 250 | C10H13F3N2O2 | 1000129-15-9 | 12   |
| 2    |    | 2,3-Bornanediol                     | 170 | C10H18O2     | 025696-56-4  | 12   |
| 3    |    | 2-Propanol, 1,1,1-trichloro-2-me... | 176 | C4H7Cl3O     | 000057-15-8  | 10   |
| 4    |    | Ethyl dodecyl ether                 | 214 | C14H30O      | 007289-37-4  | 10   |
| 5    |    | 3-Heptanol, 4-methyl-               | 130 | C8H18O       | 014979-39-6  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

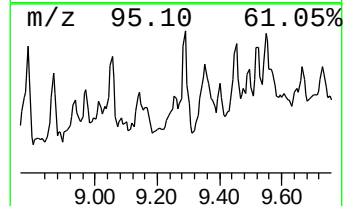
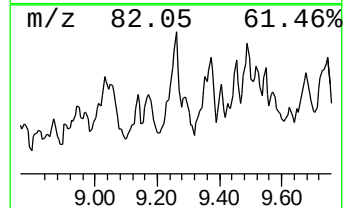
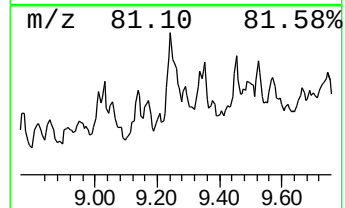
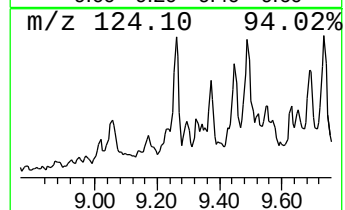
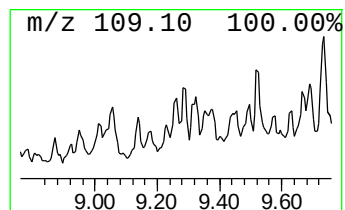
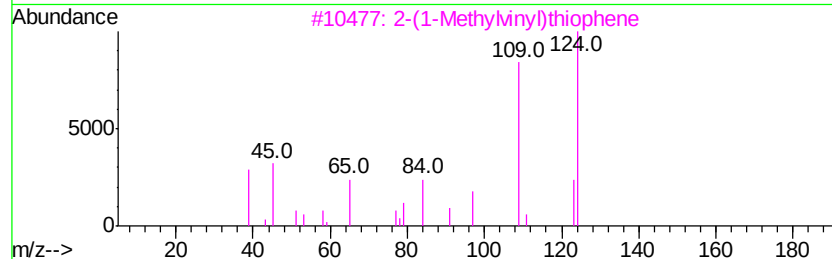
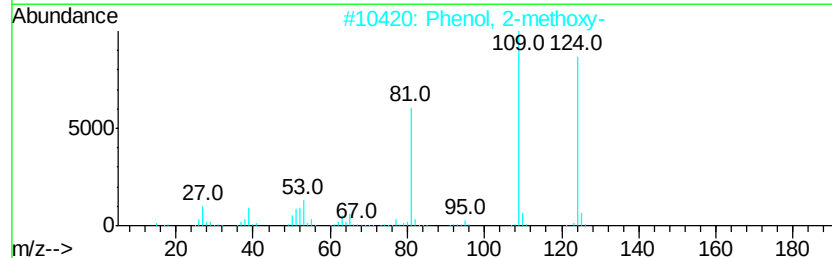
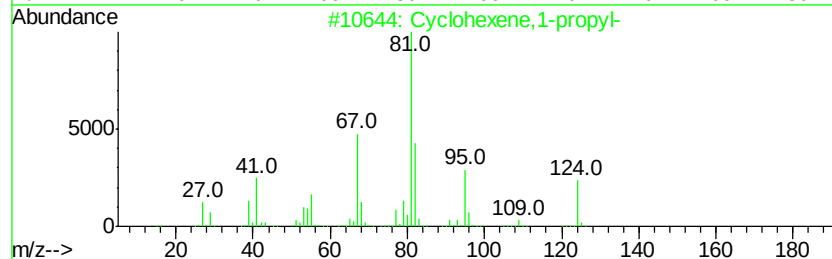
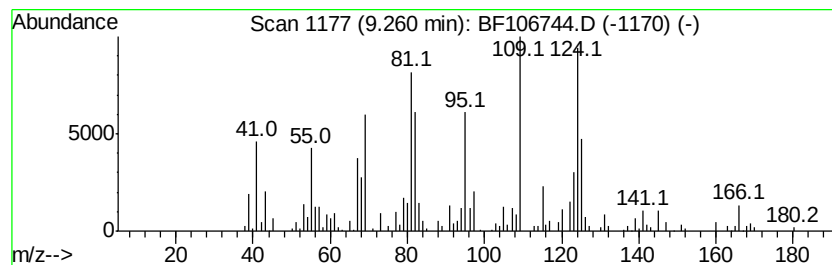
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown9.26 Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD | R.T.  |
|------|----------|--------|------------------|-------|
| 9.26 | 25.03 ng | 360275 | Acenaphthene-d10 | 10.01 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#         | Qual |
|------|----|-----------------------------------|-----|---------|--------------|------|
| 1    | 5  | Cyclohexene,1-propyl-             | 124 | C9H16   | 002539-75-5  | 46   |
| 2    |    | Phenol, 2-methoxy-                | 124 | C7H8O2  | 000090-05-1  | 43   |
| 3    |    | 2-(1-Methylvinyl)thiophene        | 124 | C7H8S   | 030616-73-0  | 38   |
| 4    |    | 2-Propanone, 1-(1H-pyrazol-1-yl)- | 124 | C6H8N2O | 1000337-51-8 | 35   |
| 5    |    | 1,3-Heptadiene, 2,3-dimethyl-     | 124 | C9H16   | 074779-65-0  | 30   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

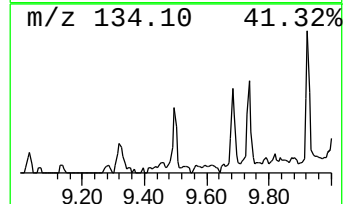
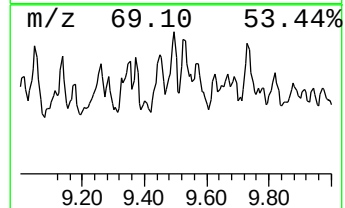
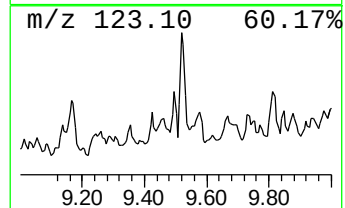
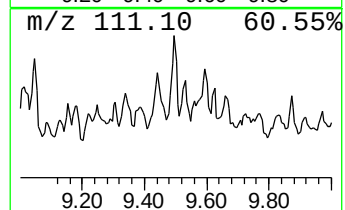
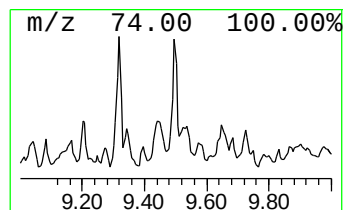
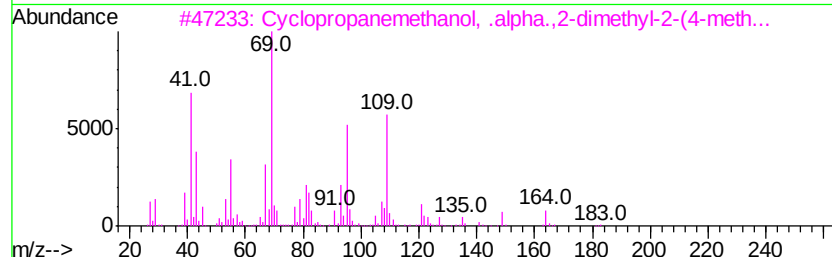
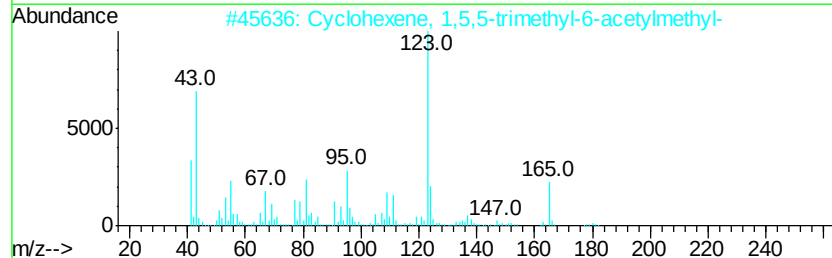
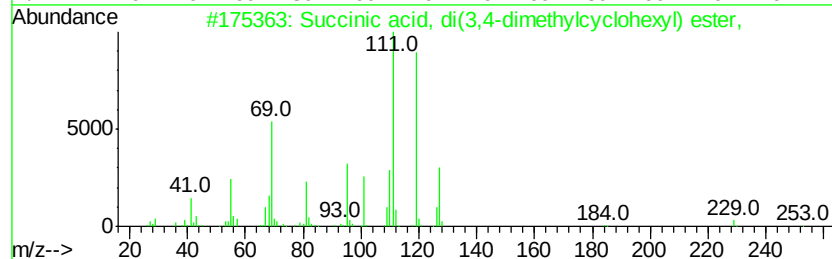
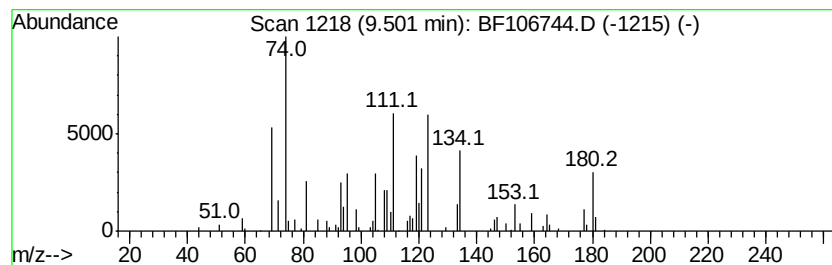
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 unknown9.50 Concentration Rank 15

| R.T. | EstConc  | Area   | Relative to ISTD | R.T.  |
|------|----------|--------|------------------|-------|
| 9.50 | 11.00 ng | 158361 | Acenaphthene-d10 | 10.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Succinic acid, di(3,4-dimethylcv... | 338 | C20H34O4 | 1000330-03-7 | 22   |
| 2    |    | Cyclohexene, 1,5,5-trimethyl-6-a... | 180 | C12H20O  | 211563-96-1  | 22   |
| 3    |    | Cyclopropanemethanol, .alpha.,2-... | 182 | C12H22O  | 121959-70-4  | 16   |
| 4    |    | 2,2-Dimethylocta-3,4-dienal         | 152 | C10H16O  | 000590-71-6  | 14   |
| 5    |    | 1,3-Hexadiene, 3-ethyl-2-methyl-... | 124 | C9H16    | 074752-97-9  | 11   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

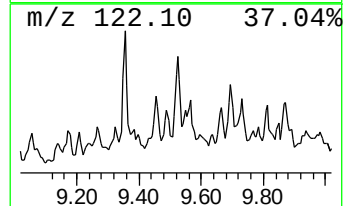
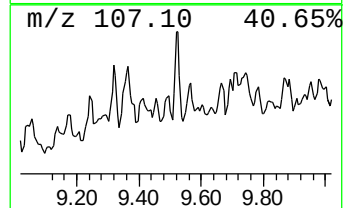
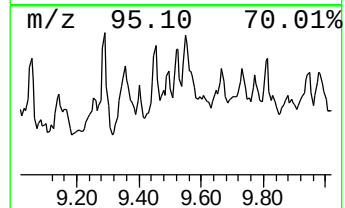
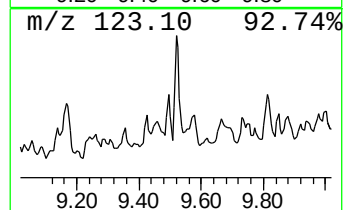
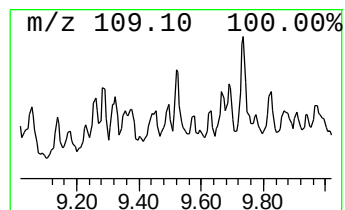
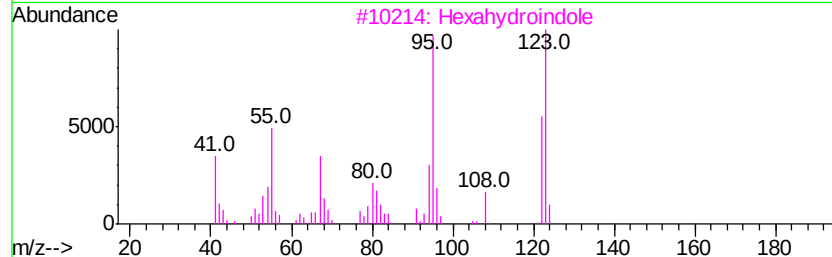
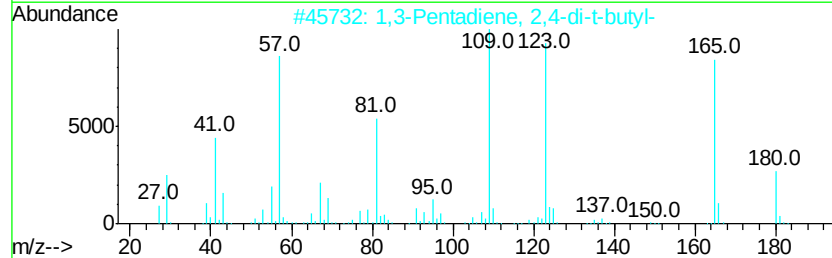
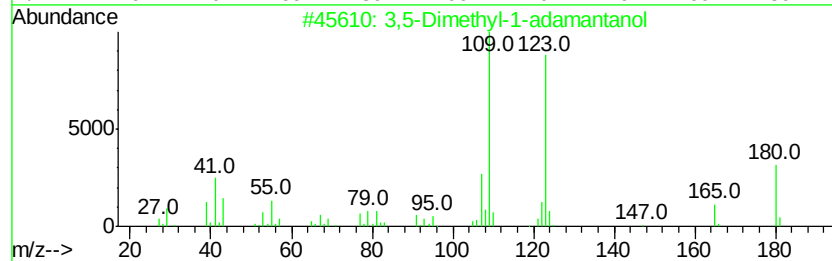
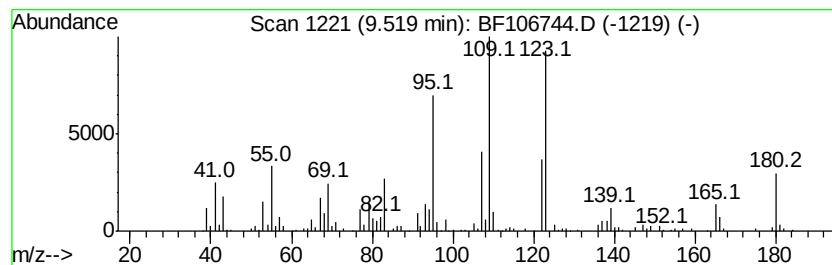
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown9.52 Concentration Rank 6

| R.T. | EstConc  | Area   | Relative to ISTD | R.T.  |
|------|----------|--------|------------------|-------|
| 9.52 | 21.36 ng | 307408 | Acenaphthene-d10 | 10.01 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-------------------------------------|---|--------------|-----|----------|-------------|------|
| 1    | 3,5-Dimethyl-1-adamantanol          |   |              | 180 | C12H20O  | 000707-37-9 | 49   |
| 2    | 1,3-Pentadiene, 2,4-di-t-butyl-     |   |              | 180 | C13H24   | 132850-21-6 | 45   |
| 3    | Hexahydroindole                     |   |              | 123 | C8H13N   | 018159-32-5 | 38   |
| 4    | Tricyclo[4.3.1.1(3,8)]undecane, ... |   |              | 180 | C12H20O  | 021898-92-0 | 38   |
| 5    | cis-1,1-Diethoxyhex-4-en-2-yne      |   |              | 168 | C10H16O2 | 093847-15-5 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

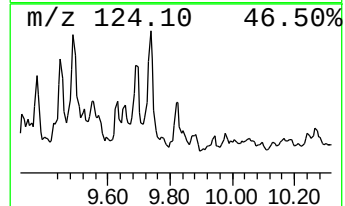
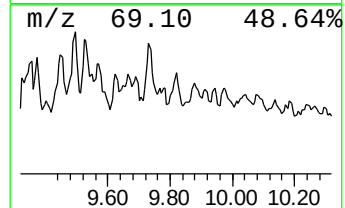
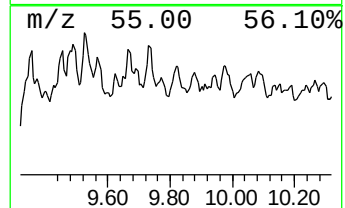
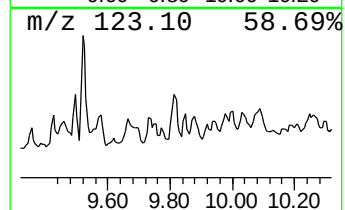
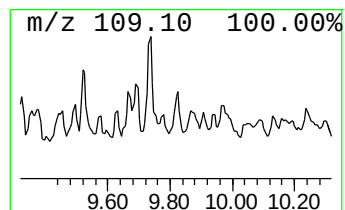
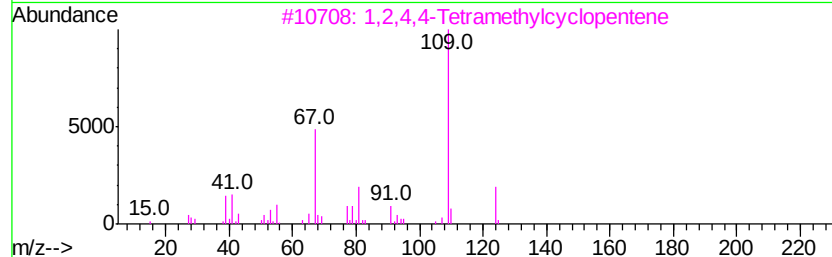
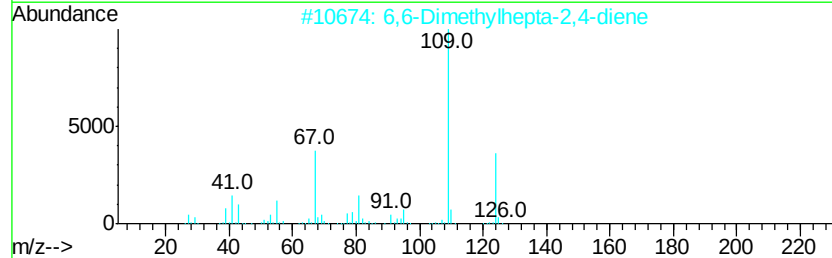
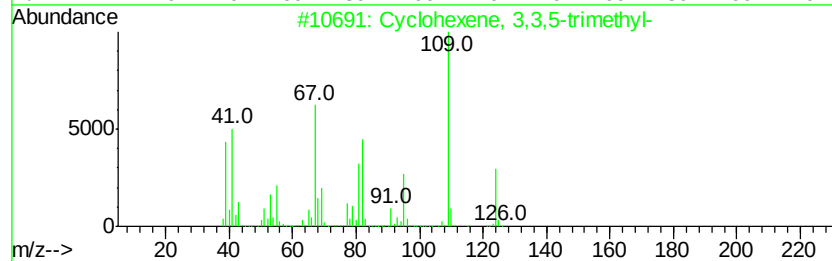
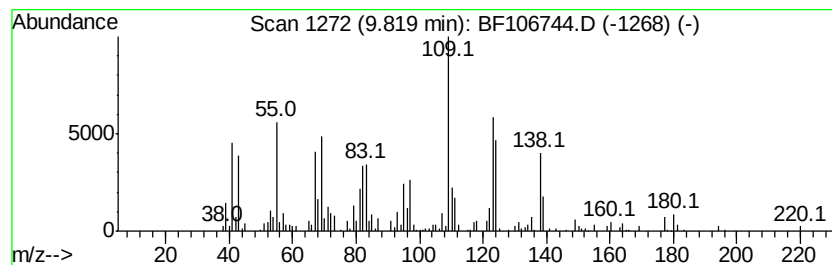
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 unknown9.82 Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD | R.T.  |
|------|----------|--------|------------------|-------|
| 9.82 | 16.49 ng | 237354 | Acenaphthene-d10 | 10.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Cyclohexene, 3,3,5-trimethyl-       | 124 | C9H16   | 000503-45-7  | 49   |
| 2    |    | 6,6-Dimethylhepta-2,4-diene         | 124 | C9H16   | 1000195-03-3 | 43   |
| 3    |    | 1,2,4,4-Tetramethylcyclopentene     | 124 | C9H16   | 065378-76-9  | 38   |
| 4    |    | 2,4-Heptadiene, 2,6-dimethyl-       | 124 | C9H16   | 004634-87-1  | 38   |
| 5    |    | Ethanone, 1-(1,3-dimethyl-3-cycl... | 152 | C10H16O | 051733-68-7  | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

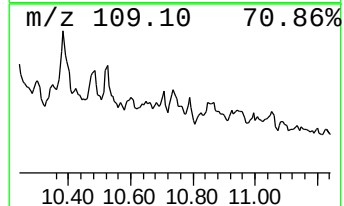
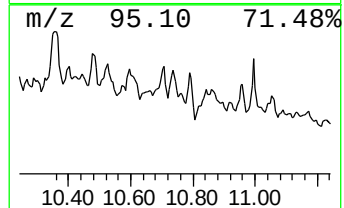
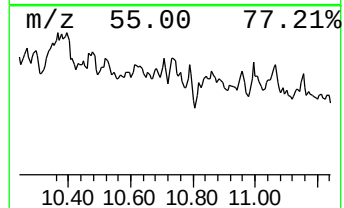
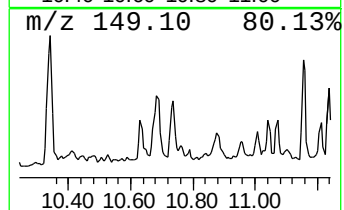
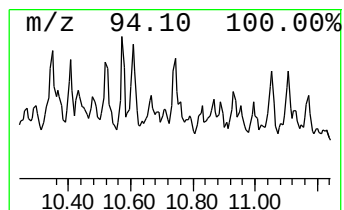
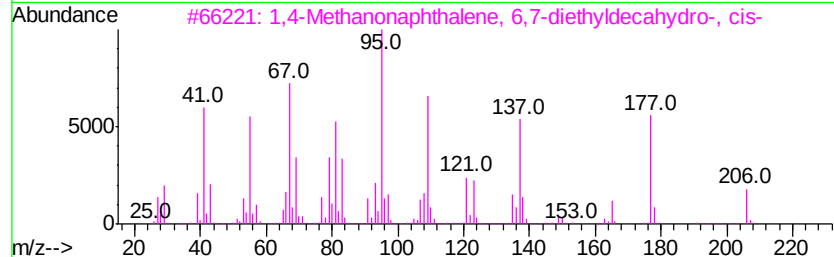
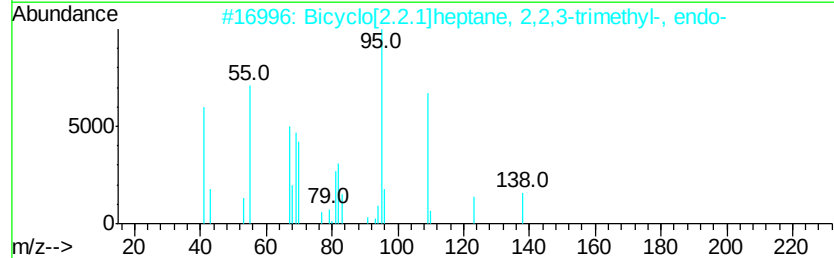
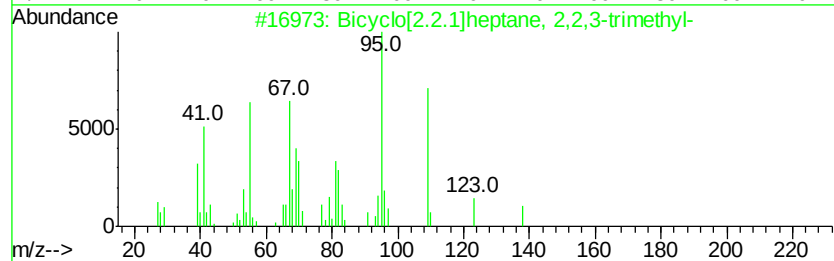
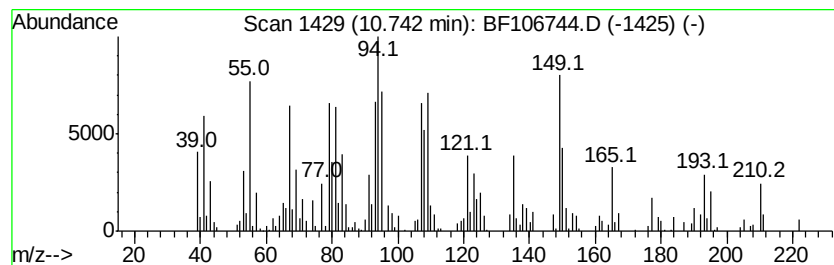
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 unknown10.74 Concentration Rank 17

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.74 | 10.31 ng | 148434 | Acenaphthene-d10 | 10.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18   | 000473-19-8  | 42   |
| 2    |    | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18   | 020536-40-7  | 38   |
| 3    |    | 1,4-Methanonaphthalene, 6,7-diet... | 206 | C15H26   | 016539-02-9  | 38   |
| 4    |    | Cyclopentanecarboxylic acid, 2-m... | 154 | C9H14O2  | 074764-24-2  | 25   |
| 5    |    | Phthalic acid, entyl undec-2-en-... | 346 | C21H30O4 | 1000315-41-3 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

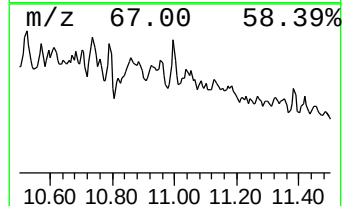
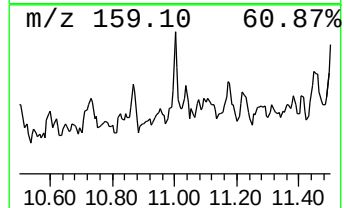
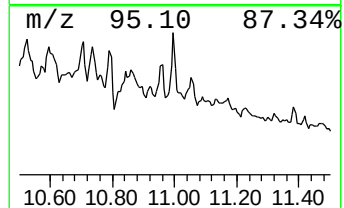
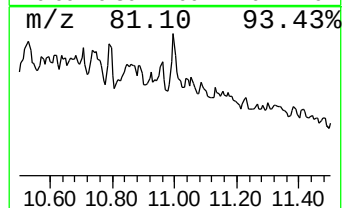
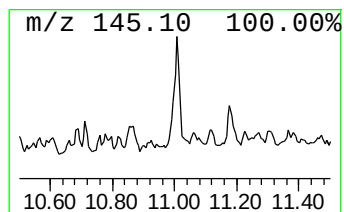
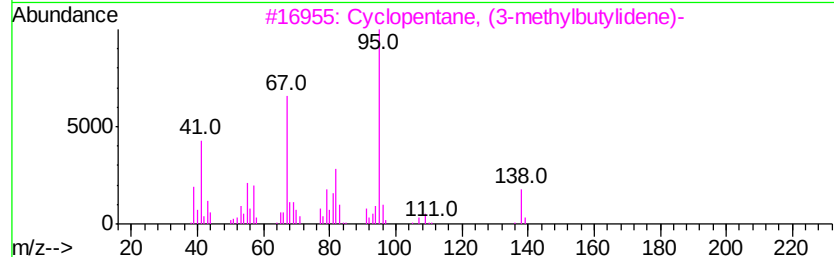
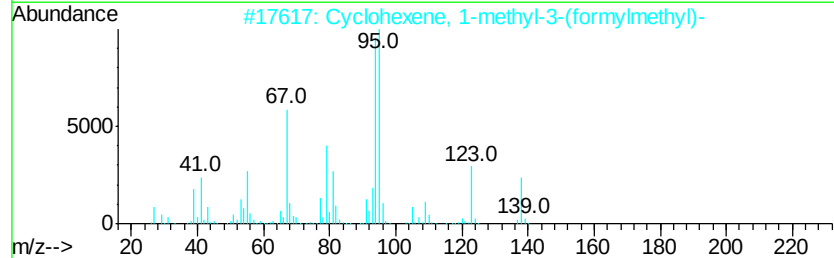
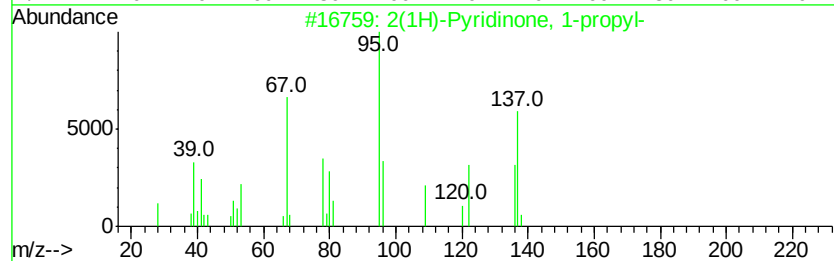
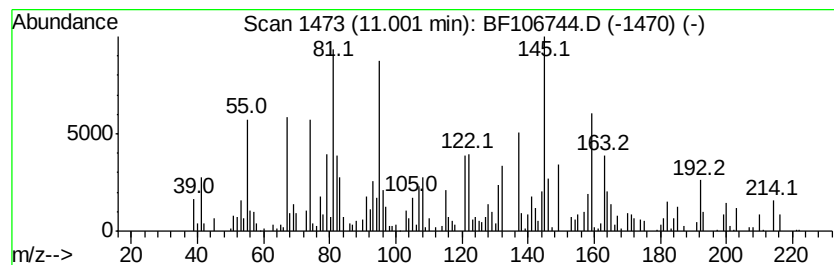
Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 19 2(1H)-Pyridinone, 1-propyl- Concentration Rank 18

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.00     | 9.93 ng                             | 129712 | Phenanthrene-d10 | 11.50       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 2(1H)-Pyridinone, 1-propyl-         | 137    | C8H11NO          | 019006-63-4 | 47   |
| 2         | Cyclohexene, 1-methyl-3-(formylm... | 138    | C9H14O           | 129993-40-4 | 46   |
| 3         | Cyclopentane, (3-methylbutylidene)- | 138    | C10H18           | 053366-51-1 | 38   |
| 4         | 1,3-Dimethyl-1-cyclohexene          | 110    | C8H14            | 002808-76-6 | 38   |
| 5         | .alpha.-Santoline alcohol           | 154    | C10H18O          | 090823-36-2 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\  
Data File : BF106744.D  
Acq On : 23 Jun 2018 7:46  
Operator : JU/SJ  
Sample : J3596-05  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

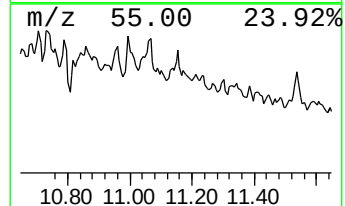
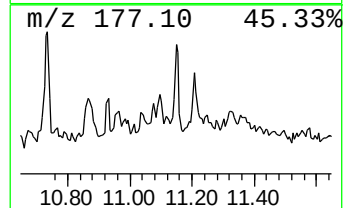
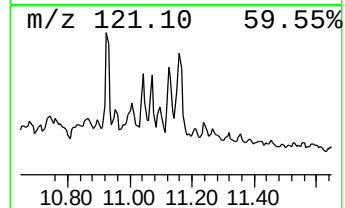
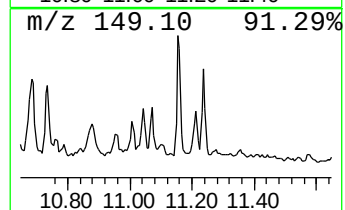
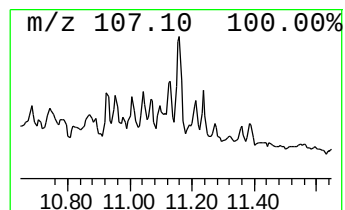
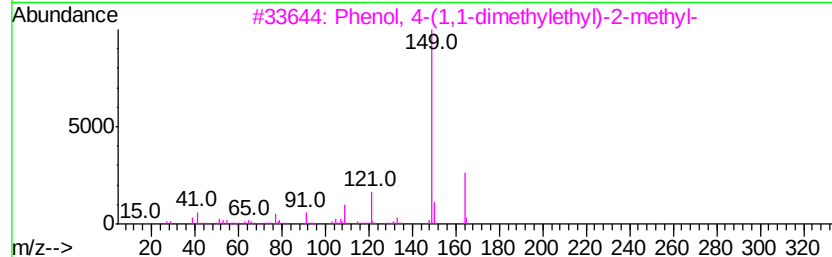
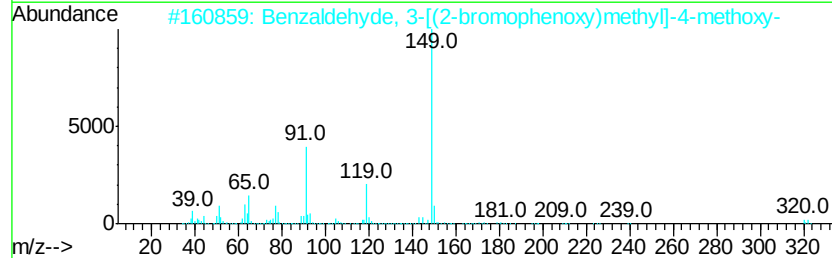
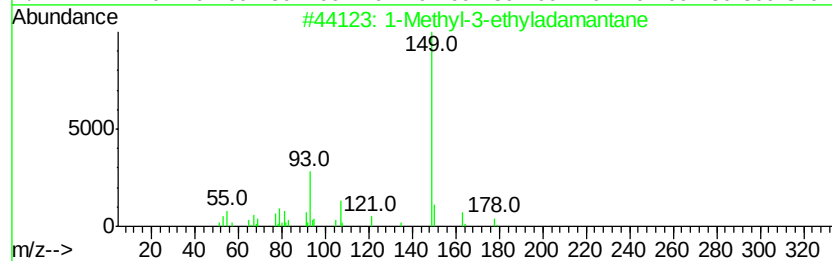
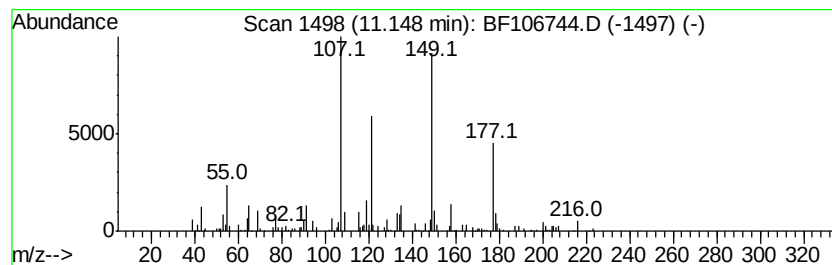
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 unknown11.15 Concentration Rank 14

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.15 | 11.12 ng | 145275 | Phenanthrene-d10 | 11.50 |

| Hit# | of | Tentative ID                                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-----------------------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 1-Methyl-3-ethyladamantane                          | 178 | C13H22     | 001687-34-9  | 47   |
| 2    |    | Benzaldehyde, 3-[(2-bromophenoxy)methyl]-4-methoxy- | 320 | C15H13BrO3 | 1000350-79-6 | 43   |
| 3    |    | Phenol, 4-(1,1-dimethylethyl)-2-methyl-             | 164 | C11H16O    | 000098-27-1  | 43   |
| 4    |    | Benzaldehyde, 3-(4-bromophenoxy)methyl-             | 320 | C15H13BrO3 | 351365-97-4  | 43   |
| 5    |    | 2,4,7-Trioxabicyclo[4.4.0]9-dec-5-ene               | 468 | C30H44O4   | 1000160-93-8 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF062218\  
 Data File : BF106744.D  
 Acq On : 23 Jun 2018 7:46  
 Operator : JU/SJ  
 Sample : J3596-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF061918.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 |
| 2-Pentanone, 4-hy...  | 5.20  | 45.3    | ng    | 603258   | 1                     | 6.96  | 266381 | 20.0 |
| Hexylene glycol       | 6.12  | 9.7     | ng    | 128653   | 1                     | 6.96  | 266381 | 20.0 |
| 1,1-Hexylenedioxy...  | 6.32  | 13.9    | ng    | 185010   | 1                     | 6.96  | 266381 | 20.0 |
| unknown8.12           | 8.12  | 9.9     | ng    | 168629   | 2                     | 8.24  | 341704 | 20.0 |
| Bicyclo[3.3.2]dec...  | 8.37  | 10.8    | ng    | 184437   | 2                     | 8.24  | 341704 | 20.0 |
| unknown8.47           | 8.47  | 13.8    | ng    | 235662   | 2                     | 8.24  | 341704 | 20.0 |
| unknown8.69           | 8.69  | 22.1    | ng    | 377727   | 2                     | 8.24  | 341704 | 20.0 |
| unknown8.77           | 8.77  | 13.3    | ng    | 227592   | 2                     | 8.24  | 341704 | 20.0 |
| unknown8.92           | 8.92  | 12.0    | ng    | 205088   | 2                     | 8.24  | 341704 | 20.0 |
| Tricyclo[4.3.1.1(...  | 9.01  | 12.2    | ng    | 209192   | 2                     | 8.24  | 341704 | 20.0 |
| unknown9.05           | 9.05  | 18.2    | ng    | 311004   | 2                     | 8.24  | 341704 | 20.0 |
| unknown9.26           | 9.26  | 25.0    | ng    | 360275   | 3                     | 10.01 | 287838 | 20.0 |
| unknown9.50           | 9.50  | 11.0    | ng    | 158361   | 3                     | 10.01 | 287838 | 20.0 |
| unknown9.52           | 9.52  | 21.4    | ng    | 307408   | 3                     | 10.01 | 287838 | 20.0 |
| unknown9.82           | 9.82  | 16.5    | ng    | 237354   | 3                     | 10.01 | 287838 | 20.0 |
| unknown10.74          | 10.74 | 10.3    | ng    | 148434   | 3                     | 10.01 | 287838 | 20.0 |
| 2(1H)-Pyridinone, ... | 11.00 | 9.9     | ng    | 129712   | 4                     | 11.50 | 261274 | 20.0 |
| unknown11.15          | 11.15 | 11.1    | ng    | 145275   | 4                     | 11.50 | 261274 | 20.0 |