

## Supporting Information

*Rec. Nat. Prod.* X:X (2021) XX-XX

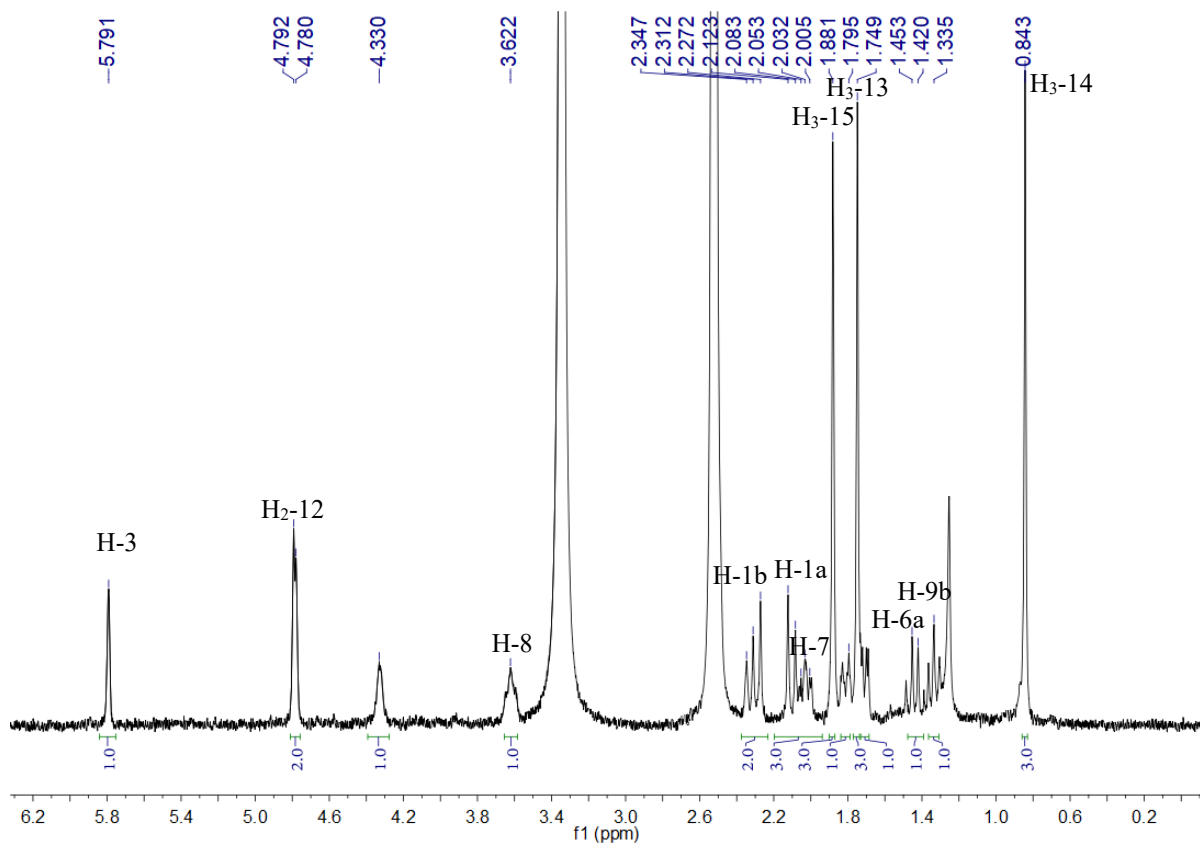
### Eudesmane Sesquiterpenoids from *Salvia plebeia*

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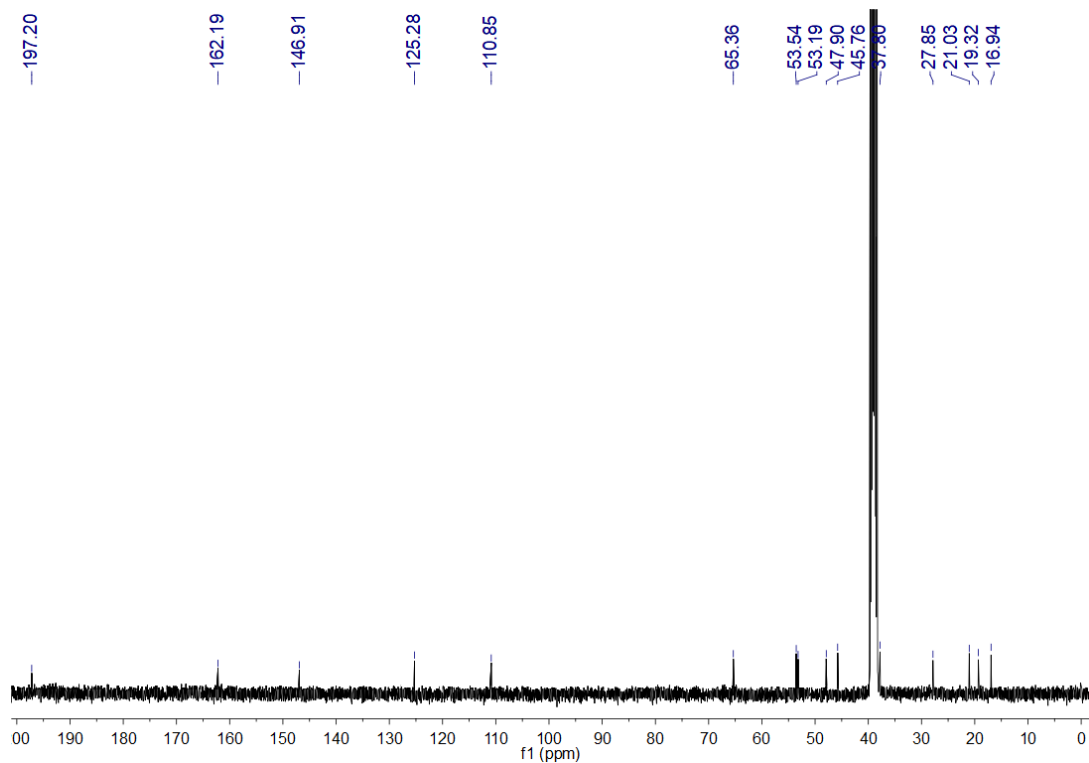
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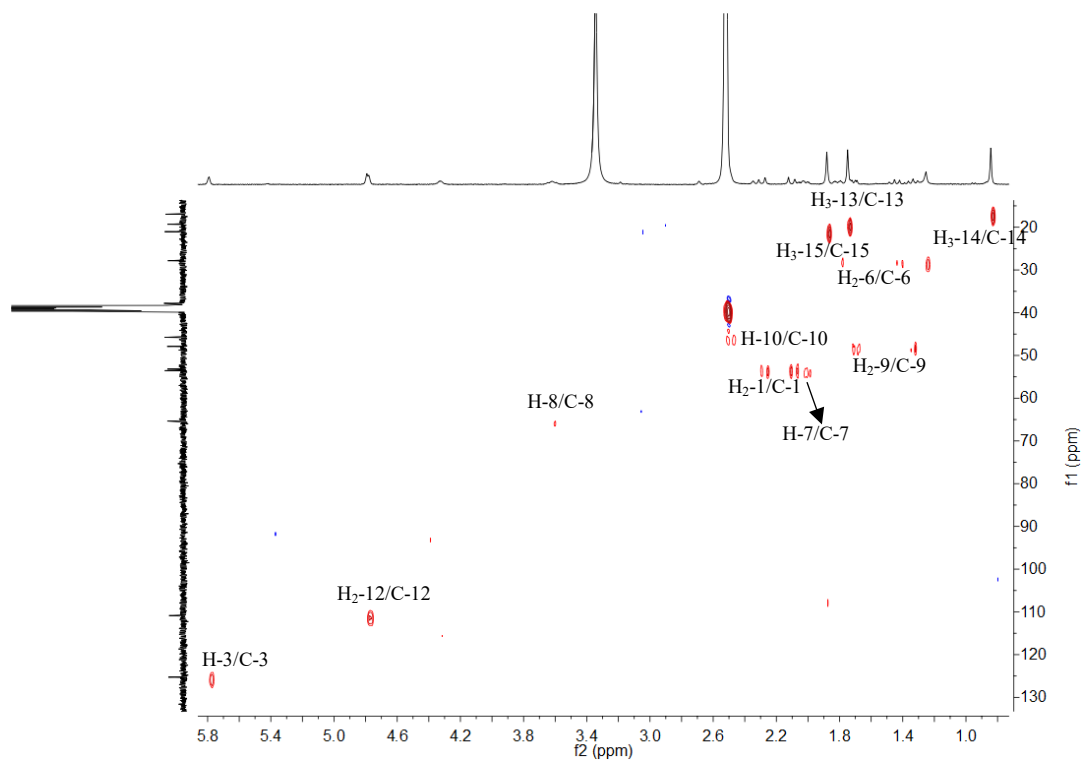
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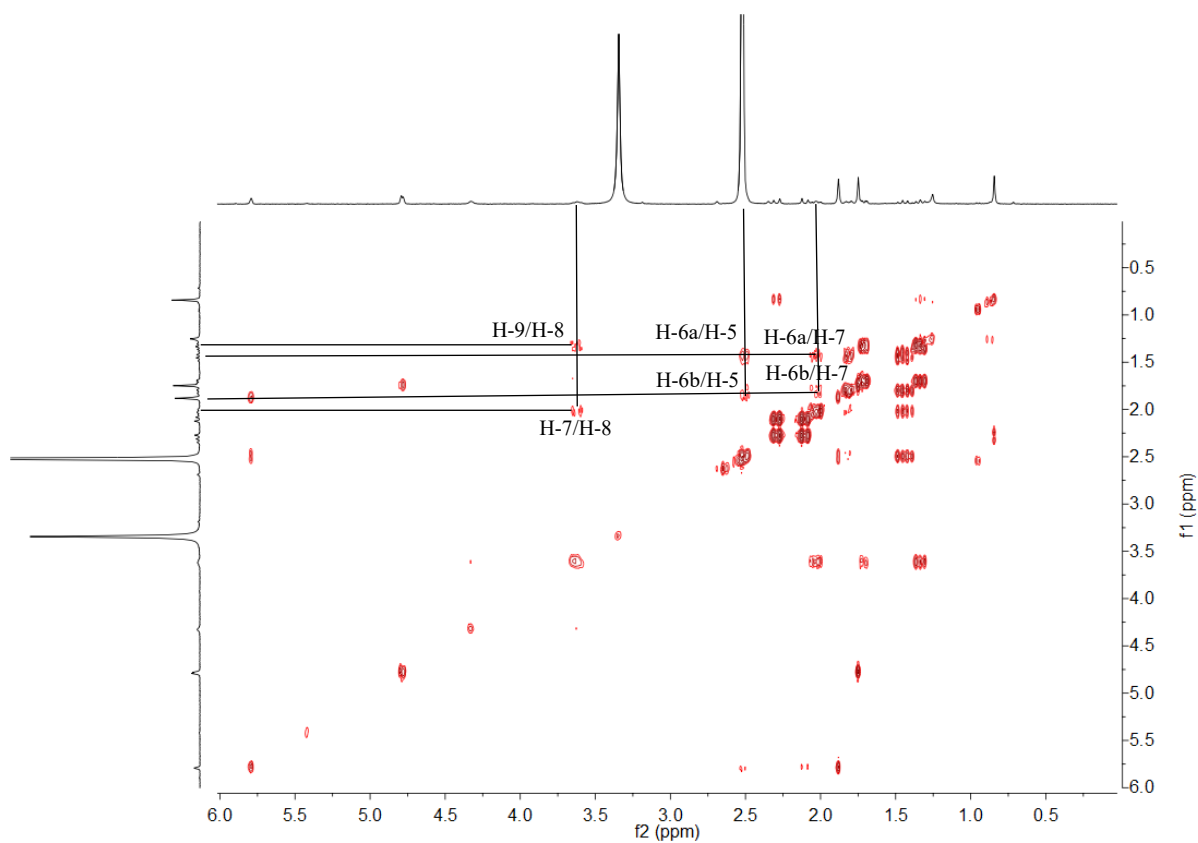
**Figure S1:** <sup>1</sup>H NMR Spectrum of **1** in DMSO-*d*<sub>6</sub> (400 MHz)



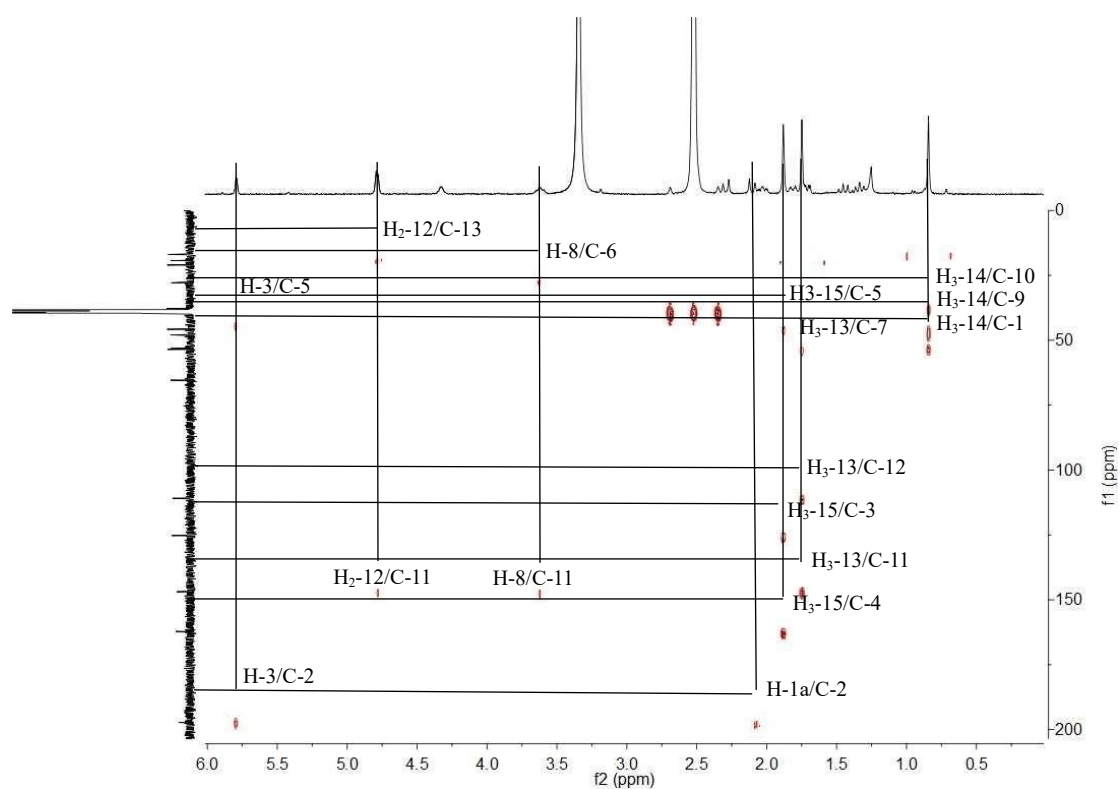
**Figure S2:** <sup>13</sup>C NMR Spectrum of **1** in DMSO-*d*<sub>6</sub> (100 MHz)



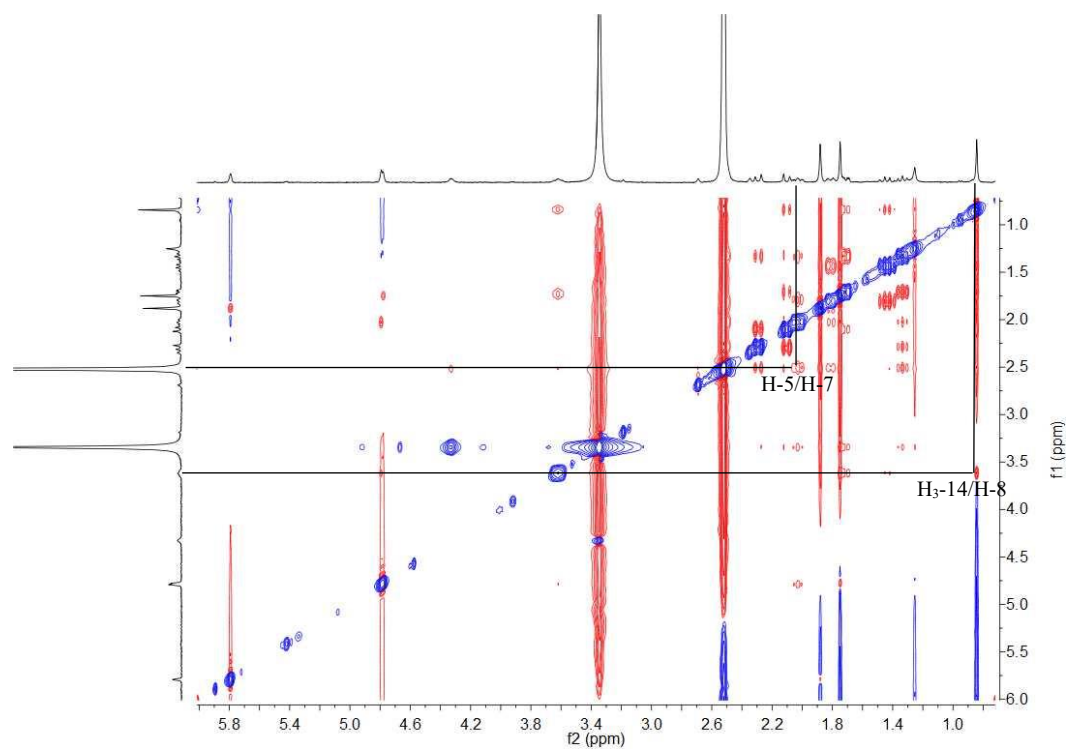
**Figure S3:** HSQC Spectrum of **1** in DMSO-*d*<sub>6</sub>



**Figure S4:** <sup>1</sup>H-<sup>1</sup>H COSY Spectrum of **1** in DMSO-*d*<sub>6</sub>



**Figure S5:** HMBC Spectrum of **1** in DMSO- $d_6$



**Figure S6:** NOESY Spectrum of **1** in DMSO- $d_6$

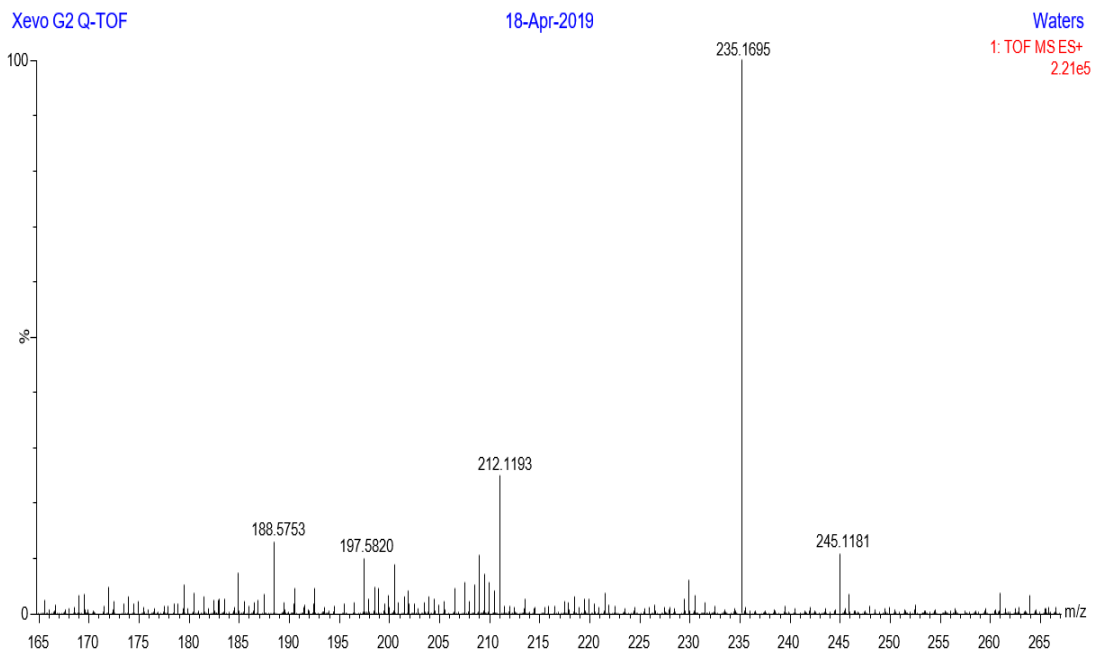
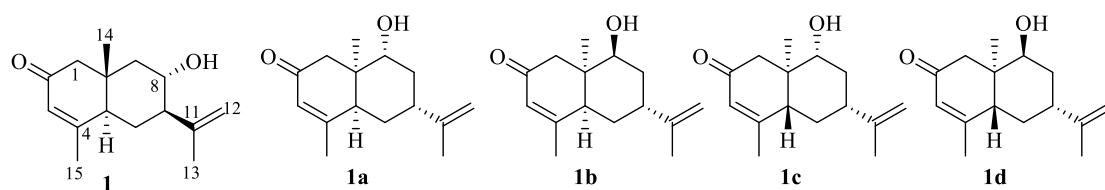


Figure S7: HRESIMS spectrum of **1**.

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<p>Score: 95</p> <p>1. <b>1007586-28-8</b></p> <p>Rotation (-), Absolute stereochemistry.</p>	<p>Score: 95</p> <p>2. <b>1007586-37-9</b></p> <p>Rotation (+), Absolute stereochemistry.</p>
<p>Score: 95</p> <p>3. <b>1007586-38-0</b></p> <p>Rotation (-), Absolute stereochemistry.</p>	<p>Score: 95</p> <p>4. <b>1007586-39-1</b></p> <p>Rotation (-), Absolute stereochemistry.</p>

Figure S8: Scifinder report for compound **1**



**Table S1.** Comparisons of  $^1\text{H}$  NMR and  $^{13}\text{C}$  NMR data of compounds **1** and **1a–1d**

No.	<b>1</b>		<b>1a</b>		<b>1b</b>		<b>1c</b>		<b>1d</b>	
	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$
1	2.09, 2.29	53.2	2.22, 2.81	48.0	2.52, 2.20		2.75, 2.16	50.5	2.86	50.1
2		197.2		199.2		199.6		198.6		203.3
3	5.79	125.3	5.97	129.0	5.85	126.5	5.91	127.1	5.84	126.7
4		162.2		161.9		163.5		161.8		167.6
5	2.50	45.8	2.60	47.1	1.98- 2.17	44.5	2.89	48.5	2.98	42.4
6	1.81, 1.45	27.9	1.71- 1.91	28.1	1.98- 2.17	29.9	1.93, 1.34	28.1	1.39, 1.93- 2.02	34.2
7	2.03	53.5	1.71- 1.91	40.3	2.45	37.5	2.12- 2.22	43.4	2.56	40.7
8	3.62	65.4	1.52	35.6	1.74	32.9	1.86 1.54	34.9	1.81, 1.71	34.2
9	1.70, 1.34	47.9	3.78	69.6	3.61	73.3	3.59	77.6	3.52	74.6
10		37.8		41.3		41.0		42.9		43.7
11		146.9		148.4		146.2		148.4		151.0
12	4.78,4.79	110.9	4.75, 4.76	109.6	4.84, 4.93	111.1	4.79	109.9	4.79, 4.76	109.8
13	1.75	19.3	1.76	20.9	1.78	22.4	1.78	20.9	1.78	21.1
14	0.84	16.9	1.14	20.4	1.09	24.9	0.90	11.2	0.89	17.7
15	1.87	21.0	1.94	22.5	1.96	23.4	1.94	22.6	1.97	22.4

Note: **1** in  $\text{DMSO-}d_6$ , **1a**, **1b**, **1c** in  $\text{CDCl}_3$ , **1d** in  $\text{methanol-}d_4$ ,