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APPENDICES

Retention time (min)	Compound	Structure
4.38	α-thujene	
4.98	tricyclene	
5.03	α-pinene	
5.48	camphene	
5.98	sylvestrene	
6.01	benzaldehyde	HC=0
6.08	sabinene	
6.09	isosylvestrene	Ť Ţ
6.11	myrcene	
		<u> </u>

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A. The chemical components of essential oil isolated from selected Lauraceous Plants

Retention time (min)	Compound	Structure
6.23	β-phellandrene	
6.37	6-methyl-5-hepten-2-one	
6.54	β-pinene	
6.83	<i>n</i> -decane	CH ₃ (CH ₂) ₈ CH ₃
7.11	α-phellandrene	
7.16	δ-3-carene	
7.44	δ-2-carene	
7.59	o-cymene	
7.78	verbenene	tr
7.88	limonene	

Retention time (min)	Compound	Structure
7.94	<i>trans</i> -sabinene hydrate acetate	O-Ac
7.98	1,8-cineole	↓
8.09	terpin-4-ol-acetate	O-Ac
8.13	(Z)-β-ocimene	
8.48	(E)-β+ocimene	E A
8.91	γ-terpinene	
9.51	<i>cis</i> -linalool oxide	н
9.61	<i>trans</i> -sabinene hydrate	HO

Retention time (min)	Compound	Structure
9.76	α-terpinene	
9.93	terpinolene	
10.13	trans-linalool oxide	ОН
10.23	2-nonanone	CH ₃ CO(CH ₂) ₆ CH ₃
10.27	1,4-cineole	
10.35	verbenone	C C C C C C C C C C C C C C C C C C C
10.44	n-heneicosane	CH ₃ (CH ₂) ₁₉ CH ₃
10.81	linalool	ОН
11.73	α-campholenal	ST H
11.81	trans-para-menth-2-en-1-ol	OH

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Retention time (min)	Compound	Structure
12.54	camphor	X
12.76	citronellal	H O
13.19	<i>cis</i> -limonene oxide	
13.25	acetophenone	
13.46	(Z)-cinnamyl alcohol	СТОН
13.66	ethyl benzoate	O-CH ₂ CH ₃
13.88	santolina alcohol	OH
13.93	isoborneol	
13.99	borneol	ОН

Retention time (min)	Compound	Structure
14.16	terpin-4-ol	
14.88	methyl chavicol	O-Me
		11
14.98	a-terpineol	1
16.24		Сон
16.34	citronellol	
. 16.63	neral	OH H OH
		\sim
16.81	(Z)-cinnamaldehyde	Z O H
17.23	cis-carveol	
		он
17.38	geraniol	1
		I OH

Retention time	Compound	Structure
(min)		
17.94	<i>(Z)</i> -methyl cinnamate	
18.04	<i>n</i> -pentyl benzoate	O-Me
18.44	bornyl acetate	O-Ac
18.46	isobornyl acetate	, MO-Ac
18.50	2-undecanone	CH ₃ CO(CH ₂) ₈ CH ₃
18.61	<i>(E)-</i> cinnamaldehyde	
18.77	geranial	O H H
18.99	<i>(Z)</i> -isosafrole	
19.79	α-terpinyl acetate	O-Ac

Retention time (min)	Compound	Structure
20.63	δ-elemene	
20.17	α-cubebene	H
21.02	(E)-isoeugenol	U-Me HO
21.13	β-cubebene	
21.21	<i>(E)</i> -methyl cinnamate	
21.29	3-thujyl acetate	H H H O-Ac
22.10	(Z)-isoeugenol	HO HO Z
22.38	α-copaene	H H H

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Retention time (min)	Compound	Structure
22.66	<i>n</i> -hexyl benzoate	
22.73	geranyl acetate	O-Ac
22.93	germacrene D	
23.01	<i>cis</i> -verbenyl acetate	Unit O-Ac
23.04	<i>(E)</i> -isosafrole	
23.09	β-elemene	
23.71	<i>cis</i> -β-guiene	
23.98	methyl eugenol	O-Me O-Me

Retention time (min)	Compound	Structure
24.10	α-bulnesene	H
24.16	tetradecanal	CH ₃ (CH ₂) ₁₂ COH
24.20	cumin aldehyde	H O
24.33	<i>(E)</i> -caryophyllene	H
24.89	<i>(Z)</i> -cinnamyl acetate	C-Ac
25.07	<i>9-epi-(E)-</i> caryophyllene	H
25.11	longifolene	\$
25.23	α-gurjunene	H
25.37	seychellene	

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Retention time (min)	Compound	Structure
25.41	(Z)-methyl-butyl-benzoate	
25.44	(Z)-α-bisabolene	
25.54	cyperene	H
25.56	<i>allo-</i> aromadendrene	H H H H H H H
25.58	<i>trans</i> -calamenene	
25.73	<i>cis</i> -β-guaiene	
25.89	α-humulene	
26.03	valencene	
26.58	δ-cadinene	H H

Retention time (min)	Compound	Structure
26.61	β-patchoulene	
26.75	γ-cadinene	
26.89	α-muurolene	H H H H
26.92	2-methyl-undecanal	CH ₃ (CH ₂) ₈ CH(CH ₃)COH
26.99	γ-muurolene	H
27.34	β-selinene	
27.38	viridiflorene	
27.43	isoledene	
27.59	bicyclogermacrene	

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Retention time (min)	Compound	Structure
27.63	α-selinene	
27.84	epi-cubebol	OH IIII H H H
27.88	β-gurjunene	H
28.11	germacrene A	
28.39	<i>cis</i> -muurola-4(14), 5-diene	I H
28.59	<i>trans</i> -β-guaiene	
28.76	α-cadinene	H
28.78	<i>cis</i> -calamenene	

Retention time (min)	Compound	Structure
28.84	eugenyl acetate	O-Ac O-Ac
28.91	<i>3</i> , <i>7(11)</i> -selinadiene	
29.09	myristicin	Me-O
29.19	cadina-1,4-diene	H ,
29.61	α-calacorene	
30.15	elemicin	O-Me Me-O I
30.19	germacrene B	
30.66	<i>(E)</i> -nerolidol	OH OH

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Retention time (min)	Compound	Structure
31.29	spathulenol	H H H
31.32	caryophyllene oxide	H
31.47	globulol	H. M.
31.59	α-eudesmol acetate	O-Ac
31.63	β-eudesmol acetate	O-Ac
31.96	hinesol acetate	O-Ac
32.26	longiborneol acetate	Ac-O Lundart
32.30	ledol	HOH

Retention time (min)	Compound	Structure
32.44	humulene epoxide II	
32.64	dodecanal	CH ₃ (CH ₂) ₁₀ COH
33.08	juniper camphor acetate	H I O-Ac
33.19	<i>1-epi-</i> cubenol	OH H
33.54	bicyclo-vetivenol	€он
33.68	cubenol	OH H H
33.93	α-muurolol	HOH
33.98	<i>epi-</i> α-muurolol	HO

Retention time (min)	Compound	Structure
34.36	himachalol	H H ,,,, OH
34.48	α-eudesmol	Н СН
34.49	αcadinol	HO
34.58	selin-11-en-4-alpha-ol	Л Н Н Н ОН
38.86	benzyl benzoate	0=c-0

Note : GC/MS Condition
Instrument model : Varian Saturn 3
Column : fused silica capillary column (30 m X 0.25 mm i.d.) coated with DB-5 (J&w), film thickness 0.25 μm
Column programming : 60-240°C rate 3.3°C/min
Injector temperature : 180°C
Helium carrier gas : 1 ml/min
Split ratio : 100 : 1
Accelerating voltage : 1700 volts
Sample size : 1 μl
Solvent : methanol (HPLC grade)













Figure 42 Mass spectra of benzaldehyde (A) and authentic benzaldehyde (B) by GC-MS



Figure 43 Mass spectra of benzyl benzoate (A) and authentic benzyl benzoate (B) by GC-MS







Figure 45 Mass spectra of bisabolene $\langle Z \rangle$ - $\alpha \rightarrow \langle A \rangle$ and authentic bisabolene $\langle Z \rangle$ - $\alpha \rightarrow \langle B \rangle$ by GC-MS











Figure 48 Mass spectra of bulnesene $\langle \alpha - \rangle$ (A) and authentic bulnesene $\langle \alpha - \rangle$ (B) by GC-MS





Mass spectra of cadina-1.4-diene (A) and authentic cadina-1.4-diene (B) by GC-MS



Figure 50 Mass spectra of cadinene $\langle \alpha \rangle$ (A) and authentic cadinene $\langle \alpha \rangle$ (B) by GC-MS



Figure 51 Mass spectra of cadinene $<\delta >$ (A) and authentic cadinene $<\delta >$ (B) by GC-MS



Figure 52 Mass spectra of cadinene $\langle \gamma \rangle$ (A) and authentic cadinene $\langle \gamma \rangle$ (B) by GC-MS



Figure 53 Mass spectra of cadinol $<\alpha ->$ (A) and authentic cadinol $<\alpha ->$ (B) by GC-MS



Figure 54 Mass spectra of calacorene $\langle \alpha \rangle$ (A) and authentic calacorene $\langle \alpha \rangle$ (B) by GC-MS



Figure 55 Mass spectra of calamenene $\langle cis \rangle$ (A) and authentic calamenene $\langle cis \rangle$ (B) by GC-MS



Figure 56 Mass spectra of calamenene < *trans->* (A) and authentic calamenene < *trans->* (B) by GC-MS



Figure 57 Mass spectra of camphene (A) and authentic camphene (B) by GC-MS



Figure 58 Mass spectra of campholenal $<\alpha ->$ (A) and authentic campholenal $<\alpha ->$ (B) by GC-MS











Figure 61 Mass spectra of carene $< \delta - 3 - >$ (A) and authentic carene $< \delta - 3 - >$ (B) by GC-MS



Figure 62 Mass spectra of carveol < cis > (A) and authentic carveol < cis > (B) by GC-MS



Figure 63 Mass spectra of caryophellene $\langle E \rangle - \rangle$ (A) and authentic caryophellene $\langle E \rangle - \rangle$ (B) by GC-MS







Figure 65 Mass spectra of caryophyllene oxide (A) and authentic caryophyllene oxide (B) by GC-MS



Figure 66 Mass spectra of cineole < 1, 4 -> (A) and authentic cineole < 1, 4 -> (B) by GC-MS



Figure 67 Mass spectra of cineole < 1, 8 -> (A) and authentic cineole < 1, 8 -> (B) by GC-MS







Figure 69 Mass spectra of cinnamaldehyde $\langle (Z) \rangle$ (A) and authentic cinnamaldehyde $\langle (Z) \rangle$ (B) by GC-MS



Figure 70 Mass spectra of cinnamyl acetate < (Z) -> (A) and authentic cinnamyl acetate < (Z) -> (B) by GC-MS



Figure 71 Mass spectra of cinnamyl alcohol < (Z) -> (A) and authentic cinnamyl alcohol < (Z) -> (B) by GC-MS



Figure 72 Mass spectra of citronellal (A) and authentic citronellal (B) by GC-MS



Figure 73 Mass spectra of citronellol (A) and authentic citronellol (B) by GC-MS



Figure 74 Mass spectra of copaene $< \alpha ->$ (A) and authentic copaene $< \alpha ->$ (B) by GC-MS



Figure 75 Mass spectra of cubebene $< \alpha ->$ (A) and authentic cubebene $< \alpha ->$ (B) by GC-MS



Figure 76 Mass spectra of cubebene $<\beta ->$ (A) and authentic cubebene $<\beta ->$ (B) by GC-MS



Figure 77 Mass spectra of cubebol < *epi->* (A) and authentic cubebol < *epi->* (B) by GC-MS



Figure 78 Mass spectra of cubenol (A) and authentic cubenol (B) by GC-MS



Figure 79 Mass spectra of cubenol <*l-epi*-> (A) and authentic cubenol <*l-epi*-> (B) by GC-MS







Figure 81 Mass spectra of cymene $\langle o \rangle$ (A) and authentic cymene $\langle o \rangle$ (B) by GC-MS



Figure 82 Mass spectra of cyperene (A) and authentic cyperene (B) by GC-MS







Figure 84 Mass spectra of dodecanal (A) and authentic dodecanal (B) by GC-MS











Figure 87 Mass spectra of elemicin (A) and authentic elemicin (B) by GC-MS











Figure 90 Mass spectra of eudesmol acetate $< \alpha ->$ (A) and authentic eudesmol acetate $< \alpha ->$ (B) by GC-MS



Figure 91 Mass spectra of eudesmol acetate $< \beta > (A)$ and authentic eudesmol acetate $< \beta > (B)$ by GC-MS



Figure 92 Mass spectra of eugenvl acetate (A) and authentic eugenvl acetate (B) by GC-MS



Figure 93 Mass spectra of geranial (A) and authentic geranial (B) by GC-MS



Figure 94 Mass spectra of geraniol (A) and authentic geraniol (B) by GC-MS







Figure 96 Mass spectra of germacrene A (A) and authentic germacrene A (B) by GC-MS















Figure 100 Mass spectra of guaiene $\langle cis-\beta \rangle$ (A) and authentic guaiene $\langle cis-\beta \rangle$ (B) by GC-MS



Figure 101 Mass spectra of guaiene $< trans-\beta > (A)$ and authentic guaiene $< trans-\beta > (B)$ by GC-MS



Figure 102 Mass spectra of gurjunene $<\beta >$ (A) and authentic gurjunene $<\beta >$ (B) by GC-MS







Figure 104 Mass spectra of heneicosane < n > (A) and authentic heneicosane < n > (B) by GC-MS



Figure 105 Mass spectra of hepten-2-one <6-methyl-5-> (A) and authentic hepten-2one <6-methyl-5-> (B) by GC-MS



Figure 106 Mass spectra of hexyl benzoate < n > (A) and authentic hexyl benzoate < n > (B) by GC-MS







Figure 108 Mass spectra of hinesol acetate (A) and authentic hinesol acetate (B) by GC-MS



Figure 109 Mass spectra of humulene $< \alpha ->$ (A) and authentic humulene $< \alpha ->$ (B) by GC-MS



Figure 110 Mass spectra of humulene epoxide II (A) and authentic humulene epoxide II (B) by GC-MS



Figure 111 Mass spectra of isoborneol (A) and authentic isoborneol (B) by GC-MS



Figure 112 Mass spectra of isobornyl acetate (A) and authentic isobornyl acetate (B) by GC-MS



Figure 113 Mass spectra of isoeugenol $\langle Z \rangle$ -> (A) and authentic isoeugenol $\langle Z \rangle$ -> (B) by GC-MS



Figure 114 Mass spectra of isoledene (A) and authentic isoledene (B) by GC-MS



Figure 115 Mass spectra of isosafrole $\langle E \rangle > \langle A \rangle$ and authentic isosafrole $\langle E \rangle > \langle B \rangle$ by GC-MS



Figure 116 Mass spectra of isosafrole $\langle Z \rangle$ -> (A) and authentic isosafrole $\langle Z \rangle$ -> (B) by GC-MS



Figure 117 Mass spectra of isosylvestrene (A) and authentic isosylvestrene (B) by GC-MS



Figure 118 Mass spectra of juniper camphor acetate (A) and authentic juniper camphor acetate (B) by GC-MS



Figure 119 Mass spectra of ledol (A) and authentic ledol (B) by GC-MS



Figure 120 Mass spectra of limonene (A) and authentic limonene (B) by GC-MS



Figure 121 Mass spectra of limonene oxide $\langle cis \rangle$ (A) and authentic limonene oxide $\langle cis \rangle$ (B) by GC-MS







Figure 123 Mass spectra of linalool oxide $\langle cis \rangle$ (A) and authentic linalool oxide $\langle cis \rangle$ (B) by GC-MS



Figure 124 Mass spectra of linalool oxide < *trans->* (A) and authentic linalool oxide < *trans->* (B) by GC-MS



Figure 125 Mass spectra of longiborneol acetate (A) and authentic longiborneol acetate (B) by GC-MS



Figure 126 Mass spectra of longifolene (A) and authentic longifolene (B) by GC-MS











Figure 129 Mass spectra of methyl cinnamate $\langle (E) \rangle$ (A) and authentic methyl cinnamate $\langle (E) \rangle$ (B) by GC-MS



Figure 130 Mass spectra of methyl cinnamate $\langle Z \rangle - \langle A \rangle$ and authentic methyl cinnamate $\langle Z \rangle - \langle B \rangle$ by GC-MS



Figure 131 Mass spectra of methyl eugenol (A) and authentic methyl eugenol (B) by GC-MS



Figure 132 Mass spectra of 6-methyl-5-hepten-2-one (A) and authentic 6-methyl-5hepten-2-one (B) by GC-MS



Figure 133 Mass spectra of methyl-butyl-benzoate < 2-> (A) and authentic methylbutyl-benzoate < 2-> (B) by GC-MS



Figure 134 Mass spectra of methyl-undecanal < 2-> (A) and authentic methylundecanal < 2-> (B) by GC-MS



Figure 135 Mass spectra of muurola-4(14), 5-diene < cis-> (A) and authentic muurola-4(14), 5-diene < cis-> (B) by GC-MS



Figure 136 Mass spectra of muurolene $< \alpha -> (A)$ and authentic muurolene $< \alpha -> (B)$ by GC-MS



Figure 137 Mass spectra of muurolene $<\gamma >$ (A) and authentic muurolene $<\gamma >$ (B) by GC-MS



Figure 138 Mass spectra of muurolol $\langle epi-\alpha - \rangle$ (A) and authentic muurolol $\langle epi-\alpha - \rangle$ (B) by GC-MS



Figure 139 Mass spectra of muurolol $< \alpha > (A)$ and authentic muurolol $< \alpha > (B)$ by GC-MS



Figure 140 Mass spectra of myrcene (A) and authentic myrcene (B) by GC-MS



Figure 141 Mass spectra of myristicin (A) and authentic myristicin (B) by GC-MS



Figure 142 Mass spectra of neral (A) and authentic neral (B) by GC-MS



Figure 143 Mass spectra of nerolidol $\langle (E) \rangle$ (A) and authentic nerolidol $\langle (E) \rangle$ (B) by GC-MS



Figure 144 Mass spectra of nonanone <2-> (A) and authentic nonanone <2-> (B) by GC-MS



Figure 145 Mass spectra of ocimene $< (E)-\beta -> (A)$ and authentic ocimene $< (E)-\beta ->$ (B) by GC-MS



Figure 146 Mass spectra of ocimene $\langle Z \rangle$ - β - \rangle (A) and authentic ocimene $\langle Z \rangle$ - β - \rangle (B) by GC-MS



Figure 147 Mass spectra of patchoulene $<\beta>$ (A) and authentic patchoulene $<\beta>$ (B) by GC-MS



Figure 148 Mass spectra of pentyl benzoate < n > (A) and authentic pentyl benzoate < n > (B) by GC-MS



Figure 149 Mass spectra of phellandrene $<\beta >$ (A) and authentic phellandrene $<\beta >$ (B) by GC-MS



Figure 150 Mass spectra of phellandrene $\langle \alpha \rangle$ (A) and authentic phellandrene $\langle \alpha \rangle$ (B) by GC-MS



Figure 151 Mass spectra of pinene $<\beta ->$ (A) and authentic pinene $<\beta ->$ (B) by GC-MS



Figure 152 Mass spectra of pinene $\langle \alpha \rangle$ (A) and authentic pinene $\langle \alpha \rangle$ (B) by GC-MS



Figure 153 Mass spectra of sabinene (A) and authentic sabinene (B) by GC-MS



Figure 154 Mass spectra of sabinene hydrate < *trans->* (A) and authentic sabinene hydrate < *trans->* (B) by GC-MS



Figure 155 Mass spectra of sabinene hydrate acetate < *trans->* (A) and authentic sabinene hydrate acetate < *trans->* (B) by GC-MS



Figure 156 Mass spectra of santolina alcohol (A) and authentic santolina alcohol (B) by GC-MS



Figure 157 Mass spectra of selin-*11-en-4-alpha-ol* (A) and authentic selin-*11-en-4-alpha-ol* (B) by GC-MS



Figure 158 Mass spectra of selinadiene $\langle 3, 7(11) \rangle$ (A) and authentic selinadiene $\langle 3, 7(11) \rangle$ (B) by GC-MS



Figure 159 Mass spectra of selinene $<\beta ->$ (A) and authentic selinene $<\beta ->$ (B) by GC-MS



Figure 160 Mass spectra of selinene $\langle \alpha - \rangle$ (A) and authentic selinene $\langle \alpha - \rangle$ (B) by GC-MS







Figure 162 Mass spectra of spathulenol (A) and authentic spathulenol (B) by GC-MS



Figure 163 Mass spectra of sylvestrene (A) and authentic sylvestrene (B) by GC-MS



Figure 164 Mass spectra of terpin-4-ol (A) and authentic terpin-4-ol (B) by GC-MS



Figure 165 Mass spectra of terpin-4-ol-acetate (A) and authentic terpin-4-ol-acetate (B) by GC-MS



Figure 166 Mass spectra of terpinene $\langle \gamma \rangle$ (A) and authentic terpinene $\langle \gamma \rangle$ (B) by GC-MS



Figure 167 Mass spectra of terpinene $< \alpha ->$ (A) and authentic terpinene $< \alpha ->$ (B) by GC-MS



Figure 168 Mass spectra of terpineol $< \alpha ->$ (A) and authentic terpineol $< \alpha ->$ (B) by GC-MS



Figure 169 Mass spectra of terpinolene (A) and authentic terpinolene (B) by GC-MS



Figure 170 Mass spectra of terpinyl acetate $\langle \alpha \rangle$ (A) and authentic terpinyl acetate $\langle \alpha \rangle$ (B) by GC-MS



Figure 171 Mass spectra of tetradecanal (A) and authentic tetradecanal (B) by GC-MS



Figure 172 Mass spectra of thujene $< \alpha ->$ (A) and authentic thujene $< \alpha ->$ (B) by GC-MS



Figure 173 Mass spectra of thyjyl acetate $\langle 3 \rangle$ (A) and authentic thyjyl acetate $\langle 3 \rangle$ (B) by GC-MS



Figure 174 Mass spectra of tricyclene (A) and authentic tricyclene (B) by GC-MS



Figure 175 Mass spectra of undecanone < 2-> (A) and authentic undecanone < 2-> (B) by GC-MS



Figure 176 Mass spectra of valencene (A) and authentic valencene (B) by GC-MS



Figure 177 Mass spectra of verbenene (A) and authentic verbenene (B) by GC-MS



Figure 178 Mass spectra of verbenone (A) and authentic verbenone (B) by GC-MS



Figure 179 Mass spectra of verbenyl acetate $\langle cis \rangle$ (A) and authentic verbenyl acetate $\langle cis \rangle$ (B) by GC-MS



Figure 180 Mass spectra of vetivenol < *bicyclo->* (A) and authentic vetivenol < *bicyclo->* (B) by GC-MS



Figure 181 Mass spectra of viridiflorene (A) and authentic viridiflorene (B) by GC-MS

VITA

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