SUPPLEMENTARY MATERIAL

A new dimeric alkylresorcinol from the stem barks of *Swintonia* floribunda (Anacardiaceae)

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ABSTRACT

From an EtOAc-soluble fraction of the stem barks of *Swintonia floribunda* (Anacardiaceae), one new dimeric alkylresorcinol named integracin E (1), together with 4 known compounds (2–5) were isolated. Their chemical structures were elucidated based on the spectroscopic data interpretation. The absolute configuration of 1 was determined by the specific rotation analysis of its acid-catalyzed hydrolysis product. Compound 1 showed potent tyrosinase inhibitory activity with an IC₅₀ value of $48.2 \mu M$.

Keywords: Swintonia floribunda, Anacardiaceae, alkylresorcinol, tyrosinase inhibitory.

Table S1. NMR Spectroscopic Data for Compound 1 in $CDCl_3$.

	Integracin E (1)	
Position	$\delta_{\rm C}$, type	$\frac{\delta_{\rm H} (J \text{ in Hz})}{\delta_{\rm H} (J \text{ in Hz})}$
1	105.5, C	OH (J III IIZ)
2	165.4, C	_
3	103.4, C 101.6, CH	- 6 20 d (2 6)
3 4	*	6.29, d (2.6)
	160.4, C	-
5	111.0, CH	6.24, d (2.6)
6	149.1, C	- 2.06 1: (12.0 7.0)
7	37.1, CH ₂	2.86, dt (12.9, 7.9)
	22 4 677	2.81, dt (12.9, 7.9)
8	32.4, CH ₂	1.53, m
9–16	29.0–29.9, CH ₂	1.22–1.45
17	32.1, CH_2	1.26, m
18	22.8, CH_2	1.29, m
19	14.3, CH ₃	0.88, t (6.9)
1'	146.1, C	_
2′/6′	108.2, CH	6.21, d (2.1)
3′/5′	156.8, C	_
4'	100.4, CH	6.18, brs
7'	35.9, CH ₂	2.44, t (7.7)
8′	30.9, CH ₂	1.53, m
9′	$30.2, CH_2$	1.35, m
10′, 11′	29.0–29.9, CH ₂	1.22-1.45
12′	25.5, CH ₂	1.36, m
13′	34.4, CH ₂	1.65, m
14′	76.0, CH	5.26, m
15′	36.8, CH ₂	1.65, m
16′	19.0, CH_2	· · · · · · · · · · · · · · · · · · ·
17'	14.1, CH ₃	0.94, t (7.3)
	171.7, C=O	_
2-OH		12.13, brs

Figure S1. Significant HMBC correlations observed for compound 1.

$$\begin{array}{c} C_{17}H_{27}O_{2}^{+}\\ m/z\ 263.1968\\ HO_{5}^{-} \\ OH_{2}^{-} \\$$

Figure S2. MS fragmentation of 1.

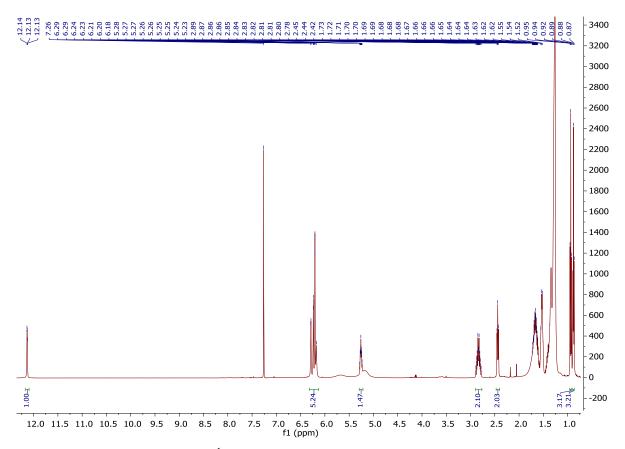


Figure S3. ¹H NMR spectrum of 1 (500 MHz, CDCl₃).

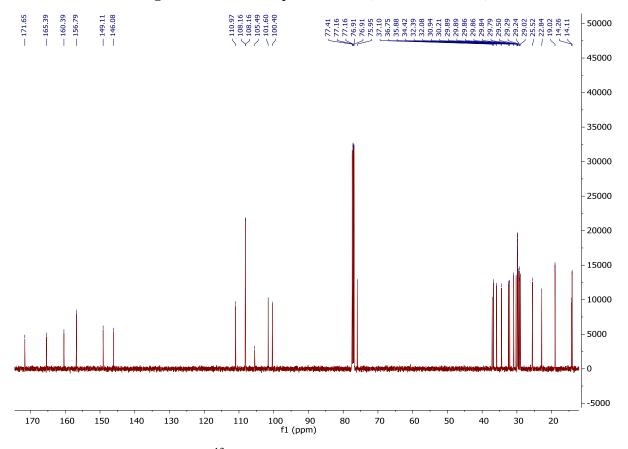


Figure S4. ¹³C NMR spectrum of 1 (125 MHz, CDCl₃).

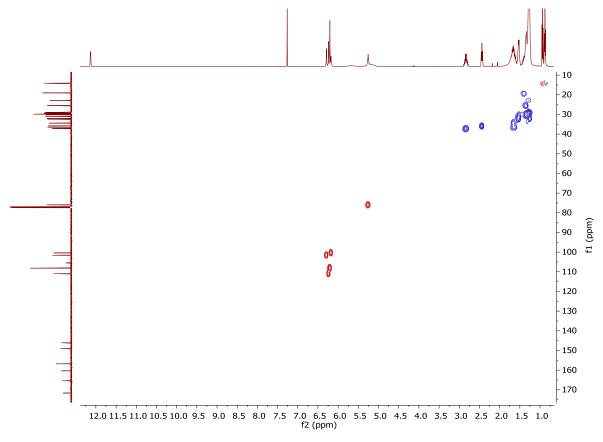


Figure S5. HSQC NMR spectrum of 1.

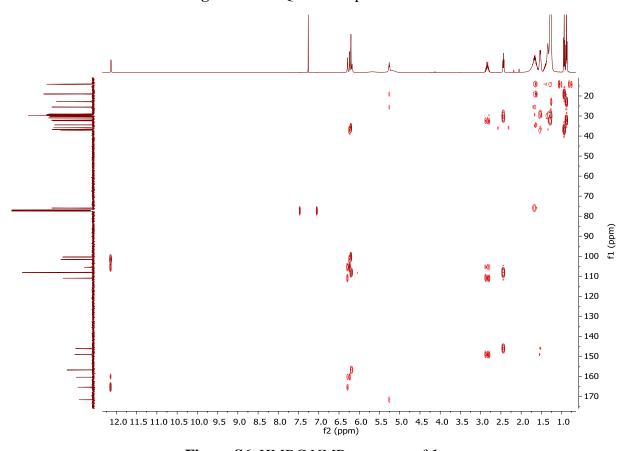


Figure S6. HMBC NMR spectrum of **1**.

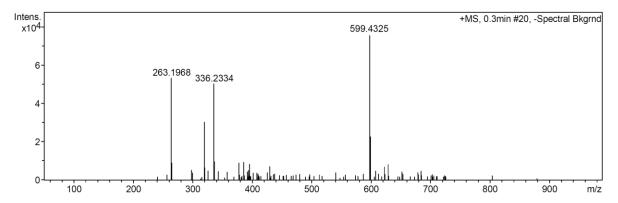


Figure S7. HRESIMS of 1.

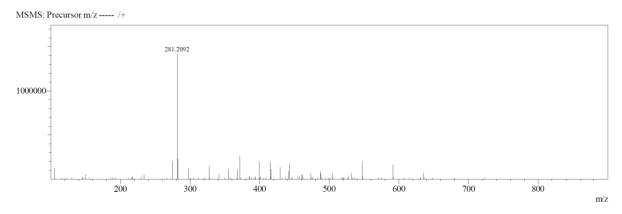


Figure S8. HRESIMS of 1a.

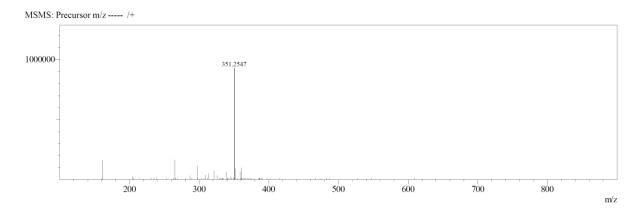


Figure S9. HRESIMS of 1b.