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## APPENDIX I

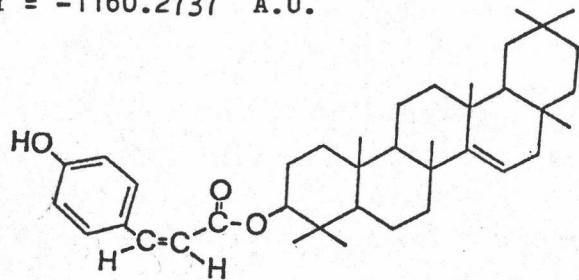
The semiempirical MO SCF were computed by CNDO/2 PROGRAMME at the IBM 3031/08 Computer Center of Chulalongkorn University.

**I. For Taraxeryl-cis-p-hydroxy-cinnamate (Compound 11)**

NUCLEAR REPULSION ENERGY = 963.2312 A.U.  
 TOTAL ENERGY = -197.0423 A.U.  
 BINDING ENERGY = -16.4969 A.U.  
 ELECTRONIC ENERGY = -1160.2737 A.U.

\*VALENCE ELECTRON DENSITY\*

1	C	3.8399				
2	C	4.0160				
3	C	4.0065				
4	C	3.9462				
5	C	3.9176	31	H	0.9939	
6	C	4.0398	32	H	0.9607	
7	O	6.2283	33	H	1.0000	
8	H	0.9649	34	H	1.0042	
9	H	0.6156	35	H	0.9937	
10	H	0.9821	36	H	1.0040	
11	H	0.9854	37	H	1.0094	
12	H	0.8619	38	H	0.9992	
13	C	4.1392	39	H	0.9463	
14	H	1.0076	40	H	1.0202	
15	C	4.2613	41	H	1.0128	
16	H	0.6306	42	H	1.0010	
17	C	4.1819	43	H	0.9701	
18	O	6.2549	44	H	1.0104	
19	O	6.2216	45	H	1.0207	
20	C	3.8049				
21	H	1.0479				
22	C	4.0519			DIPOLE MOMENTS	
23	C	3.9687				
24	C	4.0472	COMPONENTS	X	Y	Z
25	C	4.0089	DENSITIES	-0.51933	5.23556	0.16725
26	C	3.9867	S.P.	-0.63539	3.89430	-0.16232
27	H	1.0142	P.D.	0.00000	0.00000	0.00000
28	C	3.9601	TOTAL	-1.15471	9.12986	0.00494
29	C	4.0537	DIPOLE MOMENT=	9.20259	DEBYES	
30	C	3.9889				



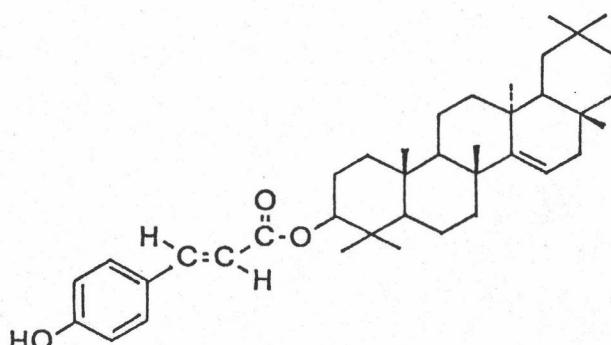
	COORDINATES		
	*X*	*Y*	*Z*
C	3.7360997	6.4709797	0.0000100
C	2.5262403	5.7725000	0.0000100
C	2.5262203	4.3754997	0.0000100
C	3.7360497	3.6769800	0.0000100
C	4.9459000	4.3754702	0.0000100
C	4.9459105	5.7724705	0.0000100
O	3.7361002	7.8309803	0.0000100
H	5.8812199	6.3124599	0.0000100
H	5.8811998	3.8354597	0.0000100
H	1.5909004	3.8355198	0.0000100
H	1.5909405	6.3125200	0.0000100
H	2.7278404	8.2180300	0.0000100
C	3.7360401	2.1569796	0.0000100
H	2.8007202	1.6169901	0.0000100
C	4.8939104	1.4894701	0.0000100
H	5.8292198	2.0284700	0.0000100
C	6.2016096	2.2434597	0.0000100
O	6.2016201	3.4764605	0.0000100
O	7.3776703	1.5644503	0.0000100
C	8.4291496	2.4159098	0.0000200
H	8.7093201	2.6427898	-1.0180397
C	9.0374899	3.7164402	0.7258500
C	9.6198196	1.7623997	0.7258500
C	9.2222900	1.4399204	2.1782904
C	10.0125999	0.4616400	0.0010500
C	10.8169093	2.7311897	0.7258000
H	11.0968800	2.9575005	-0.2931500
C	10.4258099	4.0321798	1.4504004
C	10.0271597	3.7110901	2.9028397
C	9.2343102	4.6855803	0.7258500
H	8.3831596	0.7600200	2.1799097
H	8.9472198	2.3516598	2.6876902
H	10.0572996	0.9816100	2.6873198
H	10.2921801	0.6868200	-1.0175400
H	9.1738701	-0.2187500	0.0015900
H	10.8476105	0.0033300	0.5100800
H	9.7522898	4.6233902	3.4113197
H	10.8615799	3.2526999	3.4127703
H	9.1876402	3.0316801	2.9044600
H	8.9588404	5.5973997	1.2348804
H	9.5122099	4.9124298	-0.2928300
H	7.7579098	3.4888401	1.7439003
H	7.2021904	4.1751099	0.2176300
H	11.6519203	2.2728901	1.2341099
H	11.2653303	4.7115893	1.4498596

**II. For Taraxeryl-trans-p-hydroxycinnamate (Compound 12)**

NUCLEAR REPULSION ENERGY = 886.9883 A.U.  
 TOTAL ENERGY = -199.1539 A.U.  
 BINDING ENERGY = -18.6086 A.U.  
 ELECTRONIC ENERGY = -1086.1423 A.U.

\*VALENCE ELECTRON DENSITY\*

1	C	3.8152				
2	C	4.0593				
3	C	3.9818				
4	C	3.9896				
5	C	3.9757				
6	C	4.0501				
7	C	3.2333				
8	H	0.9923				
9	H	0.9941				
10	H	1.0051				
11	H	0.9982				
12	H	0.8753				
13	C	3.9242				
14	H	1.0369				
15	C	4.4224				
16	H	0.5789				
17	C	3.6266				
18	O	6.2923				
19	O	6.2292				
20	C	3.8165				
21	H	1.0298				
22	C	4.0859				
23	C	3.9639				
24	C	4.0513				
25	C	4.0119				
26	C	3.9855				
27	H	1.0121				
28	C	3.9626				
29	C	4.0519				
30	C	3.9850				



DIPOLE MOMENTS

	COMPONENTS	X	Y	Z
DENSITIES	0.63009	-1.55317	-0.28227	
S.P.	-2.95125	-1.68784	0.03230	
P.D.	0.00000	0.00000	0.00000	
TOTAL	-2.32115	-3.24101	-0.24997	
DIPOLE MOMENT=	3.99429	DEBYES		

## COORDINATES

	*X*	*Y*	*Z*
C	3.7360897	6.4709797	0.0000100
C	2.5242403	5.7725000	0.0000100
C	2.5242203	4.3754997	0.0000100
C	3.7360497	3.6769800	0.0000100
C	4.9459000	4.3754702	0.0000100
C	4.9459105	5.7724705	0.0000100
C	3.7361002	7.8309803	0.0000100
H	5.8812199	6.3124599	0.0000100
H	5.8811993	3.8354597	0.0000100
H	1.5909004	3.6355198	0.0000100
H	1.5909405	6.3125200	0.0000100
H	2.7278404	9.2166300	0.0000100
C	3.7360401	2.1569796	0.0000100
C	2.8907202	1.6169901	0.0000100
C	4.8939104	1.4884701	0.0000100
C	4.8938999	0.4084700	0.0000100
C	4.8938999	-0.6215200	0.0000100
C	3.8246301	-0.6380100	0.0000100
C	6.0699501	-0.7005300	0.0000100
C	5.8582897	-2.0368700	0.0000100
C	5.8018904	-2.3929396	1.0180597
C	4.5361700	-2.3477300	-0.7258200
C	7.0195704	-2.7412796	-0.7258200
C	7.1000900	-2.2357702	-2.1782598
C	8.3424597	-2.4310797	-0.0016200
C	6.7790899	-4.2623796	-0.7258500
H	6.7230797	-4.6180000	0.2931800
C	5.4568501	-4.5741596	-1.4503498
C	5.5356102	-4.0493499	-2.9028101
C	4.2952499	-3.8499699	-0.7258200
H	7.2693501	-1.1691198	-2.1798897
H	6.1729603	-2.4534101	-2.6876497
H	7.9144897	-2.7297697	-2.6872902
H	8.2372200	-2.7827904	1.0173495
H	8.5123396	-1.3645296	-0.0015600
H	9.4588604	-2.9250903	-0.5100500
H	4.60930999	-4.2864704	-3.4112902
H	6.3497801	-4.5618200	-3.4127399
H	5.7042503	-3.0016203	-2.9044304
H	3.3478503	-4.0892999	-1.2348499
H	4.2377396	-4.2230501	0.2928600
H	4.5934896	-1.9919996	-1.7438693
H	3.7213001	-1.8538504	-0.2176100
H	7.5934896	-4.7563801	-1.2340803
H	5.12562099	-5.6409101	-1.4498301

**VITA**

Mr. Warinthorn Chavasiri was born on October 20, 1963 in Songkla province, Thailand. He received Bachelor Degree of Science of Chemistry at Chulalongkorn University in 1985. Since 1985, he has been a graduate student studying Organic Chemistry in Chulalongkorn University. During the study towards the Master's degree, he was awarded as a research assistance by Faculty of Science during 1985-1987 and supported a research grant for his Master degree's thesis from the Graduate School, Chulalongkorn University.

