

Supplementary Data

Oligostilbenoids from *Vatica lowii* King (Dipterocarpaceae) and their chemotaxonomic significance

Aisyah Salihah Kamarozaman^a, Jalifah Latip^{b*}, Yana M. Syah^c, Christian Paetz^d, Yoshinori Asakawa^e, Yamaguchi Kentaro^f

^a*Centre of Foundation Studies, Universiti Teknologi MARA, Cawangan Selangor, Kampus Dengkil, 43800 Dengkil, Selangor, Malaysia*

^b*Department of Chemical Sciences, Faculty of Science and Technology, Universiti Kebangsaan Malaysia, 43600 Bandar Baru Bangi, Selangor, Malaysia*

^c*Department of Chemistry, Bandung Institute of Technology (ITB), Jln. Ganesha 10, 40132 Bandung, Indonesia*

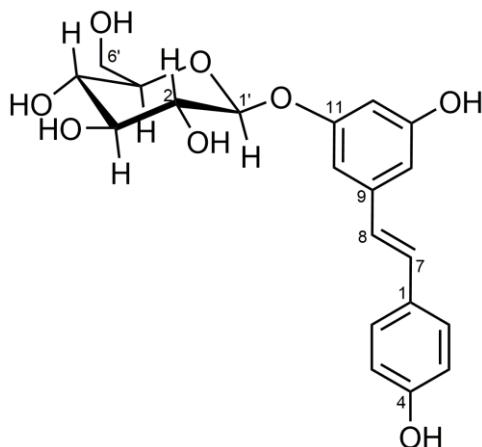
^d*Abteilung NMR, Max-Planck-Institute for Chemical Ecology, Hans-Knöll-Straße 8, 07745 Jena, Germany*

^e*Faculty of Pharmaceutical Sciences, Tokushima Bunri University, Nishihama Yamashiro-cho, Tokushima 770-8514, Japan*

^f*Faculty of Pharmaceutical Sciences, Tokushima Bunri University, Shido, Sanuki-city, Kagawa 769-2193, Japan*

*Corresponding author
E-mail: jalifah@ukm.edu.my

Piceid (1)



[α] _D	: -65° (c=0.1, MeOH)
Molecular formula	: C ₂₀ H ₂₂ O ₈
UV λ_{max} (MeOH)	: 225, 298 nm
IR ν_{max} (KBr)	: 3379 (OH); 1655 cm ⁻¹ (C=C Ar)

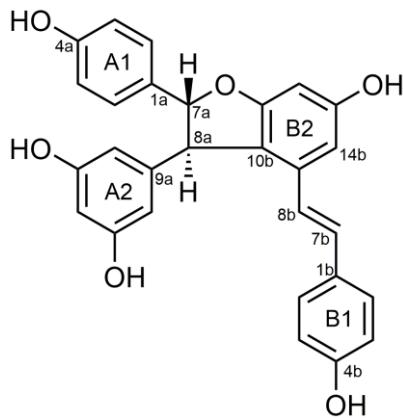
NMR data of compound 1

No	δ_{H} (mult., J in Hz)		δ_{C}	
	1*	1**	1*	1**
1	-	-	129.9	130.0
2,6	7.42 (d, 8.4)	7.39 (d, 8.6)	128.8	128.0
3,5	6.83 (d, 8.4)	6.72 (d, 8.6)	116.4	115.6
4	-	-	158.3	157.3
7	7.10 (d, 16.4)	7.05 (d, 17.2)	129.6	128.6
8	6.92 (d, 16.4)	6.85 (d, 17.2)	126.4	125.2
9	-	-	140.8	139.4
10	6.80 (t, 1.8)	6.70 (dd, 1.8, 1.8)	103.8	102.8
11	-	-	159.4	158.4
12	6.45 (t, 1.8)	6.32 (dd, 1.8, 1.8)	106.4	104.7
13	-	-	160.3	158.9
14	6.66 (t, 1.8)	6.55 (dd, 1.8, 1.8)	108.2	107.2
1'	4.93 (d, 7.7)	4.80 (d, 7.0)	102.0	100.7
2'	3.45 (m)	3.44 (m)	74.7	73.3
3'	3.51 (m)	3.50 (m)	77.9	76.7
4'	3.43 (m)	3.43 (m)	71.4	69.8
5'	3.54 (m)	3.52 (m)	78.1	77.2
6a'	3.92 (d, 11.7)	3.90 (d, 12.0)	62.7	60.7
6b'	3.72 (dd, 11.7, 1.8)	3.72 (d, 12.0, 1.8)	62.7	60.7

* measured in acetone-d₆

** measured in DMSO-d₆ (Orsini et al. 1997)

(-)- ε -viniferin (2)



$[\alpha]_D$: -63° (c=0.1, MeOH)
Molecular formula	: C ₂₈ H ₂₂ O ₆
UV λ_{max} (MeOH)	: 225, 323 nm
IR ν_{max} (KBr)	: 3366 (OH); 1604, 1444 cm ⁻¹ (C=C Ar)

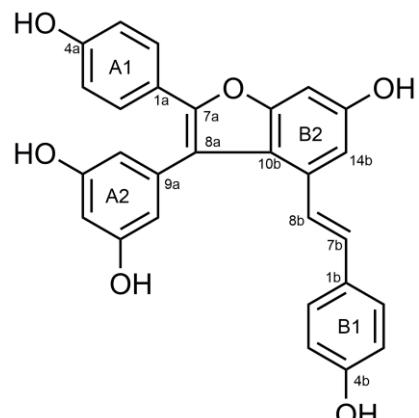
NMR data of compound 2

No	δ_{H} (mult., J in Hz)		δ_{C}	
	2'	2''	2'	2''
1a	-	-	133.7	133.8
2a,6a	7.21 (d, 8.8)	7.19 (d, 8.0)	127.9	127.8
3a,5a	6.83 (d, 8.8)	6.80 (d, 8.0)	116.1	116.0
4a	-	-	158.3	158.0
7a	5.42 (d, 5.4)	5.40 (d, 5.0)	93.9	93.8
8a	4.48 (d, 5.4)	4.44 (d, 5.0)	57.1	57.0
9a	-	-	147.4	147.3
10a,14a	6.23 (br s)	6.22 (br s)	106.9	106.9
11a	-	-	159.9	159.7
12a	6.23 (br s)	6.22 (br s)	102.1	101.9
13a	-	-	159.9	159.7
1b	-	-	129.8	130.0
2b,6b	7.18 (d, 8.8)	7.15 (d, 8.0)	128.7	128.6
3b,5b	6.72 (d, 8.8)	6.71 (d, 8.0)	116.3	116.1
4b	-	-	158.3	158.0
7b	6.92 (d, 16.4)	6.91 (d, 16.0)	123.4	123.3
8b	6.71 (d, 16.4)	6.65 (d, 16.0)	130.0	130.1
9b	-	-	136.3	136.3
10b	-	-	119.8	119.7
11b	-	-	162.4	162.3
12b	6.32 (d, 1.8)	6.30 (d, 2.0)	96.8	96.7
13b	-	-	159.6	159.4
14b	6.73 (d, 1.8)	6.70 (d, 2.0)	104.1	104.1

* measured in acetone-d₆

** measured in acetone-d₆ (Oshima et al. 1995)

Viniferifuran (3)



Molecular formula

: C₂₈H₂₀O₆

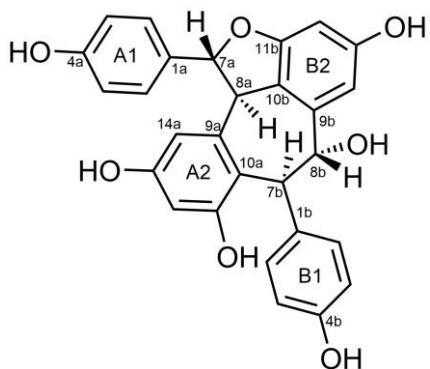
NMR data of compound 3

No	δ_{H} (mult., J in Hz)		δ_{C}	
	3*	3**	3*	3**
1a	-	-	123.8	123.8
2a,6a	7.42 (d, 8.8)	7.42 (d, 8.8)	128.7	128.4
3a,5a	6.82 (d, 8.8)	6.69 (d, 8.8)	116.2	116.2
4a	-	-	160.0	158.4
7a	-	-	154.0	150.6
8a	-	-	117.0	117.4
9a	-	-	140.8	138.7
10a	6.52 (d, 2.2)	6.40 (d, 2.2)	108.6	110.2
11a	-	-	159.7	160.6
12a	6.25 (t, 2.2)	6.47 (t, 2.2)	105.6	103.1
13a	-	-	159.7	160.6
14a	6.52 (d, 2.2)	6.40 (d, 2.2)	108.6	110.2
1b	-	-	131.9	130.7
2b,6b	7.07 (d, 8.8)	6.98 (d, 8.8)	129.1	128.8
3b,5b	6.81 (d, 8.8)	6.65 (d, 8.8)	116.4	116.3
4b	-	-	158.3	158.2
7b	6.98 (d, 16.1)	6.94 (d, 16.3)	123.3	123.3
8b	6.85 (d, 16.1)	6.85 (d, 16.3)	129.9	129.3
9b	-	-	133.1	133.3
10b	-	-	122.5	122.5
11b	-	-	158.2	156.4
12b	6.83 (d, 1.8)	6.80 (d, 2.0)	102.7	97.3
13b	-	-	158.2	156.5
14b	7.34 (d, 1.8)	6.99 (d, 2.0)	105.6	107.4

* measured in acetone-d₆

** measured in methanol-d₄ (Ito et al. 1999)

Hemsleyanol A (4)



$[\alpha]_D$: +12° (c=0.1, MeOH)
Molecular formula	: C ₂₈ H ₂₂ O ₇
UL λ_{max} (MeOH)	: 220, 232, 284 nm
IR ν_{max} (KBr)	: 3302 (OH); 1611, 1454 cm ⁻¹ (C=C Ar)

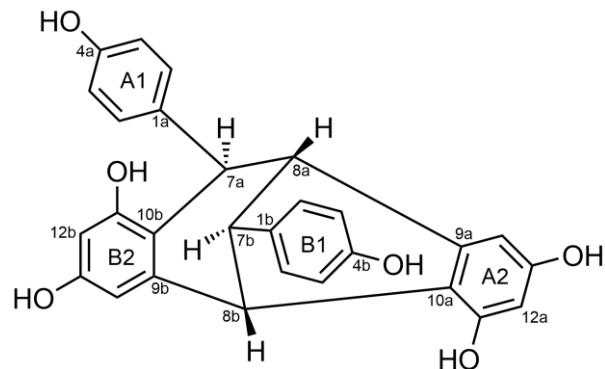
NMR data of compound 4

No	δ_{H} (mult., J in Hz)		δ_{C}	
	4*	4**	4*	4**
1a	-	-	134.0	133.8
2a,6a	7.48 (d, 8.3)	7.48 (d, 8.3)	130.7	130.4
3a,5a	6.91 (d, 8.3)	6.94 (d, 8.3)	116.7	116.4
4a	-	-	155.8	155.6
7a	5.74 (d, 9.6)	5.75 (d, 9.8)	94.3	93.3
8a	5.38 (d, 9.6)	5.41 (d, 9.8)	52.4	51.9
9a	-	-	141.8	141.4
10a	-	-	118.7	118.4
11a	-	-	158.8	158.3
12a	6.15 (br s)	6.24 (br s)	102.1	102.0
13a	-	-	156.8	156.6
14a	6.14 (br s)	6.23 (br s)	106.6	106.6
1b	-	-	135.7	135.2
2b,6b	6.66 (d, 8.3)	6.72 (d, 8.3)	130.8	130.5
3b,5b	6.44 (d, 8.3)	6.48 (d, 8.3)	115.1	114.9
4b	-	-	159.0	158.5
7b	5.04 (d, 5.8)	5.07 (d, 5.9)	49.5	49.3
8b	4.70 (d, 5.8)	4.76 (br d)	79.0	78.4
8b (OH)	4.68 (br s)	4.45 (br s)	79.0	78.4
9b	-	-	140.6	140.9
10b	-	-	118.5	118.1
11b	-	-	161.0	160.7
12b	6.00 (d, 1.9)	6.04 (d, 2.0)	96.6	96.2
13b	-	-	159.0	158.5
14b	5.67 (d, 1.9)	5.70 (d, 2.0)	108.7	108.2

* measured in methanol-d₄

** measured in acetone -d₆ (Ito et al. 2000b)

(-)-Ampelopsin F (5)



$[\alpha]_D$: -28° (c=0.1, MeOH)
Molecular formula	: C ₂₈ H ₂₂ O ₆
UL λ_{max} (MeOH)	: 220, 232, 284 nm
IR ν_{max} (KBr)	: 3302 (OH); 1611, 1454 cm ⁻¹ (C=C Ar)

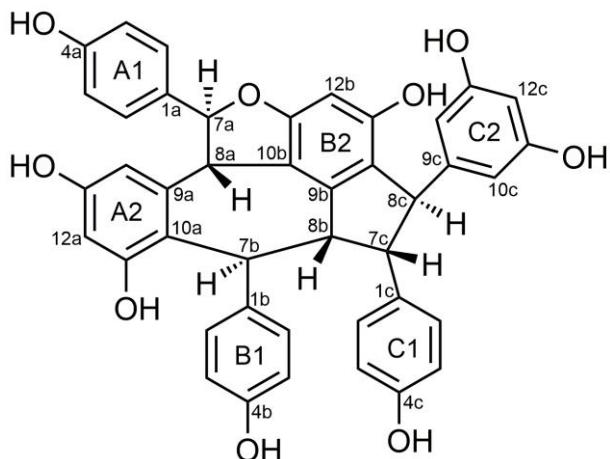
NMR data of compound 5

No	δ_{H} (mult., J in Hz)		δ_{C}	
	5'	5''	5'	5''
1a	-	-	139.1	138.4
2a,6a	7.01 (d, 8.8)	7.08 (d, 8.5)	130.1	129.8
3a,5a	6.68 (d, 8.8)	6.76 (d, 8.5)	115.7	115.5
4a	-	-	156.2	156.1
7a	4.08 (d, 1.5)	4.18 (d, 1.5)	47.6	47.1
8a	3.21 (br s)	3.34 (br s)	59.3	58.2
9a	-	-	147.5	147.2
10a	-	-	128.7	127.7
11a	-	-	153.2	153.1
12a	6.00 (d, 1.8)	6.07 (d, 1.9)	101.9	101.7
13a	-	-	158.5	158.5
14a	6.37 (d, 1.8)	6.51 (d, 1.9)	104.3	104.0
1b	-	-	136.1	135.3
2b	6.72 (d, 8.4)	6.78 (d, 8.5)	129.4	129.1
3b	6.50 (d, 8.4)	6.58 (d, 8.5)	115.6	115.4
4b	-	-	156.1	156.1
7b	3.55 (br s)	3.64 (br s)	50.9	50.4
8b	4.01 (br s)	4.12 (br s)	49.8	49.6
9b	-	-	147.8	147.5
10b	-	-	114.0	113.2
11b		-	158.1	157.7
12b	6.07 (d, 1.8)	6.17 (d, 2.4)	101.9	101.6
13b	-	-	157.1	157.0
14b	6.34 (d, 1.8)	6.44 (d, 2.4)	105.7	105.5

* measured in methanol-d₄

** measured in acetone-d₆ (Luo et al. 2001)

Vaticanol A (6)



[α]_D : -169° (c=0.1, MeOH)

Molecular formula : C₄₂H₃₂O₉

UV λ_{max} (MeOH) : 216, 236, 284 nm

IR ν_{max} (KBr) : 3369 (OH); 1612, 1451 cm⁻¹ (C=C Ar)

NMR data of compound 6

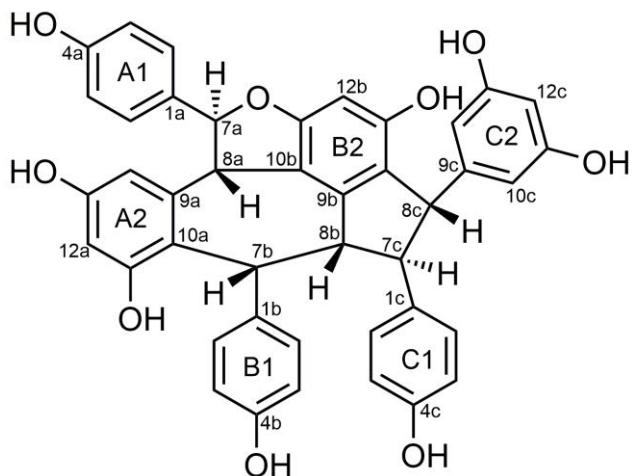
No	δ_{H} (mult., J in Hz)		δ_{C}	
	6'	6''	6'	6''
1a	-	-	134.5	134.4
2a,6a	7.27 (d, 8.8)	7.28 (d, 8.8)	128.1	128.0
3a,5a	6.82 (d, 8.8)	6.83 (d, 8.8)	116.1	116.0
4a	8.45 (br s)	8.42 (br s)	157.9	157.9
7a	6.16 (d, 4.0)	6.18 (br d, 3.9)	86.7	86.5
8a	4.47 (d, 4.0)	4.51 (d, 3.9)	50.4	50.3
9a	-	-	144.8	144.7
10a	-	-	119.4	119.3
11a	7.38 (br s)	7.31 (br s)	158.0	157.7
12a	6.07 (d, 2.2)	6.09 (d, 2.4)	101.4	101.3
13a	7.92 (br s)	7.95 (br s)	156.5	156.3
14a	6.46 (d, 2.2)	6.48 (d, 2.4)	103.4	103.3
1b	-	-	138.8	138.7
2b,6b	7.05 (d, 8.8)	7.07 (d, 8.8)	129.4	129.2
3b,5b	6.58 (d, 8.8)	6.60 (d, 8.8)	115.4	115.4
4b	8.03 (br s)	8.00 (br s)	155.9	155.7
7b	5.14 (br s)	5.17 (br s)	36.1	36.0
8b	4.49 (d, 7.3)	4.52 (d, 7.3)	48.7	48.6
9b	-	-	145.0	144.9
10b	-	-	118.7	118.6
11b	-	-	160.0	159.9
12b	6.20 (br s)	6.22 (s)	95.4	95.3
13b	8.02 (br s)	7.97 (br s)	155.6	155.4

14b	-	-	122.3	122.2
1c	-	-	135.9	135.8
2c,6c	6.53 (d, 8.8)	6.55 (d, 8.8)	129.7	129.6
3c,5c	6.35 (d, 8.8)	6.37 (d, 8.8)	115.9	114.9
4c	7.90 (br s)	7.89 (br s)	156.6	156.4
7c	3.62 (d, 7.3)	3.65 (d, 7.3)	64.5	64.3
8c	4.16 (br s)	4.20 (br s)	57.6	57.5
9c	-	-	147.6	147.5
10c,14c	6.24 (d, 2.2)	6.27 (d, 2.0)	106.8	106.7
11c,13c	8.10 (br s)	8.09 (br s)	159.4	159.2
12c	6.18 (t, 2.2)	6.21 (t, 2.0)	101.3	101.2

* measured in acetone-*d*₆

** measured in acetone-*d*₆ (Tanaka et al. 2000a)

Vaticanol E (7)



[α]_D : +191° (c=0.1, MeOH)

Molecular formula : C₄₂H₃₂O₉

UV λ_{max} (MeOH) : 216, 292 nm

IR ν_{max} (KBr) : 3295 (OH); 1606, 1453 cm⁻¹ (C=C Ar)

NMR data of compound 7

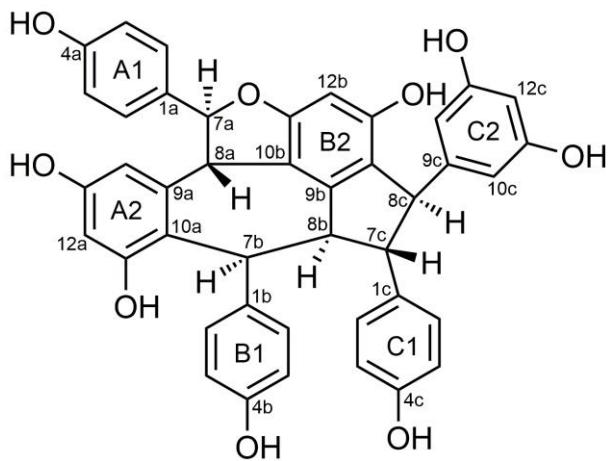
No	δ_{H} (mult., J in Hz)		δ_{C}	
	7'	7''	7'	7''
1a	-	-	133.6	133.5
2a,6a	7.50 (d, 8.3)	7.77 (d, 8.3)	130.9	130.6
3a,5a	6.93 (d, 8.3)	6.99 (d, 8.3)	116.7	116.4
4a	-	-	159.0	158.6
7a	5.63 (d, 9.3)	5.66 (d, 9.3)	96.4	95.6
8a	5.38 (d, 9.3)	5.35 (d, 9.3)	52.6	52.3
9a	-	-	140.6	140.4
10a	-	-	124.2	123.5
11a	-	-	158.5	157.9

12a	6.03 (d, 2.3)	6.15 (d, 2.0)	102.0	102.0
13a	-	-	156.2	156.3
14a	6.02 (d, 2.3)	6.12 (br d, 2.0)	108.0	108.0
1b	-	-	133.5	132.8
2b,6b	5.97 (d, 8.5)	6.05 (d, 8.6)	133.3	132.9
3b,5b	6.27 (d, 8.5)	6.32 (d, 8.6)	113.8	113.7
4b	-	-	155.8	155.9
7b	4.79 (d, 2.8)	4.83 (d, 2.9)	42.9	42.4
8b	4.28 (dd, 9.8, 2.8)	4.30 (dd, 8.3, 2.9)	51.7	51.2
9b	-	-	144.7	144.1
10b	6.19 (d, 2.0)	6.30 (d, 2.0)	116.3	116.0
11b	-	-	160.4	160.4
12b	6.05 (s)	6.08 (s)	96.5	96.3
13b	-	-	155.8	155.6
14b	-	-	124.7	123.8
1c	-	-	130.9	130.3
2c,6c	7.14 (d, 8.3)	7.20 (d, 8.6)	131.6	131.3
3c,5c	6.79 (d, 8.3)	6.85 (d, 8.6)	115.5	115.4
4c	-	-	156.2	156.3
7c	3.94 (t, 9.8)	4.07 (dd, 10.3, 8.3)	58.6	58.1
8c	3.51 (t, 9.8)	3.59 (d, 10.3)	51.6	51.2
9c	-	-	148.9	148.1
10c,14c	6.19 (d, 2.0)	6.30 (d, 2.0)	108.0	107.6
11c,13c	-	-	159.0	159.3
12c	6.03 (t, 2.3)	6.13 (t, 2.0)	101.3	101.7

* measured in methanol-*d*₄

** measured in acetone-*d*₆ (Ito et al. 2000a)

Pauciflorol B (8)



[α]_D : +20° (c=0.1, MeOH)

Molecular formula : C₄₂H₃₂O₉

UV λ_{max} (MeOH) : 216, 296 nm

IR ν_{max} (KBr) : 3181 (OH); 1598, 1451 cm⁻¹ (C=C Ar)

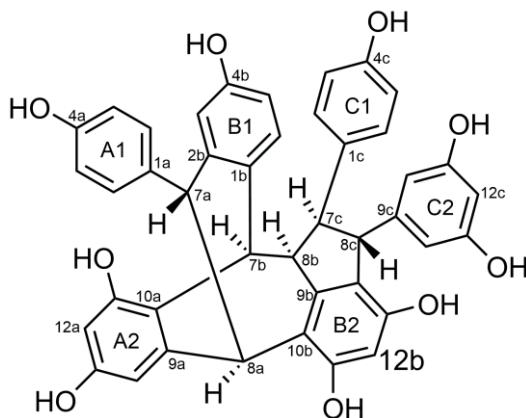
NMR data of compound **8**

No	δ_H (mult., J in Hz)		δ_C	
	8[*]	8^{**}	8[*]	8^{**}
1a	-	-	130.9	130.1
2a,6a	7.16 (d, 8.5)	7.26 (d, 8.6)	130.5	129.5
3a,5a	6.74 (d, 8.5)	6.81 (d, 8.6)	116.5	115.4
4a	-	-	158.9	157.9
7a	5.75 (d, 11.6)	5.84 (d, 11.7)	91.2	89.9
8a	4.30 (d, 11.6)	4.47 (d, 11.9)	~48.6	48.1
9a	-	-	141.9	141.0
10a	-	-	125.7	124.1
11a	-	-	156.3	155.1
12a	6.22 (d, 2.0)	6.35 (d, 2.0)	101.4	101.0
13a	-	-	156.8	156.1
14a	5.98 (br d, 2.0)	6.17 (br d, 2.2)	105.7	105.3
1b	-	-	134.0	132.7
2b,6b	7.13 (d, 8.0)	7.19 (d, 8.5)	130.6	129.8
3b,5b	6.64 (d, 8.0)	6.69 (d, 8.5)	115.5	114.8
4b	-	-	155.8	155.1
7b	5.16 (d, 3.3)	5.28 (d, 3.3)	37.8	36.8
8b	3.65 (br d, 11.6)	3.66 (m)	52.8	51.8
9b	-	-	144.8	143.4
10b	-	-	116.2	115.3
11b	-	-	159.5	158.8
12b	6.08 (s)	6.15 (s)	96.7	96.0
13b	-	-	154.9	154.0
14b	-	-	121.5	120.3
1c	-	-	133.4	132.1
2c,6c	6.96 (d, 8.0)	7.02 (d, 8.5)	130.3	129.3
3c,5c	6.69 (d, 8.0)	6.74 (d, 8.5)	115.9	115.2
4c	-	-	156.8	156.1
7c	3.70 (dd, 11.6, 9.3)	3.76 (dd, 11.6, 9.3)	62.9	61.4
8c	4.17 (d, 9.3)	4.25 (d, 9.3)	58.4	56.7
9c	-	-	147.5	146.1
10c,14c	6.12 (s)	6.20 (s)	106.5	106.7
11c,13c	-	-	159.0	158.5
12c	6.10 (t, 2.0)	6.19 (s)	101.6	101.1

* measured in methanol-*d*₄

** measured in acetone-*d*₆ (Ito et al. 2003)

Vaticanol G (9)



[α]_D : +149° (c=0.1, MeOH)

Molecular formula : C₄₂H₃₂O₉

UV λ_{max} (MeOH) : 216, 236, 280 nm

IR ν_{max} (KBr) : 3368 (OH); 1611, 1449 cm⁻¹ (C=C Ar)

NMR data of compound 9

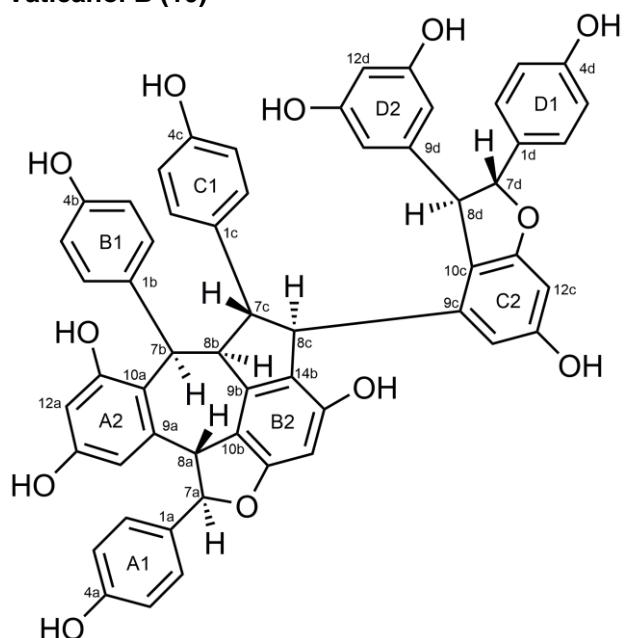
No	δ_{H} (mult., J in Hz)		δ_{C}	
	9*	9**	9*	9**
1a	-	-	140.0	140.0
2a,6a	6.42 (br s)	6.42 (br s)	130.3	130.0
3a,5a	6.45 (br s)	6.45 (br s)	114.8	114.8
4a	8.00 (br s)	7.93 (br s)	158.6	155.6
7a	4.52 (d, 3.7)	4.55 (d, 3.4)	57.2	57.2
8a	4.60 (d, 3.7)	4.63 (d, 3.4)	50.4	50.3
9a	-	-	142.0	142.0
10a	-	-	126.1	126.2
11a	8.16 (br s)	8.08 (br s)	153.3	153.3
12a	6.17 (d, 2.2)	6.20 (d, 2.4)	101.7	101.8
13a	7.72 (br s)	7.68 (br s)	156.0	155.9
14a	5.64 (d, 2.6)	5.67 (d, 2.4)	111.5	111.5
1b	-	-	129.3	129.3
2b	-	-	141.9	141.8
3b	6.05 (d, 2.6)	6.08 (d, 2.4)	119.5	119.5
4b	8.00 (br s)	7.49 (br s)	154.9	154.9
5b	5.74 (dd, 8.4, 2.9)	5.76 (dd, 8.3, 2.4)	112.9	112.9
6b	5.99 (d, 8.8)	6.02 (d, 8.3)	135.1	135.1
7b	4.86 (d, 2.6)	4.89 (d, 2.0)	42.8	42.8
8b	3.83 (dd, 7.0, 2.2)	3.86 (dd, 7.0, 2.0)	53.9	53.9
9b	-	-	147.1	147.1
10b	-	-	117.6	117.7
11b	8.63 (br s)	8.54 (br s)	155.1	155.1
12b	6.45 (br s)	6.47 (s)	102.0	102.0
13b	7.71 (br s)	7.67 (br s)	153.1	153.1

14b	-	-	121.9	122.0
1c	-	-	137.2	137.2
2c	5.92 (br s)	5.92 (br s)	127.6	127.5
3c	5.97 (br s)	5.98 (br s)	114.3	114.3
4c	8.00 (br s)	7.94 (br s)	156.6	156.5
5c	6.66 (br s)	6.68 (br s)	116.2	116.2
6c	7.13 (br s)	7.12 (br s)	130.7	130.7
7c	3.49 (d, 7.0)	3.52 (d, 7.0)	63.1	63.0
8c	4.07 (br s)	4.11 (s)	57.1	57.1
9c	-	-	147.8	147.7
10c,14c	5.93 (d, 2.2)	5.97 (d, 2.4)	106.5	106.5
11c,13c	8.08 (br s)	8.04 (br s)	159.2	159.1
12c	6.08 (t, 2.2)	6.12 (t, 2.4)	101.1	101.1

* measured in acetone-*d*₆

** measured in acetone-*d*₆ (Ito et al. 2001)

Vaticanol B (10)



$[\alpha]_D$: -28° (c=0.1, MeOH)

Molecular formula : C₅₆H₄₂O₁₂

UV λ_{max} (MeOH) : 216, 236, 284 nm

IR ν_{max} (KBr) : 3368 (OH); 1614, 1450 cm⁻¹ (C=C Ar)

NMR data of compound 10

No	δ_{H} (mult., J in Hz)		δ_{C}	
	10*	10**	10*	10**
1a	-	-	130.8	130.8
2a,6a	7.22 (d, 8.8)	7.23 (d, 8.8)	130.2	130.2
3a,5a	6.79 (d, 8.8)	6.79 (d, 8.8)	115.4	116.0

4a-OH	8.54 (s)	8.48 (s)	158.6	158.5
7a	5.76 (d, 11.7)	5.77 (d, 11.7)	90.5	90.4
8a	4.42 (d, 11.0)	4.44 (d, 11.7)	48.9	48.8
9a	-	-	141.8	141.8
10a	-	-	124.5	124.5
11a-OH	8.21 (s)	8.17 (s)	155.7	155.7
12a	6.27 (d, 2.2)	6.29 (d, 2.0)	101.6	101.6
13a-OH	8.08 (s)	8.06 (s)	156.8	156.7
14a	6.16 (d, 2.2)	6.12 (d, 2.0)	105.8	105.8
1b	-	-	133.5	133.5
2b,6b	7.15 (d, 8.4)	7.17 (d, 8.8)	130.7	130.7
3b,5b	6.68 (d, 8.4)	6.70 (d, 8.8)	115.5	115.5
4b-OH	8.21 (s)	8.16 (s)	155.9	155.9
7b	5.18 (d, 3.3)	5.21 (d, 3.9)	37.1	37.1
8b	3.09 (dd, 11.3, 3.3)	3.13 (dd, 11.3, 3.9)	53.2	53.1
9b	-	-	143.3	143.2
10b	-	-	115.9	115.7
11b	-	-	158.8	158.8
12b	6.02 (s)	6.05 (s)	96.5	96.5
13b-OH	7.58 (s)	7.45 (s)	155.0	154.9
14b	-	-	122.2	122.1
1c	-	-	131.4	131.4
2c,6c	6.37 (d, 8.4)	6.42 (d, 8.8)	129.3	129.2
3c,5c	6.49 (d, 8.8)	6.52 (d, 8.8)	116.0	115.8
4c-OH	7.94 (s)	7.90 (s)	156.4	156.3
7c	4.09 (dd, 11.3, 10.1)	4.10 (dd, 11.3, 10.7)	57.6	57.6
8c	4.54 (d, 10.1)	4.55 (d, 10.7)	49.2	49.3
9c	-	-	141.7	141.6
10c	-	-	123.5	123.3
11c	-	-	161.7	161.7
12c	6.25 (d, 2.2)	6.21 (d, 2.0)	95.6	95.6
13c-OH	8.22 (s)	8.19 (s)	159.4	159.4
14c	6.45 (d, 2.2)	6.49 (d, 2.0)	107.0	107.0
1d	-	-	134.7	134.7
2d,6d	7.18 (d, 8.8)	7.19 (d, 8.8)	128.3	128.2
3d,5d	6.76 (d, 8.8)	6.78 (d, 8.8)	116.1	115.9
4d-OH	8.42 (s)	8.37 (s)	158.0	157.9
7d	5.35 (d, 5.1)	5.38 (d, 4.7)	94.7	94.6
8d	4.67 (d, 5.1)	4.68 (d, 4.7)	57.6	57.5
9d	-	-	148.1	147.9
10d,14d	6.09 (d, 2.5)	6.10 (d, 2.4)	107.6	107.5
11d,13d-OH	8.05 (s)	8.01 (s)	159.9	159.8
12d	6.26 (t, 2.2)	6.30 (t, 2.4)	102.2	102.2

* measured in acetone- d_6

** measured in acetone- d_6 (Tanaka et al. 2000a)