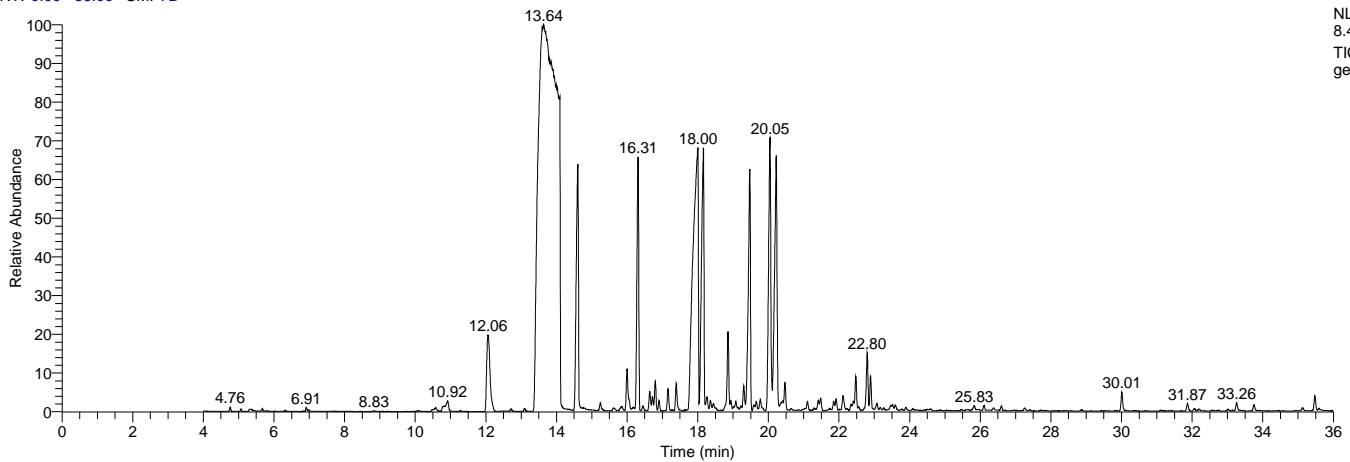


# My GC-MS Report

RT: 0.00 - 36.00 SM: 7B

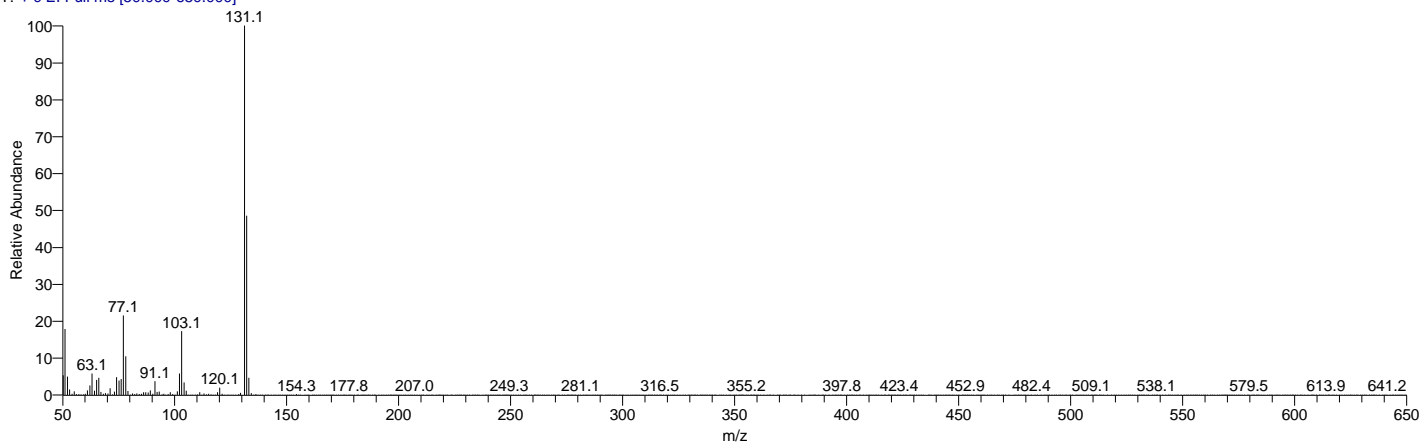


NL:  
8.47E8  
TIC MS  
gerfa\_EA

RT	Area %	Peak Area
10.92	0.24	128173820.51
12.06	1.94	1035114737.02
13.64	53.81	28642212603.14
14.60	4.27	2271905049.14
16.00	0.62	331863513.20
16.31	3.68	1956479496.62
16.64	0.24	127515039.35
16.72	0.16	84334567.09
16.80	0.35	188369875.93
16.90	0.11	57520312.66
17.16	0.28	146669995.78
17.39	0.38	200364399.84
18.00	9.92	5280783018.22
18.16	4.83	2570696982.40
18.26	0.11	59233142.90
18.86	1.05	557328644.63
19.30	0.26	137888084.26
19.47	4.14	2201097941.66
19.66	0.12	63533698.33
19.77	0.18	96526750.45
20.04	4.78	2545719837.35
20.22	4.91	2615571936.08
20.47	0.43	228816648.36
21.11	0.14	75018362.93
21.41	0.14	76347365.03
21.48	0.16	86057337.89
21.85	0.11	60872365.51
21.92	0.14	75870020.41
22.12	0.22	118894434.31
22.48	0.53	283703031.75
22.80	0.77	411968559.79
22.89	0.37	199251685.20
30.01	0.25	135159887.79
33.26	0.12	64939646.82
35.47	0.21	109604833.07

# My GC-MS Report

gerfa\_EA #2064 RT: 10.92 AV: 1 NL: 8.04E6  
T: + c EI Full ms [50.000-650.000]

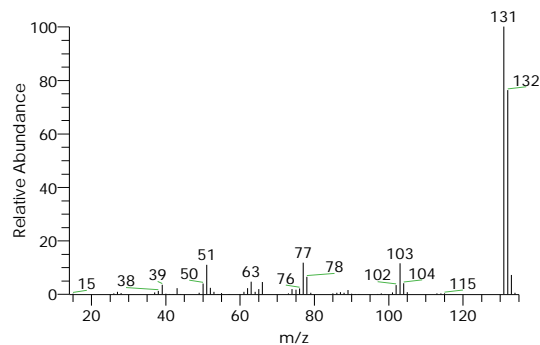
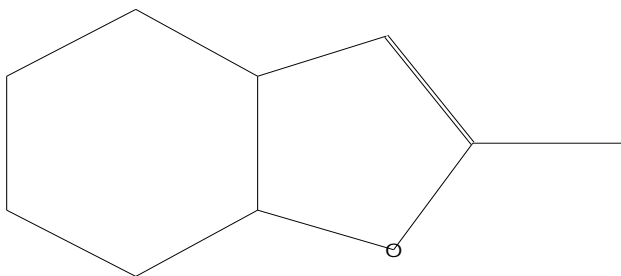


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
10.92	Benzofuran, 2-methyl-	0.24	C9H8O	132	4265-25-2	901	mainlib
10.92	Benzofuran, 2-methyl-	0.24	C9H8O	132	4265-25-2	905	replib
10.92	BENZOFURAN, 2-METHYL-	0.24	C9H8O	132	4265-25-2	905	WileyRegistry8e

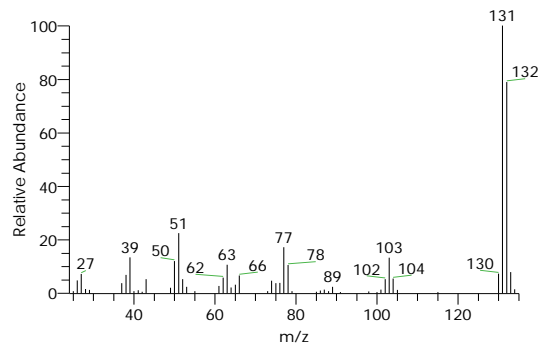
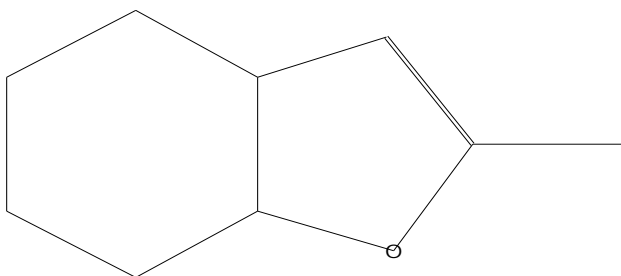
Compound Structure

Hit Spectrum

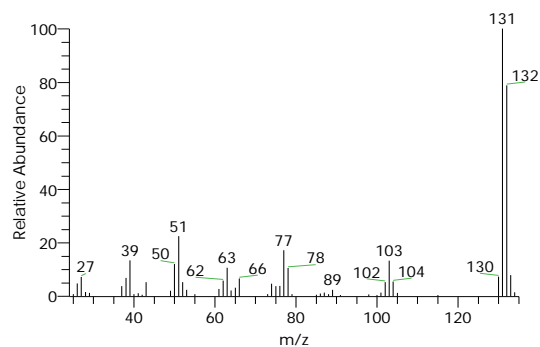
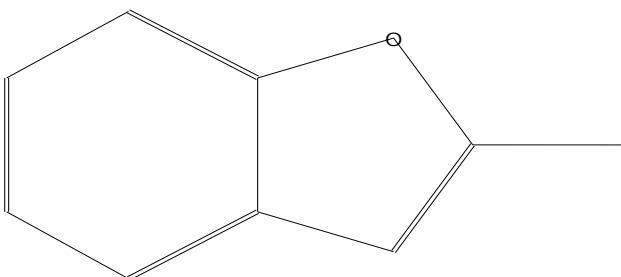
Benzofuran, 2-methyl-  
Formula C9H8O, MW 132, CAS# 4265-25-2, Entry# 115872  
2-Methylbenzofuran



Benzofuran, 2-methyl-  
Formula C9H8O, MW 132, CAS# 4265-25-2, Entry# 20943  
2-Methylbenzofuran

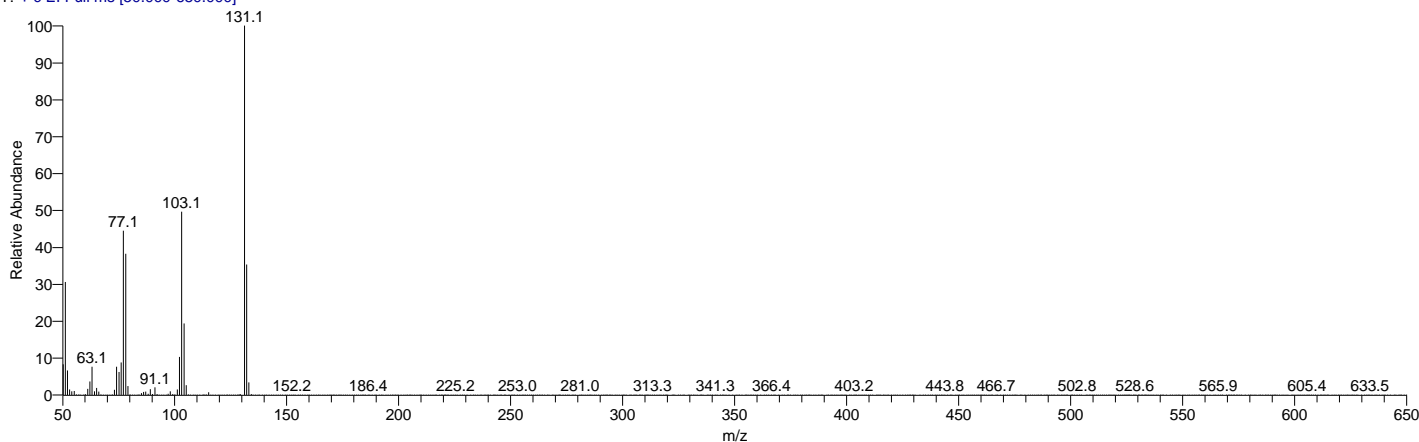


BENZOFURAN, 2-METHYL-  
Formula C9H8O, MW 132, CAS# 4265-25-2, Entry# 20576  
2-METHYLBENZOFURAN



# My GC-MS Report

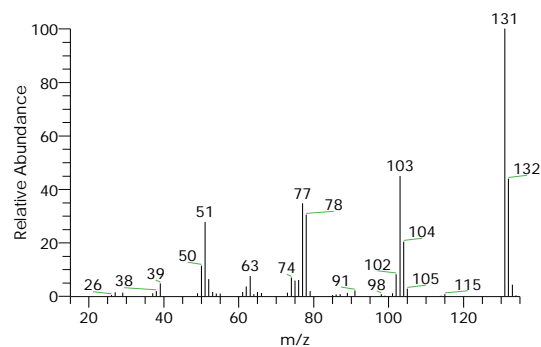
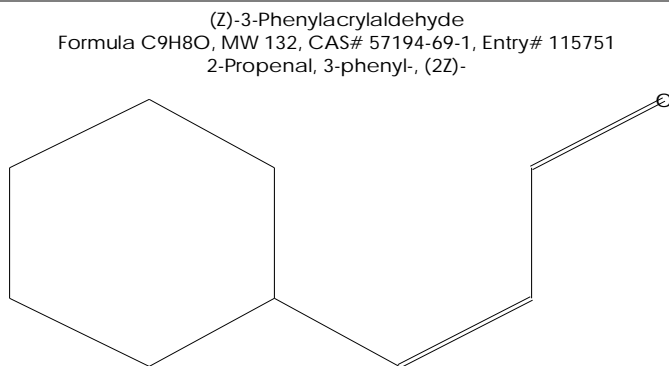
gerfa\_EA #2405 RT: 12.06 AV: 1 NL: 4.21E7  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
12.06	(Z)-3-Phenylacrylaldehyde	1.94	C9H8O	132	57194-69-1	960	mainlib
12.06	2-Propenal, 3-phenyl-	1.94	C9H8O	132	104-55-2	949	replib
12.06	Cinnamaldehyde, (E)-	1.94	C9H8O	132	14371-10-9	949	mainlib

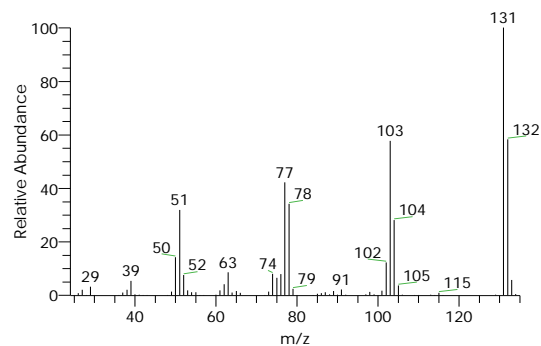
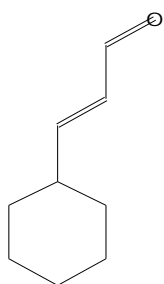
## Compound Structure

## Hit Spectrum



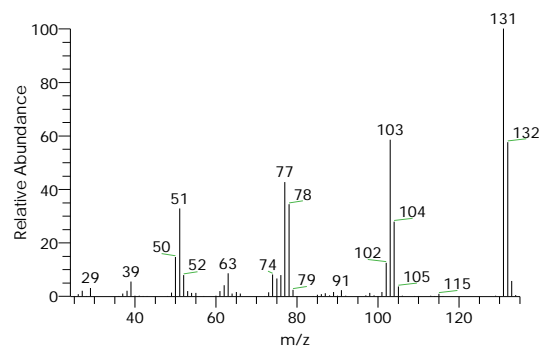
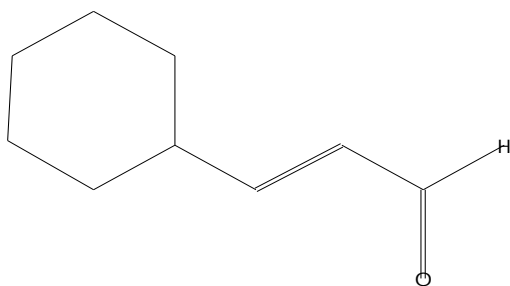
2-Propenal, 3-phenyl-  
Formula C9H8O, MW 132, CAS# 104-55-2, Entry# 20950

Cinnamaldehyde



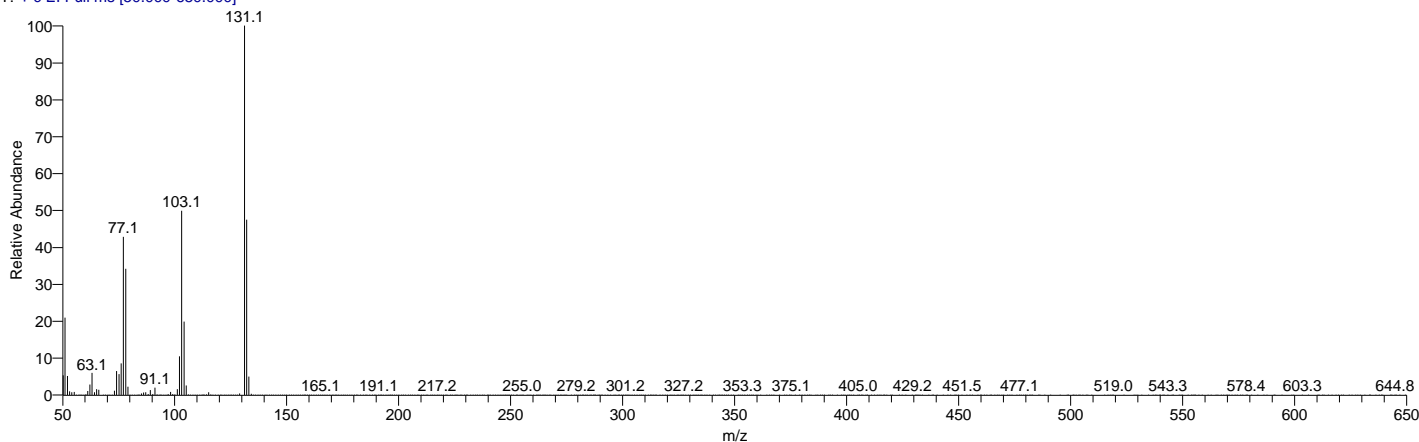
Cinnamaldehyde, (E)-  
Formula C9H8O, MW 132, CAS# 14371-10-9, Entry# 115750

(E)-Cinnamaldehyde



# My GC-MS Report

gerfa\_EA #2875 RT: 13.64 AV: 1 NL: 2.16E8  
T: + c EI Full ms [50.000-650.000]

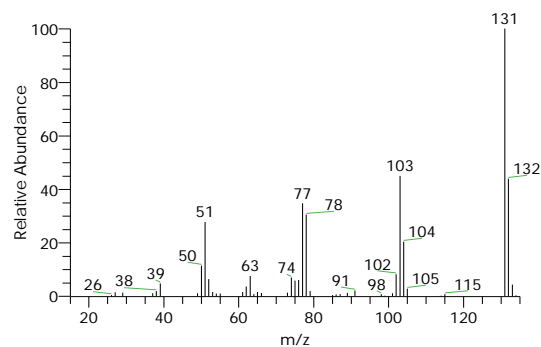
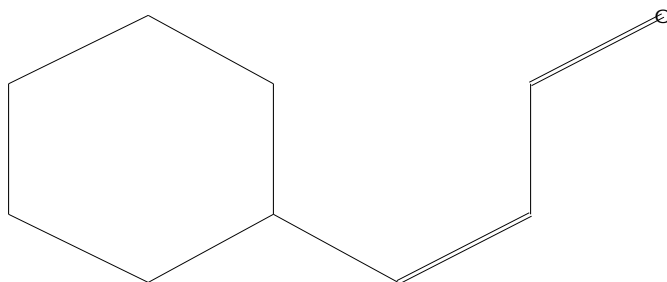


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
13.64	(Z)-3-Phenylacrylaldehyde	53.81	C9H8O	132	57194-69-1	959	mainlib
13.64	2-Propenal, 3-phenyl-	53.81	C9H8O	132	104-55-2	943	replib
13.64	Cinnamaldehyde, (E)-	53.81	C9H8O	132	14371-10-9	943	mainlib

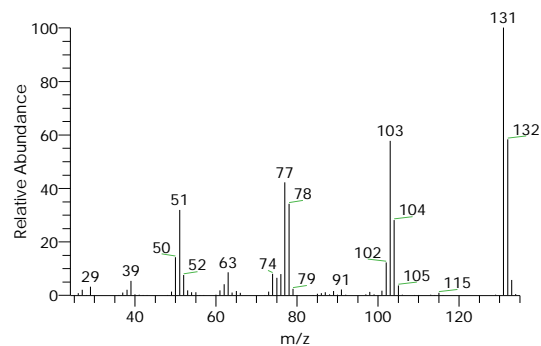
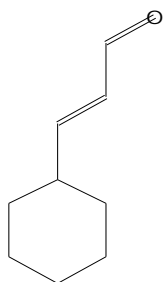
## Compound Structure

## Hit Spectrum

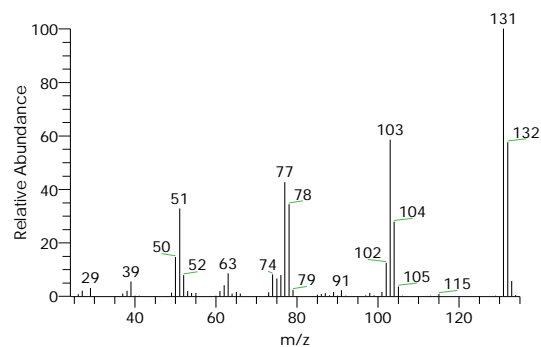
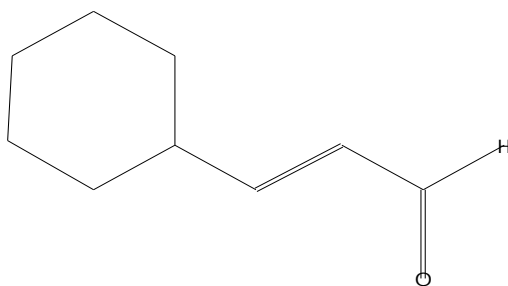
(Z)-3-Phenylacrylaldehyde  
Formula C9H8O, MW 132, CAS# 57194-69-1, Entry# 115751  
2-Propenal, 3-phenyl-, (Z)-



2-Propenal, 3-phenyl-  
Formula C9H8O, MW 132, CAS# 104-55-2, Entry# 20950  
Cinnamaldehyde

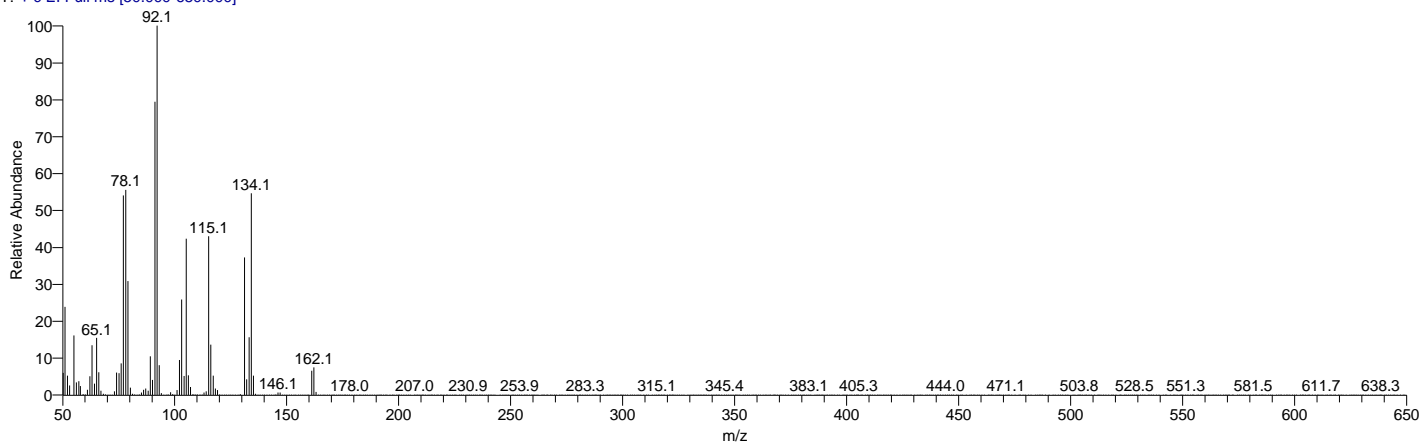


Cinnamaldehyde, (E)-  
Formula C9H8O, MW 132, CAS# 14371-10-9, Entry# 115750  
(E)-Cinnamaldehyde



# My GC-MS Report

gerfa\_EA #3162 RT: 14.60 AV: 1 NL: 6.98E7  
T: + c EI Full ms [50.000-650.000]

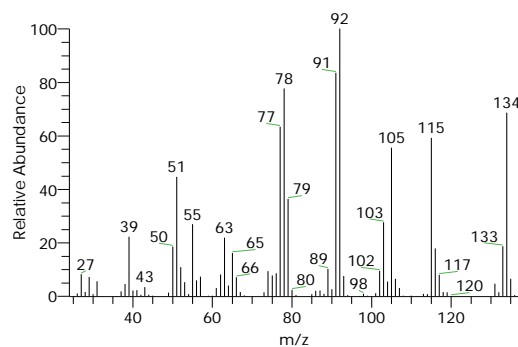
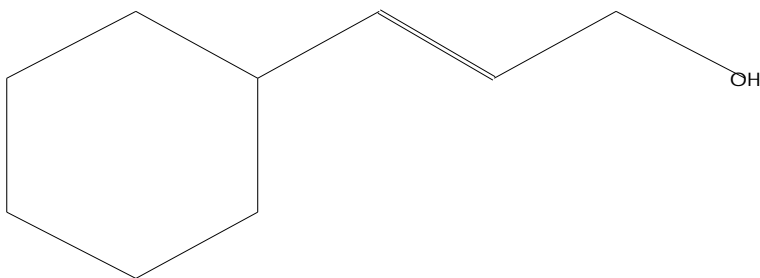


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
14.60	2-Propen-1-ol, 3-phenyl-, (E)-	4.27	C9H10O	134	4407-36-7	926	mainlib
14.60	2-Propen-1-ol, 3-phenyl-	4.27	C9H10O	134	104-54-1	914	mainlib
14.60	2-Propen-1-ol, 3-phenyl-	4.27	C9H10O	134	104-54-1	908	replib

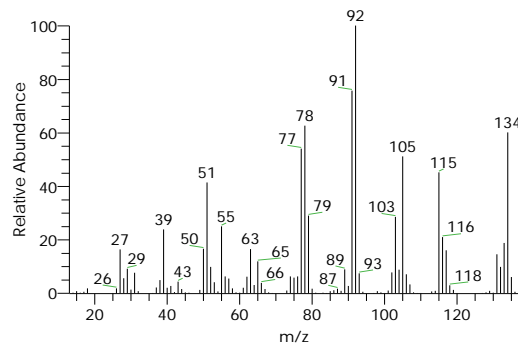
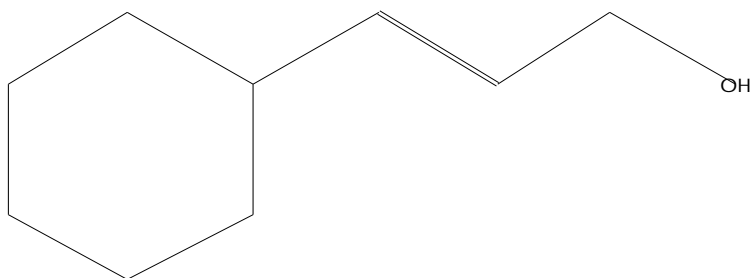
Compound Structure

Hit Spectrum

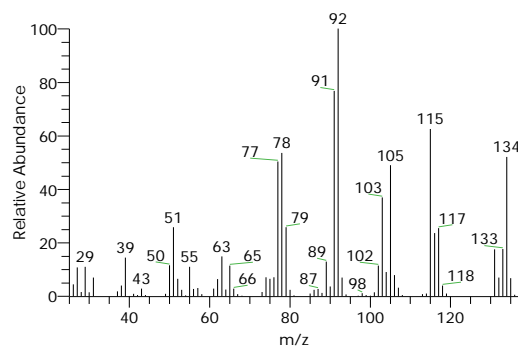
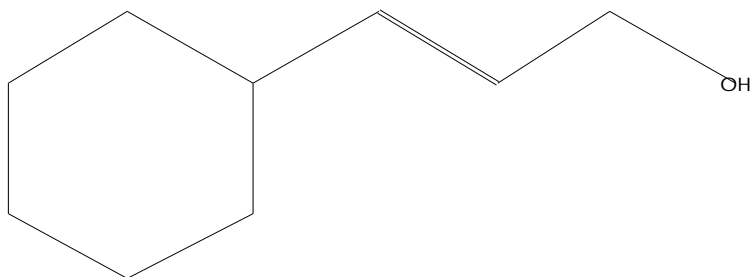
2-Propen-1-ol, 3-phenyl-, (E)-  
Formula C9H10O, MW 134, CAS# 4407-36-7, Entry# 65164  
\$:28OOCDEMITAIZTP-QPJXVBHSA-N



2-Propen-1-ol, 3-phenyl-  
Formula C9H10O, MW 134, CAS# 104-54-1, Entry# 65165  
Cinnamyl alcohol

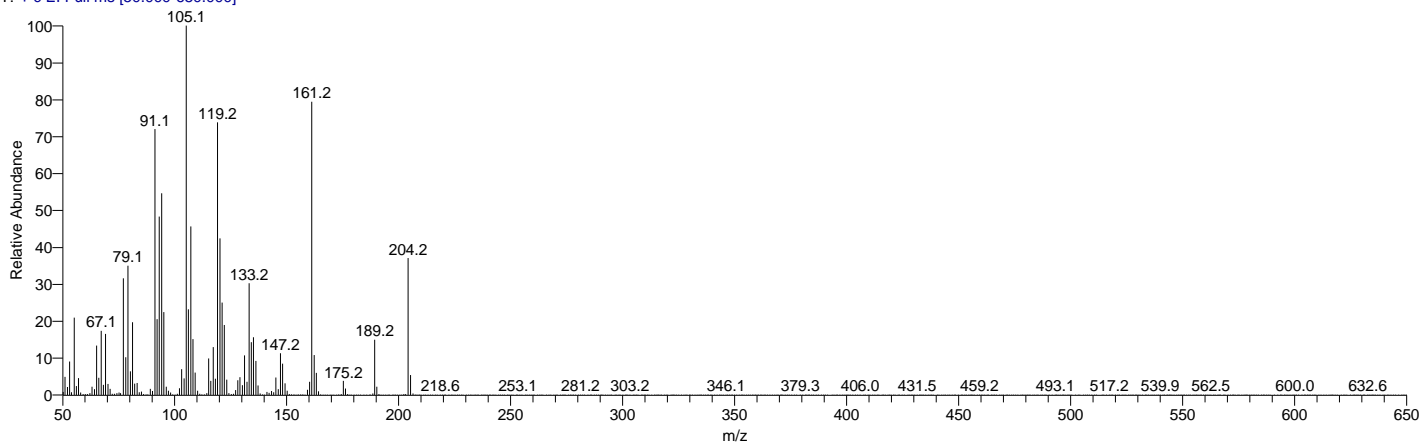


2-Propen-1-ol, 3-phenyl-  
Formula C9H10O, MW 134, CAS# 104-54-1, Entry# 13932  
Cinnamyl alcohol



# My GC-MS Report

gerfa\_EA #3579 RT: 16.00 AV: 1 NL: 8.57E6  
T: + c EI Full ms [50.000-650.000]

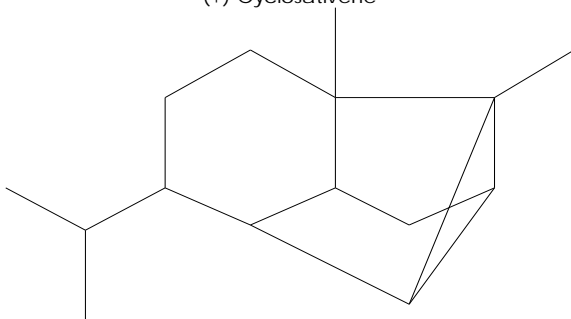


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.00	1,2,4-Metheno-1H-indene, octahydro-1,7a-dimethyl-5-(1-methylethyl)-, [1S-(1à,2à,3aà,4à,5à,7aà,8S*)]-	0.62	C15H24	204	22469-52-9	935	replib
16.00	(+)-CYCLOISOSATIVENE	0.62	C15H24	204	NA	915	WileyRegistry8e
16.00	1,2,4-Metheno-1H-indene, octahydro-1,7a-dimethyl-5-(1-methylethyl)-, [1S-(1à,2à,3aà,4à,5à,7aà,8S*)]-	0.62	C15H24	204	22469-52-9	911	replib

Compound Structure

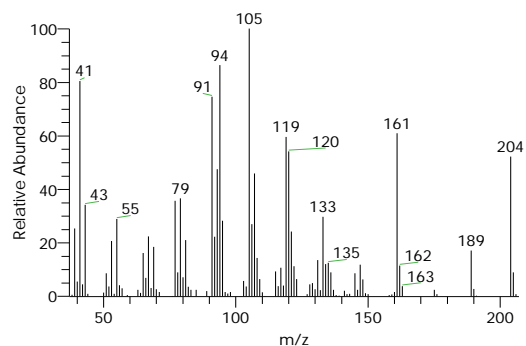
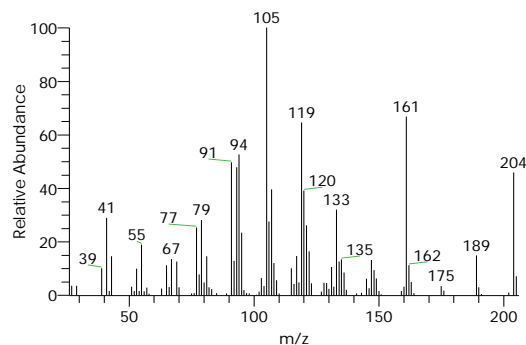
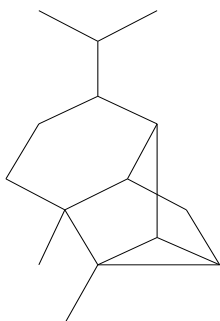
Hit Spectrum

Formula C15H24, MW 204, CAS# 22469-52-9, Entry# 16426  
(+)-Cyclosativene



(+)-CYCLOISOSATIVENE

Formula C15H24, MW 204, CAS# NA, Entry# 89651



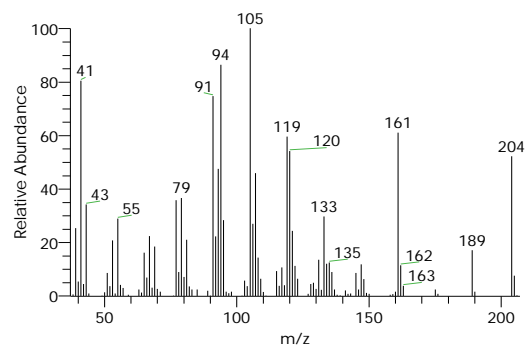
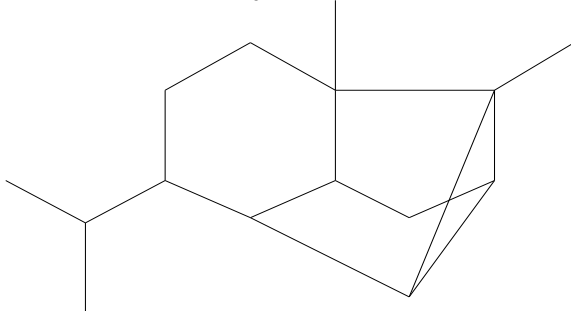
# My GC-MS Report

Compound Structure

Hit Spectrum

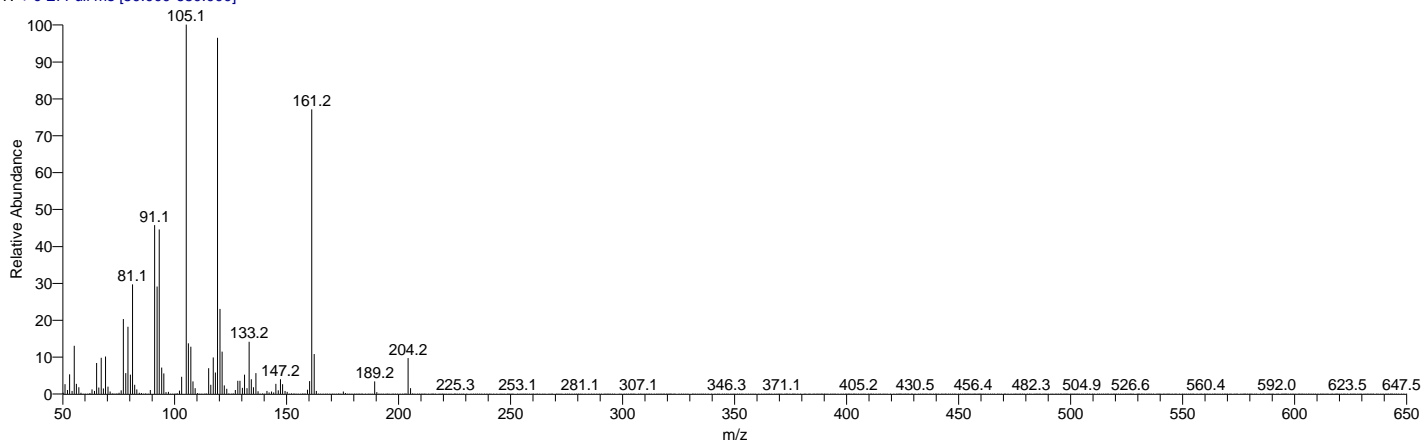
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 22469-52-9, Entry# 16246

(+)-Cyclosativene



gerfa\_EA #3671 RT: 16.31 AV: 1 NL: 7.59E7

T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.31	TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, ST	3.68	C <sub>15</sub> H <sub>24</sub>	204	3856-2 5-5	951	WileyRegi
16.31	.alfa.-Copaene	3.68	C <sub>15</sub> H <sub>24</sub>	204	NA	930	mainlib
16.31	TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, ST	3.68	C <sub>15</sub> H <sub>24</sub>	204	3856-2 5-5	952	WileyRegi

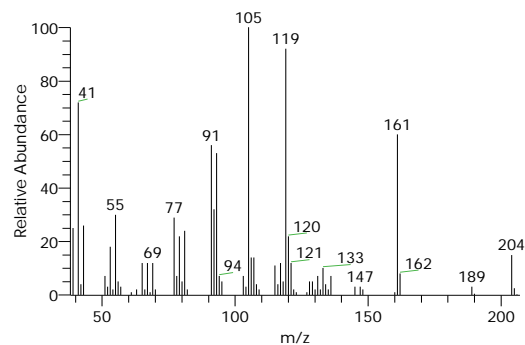
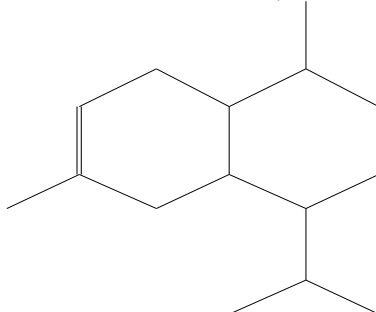
Compound Structure

Hit Spectrum

TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, ST

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3856-25-5, Entry# 89462

TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, STEREOISOMER

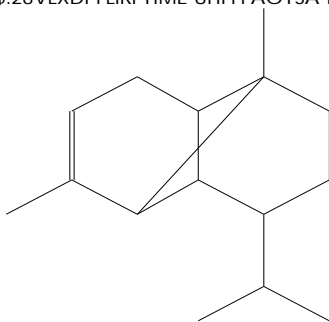


# My GC-MS Report

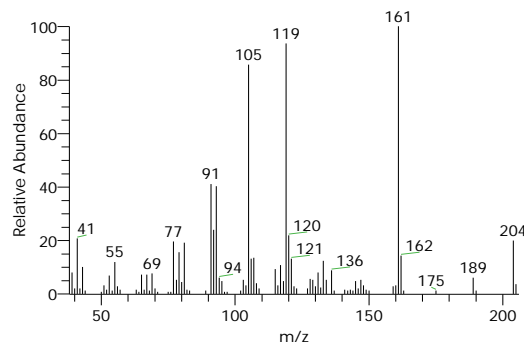
Compound Structure

Hit Spectrum

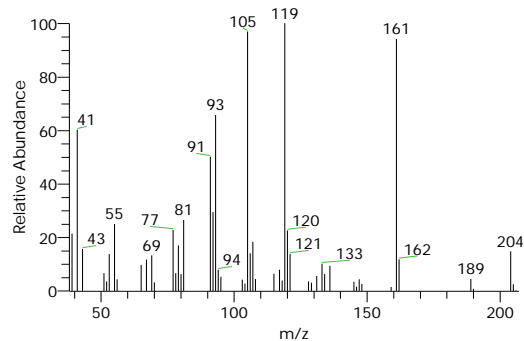
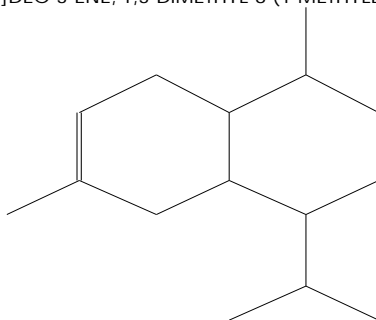
.alpha.-Copaene  
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# NA, Entry# 150578  
\$:28VLXDPFLIRFYIME-UHFFFAOYSA-N



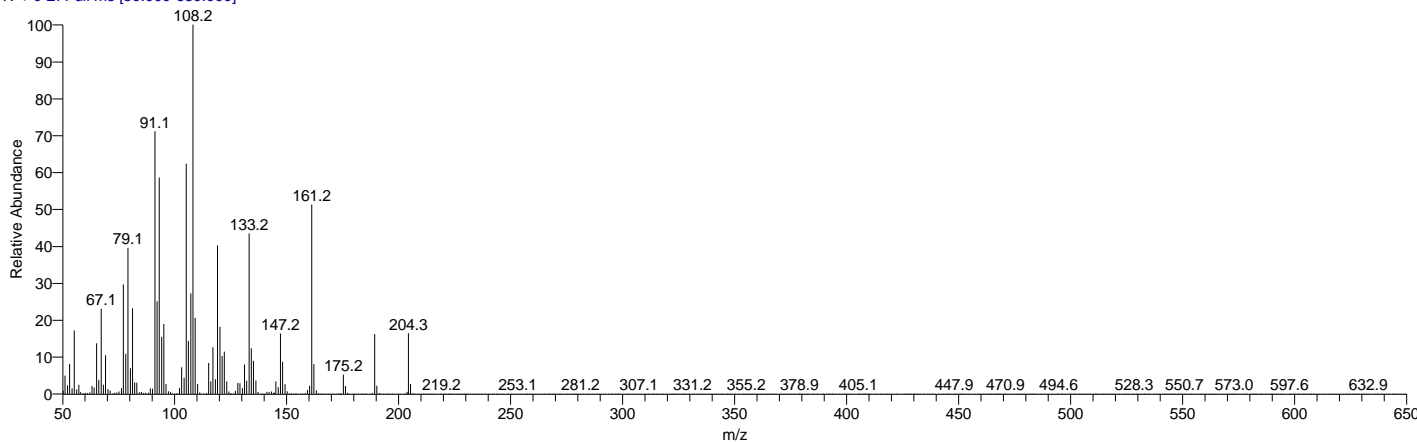
SI 928, RSI 930, mainlib, Entry# 150578, CAS# NA, .alpha.-Copaene



TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, ST  
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3856-25-5, Entry# 89463  
TRICYCLO[4.4.0.0(2,7)]DEC-3-ENE, 1,3-DIMETHYL-8-(1-METHYLETHYL)-, STEREOISOMER



gerfa\_EA #3770 RT: 16.64 AV: 1 NL: 4.70E6  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.64	1,4-Methano-1H-indene, octahydro-4-methyl-8-methylene-7-(1-methylethyl)-, [1S-(1à,3aá,4à,7à,7aá)]-	0.24	C <sub>15</sub> H <sub>24</sub>	204	3650-28-0	915	mainlib
16.64	(+)-SATIVEN	0.24	C <sub>15</sub> H <sub>24</sub>	204	3650-28-0	915	WileyRegistry8e
16.64	1,4-METHANO-1H-INDENE, OCTAHYDRO-4-METHYL-8-METHYLENE-7-(1-METHYLETHYL)-, [1S-(1à,3Aá,4à,7à,7Aá)]-	0.24	C <sub>15</sub> H <sub>24</sub>	204	3650-28-0	915	WileyRegistry8e

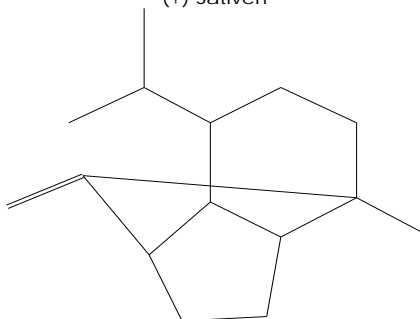


# My GC-MS Report

Compound Structure

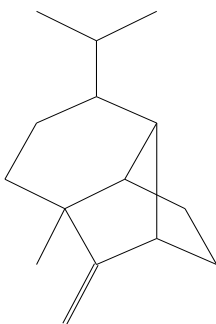
Hit Spectrum

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3650-28-0, Entry# 84625  
(+)-Sativen



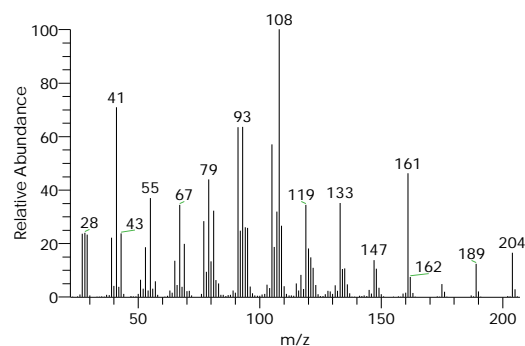
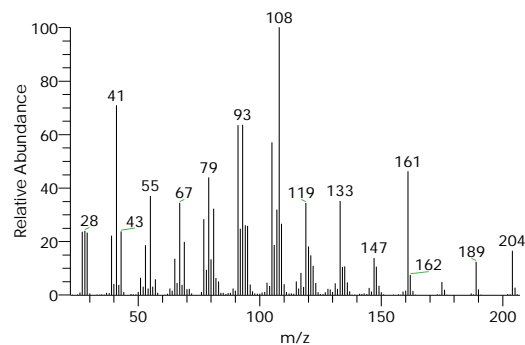
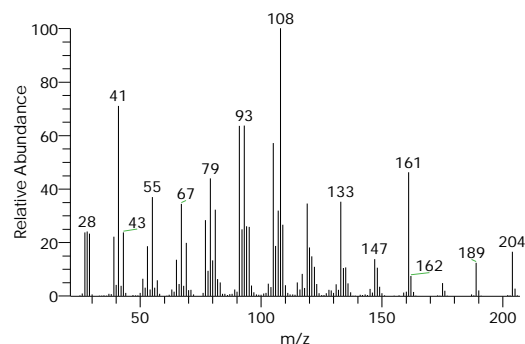
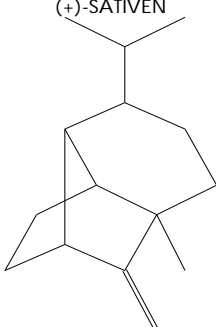
(+)-SATIVEN

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3650-28-0, Entry# 392536

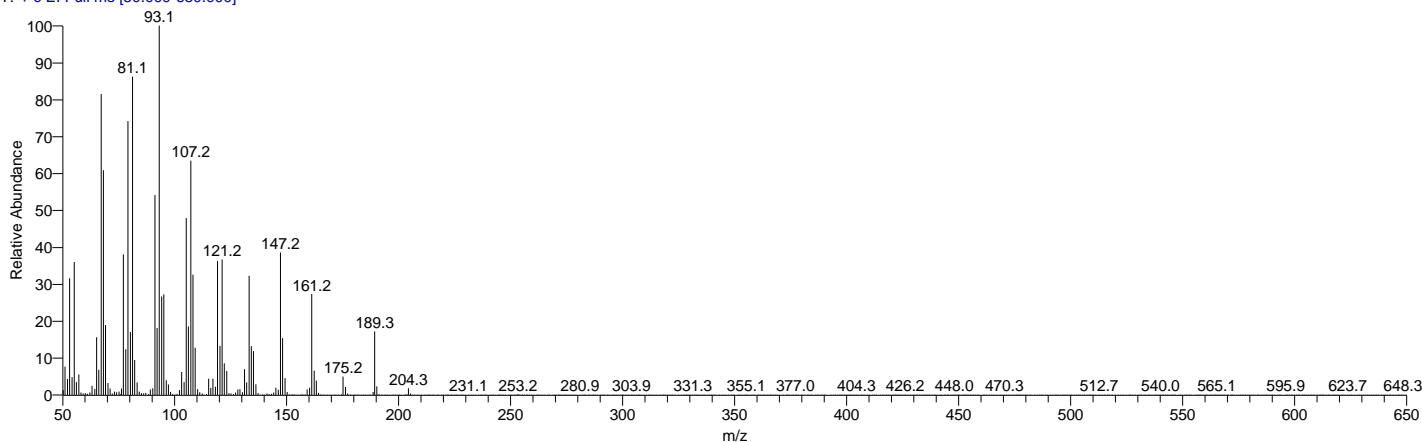


(+)-SATIVEN

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3650-28-0, Entry# 89185



gerfa\_EA #3793 RT: 16.72 AV: 1 NL: 2.71E6  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.72	CYCLOHEXANE, 1-ETHENYL-1-METHYL-2,4-BIS(1-METHYLETHENYL)-, [1S-(1à,2á,4á)]-	0.16	C <sub>15</sub> H <sub>24</sub>	204	515-13-9	949	WileyRegistry8e

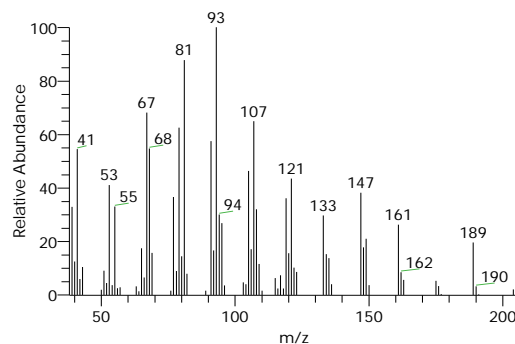
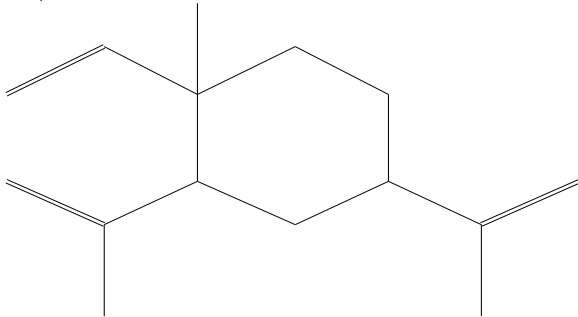
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.72	CYCLOHEXANE, 1-ETHENYL-1-METHYL-2,4-BIS(1-METHYLETHENYL)-, [1S-(1à,2á,4á)]-	0.16	C15H24	204	515-13-9	936	WileyRegistry8e
16.72	Cyclohexane, 1-ethenyl-1-methyl-2,4-bis(1-methylethenyl)-, [1S-(1à,2á,4á)]-	0.16	C15H24	204	515-13-9	914	replib

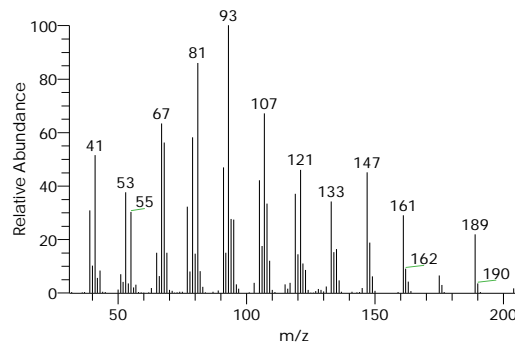
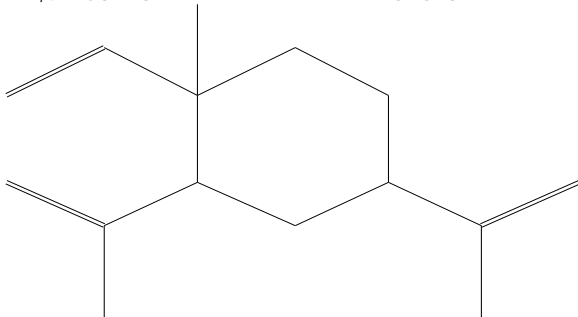
Compound Structure

Hit Spectrum

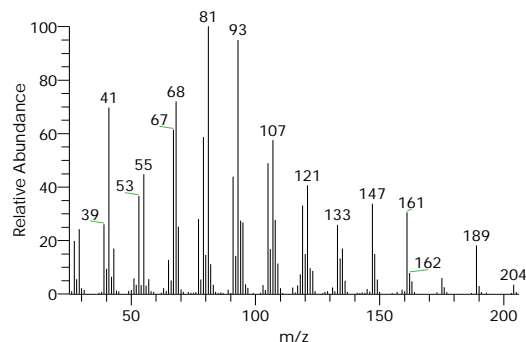
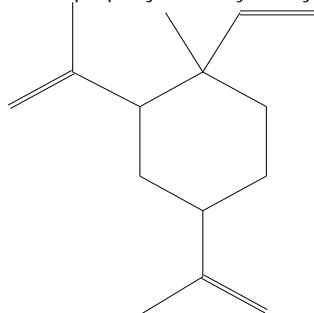
CYCLOHEXANE, 1-ETHENYL-1-METHYL-2,4-BIS(1-METHYLETHENYL)-, [1S-(1à,2á,4á)]-  
Formula C15H24, MW 204, CAS# 515-13-9, Entry# 89511  
2,4-DIISOPROPENYL-1-METHYL-1-VINYLCYCLOHEXANE #



CYCLOHEXANE, 1-ETHENYL-1-METHYL-2,4-BIS(1-METHYLETHENYL)-, [1S-(1à,2á,4á)]-  
Formula C15H24, MW 204, CAS# 515-13-9, Entry# 89566  
2,4-DIISOPROPENYL-1-METHYL-1-VINYLCYCLOHEXANE #

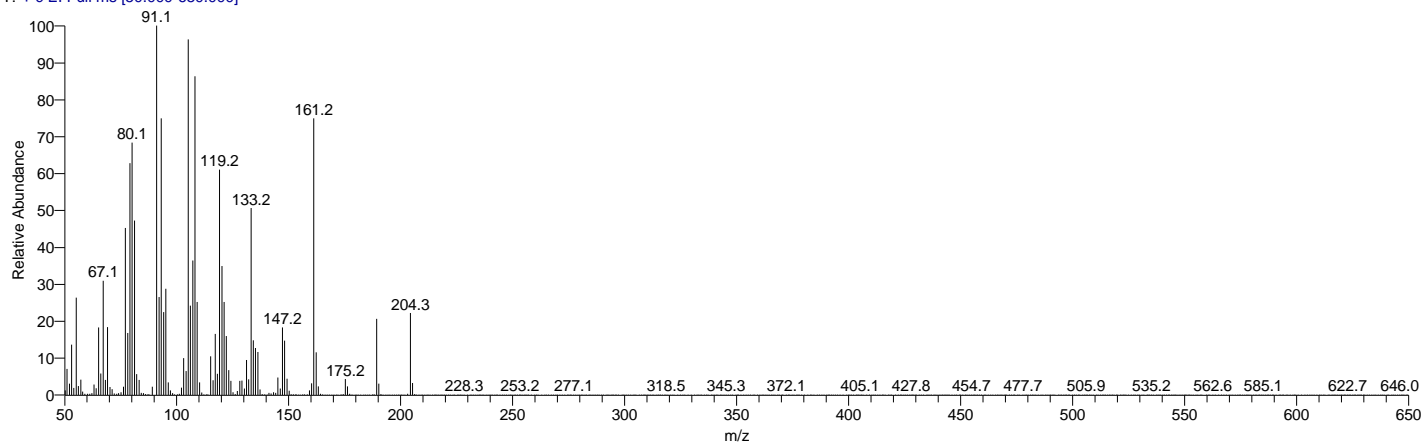


Cyclohexane, 1-ethenyl-1-methyl-2,4-bis(1-methylethenyl)-, [1S-(1à,2á,4á)]-  
Formula C15H24, MW 204, CAS# 515-13-9, Entry# 11582  
Cyclohexane, 2,4-diisopropenyl-1-methyl-1-vinyl-, (1S,2R,4R)- (-)-



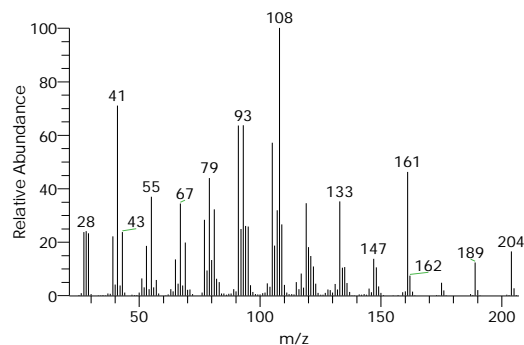
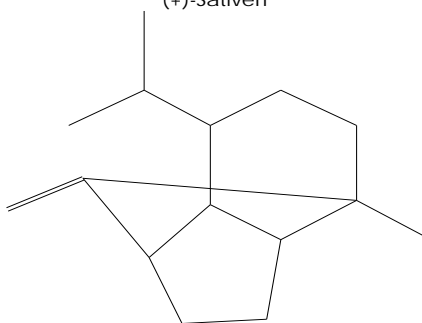
# My GC-MS Report

gerfa\_EA #3817 RT: 16.80 AV: 1 NL: 4.87E6  
T: + c EI Full ms [50.000-650.000]

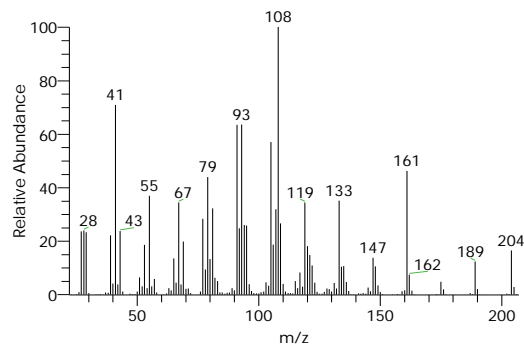
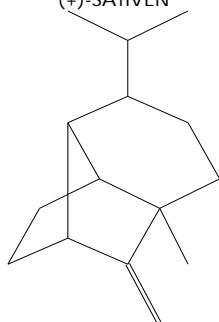


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.80	1,4-Methano-1H-indene, octahydro-4-methyl-8-methylene-7-(1-methylethyl)-, [1S-(1à,3aá,4à,7à,7aá)]-	0.35	C15H24	204	3650-28-0	908	mainlib
16.80	1,4-METHANO-1H-INDENE, OCTAHYDRO-4-METHYL-8-METHYLENE-7-(1-METHYLETHYL)-, [1S-(1à,3Aá,4à,7à,7Aá)]-	0.35	C15H24	204	3650-28-0	908	WileyRegistry8e
16.80	(+)-SATIVEN	0.35	C15H24	204	3650-28-0	908	WileyRegistry8e
Compound Structure							Hit Spectrum

Formula C15H24, MW 204, CAS# 3650-28-0, Entry# 84625  
(+)-Sativen



Formula C15H24, MW 204, CAS# 3650-28-0, Entry# 89185  
(+)-SATIVEN



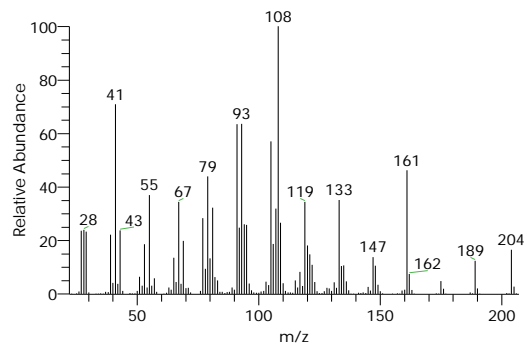
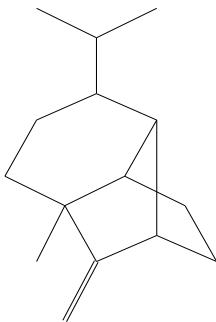
# My GC-MS Report

Compound Structure

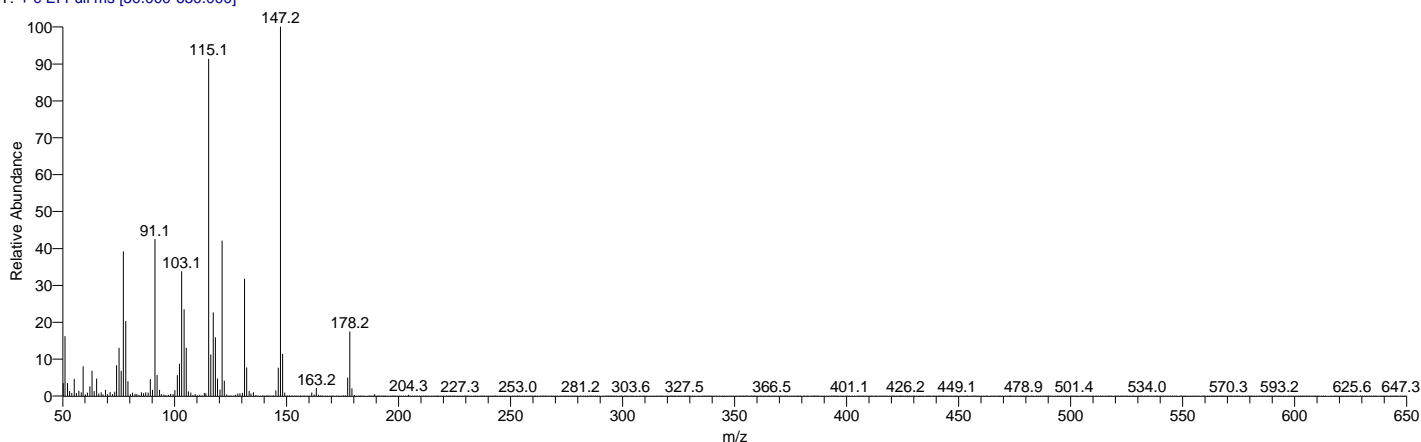
Hit Spectrum

(+)-SATIVEN

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 3650-28-0, Entry# 392536



gerfa\_EA #3849 RT: 16.90 AV: 1 NL: 3.66E6  
T: + c EI Full ms [50.000-650.000]



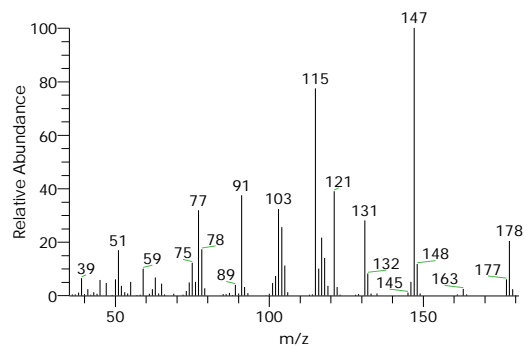
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
16.90	CINNAMALDEHYDE DIMETHYL ACETAL	0.11	C <sub>11</sub> H <sub>14</sub> O <sub>2</sub>	178	NA	921	WileyRegi stry8e
16.90	9-Methoxybicyclo[6.1.0]nona-2,4,6-triene	0.11	C <sub>10</sub> H <sub>12</sub> O	148	826-14-2	774	mainlib
16.90	9-METHOXYBICYCLO[6.1.0]NONA-2,4,6-TRIENE	0.11	C <sub>10</sub> H <sub>12</sub> O	148	NA	774	WileyRegi stry8e

Compound Structure

Hit Spectrum

CINNAMALDEHYDE DIMETHYL ACETAL

Formula C<sub>11</sub>H<sub>14</sub>O<sub>2</sub>, MW 178, CAS# NA, Entry# 60971

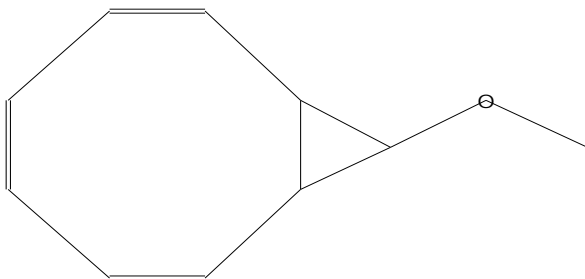


# My GC-MS Report

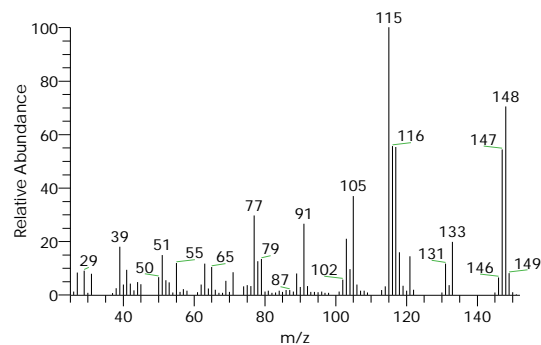
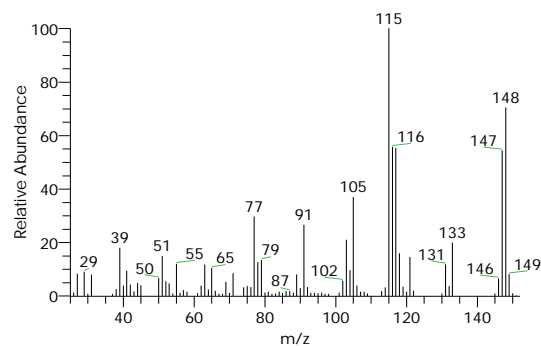
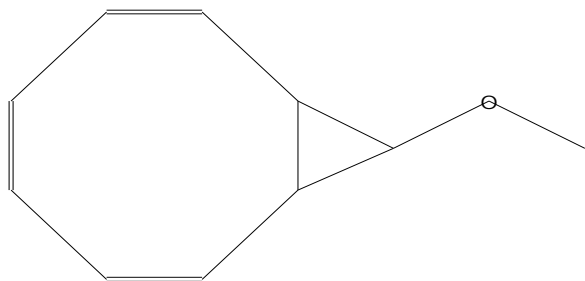
Compound Structure

Hit Spectrum

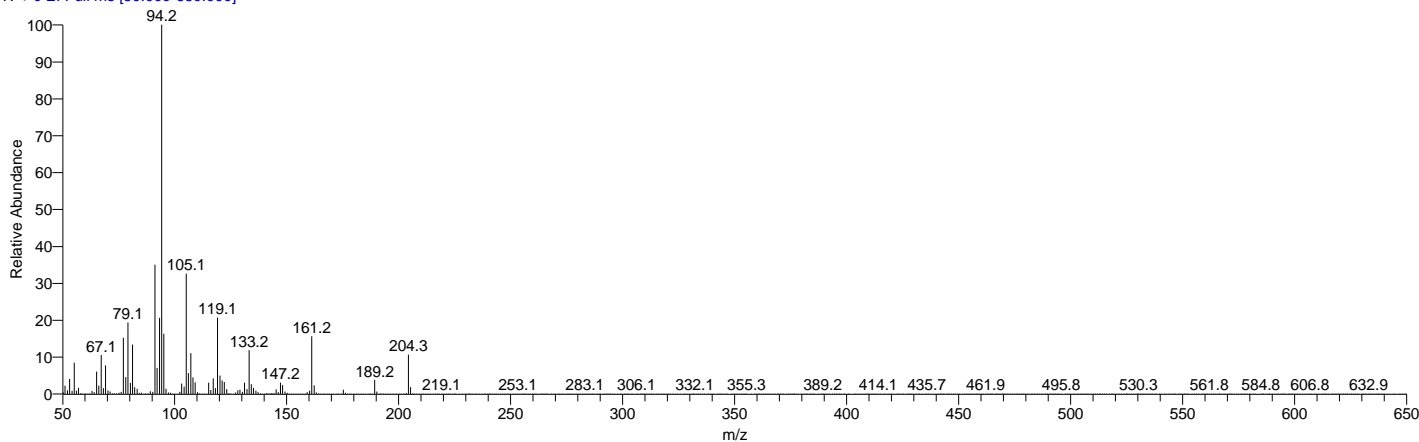
9-Methoxybicyclo[6.1.0]nona-2,4,6-triene  
Formula C<sub>10</sub>H<sub>12</sub>O, MW 148, CAS# 826-14-2, Entry# 94587  
Bicyclo[6.1.0]nona-2,4,6-trien-9-yl methyl ether #



9-METHOXYBICYCLO[6.1.0]NONA-2,4,6-TRIENE  
Formula C<sub>10</sub>H<sub>12</sub>O, MW 148, CAS# NA, Entry# 370644  
9-METHOXY-BICYCLO[6.1.0]NONA-2,4,6-TRIENE



gerfa\_EA #3925 RT: 17.16 AV: 1 NL: 1.13E7  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
17.16	Isosativene	0.28	C <sub>15</sub> H <sub>24</sub>	204	24959-83-9	958	replib
17.16	(-)-ISOSATIVENE	0.28	C <sub>15</sub> H <sub>24</sub>	204	24959-83-9	929	WileyRegistry8e
17.16	Isosativene	0.28	C <sub>15</sub> H <sub>24</sub>	204	24959-83-9	923	replib

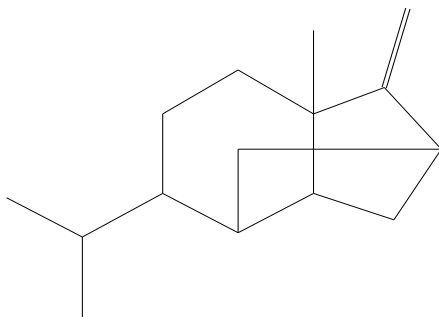
# My GC-MS Report

Compound Structure

Hit Spectrum

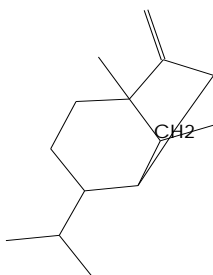
Isosativene

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 24959-83-9, Entry# 14356



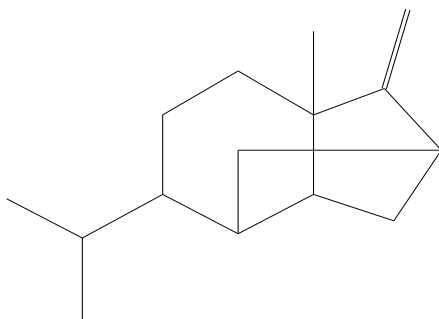
(-)-ISOSATIVENE

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 24959-83-9, Entry# 89650

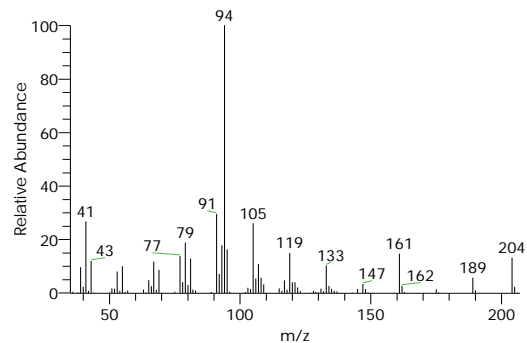
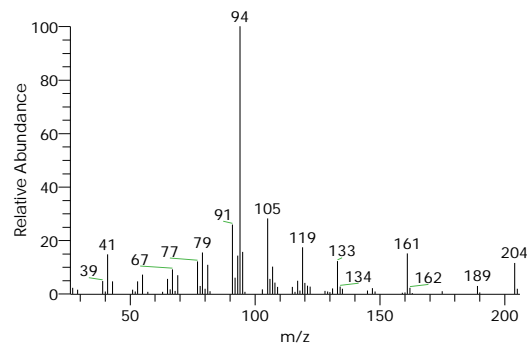


Isosativene

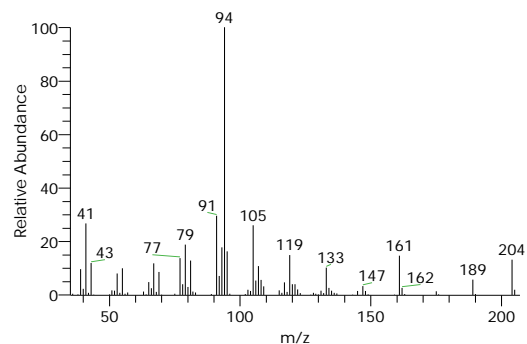
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 24959-83-9, Entry# 14330



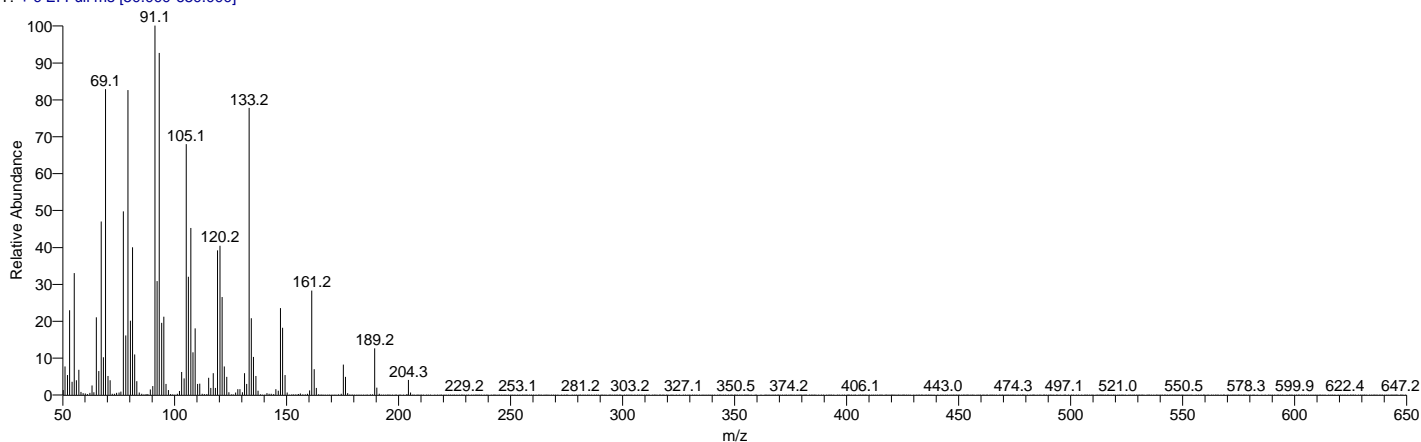
SI 938, RSI 958, replib, Entry# 14356, CAS# 24959-83-9, Isosativene



SI 917, RSI 923, replib, Entry# 14330, CAS# 24959-83-9, Isosativene



gerfa\_EA #3994 RT: 17.39 AV: 1 NL: 5.00E6  
T: + c EI Full ms [50.000-650.000]

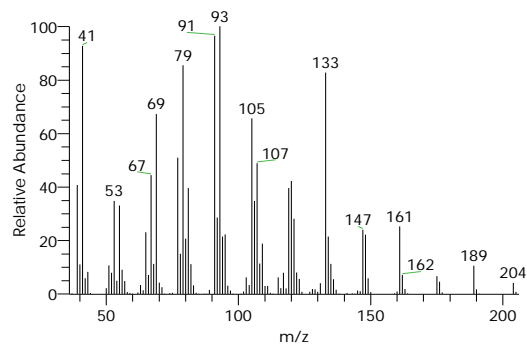
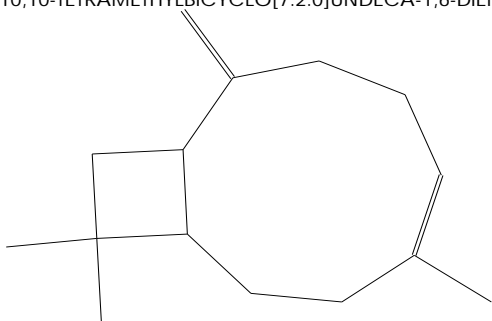


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
17.39	BICYCLO[7.2.0]UNDEC-4-ENE, 4,11,11-TRIMETHYL-8-METHYLENE- [1R-(1R*,4E,9S*)]-	0.38	C <sub>15</sub> H <sub>24</sub>	204	87-44-5	951	WileyRegi stry8e

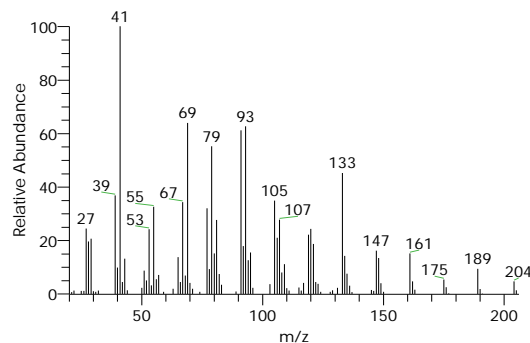
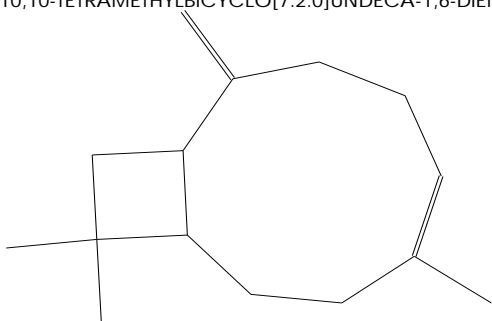
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
17.39	BICYCLO[7.2.0]UNDEC-4-ENE, 4,11,11-TRIMETHYL-8-METHYLENE-, [1R-(1R*,4E,9S*)]-	0.38	C15H24	204	87-44-5	948	WileyRegi stry8e
17.39	Caryophyllene	0.38	C15H24	204	87-44-5	934	mainlib
Compound Structure				Hit Spectrum			

BICYCLO[7.2.0]UNDEC-4-ENE, 4,11,11-TRIMETHYL-8-METHYLENE-, [1R-(1R\*,4E,9S\*)]-  
Formula C15H24, MW 204, CAS# 87-44-5, Entry# 89162  
2,6,10,10-TETRAMETHYLBICYCLO[7.2.0]UNDECA-1,6-DIENE #



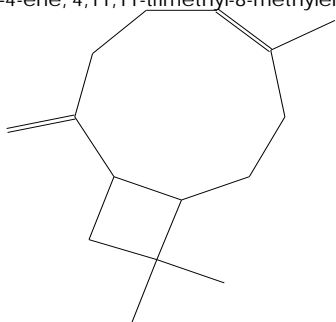
BICYCLO[7.2.0]UNDEC-4-ENE, 4,11,11-TRIMETHYL-8-METHYLENE-, [1R-(1R\*,4E,9S\*)]-  
Formula C15H24, MW 204, CAS# 87-44-5, Entry# 89170  
2,6,10,10-TETRAMETHYLBICYCLO[7.2.0]UNDECA-1,6-DIENE #



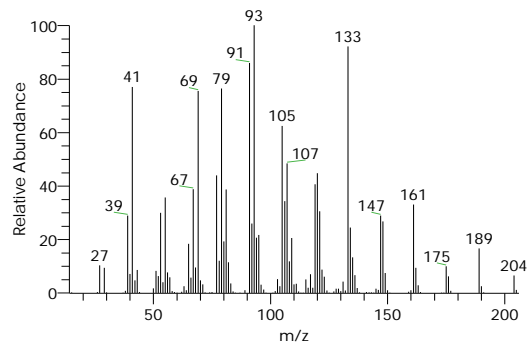
Caryophyllene

Formula C15H24, MW 204, CAS# 87-44-5, Entry# 66572

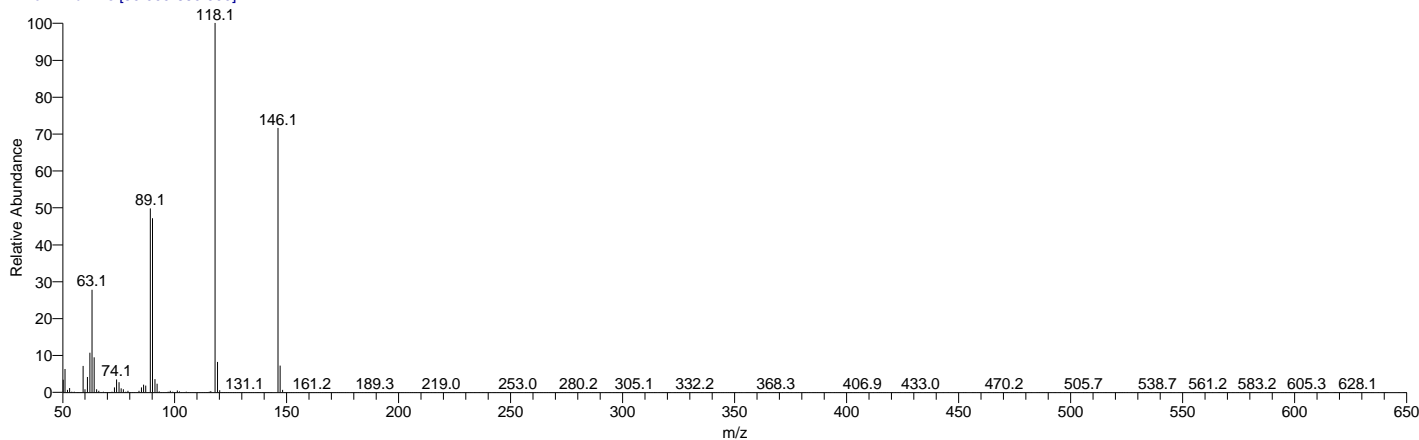
Bicyclo[7.2.0]undec-4-ene, 4,11,11-trimethyl-8-methylene-, [1R-(1R\*,4E,9S\*)]-



SI 932, RSI 934, mainlib, Entry# 66572, CAS# 87-44-5, Caryophyllene



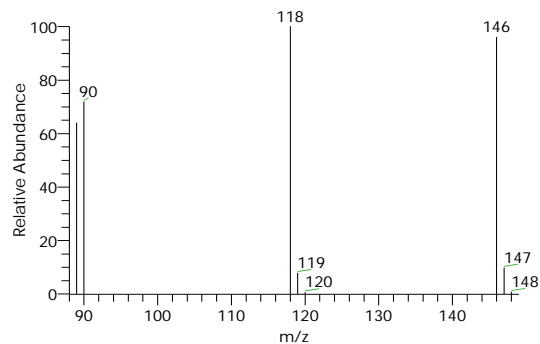
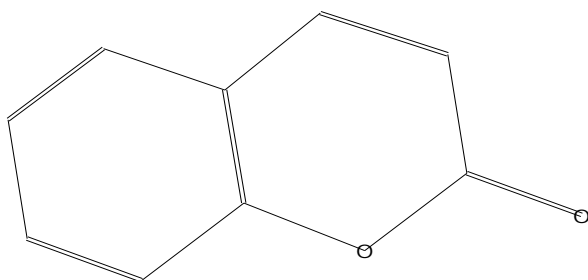
gerfa\_EA #4175 RT: 18.00 AV: 1 NL: 1.51E8  
T: + c EI Full ms [50.000-650.000]



# My GC-MS Report

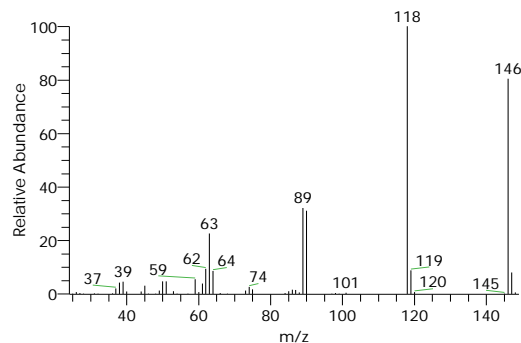
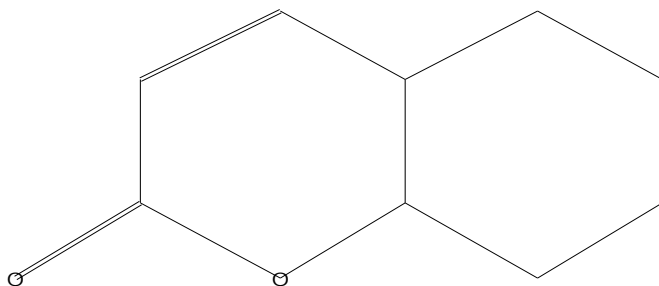
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
18.00	2H-1-BENZOPYRAN-2-ONE	9.92	C <sub>9</sub> H <sub>6</sub> O <sub>2</sub>	146	91-64-5	972	WileyRegi
18.00	Coumarin	9.92	C <sub>9</sub> H <sub>6</sub> O <sub>2</sub>	146	91-64-5	972	stry8e
18.00	Coumarin	9.92	C <sub>9</sub> H <sub>6</sub> O <sub>2</sub>	146	91-64-5	946	replib
Compound Structure							Hit Spectrum

2H-1-BENZOPYRAN-2-ONE  
Formula C<sub>9</sub>H<sub>6</sub>O<sub>2</sub>, MW 146, CAS# 91-64-5, Entry# 30896  
CHROMEN-2-ONE



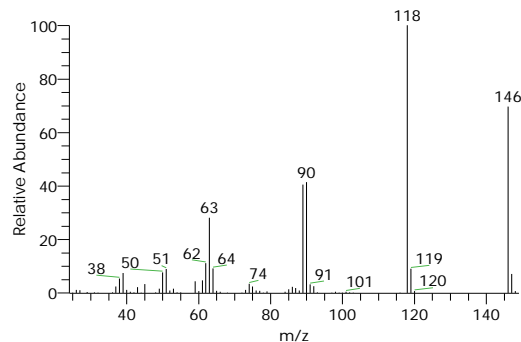
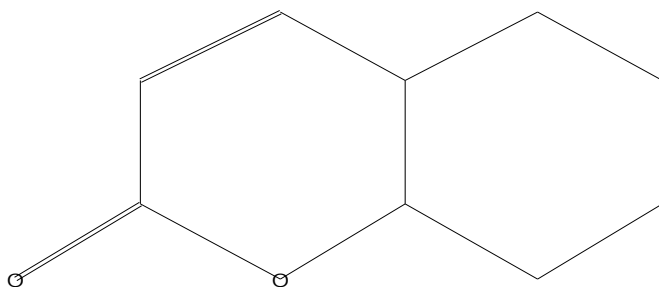
SI 960, RSI 972, replib, Entry# 18483, CAS# 91-64-5, Coumarin

Coumarin  
Formula C<sub>9</sub>H<sub>6</sub>O<sub>2</sub>, MW 146, CAS# 91-64-5, Entry# 18483  
2H-1-Benzopyran-2-one

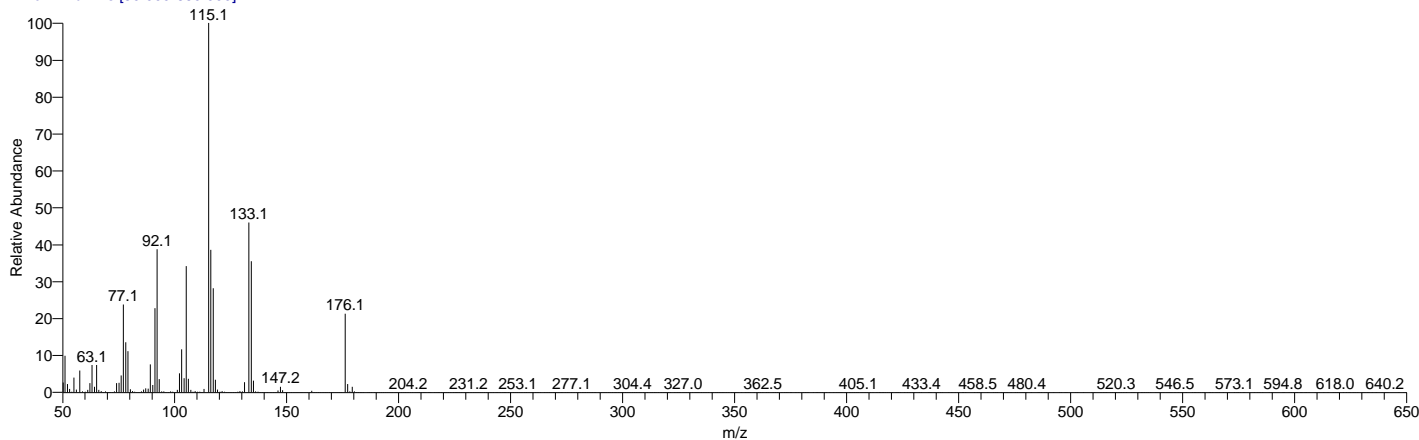


SI 946, RSI 946, replib, Entry# 18488, CAS# 91-64-5, Coumarin

Coumarin  
Formula C<sub>9</sub>H<sub>6</sub>O<sub>2</sub>, MW 146, CAS# 91-64-5, Entry# 18488  
2H-1-Benzopyran-2-one



gerfa\_EA #4223 RT: 18.16 AV: 1 NL: 1.10E8  
T: + c EI Full ms [50.000-650.000]





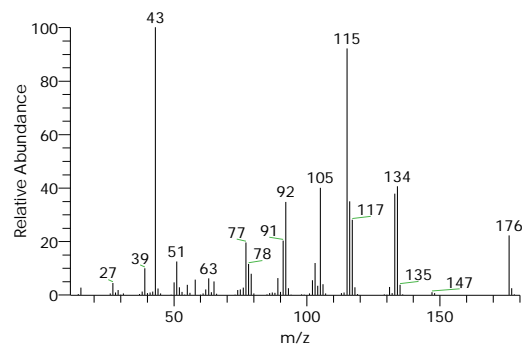
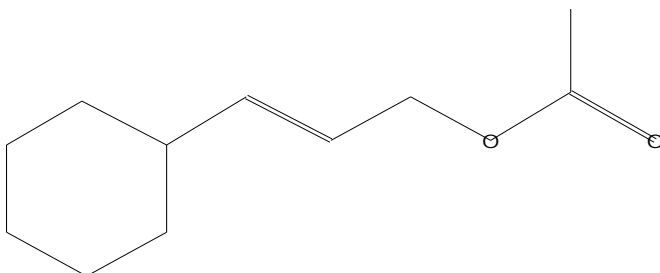
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
18.16	Acetic acid, cinnamyl ester	4.83	C11H12O2	176	103-54-8	955	mainlib
18.16	2-PROPEN-1-OL, 3-PHENYL-, ACETATE	4.83	C11H12O2	176	103-54-8	943	WileyRegi stry8e
18.16	CINNAMYL ALCOHOL, ACETATE, (E)-	4.83	C11H12O2	176	21040-45-9	968	WileyRegi stry8e

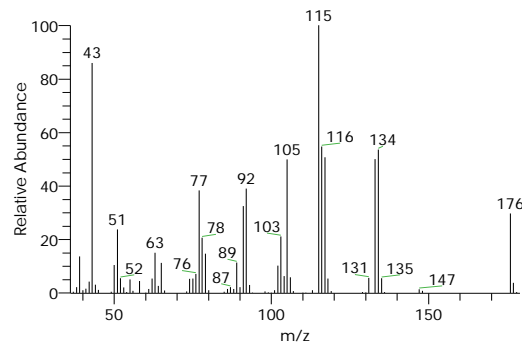
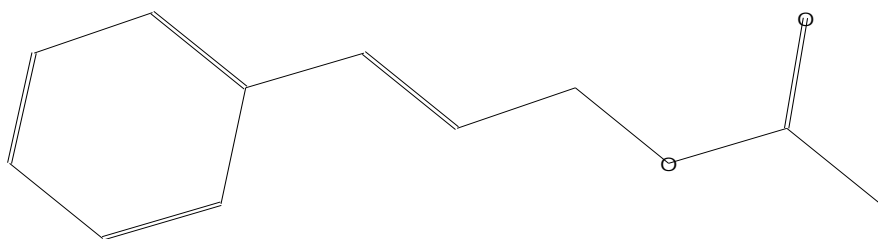
## Compound Structure

## Hit Spectrum

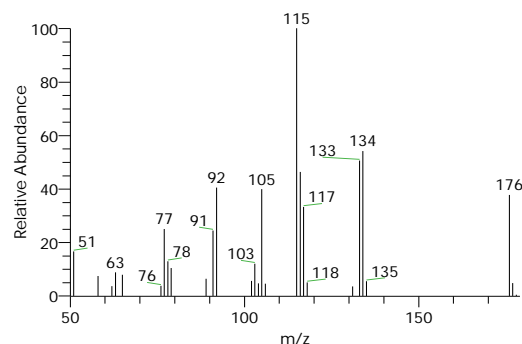
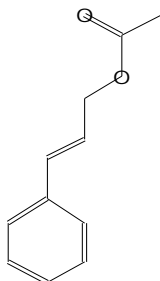
Acetic acid, cinnamyl ester  
Formula C11H12O2, MW 176, CAS# 103-54-8, Entry# 11373  
2-Propen-1-ol, 3-phenyl-, acetate



2-PROPEN-1-OL, 3-PHENYL-, ACETATE  
Formula C11H12O2, MW 176, CAS# 103-54-8, Entry# 58949  
ACETIC ACID CINNAMYL ESTER

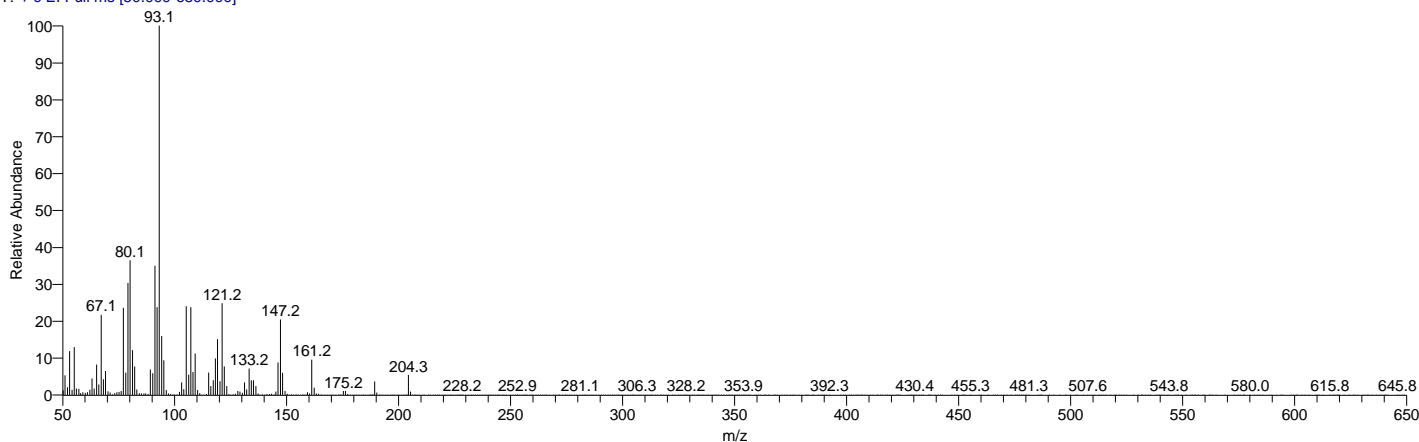


CINNAMYL ALCOHOL, ACETATE, (E)-  
Formula C11H12O2, MW 176, CAS# 21040-45-9, Entry# 59025  
(2E)-3-PHENYL-2-PROPENYL ACETATE #



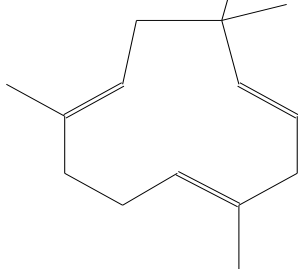
# My GC-MS Report

gerfa\_EA #4254 RT: 18.26 AV: 1 NL: 5.10E6  
T: + c EI Full ms [50.000-650.000]

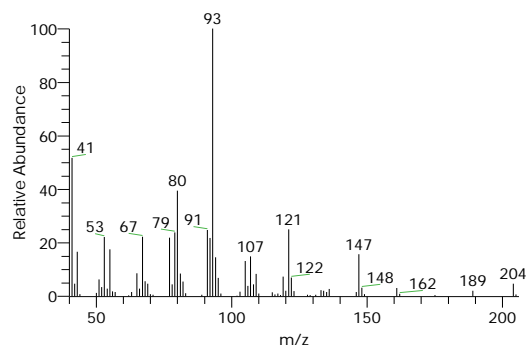


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
18.26	Humulene	0.11	C15H24	204	6753-9	907	replib
18.26	Humulene	0.11	C15H24	204	6753-9	895	replib
18.26	1,4,8-CYCLOUNDECATRIENE, 2,6,6,9-TETRAMETHYL-, (E,E,E)-	0.11	C15H24	204	6753-9	871	WileyRegistry8e
Compound Structure							Hit Spectrum

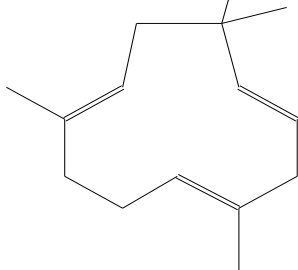
Humulene  
Formula C15H24, MW 204, CAS# 6753-98-6, Entry# 14003  
à-Caryophyllene



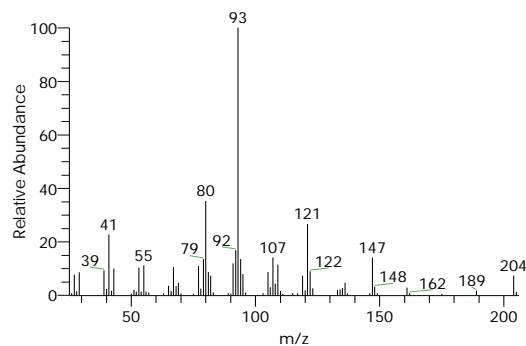
SI 887, RSI 907, replib, Entry# 14003, CAS# 6753-98-6, Humulene



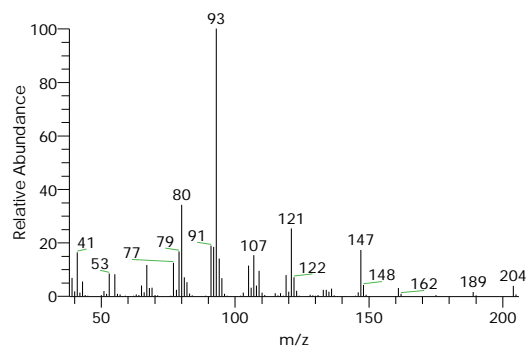
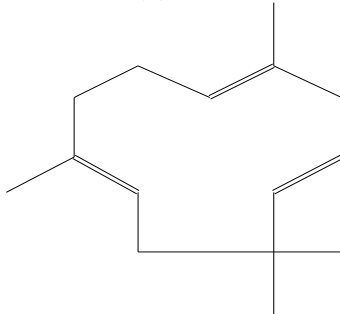
Humulene  
Formula C15H24, MW 204, CAS# 6753-98-6, Entry# 14104  
à-Caryophyllene



SI 861, RSI 895, replib, Entry# 14104, CAS# 6753-98-6, Humulene

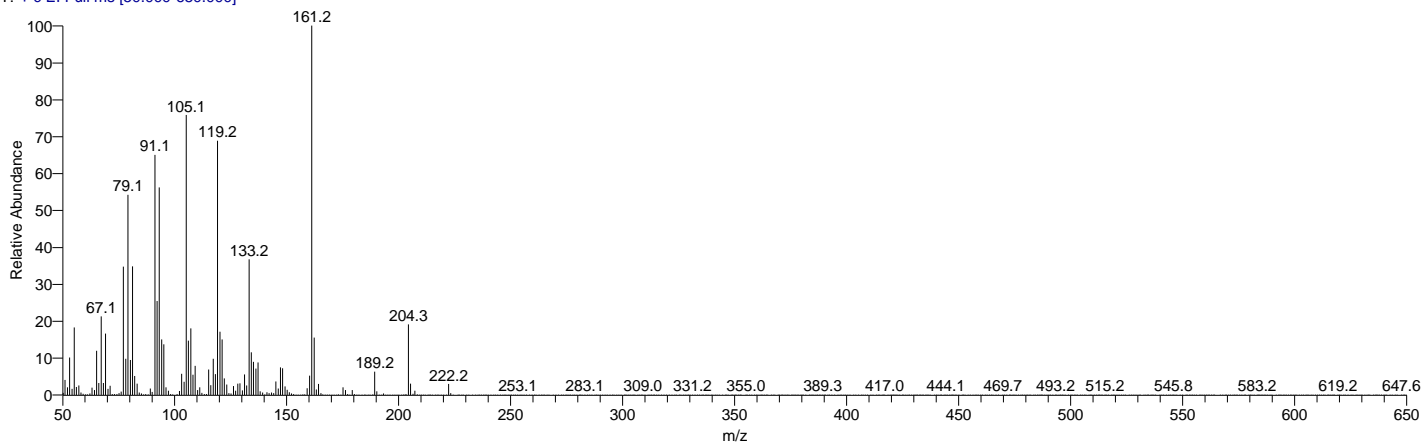


1,4,8-CYCLOUNDECATRIENE, 2,6,6,9-TETRAMETHYL-, (E,E,E)-  
Formula C15H24, MW 204, CAS# 6753-98-6, Entry# 89108  
2,6,6,9-TETRAMETHYL-1,4,8-CYCLOUNDECATRIENE #



# My GC-MS Report

gerfa\_EA #4431 RT: 18.86 AV: 1 NL: 1.84E7  
T: + c EI Full ms [50.000-650.000]



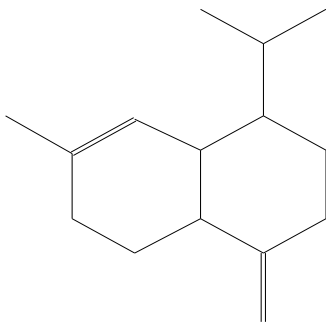
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
18.86	$\zeta$ -Muurolene	1.05	C15H24	204	30021-74-0	938	replib
18.86	Naphthalene, 1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-(1-methylethyl)-	1.05	C15H24	204	483-75-0	954	replib
18.86	$\zeta$ -Muurolene	1.05	C15H24	204	30021-74-0	929	replib

## Compound Structure

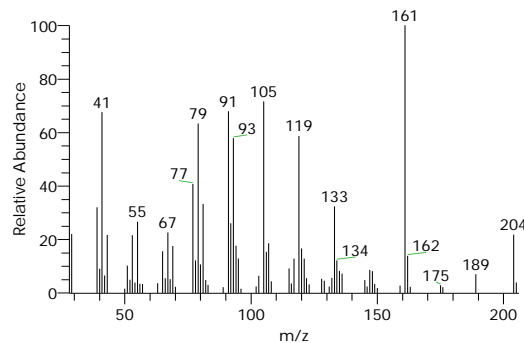
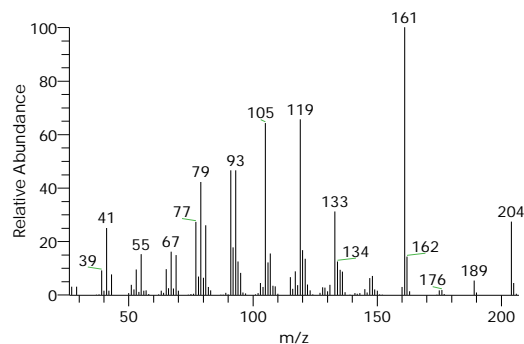
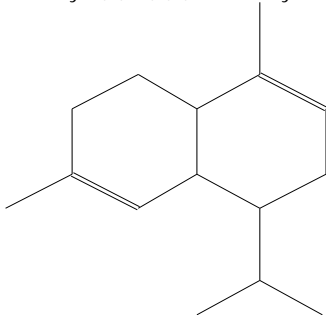
## Hit Spectrum

$\zeta$ -Muurolene

Formula C15H24, MW 204, CAS# 30021-74-0, Entry# 24906



Naphthalene, 1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-(1-methylethyl)-  
Formula C15H24, MW 204, CAS# 483-75-0, Entry# 24889  
1-Isopropyl-4,7-dimethyl-1,2,4a,5,6,8a-hexahydronaphthalene #



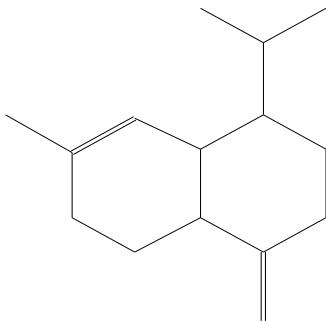
# My GC-MS Report

## Compound Structure

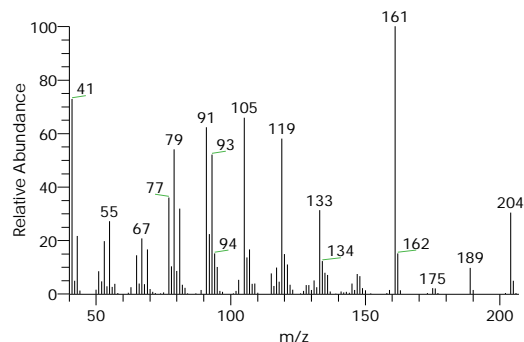
## Hit Spectrum

ç-Muurolene

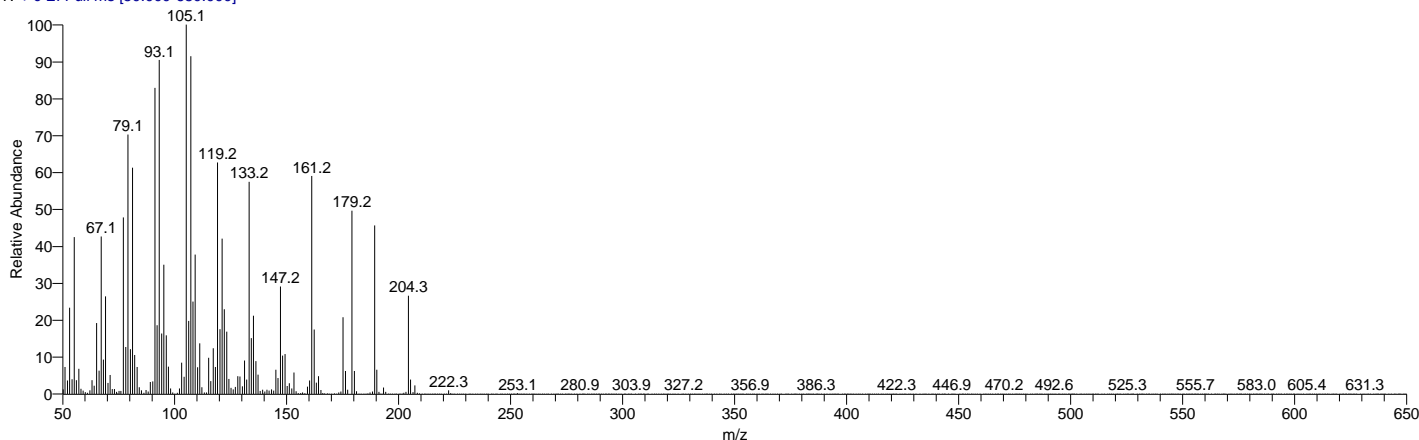
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 30021-74-0, Entry# 24843



SI 909, RSI 929, replib, Entry# 24843, CAS# 30021-74-0, ç-Muurolene



gerfa\_EA #4564 RT: 19.30 AV: 1 NL: 3.51E6  
T: + c EI Full ms [50.000-650.000]

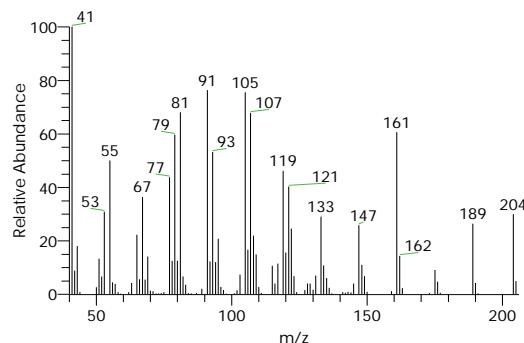
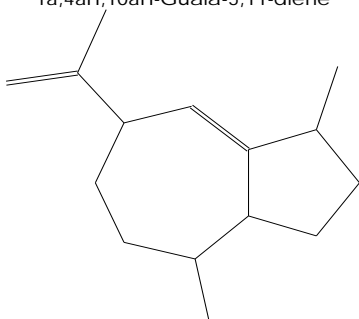


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
19.30	Azulene, 1,2,3,3a,4,5,6,7-octahydro-1,4-dimethyl-7-(1-methylethenyl)-, [1R-(1à,3aà,4à,7á)]-	0.26	C <sub>15</sub> H <sub>24</sub>	204	22567-17-5	918	replib
19.30	Aromandendrene	0.26	C <sub>15</sub> H <sub>24</sub>	204	489-39-4	909	replib
19.30	Naphthalene, 1,2,3,5,6,7,8,8a-octahydro-1,8a-dimethyl-7-(1-methylethenyl)-, [1S-(1à,7à,8aà)]-	0.26	C <sub>15</sub> H <sub>24</sub>	204	10219-75-7	898	mainlib

## Compound Structure

## Hit Spectrum

Azulene, 1,2,3,3a,4,5,6,7-octahydro-1,4-dimethyl-7-(1-methylethenyl)-, [1R-(1à,3aà,4à,7á)]-  
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 22567-17-5, Entry# 1321  
1á,4áH,10áH-Guaia-5,11-diene

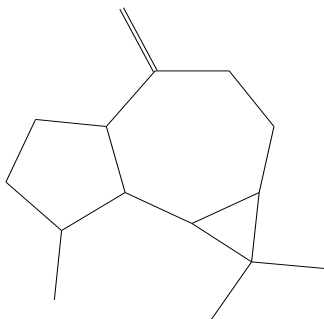


# My GC-MS Report

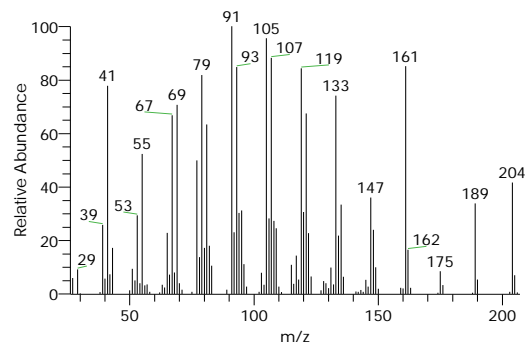
## Compound Structure

## Hit Spectrum

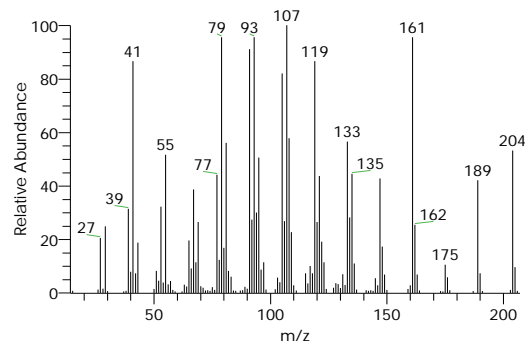
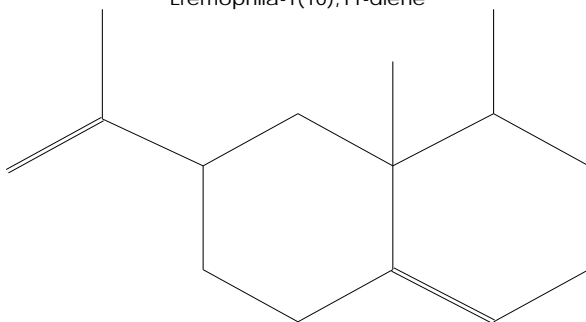
Aromandendrene  
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 489-39-4, Entry# 13310



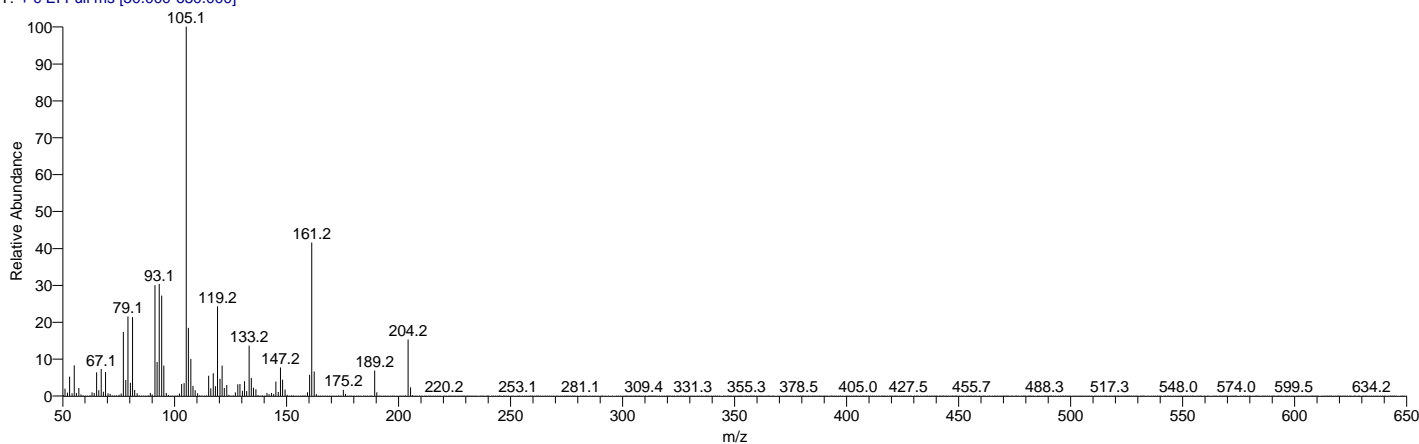
SI 860, RSI 909, replib, Entry# 13310, CAS# 489-39-4, Aromandendrene



Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 10219-75-7, Entry# 83645  
Eremophila-1(10),11-diene



gerfa\_EA #4615 RT: 19.47 AV: 1 NL: 9.72E7  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
19.47	à-Muurolene	4.14	C <sub>15</sub> H <sub>24</sub>	204	31983-2 2-9	960	replib
19.47	à-Muurolene	4.14	C <sub>15</sub> H <sub>24</sub>	204	10208-8 0-7	945	replib
19.47	NAPHTHALENE, 1,2,4A,5,6,8A-HEXAHYDRO-4,7-DIMETHYL -1-(1-METHYLETHYL)-, (1à,4Aà,8Aà)-	4.14	C <sub>15</sub> H <sub>24</sub>	204	31983-2 2-9	926	WileyRegi stry8e

# My GC-MS Report

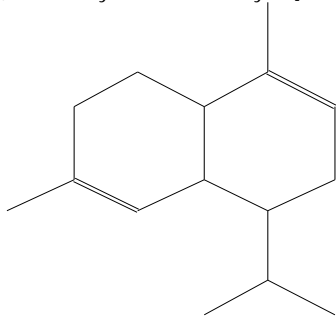
Compound Structure

Hit Spectrum

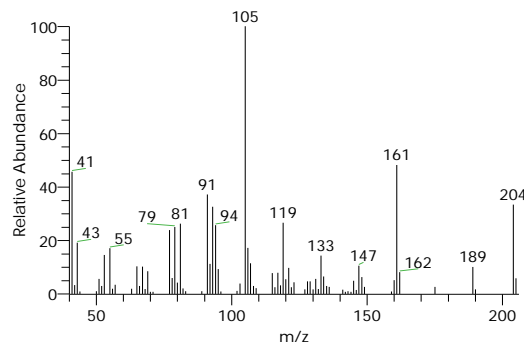
à-Muurolene

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 31983-22-9, Entry# 16421

[1à,4aà,8aà]-1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-[1-methylethyl]naphthalene



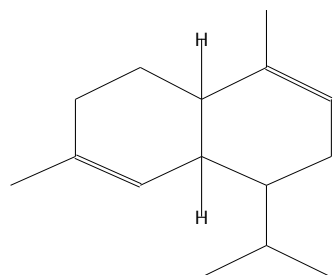
SI 952, RSI 960, replib, Entry# 16421, CAS# 31983-22-9, à-Muurolene



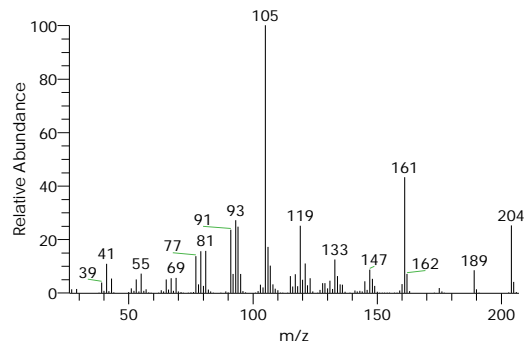
à-Muurolene

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 10208-80-7, Entry# 16423

Naphthalene, 1,2,4a,5,6,8a-hexahydro-4,7-dimethyl-1-(1-methylethyl)-, (1S,4aS,8aR)-



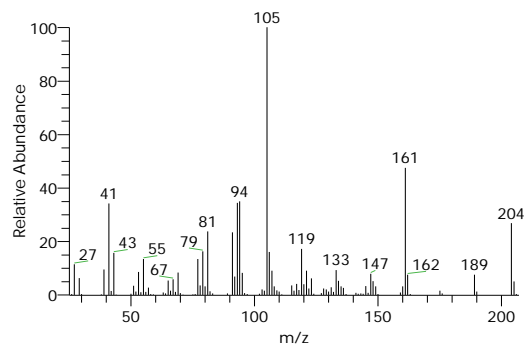
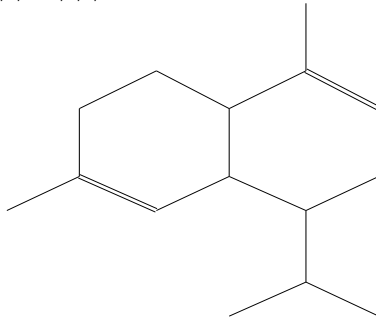
SI 945, RSI 945, replib, Entry# 16423, CAS# 10208-80-7, à-Muurolene



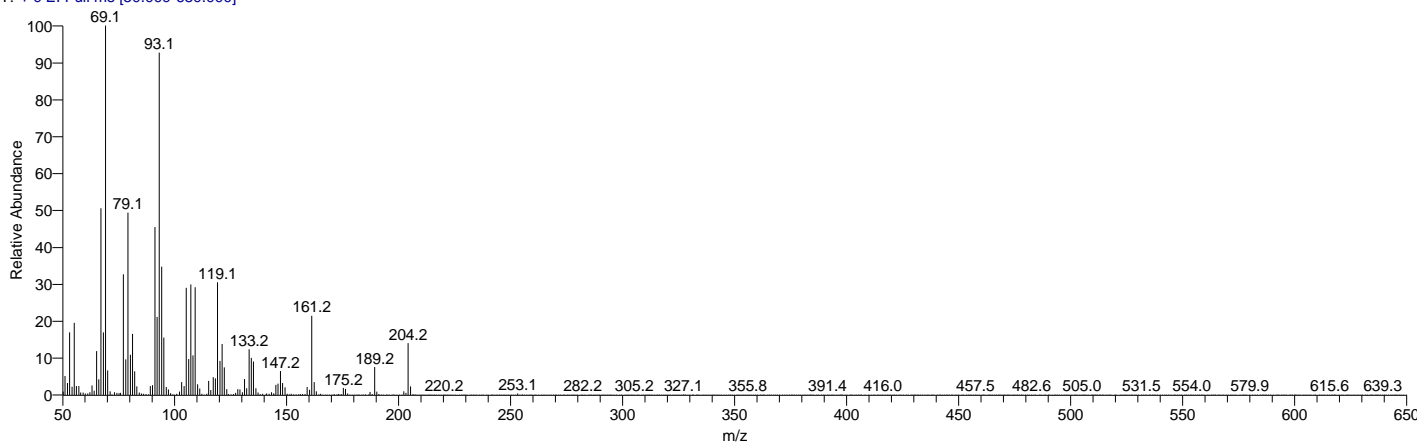
NAPHTHALENE, 1,2,4A,5,6,8A-HEXAHYDRO-4,7-DIMETHYL-1-(1-METHYLETHYL)-, (1à,4Aà,8Aà)-

Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 31983-22-9, Entry# 89350

NAPHTHALENE, 1,2,4Aa,5,6,8Aa-HEXAHYDRO-1a-ISOPROPYL-4,7-DIMETHYL-



gerfa\_EA #4670 RT: 19.66 AV: 1 NL: 2.58E6  
T: + c EI Full ms [50.000-650.000]

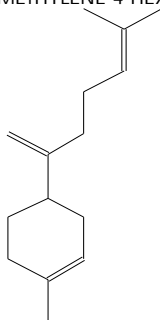


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
19.66	CYCLOHEXENE, 1-METHYL-4-(5-METHYL-1-METHYLENE-4-HEXENYL)-, (S)-	0.12	C <sub>15</sub> H <sub>24</sub>	204	495-61-4	933	WileyRegistry8e
19.66	à-Bisabolene	0.12	C <sub>15</sub> H <sub>24</sub>	204	495-61-4	896	replib

# My GC-MS Report

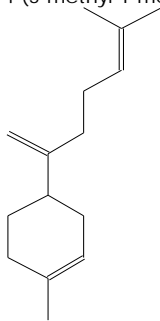
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
19.66	1-METHYL-4-(5-METHYL-1-METHYLENE-4-HEXENYL)-1-CYCLOHEXENE	0.12	C15H24	204	NA	896	WileyRegistry8e
Compound Structure				Hit Spectrum			

CYCLOHEXENE, 1-METHYL-4-(5-METHYL-1-METHYLENE-4-HEXENYL)-, (S)-  
Formula C15H24, MW 204, CAS# 495-61-4, Entry# 89082  
1-METHYL-4-(5-METHYL-1-METHYLENE-4-HEXENYL)-1-CYCLOHEXENE #

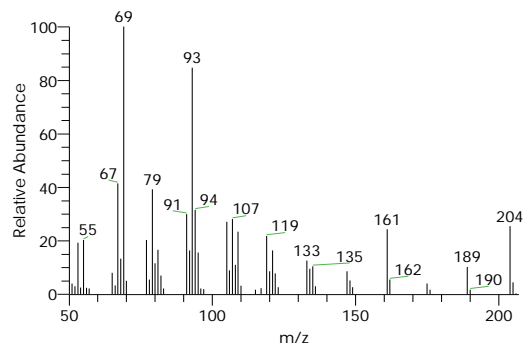
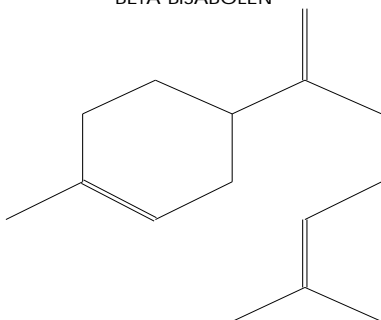


α-Bisabolene

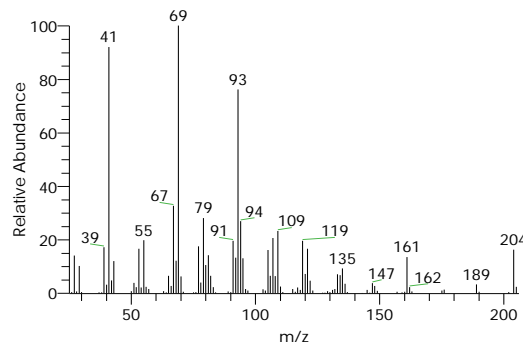
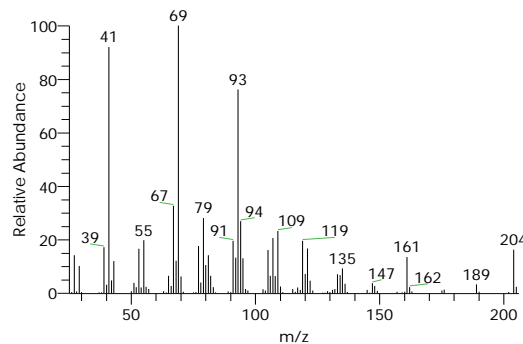
Formula C15H24, MW 204, CAS# 495-61-4, Entry# 8536  
Cyclohexene, 1-methyl-4-(5-methyl-1-methylene-4-hexenyl)-, (S)-



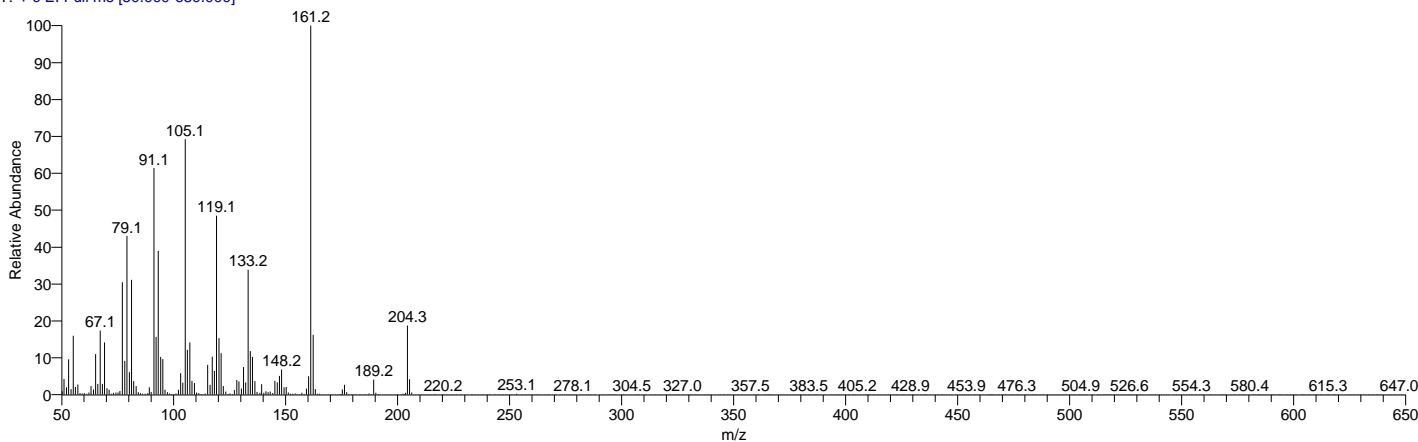
1-METHYL-4-(5-METHYL-1-METHYLENE-4-HEXENYL)-1-CYCLOHEXENE  
Formula C15H24, MW 204, CAS# NA, Entry# 388678  
BETA-BISABOLEN



SI 889, RSI 896, replib, Entry# 8536, CAS# 495-61-4, α-Bisabolene



gerfa\_EA #4704 RT: 19.77 AV: 1 NL: 3.33E6  
T: + c EI Full ms [50.000-650.000]



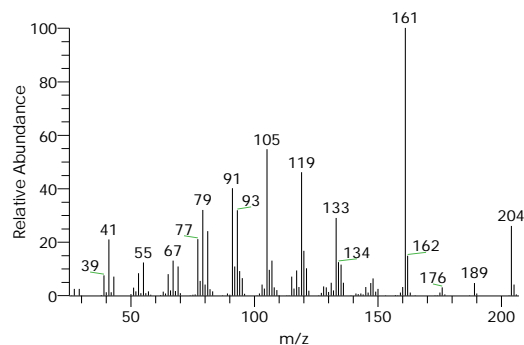
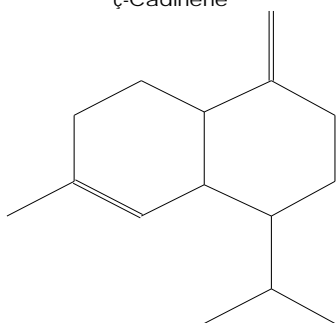
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
19.77	Naphthalene, 1,2,3,4,4a,5,6,8a-octahydro-7-methyl-4-methyle ne-1-(1-methylethyl)-, (1à,4aá,8aà)-	0.18	C15H24	204	39029-4 1-9	932	replib
19.77	Naphthalene, 1,2,3,4,4a,5,6,8a-octahydro-7-methyl-4-methyle ne-1-(1-methylethyl)-, (1à,4aá,8aà)-	0.18	C15H24	204	39029-4 1-9	932	replib
19.77	ç-Muurolene	0.18	C15H24	204	30021-7 4-0	929	replib

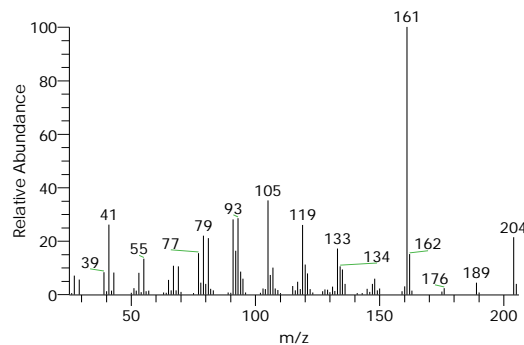
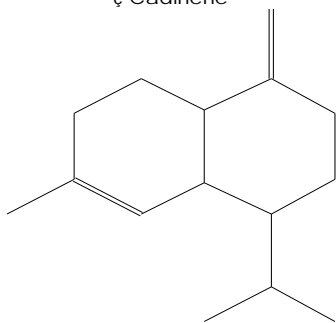
Compound Structure

Hit Spectrum

Formula C15H24, MW 204, CAS# 39029-41-9, Entry# 24895  
ç-Cadinene

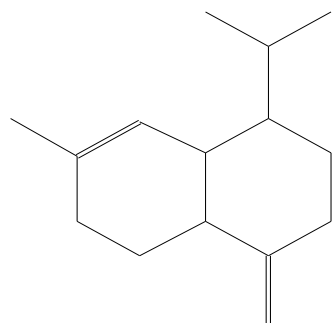


Formula C15H24, MW 204, CAS# 39029-41-9, Entry# 24890  
ç-Cadinene

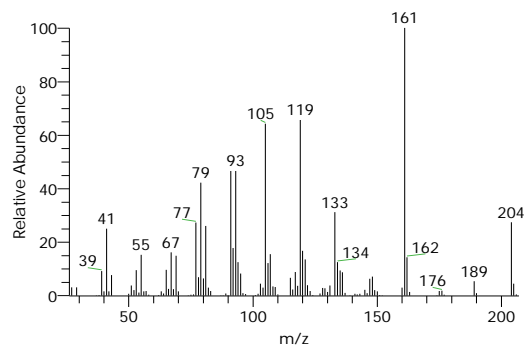


ç-Muurolene

Formula C15H24, MW 204, CAS# 30021-74-0, Entry# 24906



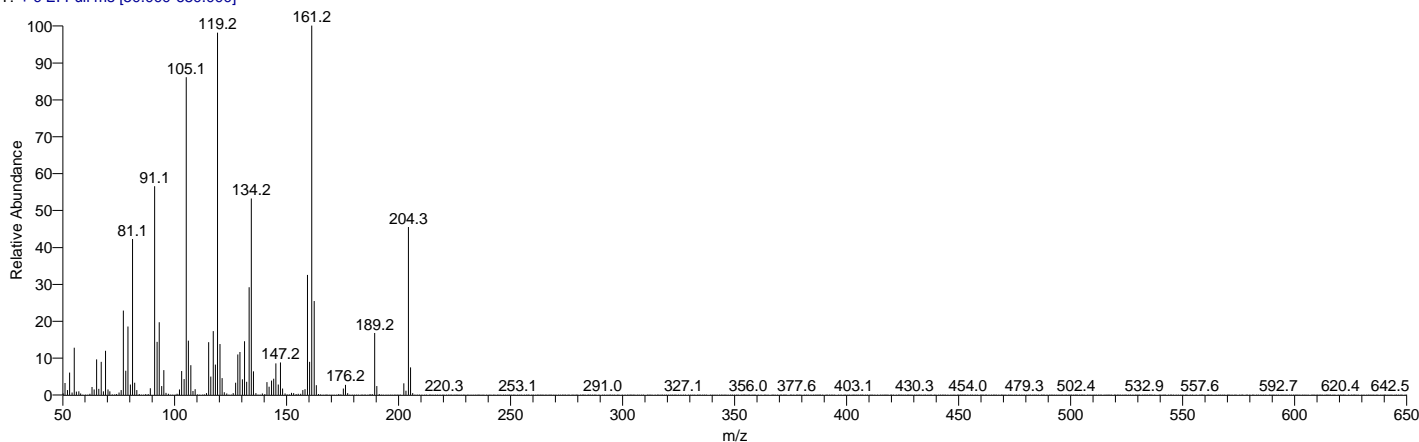
SI 921, RSI 929, replib, Entry# 24906, CAS# 30021-74-0, ç-Muurolene





# My GC-MS Report

gerfa\_EA #4785 RT: 20.04 AV: 1 NL: 6.19E7  
T: + c EI Full ms [50.000-650.000]

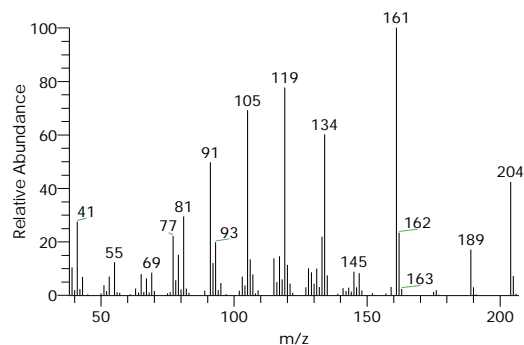
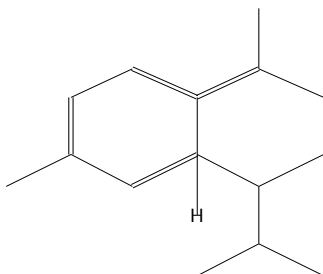


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
20.04	NAPHTHALENE, 1,2,3,5,6,8A-HEXAHYDRO-4,7-DIMETHYL-1-(1-METHYLETHYL)-, (1S-CIS)-	4.78	C15H24	204	483-76-1	942	WileyRegistry8e
20.04	NAPHTHALENE, 1,2,3,5,6,8A-HEXAHYDRO-4,7-DIMETHYL-1-(1-METHYLETHYL)-, (1S-CIS)-	4.78	C15H24	204	483-76-1	900	WileyRegistry8e
20.04	1-Isopropyl-4,7-dimethyl-1,2,3,5,6,8a-hexahydronaphthalene	4.78	C15H24	204	16729-01-4	921	mainlib

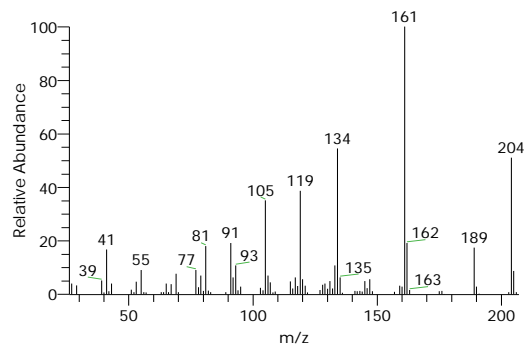
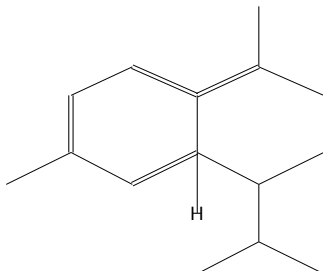
## Compound Structure

## Hit Spectrum

NAPHTHALENE, 1,2,3,5,6,8A-HEXAHYDRO-4,7-DIMETHYL-1-(1-METHYLETHYL)-, (1S-CIS)-  
Formula C15H24, MW 204, CAS# 483-76-1, Entry# 89277  
1-ISOPROPYL-4,7-DIMETHYL-1,2,3,5,6,8A-HEXAHYDRONAPHTHALENE #



NAPHTHALENE, 1,2,3,5,6,8A-HEXAHYDRO-4,7-DIMETHYL-1-(1-METHYLETHYL)-, (1S-CIS)-  
Formula C15H24, MW 204, CAS# 483-76-1, Entry# 89273  
1-ISOPROPYL-4,7-DIMETHYL-1,2,3,5,6,8A-HEXAHYDRONAPHTHALENE #

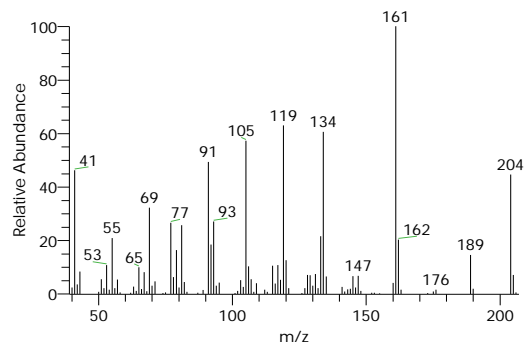
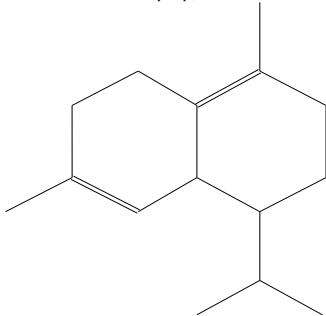


# My GC-MS Report

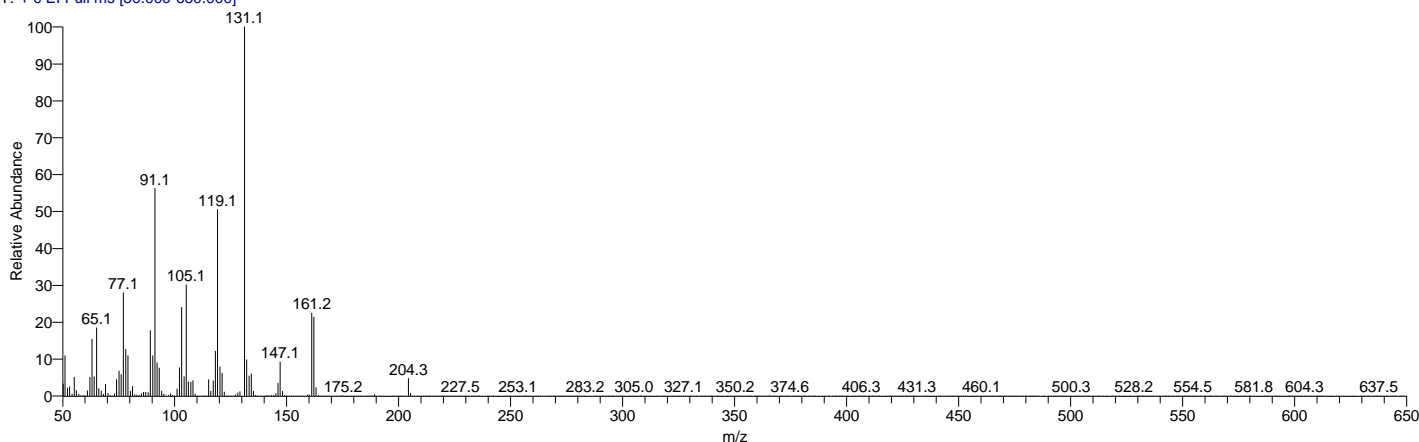
Compound Structure

Hit Spectrum

1-Isopropyl-4,7-dimethyl-1,2,3,5,6,8a-hexahydronaphthalene  
Formula C<sub>15</sub>H<sub>24</sub>, MW 204, CAS# 16729-01-4, Entry# 150583  
Cadina-1(10),4-diene



gerfa\_EA #4838 RT: 20.22 AV: 1 NL: 9.18E7  
T: + c EI Full ms [50.000-650.000]

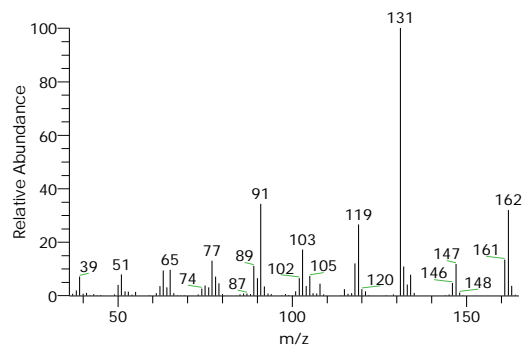
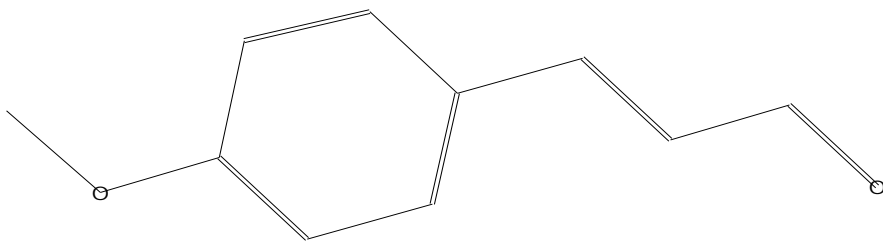


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
20.22	2-PROPENAL, 3-(4-METHOXYPHENYL)-	4.91	C <sub>10</sub> H <sub>10</sub> O <sub>2</sub>	162	1963-3	883	WileyRegi
20.22	(2E)-3-(2-METHOXYPHENYL)-2-PROPENAL #	4.91	C <sub>10</sub> H <sub>10</sub> O <sub>2</sub>	162	1504-7	891	WileyRegi
20.22	2-Propenal, 3-(2-methoxyphenyl)-	4.91	C <sub>10</sub> H <sub>10</sub> O <sub>2</sub>	162	1504-7	876	mainlib

Compound Structure

Hit Spectrum

2-PROPENAL, 3-(4-METHOXYPHENYL)-  
Formula C<sub>10</sub>H<sub>10</sub>O<sub>2</sub>, MW 162, CAS# 1963-36-6, Entry# 45236  
CINNAMALDEHYDE, P-METHOXY-

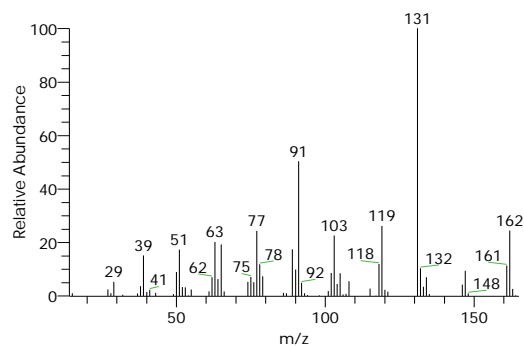
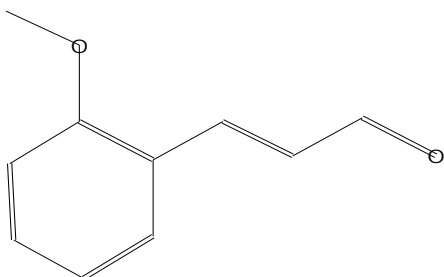


# My GC-MS Report

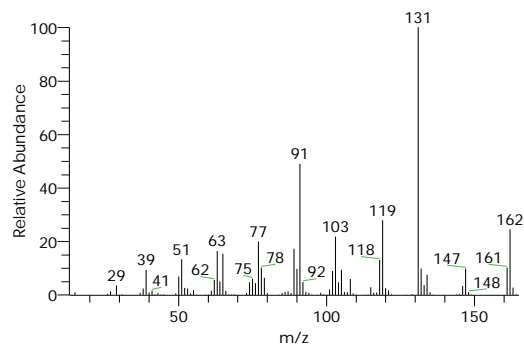
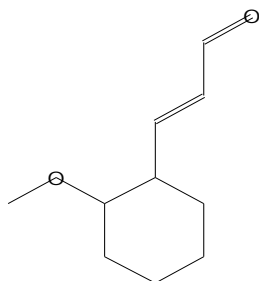
Compound Structure

Hit Spectrum

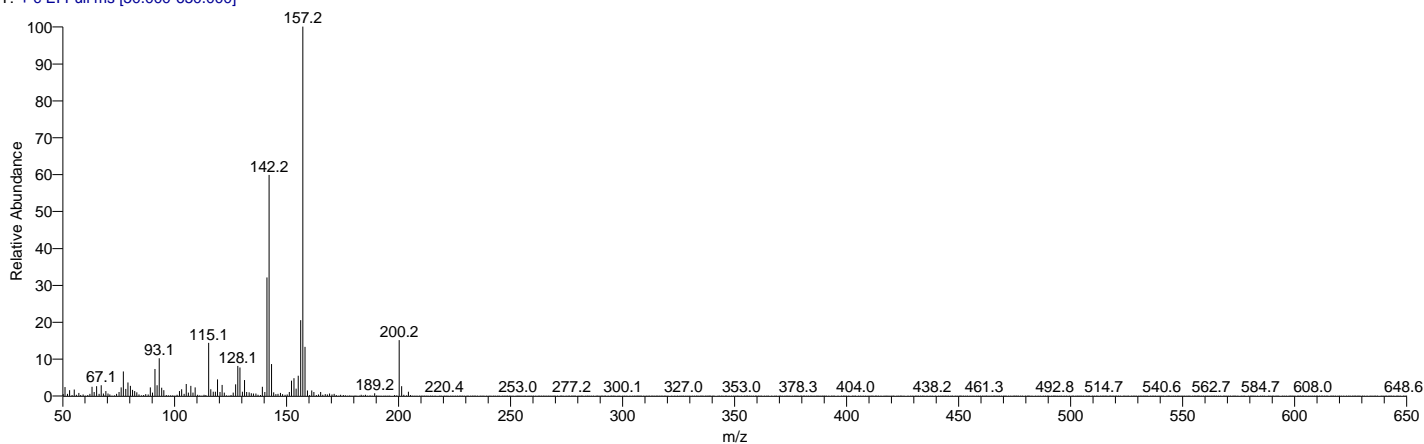
(2E)-3-(2-METHOXYPHENYL)-2-PROPENAL #  
Formula C<sub>10</sub>H<sub>10</sub>O<sub>2</sub>, MW 162, CAS# 1504-74-1, Entry# 45156  
(2E)-3-(2-METHOXYPHENYL)-2-PROPENAL



2-Propenal, 3-(2-methoxyphenyl)-  
Formula C<sub>10</sub>H<sub>10</sub>O<sub>2</sub>, MW 162, CAS# 1504-74-1, Entry# 115637  
Cinnamaldehyde, o-methoxy-



gerfa\_EA #4912 RT: 20.47 AV: 1 NL: 1.55E7  
T: + c EI Full ms [50.000-650.000]



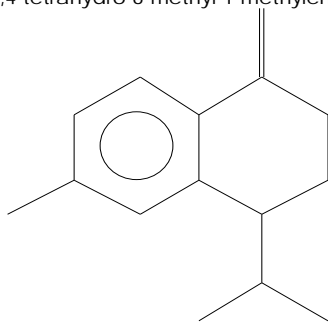
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
20.47	4-Isopropyl-6-methyl-1-methylene-1,2,3,4-tetrahydronaphthalene	0.43	C <sub>15</sub> H <sub>20</sub>	200	50277-34-4	934	mainlib
20.47	1,1,6-TRIMETHYL-1,2-DIHYDRO NAPHTHALENE	0.43	C <sub>13</sub> H <sub>16</sub>	172	NA	956	WileyRegistry8e
20.47	à-Calacorene	0.43	C <sub>15</sub> H <sub>20</sub>	200	21391-99-1	940	mainlib

# My GC-MS Report

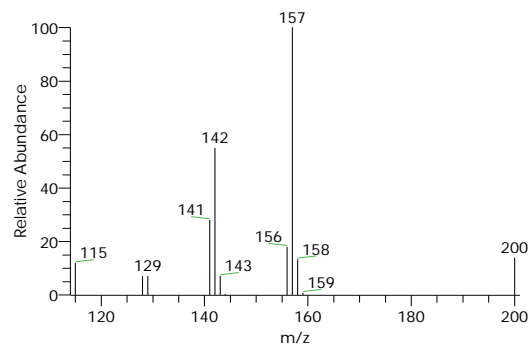
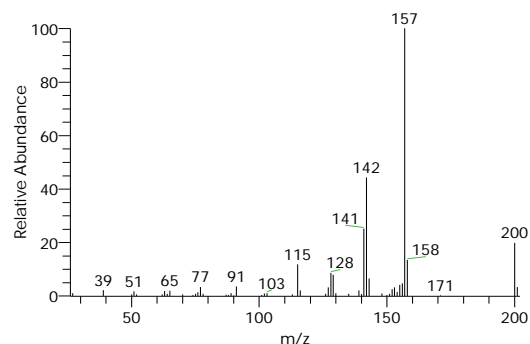
Compound Structure

Hit Spectrum

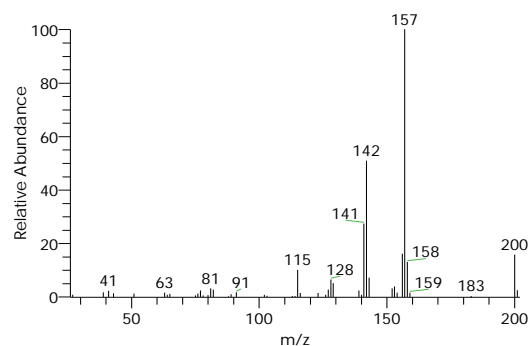
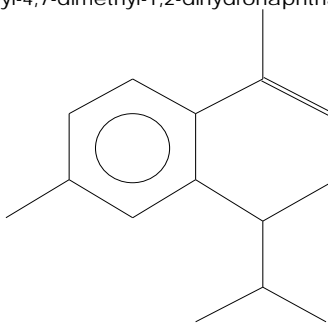
4-Isopropyl-6-methyl-1-methylene-1,2,3,4-tetrahydronaphthalene  
Formula C<sub>15</sub>H<sub>20</sub>, MW 200, CAS# 50277-34-4, Entry# 146807  
Naphthalene, 1,2,3,4-tetrahydro-6-methyl-1-methylene-4-(1-methylethyl)-



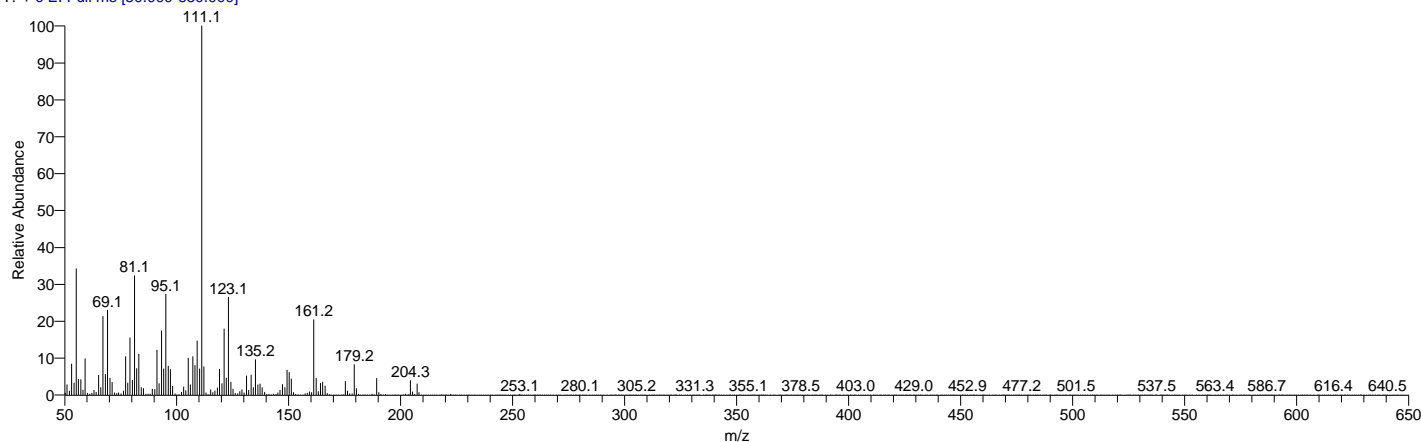
1,1,6-TRIMETHYL-1,2-DIHYDRO NAPHTHALENE  
Formula C<sub>13</sub>H<sub>16</sub>, MW 172, CAS# NA, Entry# 56092  
CALACORENE



à-Calacorene  
Formula C<sub>15</sub>H<sub>20</sub>, MW 200, CAS# 21391-99-1, Entry# 146806  
1-Isopropyl-4,7-dimethyl-1,2-dihydronaphthalene, (S)-



gerfa\_EA #5102 RT: 21.11 AV: 1 NL: 3.41E6  
T: + c EI Full ms [50.000-650.000]

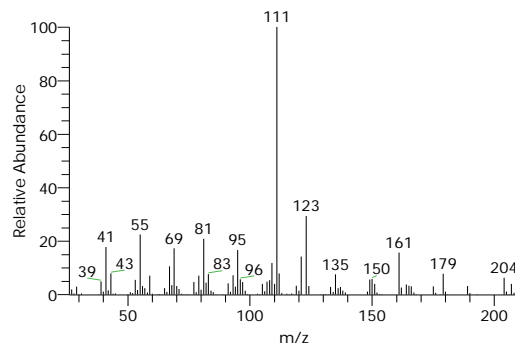
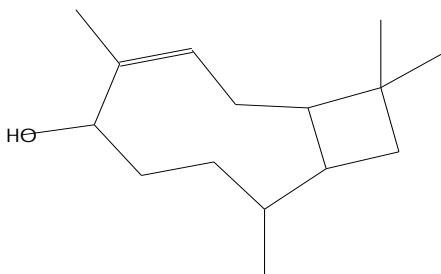


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.11	Caryophyllenyl alcohol	0.14	C <sub>15</sub> H <sub>26</sub> O	222	913176-41-7	901	mainlib
21.11	Caryophyllenyl alcohol	0.14	C <sub>15</sub> H <sub>26</sub> O	222	913176-41-7	856	replib

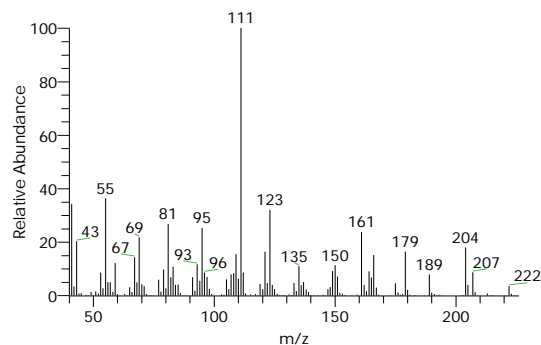
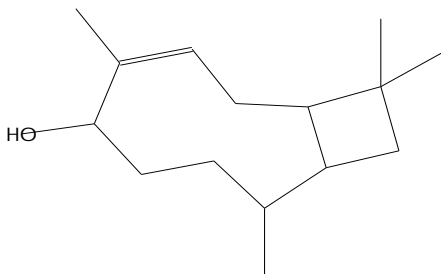
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.11	4aH-Cycloprop[e]azulen-4a-ol, decahydro-1,1,4,7-tetramethyl-, [1aR-(1aà,4á,4á,7à,7aá,7bà)]-	0.14	C15H26O	222	5986-49-2	804	replib
Compound Structure				Hit Spectrum			

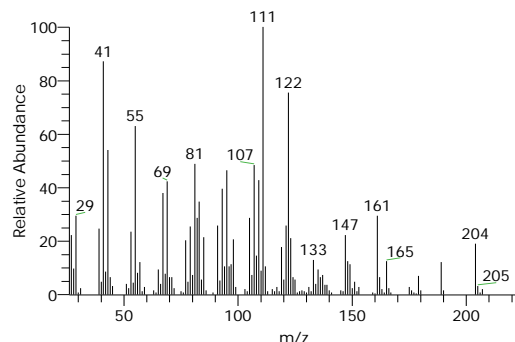
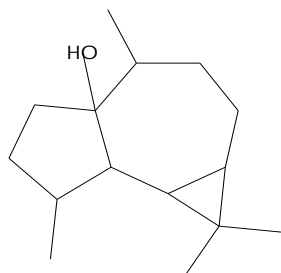
Caryophyllenyl alcohol  
Formula C15H26O, MW 222, CAS# 913176-41-7, Entry# 88498  
4,8,11,11-Tetramethylbicyclo[7.2.0]undec-3-en-5-ol #



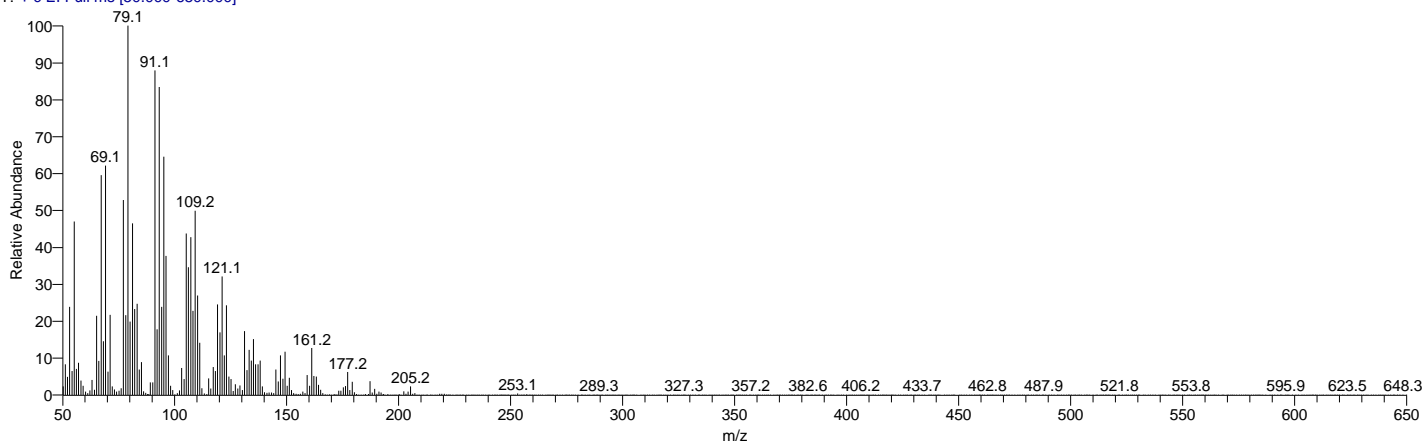
Caryophyllenyl alcohol  
Formula C15H26O, MW 222, CAS# 913176-41-7, Entry# 17579  
4,8,11,11-Tetramethylbicyclo[7.2.0]undec-3-en-5-ol #



Formula C15H26O, MW 222, CAS# 5986-49-2, Entry# 17563  
(1aR,4S,4aS,7R,7aS,7bR)-1,1,4,7-Tetramethyldecahydro-1H-cyclopropa[e]azulen-4a-ol



gerfa\_EA #5193 RT: 21.41 AV: 1 NL: 1.68E6  
T: + c EI Full ms [50.000-650.000]



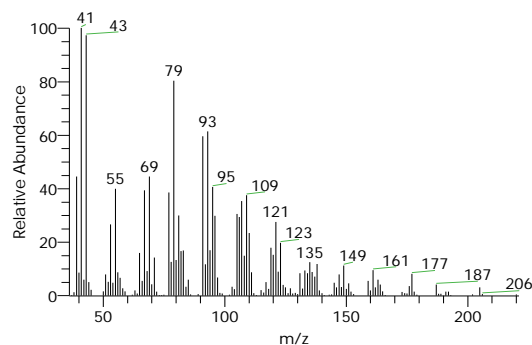
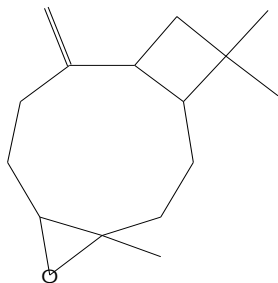
# My GC-MS Report

RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.41	(-)-5-OXATRICYCLO[8.2.0.0(4,6)]DODECANE,,12-TRIMETHYL-9-METHYLENE-, [1R-(1R*,4R*,6R*,10S*)]-	0.14	C15H24O	220	1139-30-6	921	WileyRegistry8e
21.41	(-)-5-OXATRICYCLO[8.2.0.0(4,6)]DODECANE,,12-TRIMETHYL-9-METHYLENE-, [1R-(1R*,4R*,6R*,10S*)]-	0.14	C15H24O	220	1139-30-6	925	WileyRegistry8e
21.41	Caryophyllene oxide	0.14	C15H24O	220	1139-30-6	913	mainlib

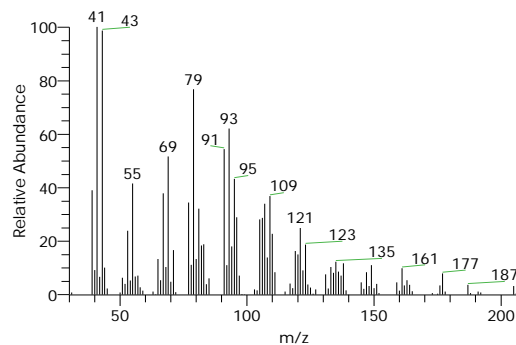
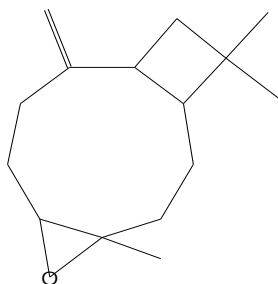
Compound Structure

Hit Spectrum

Formula C15H24O, MW 220, CAS# 1139-30-6, Entry# 107245  
5-OXATRICYCLO[8.2.0.0(4,6)]DODECANE, 4,12,12-TRIMETHYL-9-METHYLENE-, (1R,4R,6R,10S)-

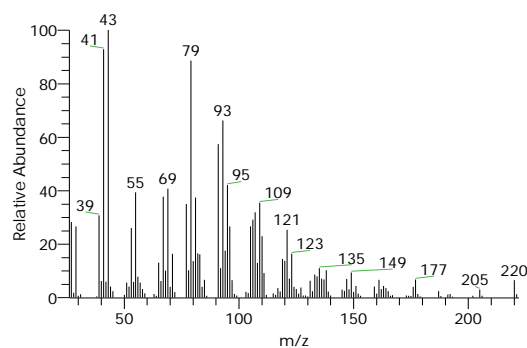
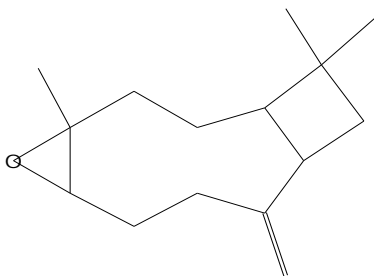


Formula C15H24O, MW 220, CAS# 1139-30-6, Entry# 107244  
5-OXATRICYCLO[8.2.0.0(4,6)]DODECANE, 4,12,12-TRIMETHYL-9-METHYLENE-, (1R,4R,6R,10S)-



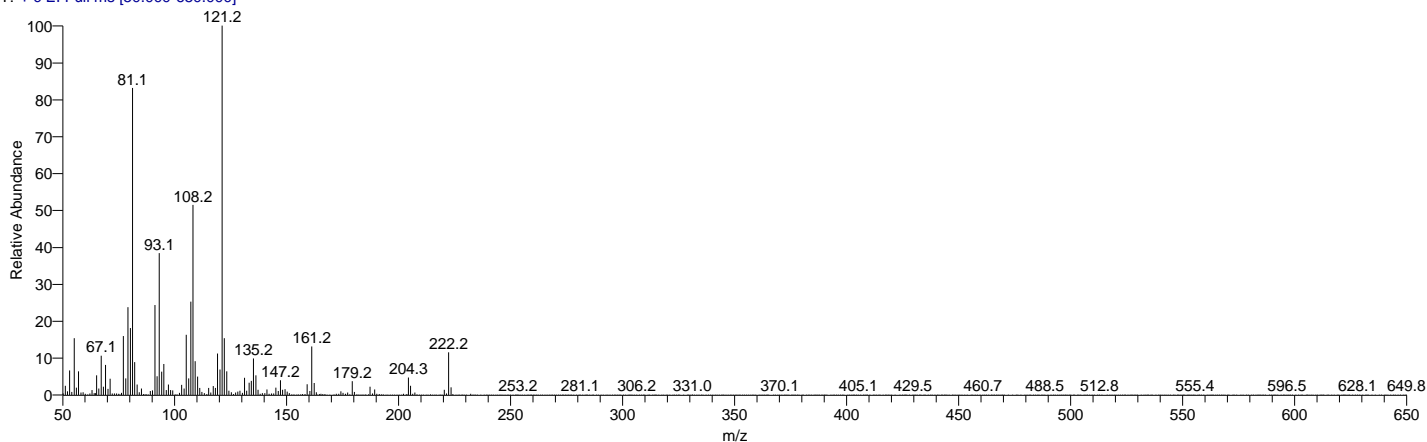
Caryophyllene oxide

Formula C15H24O, MW 220, CAS# 1139-30-6, Entry# 6247



# My GC-MS Report

gerfa\_EA #5213 RT: 21.48 AV: 1 NL: 4.55E6  
T: + c EI Full ms [50.000-650.000]

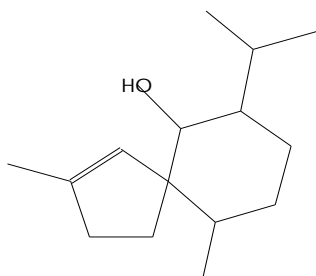


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.48	(5S,6R,7S,10R)-7-Isopropyl-2,10-dimethylspiro[4.5]dec-1-en-6-ol	0.16	C15H26O	222	72203-9	905	mainlib
21.48	gleenol	0.16	C15H26O	222	NA	877	mainlib
21.48	ç-Elemene	0.16	C15H24	204	29873-9	820	mainlib

## Compound Structure

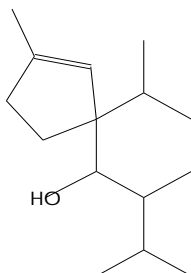
## Hit Spectrum

(5S,6R,7S,10R)-7-Isopropyl-2,10-dimethylspiro[4.5]dec-1-en-6-ol  
Formula C15H26O, MW 222, CAS# 72203-99-7, Entry# 102652  
Spiro[4.5]dec-1-en-6-ol, 2,10-dimethyl-7-(1-methylethyl)-, (5S,6R,7S,10R)-



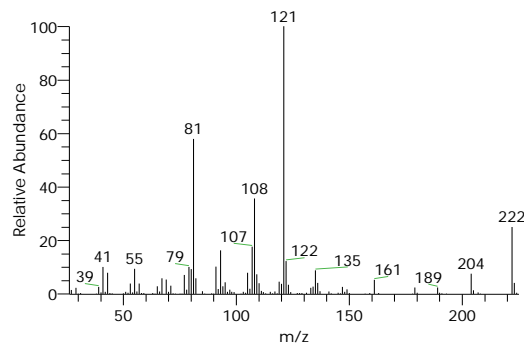
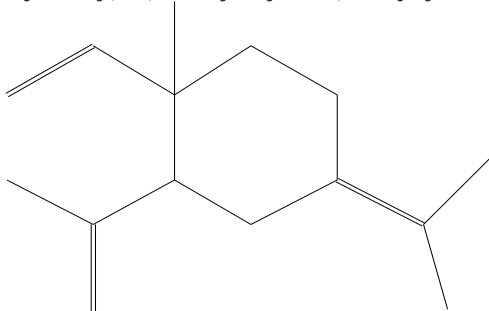
gleenol

Formula C15H26O, MW 222, CAS# NA, Entry# 102651  
\$:28MYLXGCVCCZCOHU-UHFFFAOYSA-N

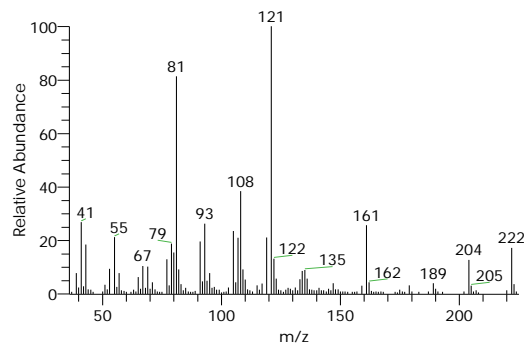


ç-Elemene

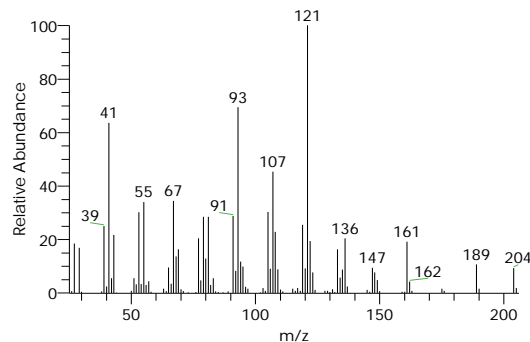
Formula C15H24, MW 204, CAS# 29873-99-2, Entry# 102784  
1-Methyl-2-(1-methylethenyl)-4-(1-methylethylidene)-1-vinylcyclohexane, (1R-trans)-



SI 864, RSI 877, mainlib, Entry# 102651, CAS# NA, gleenol

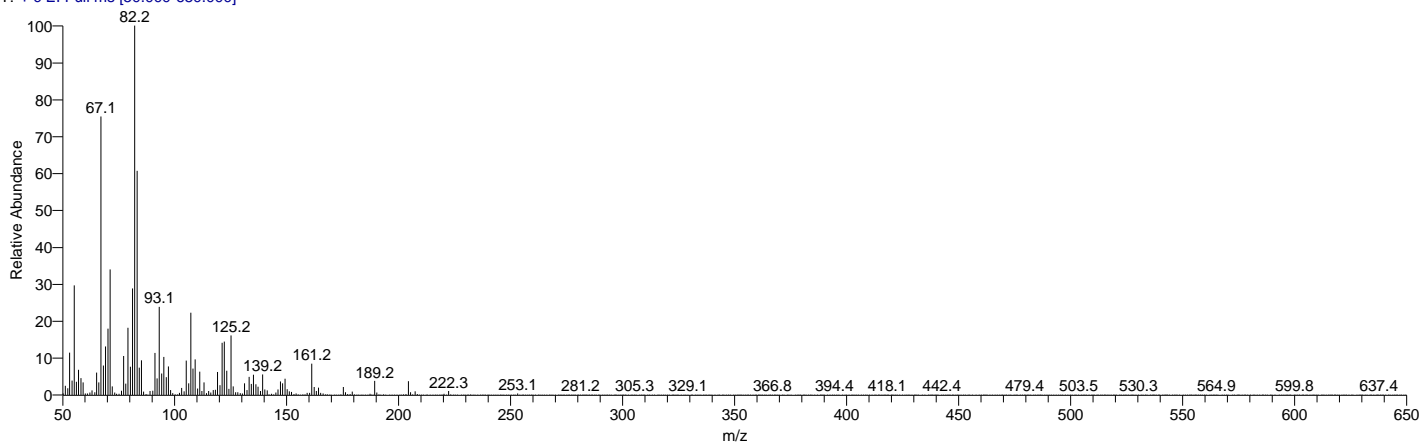


SI 775, RSI 820, mainlib, Entry# 102784, CAS# 29873-99-2, ç-Elemene



# My GC-MS Report

gerfa\_EA #5324 RT: 21.85 AV: 1 NL: 3.15E6  
T: + c EI Full ms [50.000-650.000]

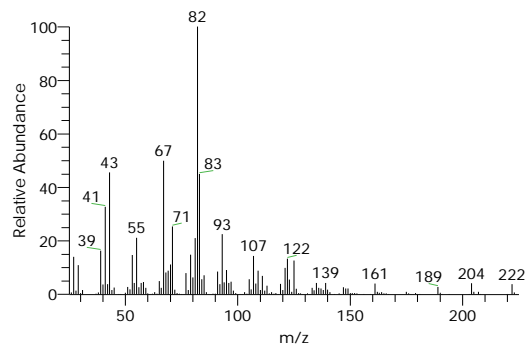
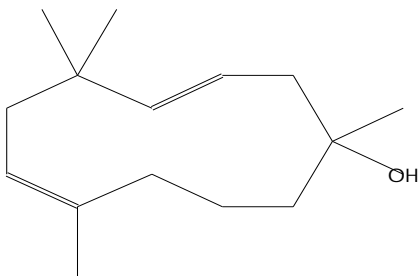


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.85	3,7-Cycloundecadien-1-ol, 1,5,5,8-tetramethyl-	0.11	C15H26O	222	118014-38-3	889	mainlib
21.85	1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL #	0.11	C15H26O	222	118014-38-3	889	WileyRegistry8e
21.85	1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL #	0.11	C15H26O	222	118014-38-3	889	WileyRegistry8e

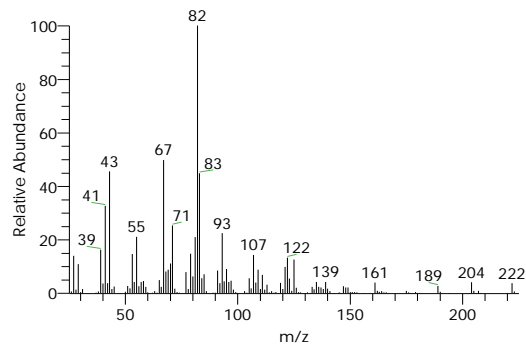
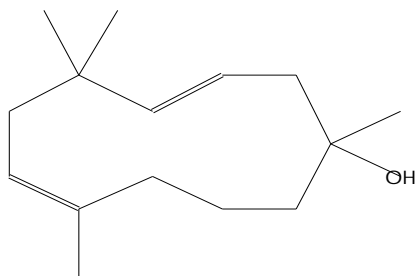
Compound Structure

Hit Spectrum

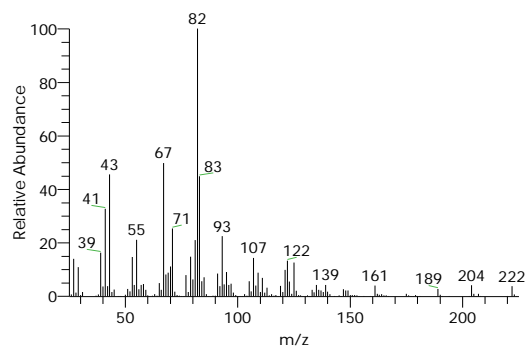
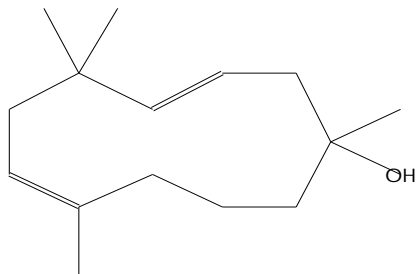
3,7-Cycloundecadien-1-ol, 1,5,5,8-tetramethyl-  
Formula C15H26O, MW 222, CAS# 118014-38-3, Entry# 51254  
1,5,5,8-Tetramethyl-3,7-cycloundecadien-1-ol #



1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL #  
Formula C15H26O, MW 222, CAS# 118014-38-3, Entry# 109756  
1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL



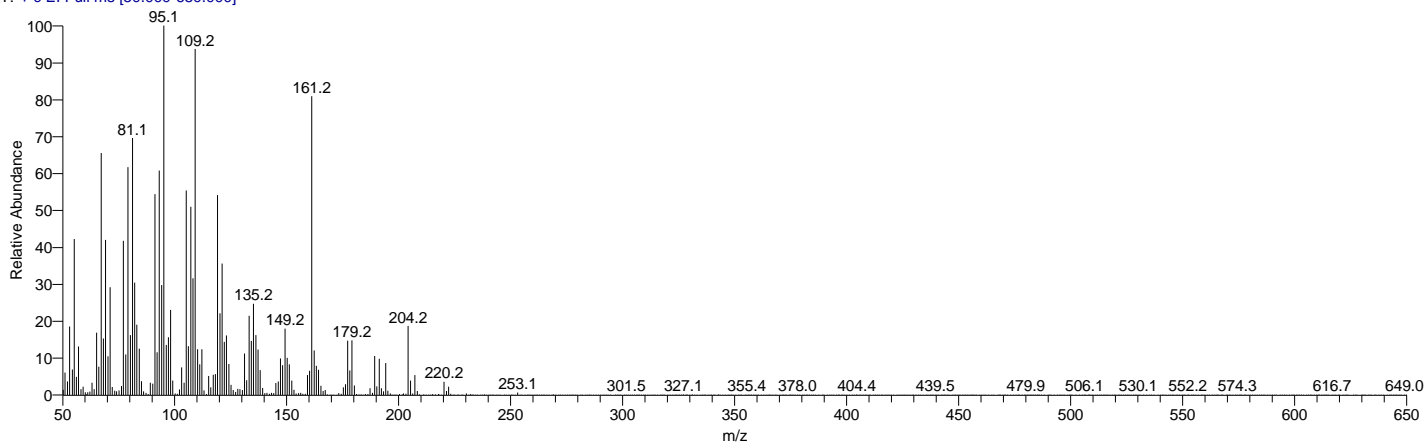
1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL #  
Formula C15H26O, MW 222, CAS# 118014-38-3, Entry# 385407  
1,5,5,8-TETRAMETHYL-3,7-CYCLOUNDECADIEN-1-OL





# My GC-MS Report

gerfa\_EA #5344 RT: 21.92 AV: 1 NL: 1.69E6  
T: + c EI Full ms [50.000-650.000]

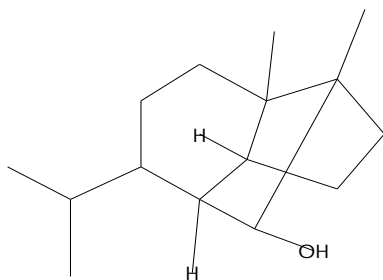


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
21.92	(1S,3aS,4S,5S,7aR,8R)-5-Isopropyl-1,7a-dimethyloctahydro-1H-1,4-methaninden-8-ol	0.14	C15H26O	222	21966-93-8	896	mainlib
21.92	Globulol	0.14	C15H26O	222	51371-47-2	844	mainlib
21.92	1,1,4,7-TETRAMETHYLDECAHYDRO-1H-CYCLOPROPA[E]AZULEN-4-OL #	0.14	C15H26O	222	51371-47-2	844	WileyRegistry8e

Compound Structure

Hit Spectrum

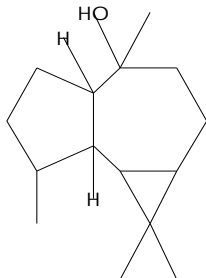
(1S,3aS,4S,5S,7aR,8R)-5-Isopropyl-1,7a-dimethyloctahydro-1H-1,4-methaninden-8-ol  
Formula C15H26O, MW 222, CAS# 21966-93-8, Entry# 86038  
Copaborneol



Globulol

Formula C15H26O, MW 222, CAS# 51371-47-2, Entry# 6178

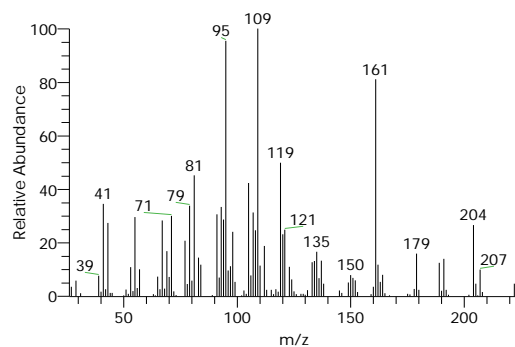
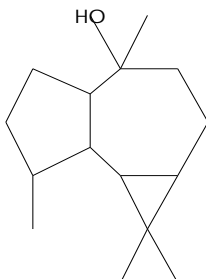
1,1,4,7-Tetramethyldecahydro-1H-cyclopropa[e]azulen-4-ol, (1a,4a,4a,7a,7a,7b)-



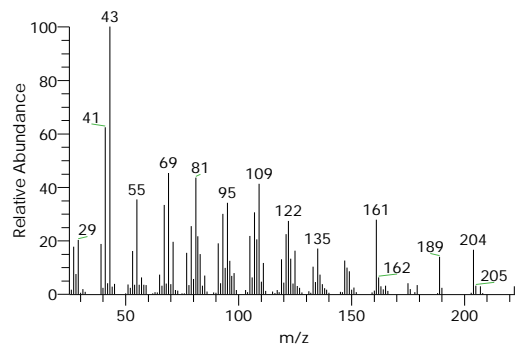
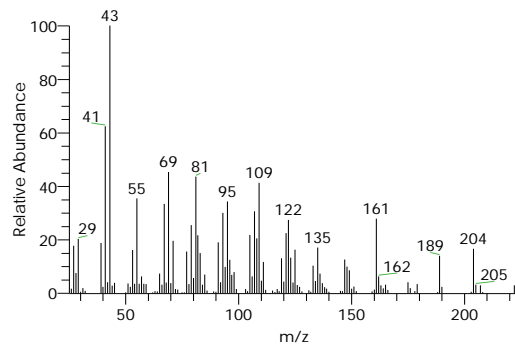
1,1,4,7-TETRAMETHYLDECAHYDRO-1H-CYCLOPROPA[E]AZULEN-4-OL #

Formula C15H26O, MW 222, CAS# 51371-47-2, Entry# 390796

1,1,4,7-TETRAMETHYLDECAHYDRO-1H-CYCLOPROPA[E]AZULEN-4-OL

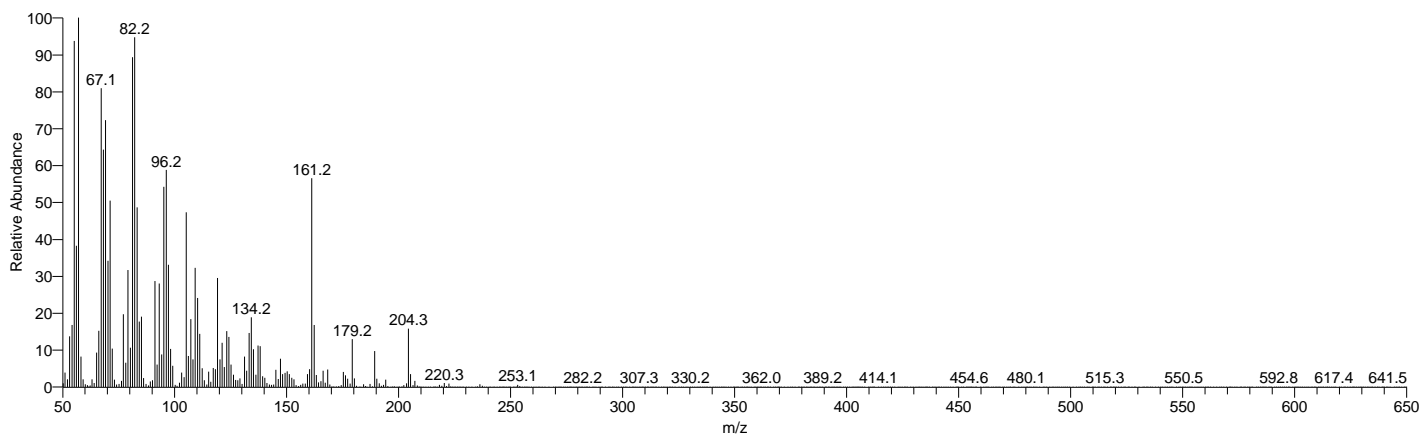


SI 816, RSI 844, mainlib, Entry# 6178, CAS# 51371-47-2, Globulol



# My GC-MS Report

gerfa\_EA #5403 RT: 22.12 AV: 1 NL: 2.07E6  
T: + c EI Full ms [50.000-650.000]

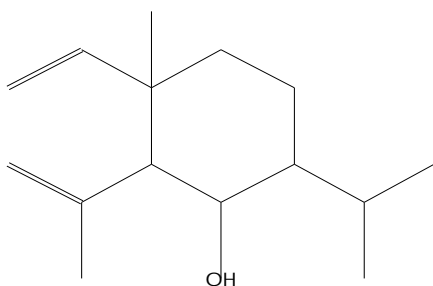


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
22.12	6-epi-shyobunol	0.22	C15H26O	222	NA	806	mainlib
22.12	Carotol	0.22	C15H26O	222	465-28-1	813	replib
22.12	3A(1H)-AZULENOL, 2,3,4,5,8,8A-HEXAHYDRO-6,8A-DIMETHYL-3-(1-METHYLETHYL)-, [3R-(3à,3Aà,8Aà)]-	0.22	C15H26O	222	465-28-1	808	WileyRegistry8e

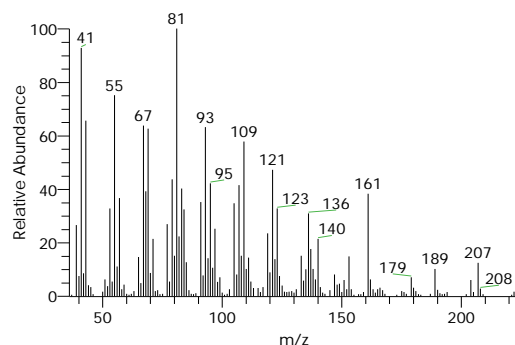
## Compound Structure

## Hit Spectrum

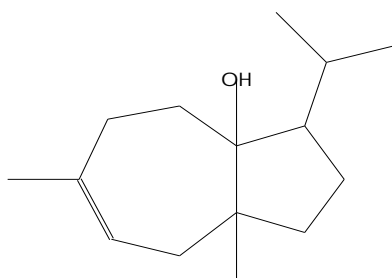
6-epi-shyobunol  
Formula C15H26O, MW 222, CAS# NA, Entry# 49759  
elema-1,3-dien-6à-ol



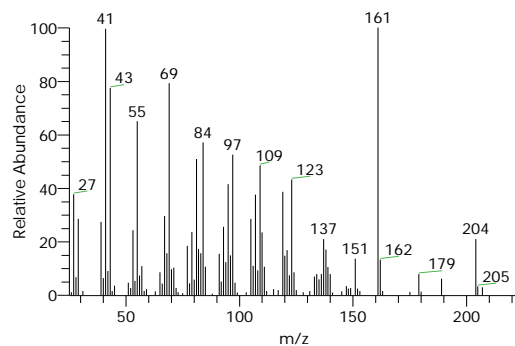
SI 794, RSI 806, mainlib, Entry# 49759, CAS# NA, 6-epi-shyobunol



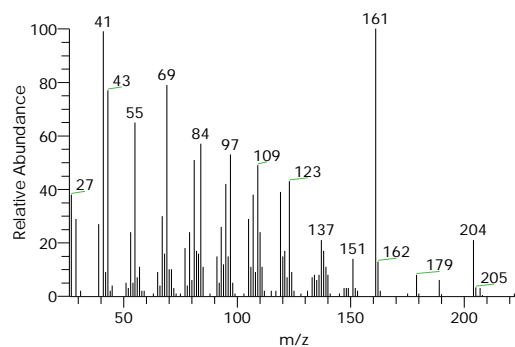
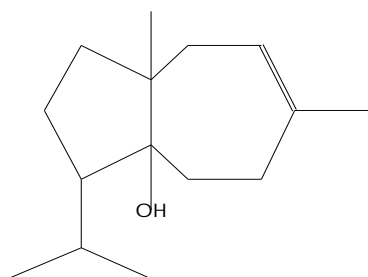
Carotol  
Formula C15H26O, MW 222, CAS# 465-28-1, Entry# 24838  
3a(1H)-Azulenol, 2,3,4,5,8,8a-hexahydro-6,8a-dimethyl-3-(1-methylethyl)-, [3R-(3à,3Aà,8Aà)]-



SI 774, RSI 813, replib, Entry# 24838, CAS# 465-28-1, Carotol

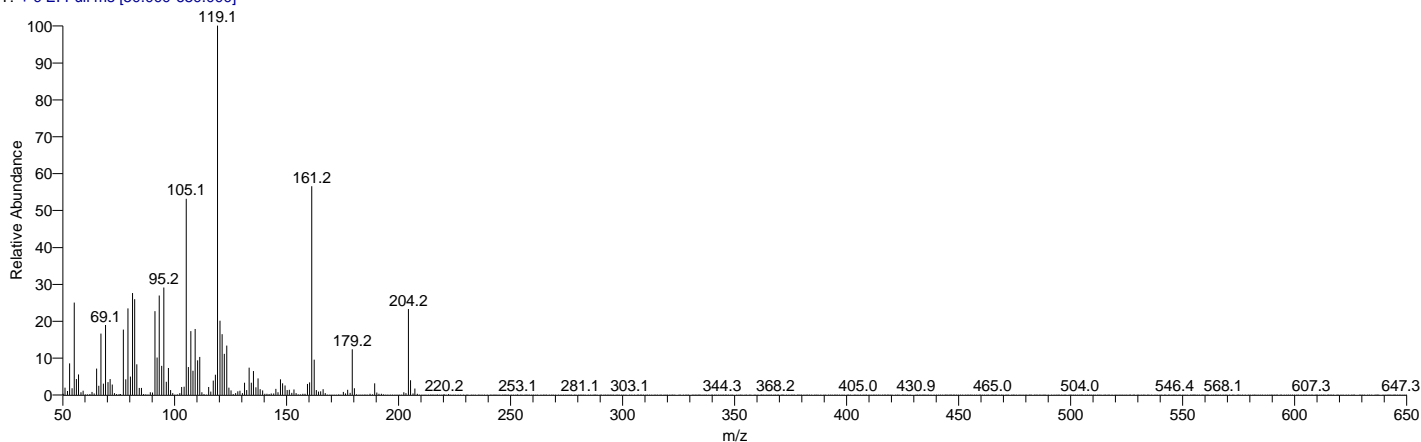


Formula C15H26O, MW 222, CAS# 465-28-1, Entry# 109801  
3-ISOPROPYL-6,8A-DIMETHYL-2,3,4,5,8,8A-HEXAHYDRO-3A(1H)-AZULENOL #



# My GC-MS Report

gerfa\_EA #5510 RT: 22.48 AV: 1 NL: 9.95E6  
T: + c EI Full ms [50.000-650.000]



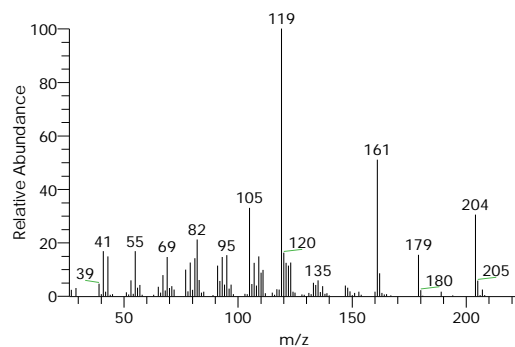
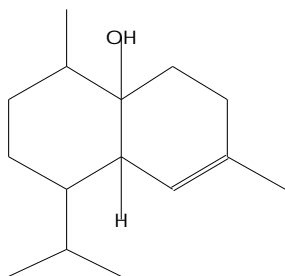
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
22.48	Di-epi-1,10-cubenol	0.53	C15H26O	222	73365-7	939	mainlib
22.48	Epicubenol	0.53	C15H26O	222	19912-6	919	mainlib
22.48	Cubenol	0.53	C15H26O	222	21284-2	914	mainlib

## Compound Structure

## Hit Spectrum

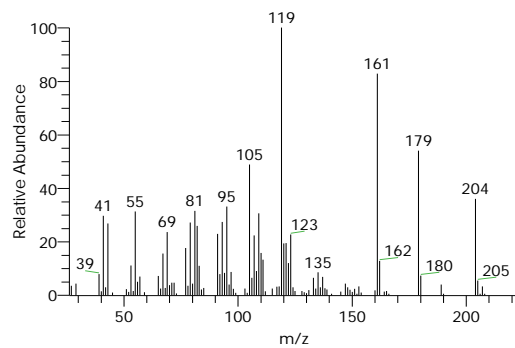
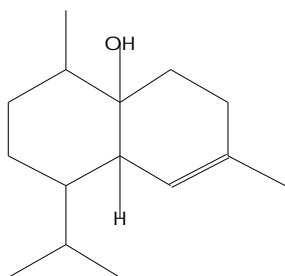
### Di-epi-1,10-cubenol

Formula C15H26O, MW 222, CAS# 73365-77-2, Entry# 100531  
(1S,4S,4aS,8aR)-1-Isopropyl-4,7-dimethyl-1,3,4,5,6,8a-hexahydronaphthalen-4a(2H)-ol



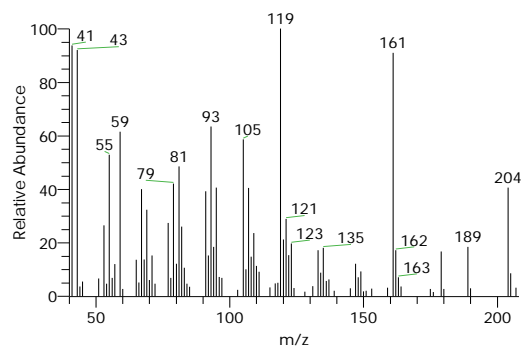
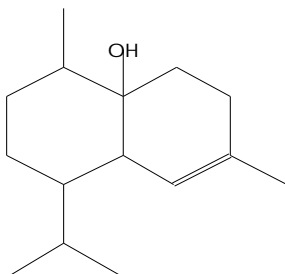
### Epicubenol

Formula C15H26O, MW 222, CAS# 19912-67-5, Entry# 100535  
(1S,4R,4aS,8aR)-1-Isopropyl-4,7-dimethyl-1,2,3,4,4a,5,6,8a-octahydronaphthalen-4a-ol



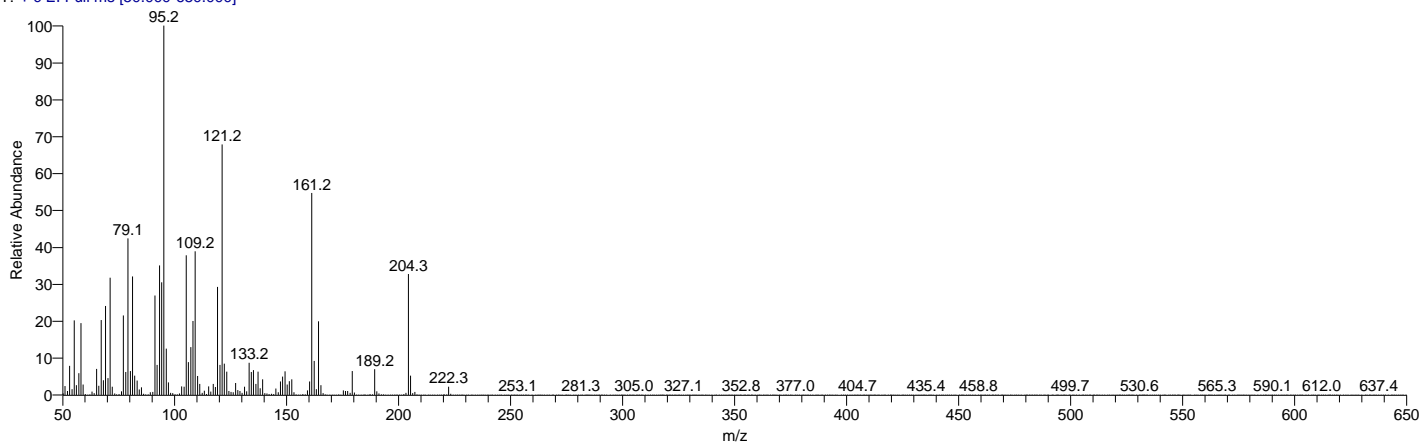
### Cubenol

Formula C15H26O, MW 222, CAS# 21284-22-0, Entry# 98767  
1-Isopropyl-4,7-dimethyl-1,3,4,5,6,8a-hexahydro-4a(2H)-naphthalenol-, [1S-(1à,4á,4aâ,8aâ)]-



# My GC-MS Report

gerfa\_EA #5606 RT: 22.80 AV: 1 NL: 1.35E7  
T: + c EI Full ms [50.000-650.000]



RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
22.80	.tau.-Muurolol	0.77	C15H26O	222	19912-62-0	949	replib
22.80	à-Cadinol	0.77	C15H26O	222	481-34-5	934	replib
22.80	.tau.-Muurolol	0.77	C15H26O	222	19912-62-0	923	mainlib

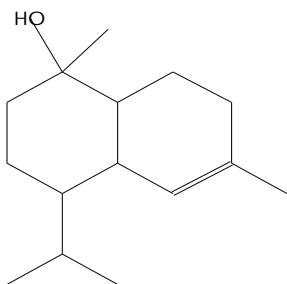
## Compound Structure

## Hit Spectrum

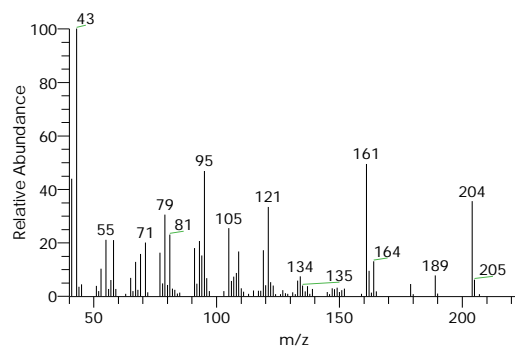
.tau.-Muurolol

Formula C15H26O, MW 222, CAS# 19912-62-0, Entry# 3342

4-Isopropyl-1,6-dimethyl-1,2,3,4,4a,7,8,8a-octahydro-1-naphthalenol-, [1S-(1à,4à,4aà,8aà)]-



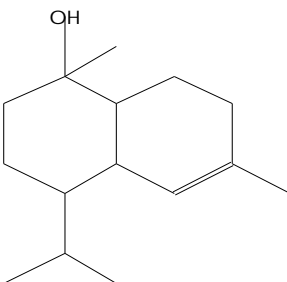
SI 932, RSI 949, replib, Entry# 3342, CAS# 19912-62-0, .tau.-Muurolol



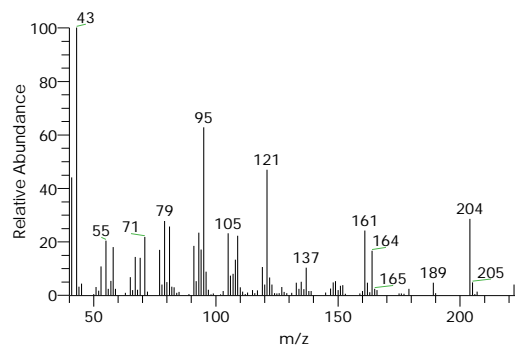
à-Cadinol

Formula C15H26O, MW 222, CAS# 481-34-5, Entry# 3029

4-Isopropyl-1,6-dimethyl-1,2,3,4,4a,7,8,8a-octahydro-1-naphthalenol #



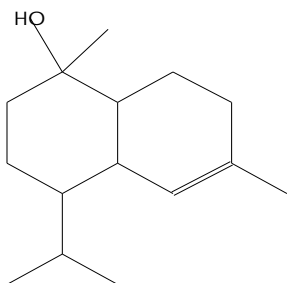
SI 924, RSI 934, replib, Entry# 3029, CAS# 481-34-5, à-Cadinol



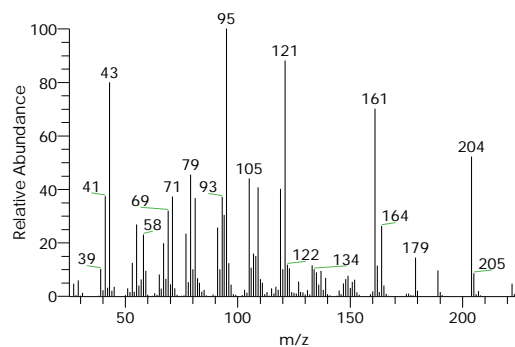
.tau.-Muurolol

Formula C15H26O, MW 222, CAS# 19912-62-0, Entry# 68534

4-Isopropyl-1,6-dimethyl-1,2,3,4,4a,7,8,8a-octahydro-1-naphthalenol-, [1S-(1à,4à,4aà,8aà)]-

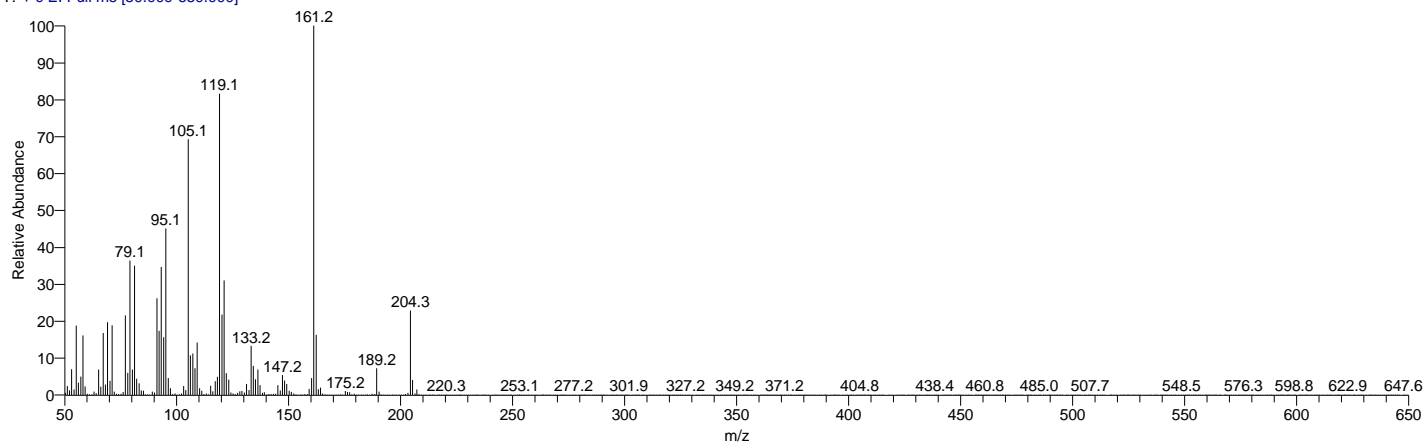


SI 923, RSI 923, mainlib, Entry# 68534, CAS# 19912-62-0, .tau.-Muurolol



# My GC-MS Report

gerfa\_EA #5635 RT: 22.89 AV: 1 NL: 9.19E6  
T: + c EI Full ms [50.000-650.000]

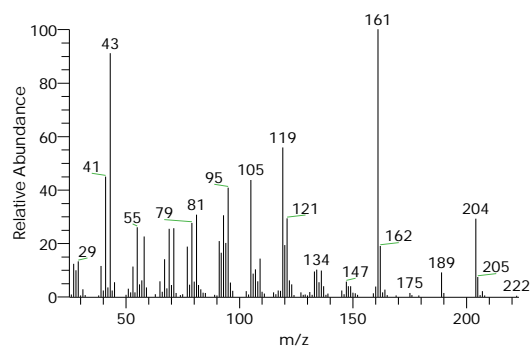
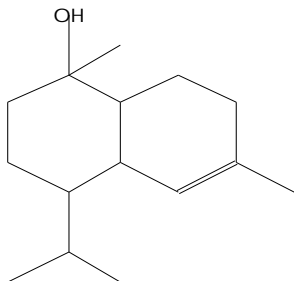


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
22.89	1-Naphthalenol, 1,2,3,4,4a,7,8,8a-octahydro-1,6-dimethyl-4-(1-m ethylethyl)-, [1R-(1à,4á,4aá,8aá)]-	0.37	C15H26O	222	19435-9 7-3	947	mainlib
22.89	1-NAPHTHALENOL, 1,2,3,4,4A,7,8,8A-OCTAHYDRO-1,6-DIMET HYL-4-(1-METHYLETHYL)-, [1R-(1à,4á,4Aá,8Aá)]-	0.37	C15H26O	222	19435-9 7-3	947	WileyRegi stry8e
22.89	1-Naphthalenol, 1,2,3,4,4a,7,8,8a-octahydro-1,6-dimethyl-4-(1-m ethylethyl)-, [1R-(1à,4á,4aá,8aá)]-	0.37	C15H26O	222	19435-9 7-3	934	replib

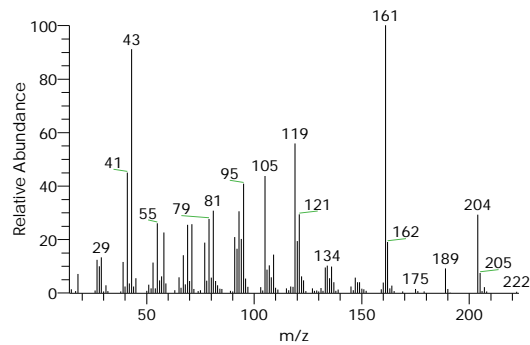
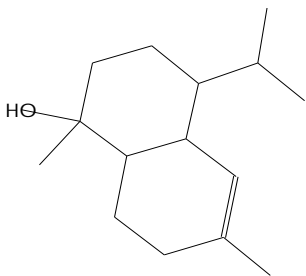
Compound Structure

Hit Spectrum

Formula C15H26O, MW 222, CAS# 19435-97-3, Entry# 150252  
(1R,4S,4aR,8aS)-4-Isopropyl-1,6-dimethyl-1,2,3,4,4a,7,8,8a-octahydronaphthalen-1-ol



Formula C15H26O, MW 222, CAS# 19435-97-3, Entry# 109844  
4-ISOPROPYL-1,6-DIMETHYL-1,2,3,4,4A,7,8,8A-OCTAHYDRO-1-NAPHTHALENOL #

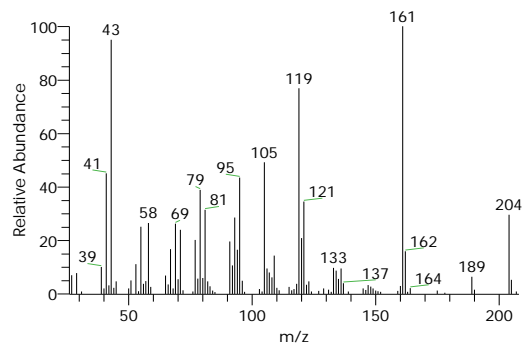
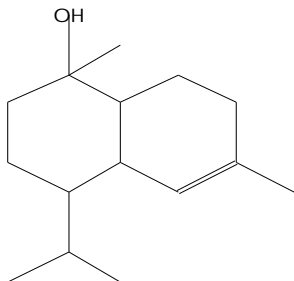


# My GC-MS Report

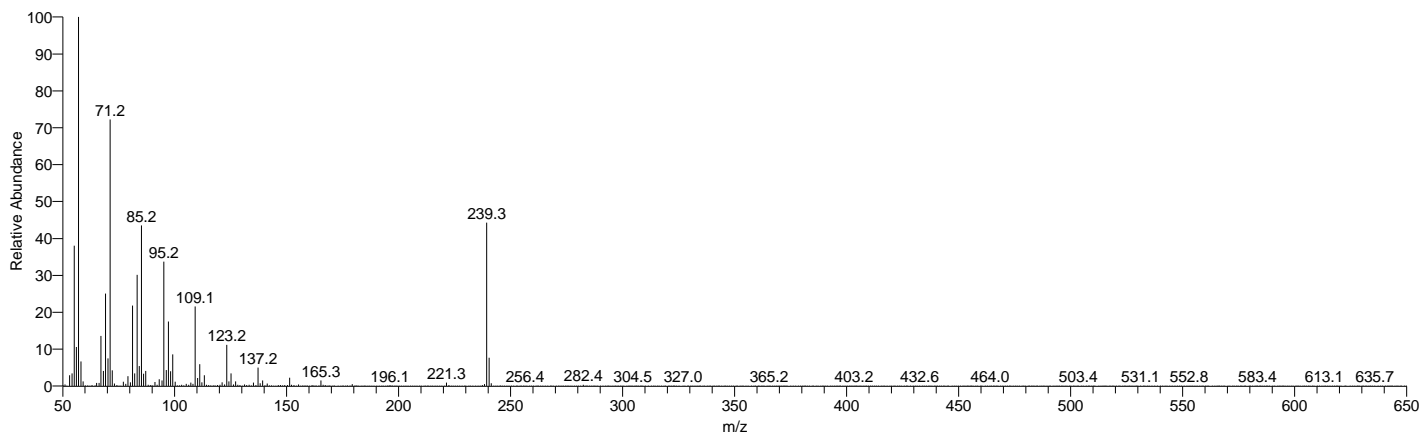
Compound Structure

Hit Spectrum

Formula C<sub>15</sub>H<sub>26</sub>O, MW 222, CAS# 19435-97-3, Entry# 24849  
(1R,4S,4aR,8aS)-4-Isopropyl-1,6-dimethyl-1,2,3,4,4a,7,8,8a-octahydronaphthalen-1-ol



gerfa\_EA #7757 RT: 30.01 AV: 1 NL: 7.40E6  
T: + c EI Full ms [50.000-650.000]

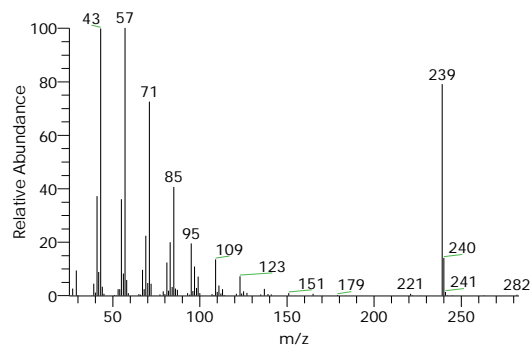
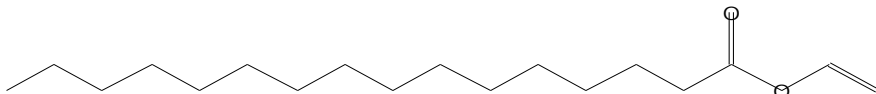


RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
30.01	Palmitic acid vinyl ester	0.25	C <sub>18</sub> H <sub>34</sub> O <sub>2</sub>	282	693-38-9	919	replib
30.01	Palmitic acid vinyl ester	0.25	C <sub>18</sub> H <sub>34</sub> O <sub>2</sub>	282	693-38-9	862	mainlib
30.01	Hexadecanoic acid, 4-nitrophenyl ester	0.25	C <sub>22</sub> H <sub>35</sub> NO <sub>4</sub>	377	1492-30-4	823	mainlib

Compound Structure

Hit Spectrum

Palmitic acid vinyl ester  
Formula C<sub>18</sub>H<sub>34</sub>O<sub>2</sub>, MW 282, CAS# 693-38-9, Entry# 6202  
Hexadecanoic acid, ethenyl ester

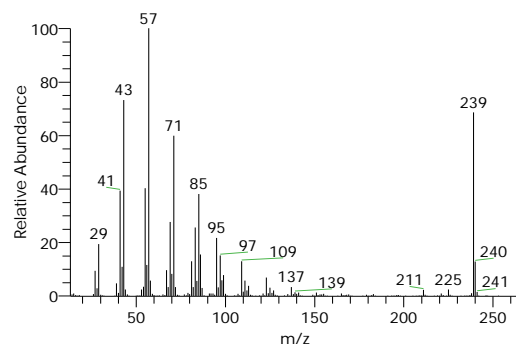
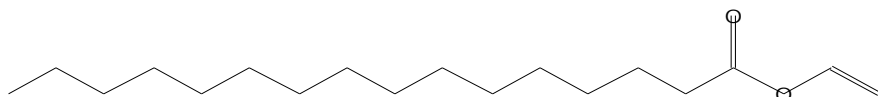


# My GC-MS Report

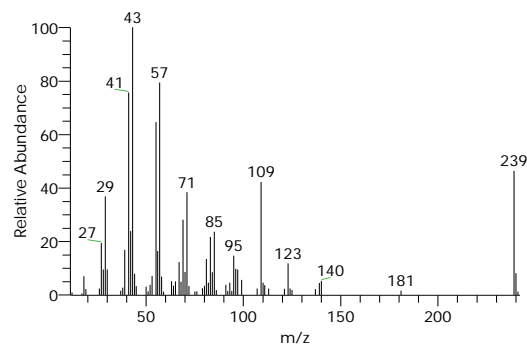
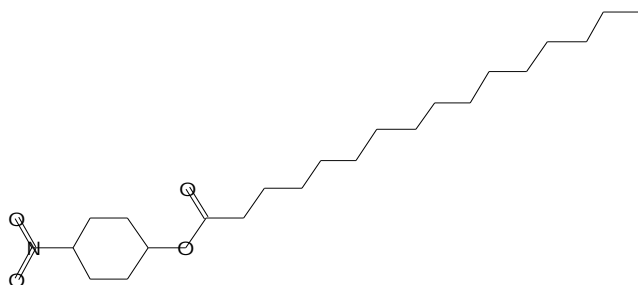
Compound Structure

Hit Spectrum

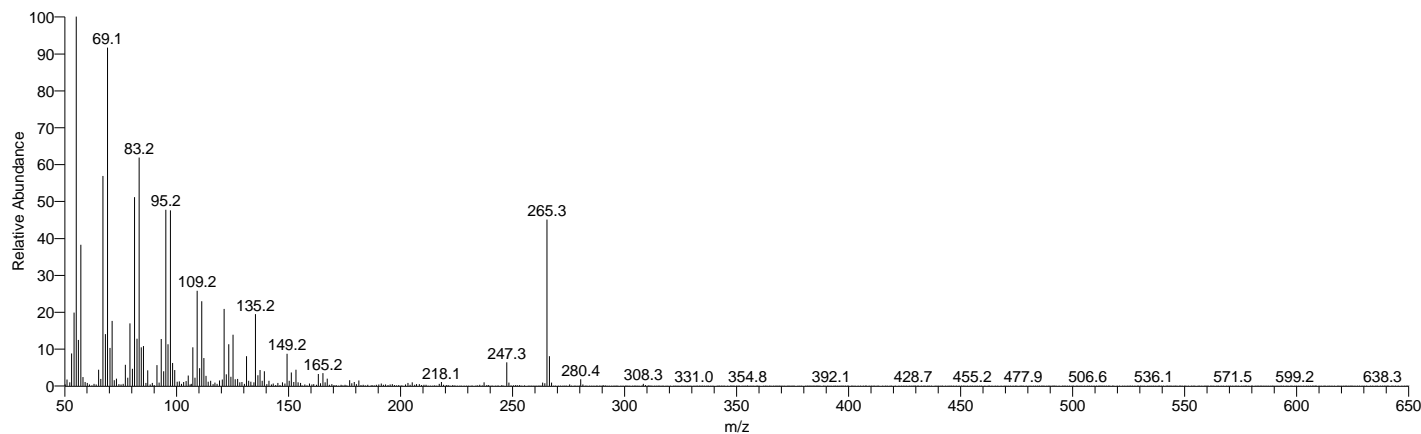
Palmitic acid vinyl ester  
Formula C18H34O2, MW 282, CAS# 693-38-9, Entry# 24552  
Hexadecanoic acid, ethenyl ester



Hexadecanoic acid, 4-nitrophenyl ester  
Formula C22H35NO4, MW 377, CAS# 1492-30-4, Entry# 7638  
4-Nitrophenyl palmitate #



gerfa\_EA #8726 RT: 33.26 AV: 1 NL: 2.03E6  
T: + c EI Full ms [50.000-650.000]



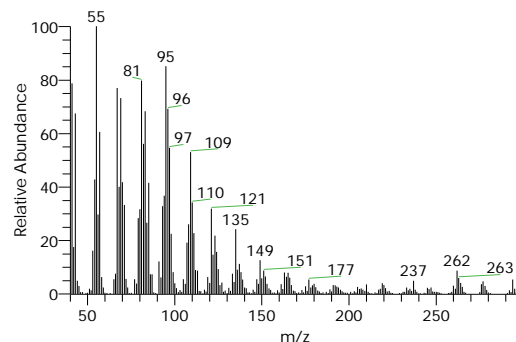
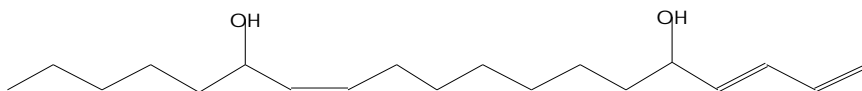
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
33.26	E,E,Z-1,3,12-Nonadecatriene-5,14-diol	0.12	C19H34O2	294	NA	720	mainlib
33.26	2,5-FURANDIONE,	0.12	C16H26O3	266	25377-7	797	WileyRegi
	3-(DODECENYL)DIHYDRO-				3-5		stry8e
33.26	2-Dodecen-1-yl(-)succinic anhydride	0.12	C16H26O3	266	19780-1	794	mainlib
					1-1		

# My GC-MS Report

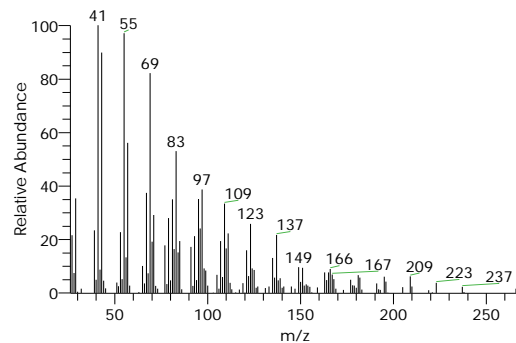
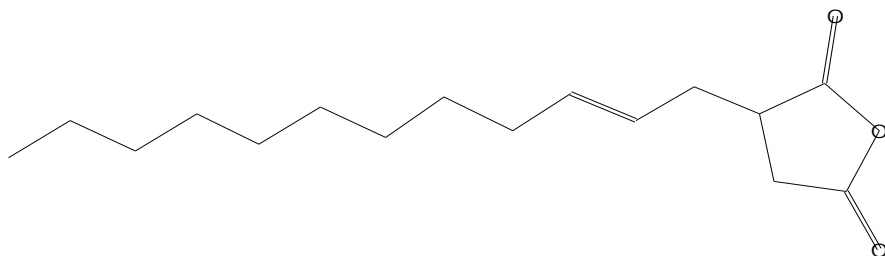
Compound Structure

Hit Spectrum

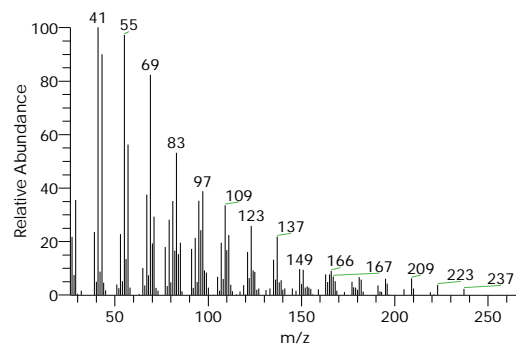
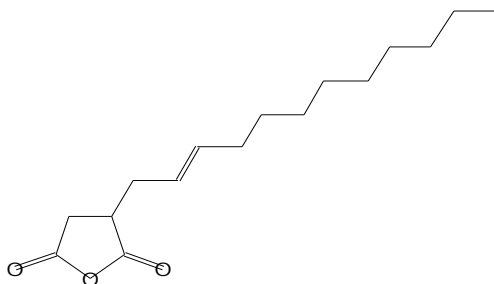
E,E,Z-1,3,12-Nonadecatriene-5,14-diol  
Formula C<sub>19</sub>H<sub>34</sub>O<sub>2</sub>, MW 294, CAS# NA, Entry# 21026  
(3E,12Z)-1,3,12-Nonadecatriene-5,14-diol #



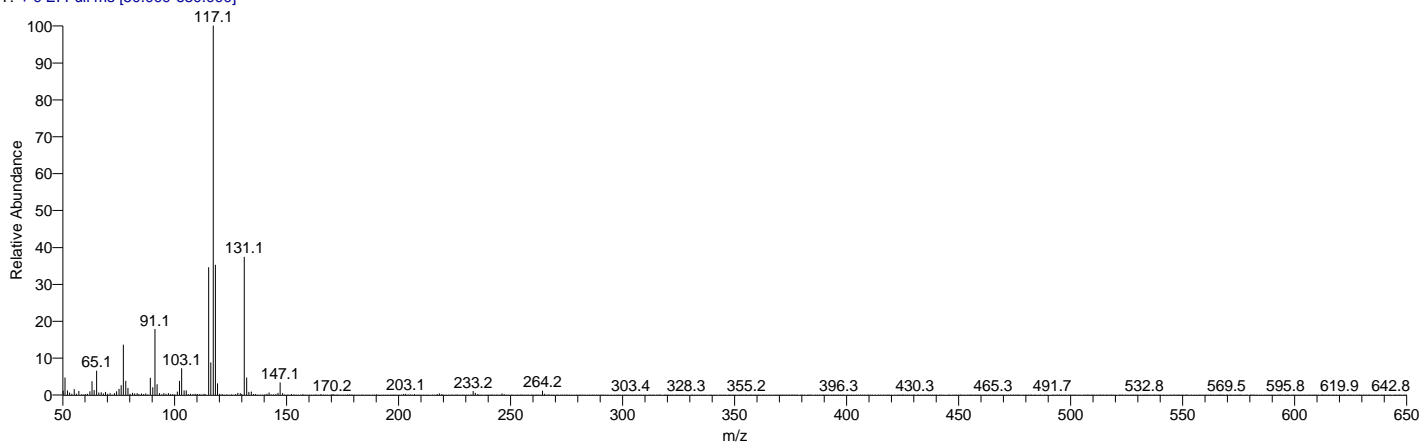
2,5-FURANDIONE, 3-(DODECENYL)DIHYDRO-  
Formula C<sub>16</sub>H<sub>26</sub>O<sub>3</sub>, MW 266, CAS# 25377-73-5, Entry# 156913  
2,5-FURANDIONE, 3-DODECYL-



2-Dodecen-1-yl(-)succinic anhydride  
Formula C<sub>16</sub>H<sub>26</sub>O<sub>3</sub>, MW 266, CAS# 19780-11-1, Entry# 2686  
2,5-Furandione, 3-dodecenyl-



gerfa\_EA #9386 RT: 35.47 AV: 1 NL: 1.07E7  
T: + c EI Full ms [50.000-650.000]



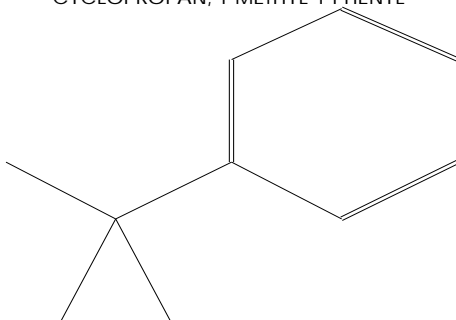
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
35.47	(1-METHYLCYCLOPROPYL)BENZENE	0.21	C <sub>10</sub> H <sub>12</sub>	132	NA	814	WileyRegi stry8e
35.47	DELTACYCLENE	0.21	C <sub>9</sub> H <sub>10</sub>	118	7785-1 0-6	897	WileyRegi stry8e



# My GC-MS Report

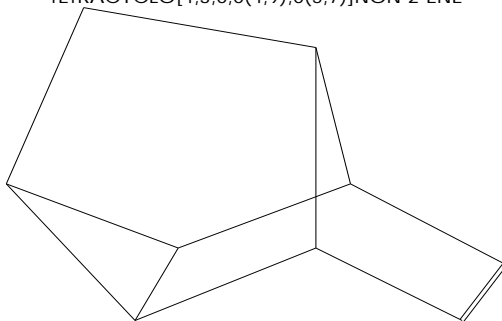
RT	Compound Name	Area %	Molecular Formula	Molecular Weight	Cas #	MF	Library
35.47	Deltacyclene	0.21	C <sub>9</sub> H <sub>10</sub>	118	7785-10-6	896	mainlib
Compound Structure				Hit Spectrum			

(1-METHYLCYCLOPROPYL)BENZENE  
Formula C<sub>10</sub>H<sub>12</sub>, MW 132, CAS# NA, Entry# 380177  
CYCLOPROPAN, 1-METHYL-1-PHENYL-



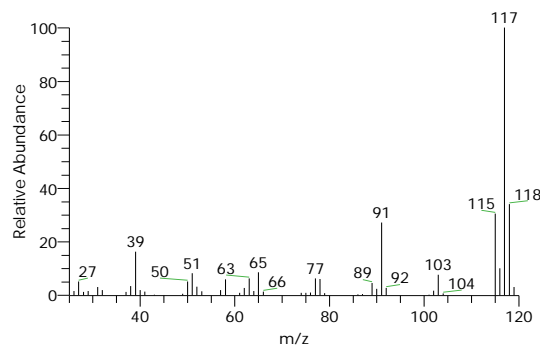
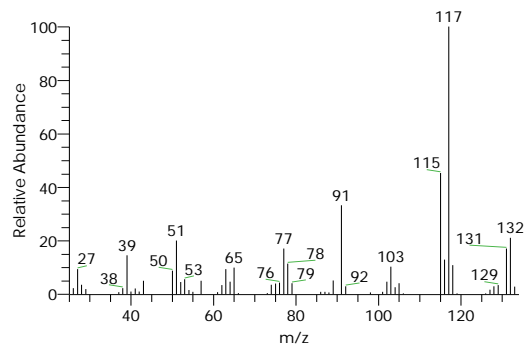
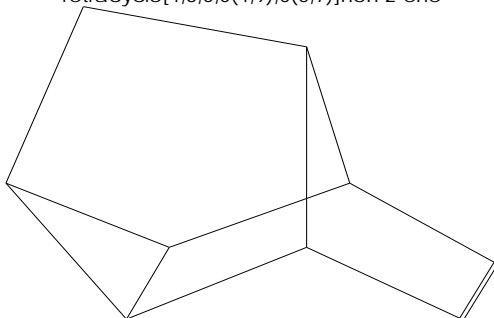
DELTACYCLENE

Formula C<sub>9</sub>H<sub>10</sub>, MW 118, CAS# 7785-10-6, Entry# 358032  
TETRACYCLO[4,3,0,0(4,9),0(5,7)]NON-2-ENE



Deltacyclene

Formula C<sub>9</sub>H<sub>10</sub>, MW 118, CAS# 7785-10-6, Entry# 97452  
Tetracyclo[4,3,0,0(4,9),0(5,7)]non-2-ene



SI 769, RSI 896, mainlib, Entry# 97452, CAS# 7785-10-6, Deltacyclene

