

# Supplementary Materials: Analysis of Indole Alkaloids from *Rhazya stricta* Hairy Roots by Ultra-Performance Liquid Chromatography-Mass Spectrometry

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Table S1. Alkaloids detected in *R. stricta* hairy root extracts by LC-MS or GC-MS.

No	Alkaloids	Identified with UPLC-MS	Identified with GC-MS	Alkaloid Type
1	Strictosidine lactam	+	-	strictosidine
2	Rhazine *	+	+	sarpagine
3	Strictosidine	+	-	strictosidine
4	Leepacine isomer I	+	-	ajmaline
5	Leepacine isomer II	+	-	ajmaline
6	Vincanine *	+	+	strychnos
7	Yohimbine *	+	+	yohimbinoïd
8	Yohimbine isomer *	+	+	yohimbinoïd
9	Serpentine isomer I	+	-	heteroyohimbine
10	Serpentine isomer II	+	-	heteroyohimbine
11	Vincamine *	+	+	eburnamine
12	Vincadifformine *	+	+	aspidosperma
13	Vallesiachotamine isomer I *	+	+	hunterburine
14	Eburenine *	+	+	aspidosperma
15	Vallesiachotamine isomer II *	+	+	hunterburine
16	Tetrahydrosecodinol	+	-	secodine
17	Tetrahydroalstonine *	+	-	heteroyohimbine
18	Tetrahydrosecodine	+	-	secodine
19	Dihydrosecodine	+	-	secodine
20	Tabersonine	+	-	aspidosperma
21	Decarbomethoxytabersonine	-	+	aspidosperma
22	Aspidospermidine	-	+	aspidosperma
23	Eburnamenine	-	+	eburnamine
24	Quebrachamine	-	+	aspidosperma
25	Tubotaiwine	-	+	aspidospermatin
26	Condylcarpine	-	+	aspidospermatin
27	Dihydrovincadifformine	-	+	aspidosperma
28	Vincadine	-	+	aspidosperma
29	Pleiocarpamine	-	+	pleiocarpaman
30	Fluorocarpamine	-	+	pleiocarpaman
31	Ajmalicine	-	+	heteroyohimbine

+: Alkaloid was detected with the method. -: alkaloid was not detected with the method. \*: Alkaloid which was detected by both LC-MS and GC-MS methods.

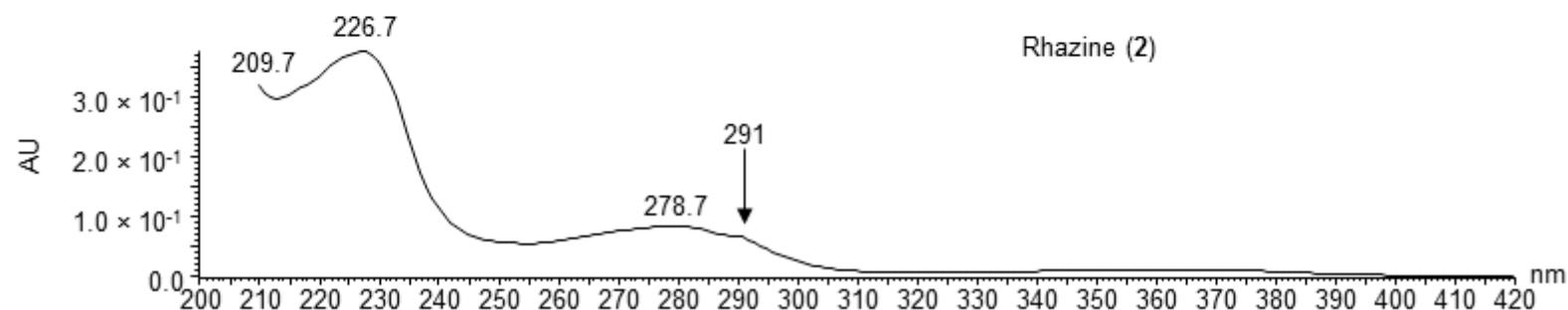
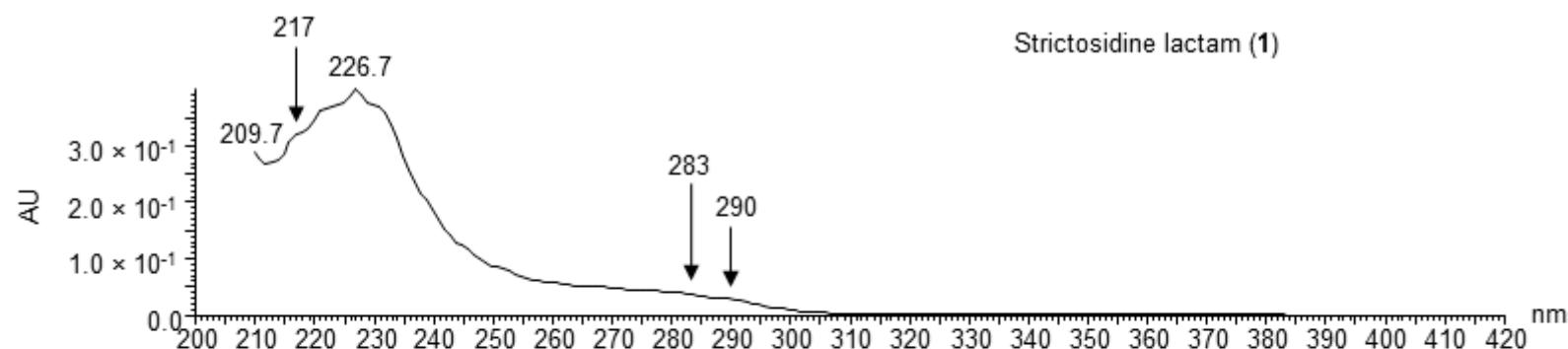


Figure S1. *Cont.*

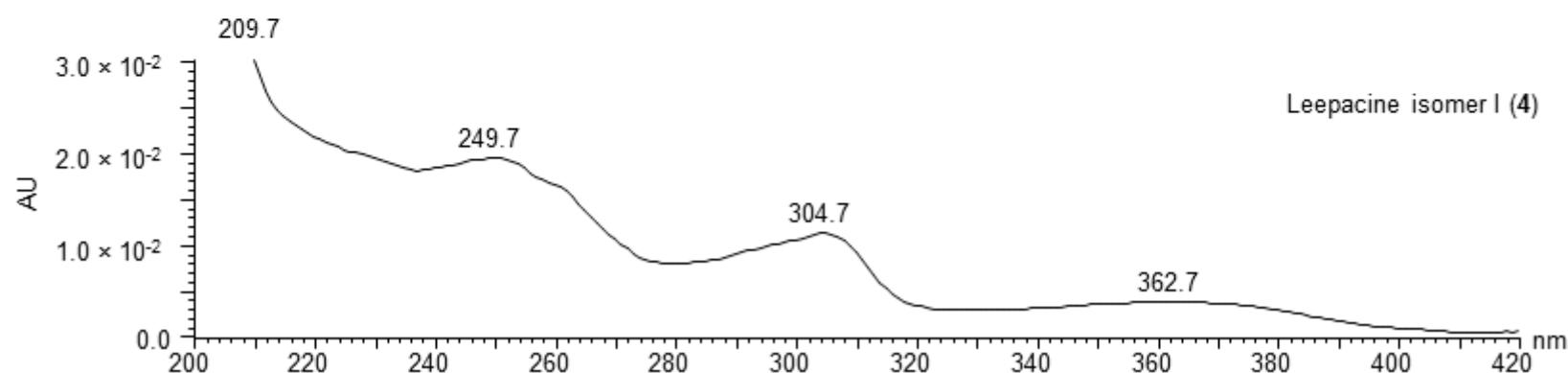
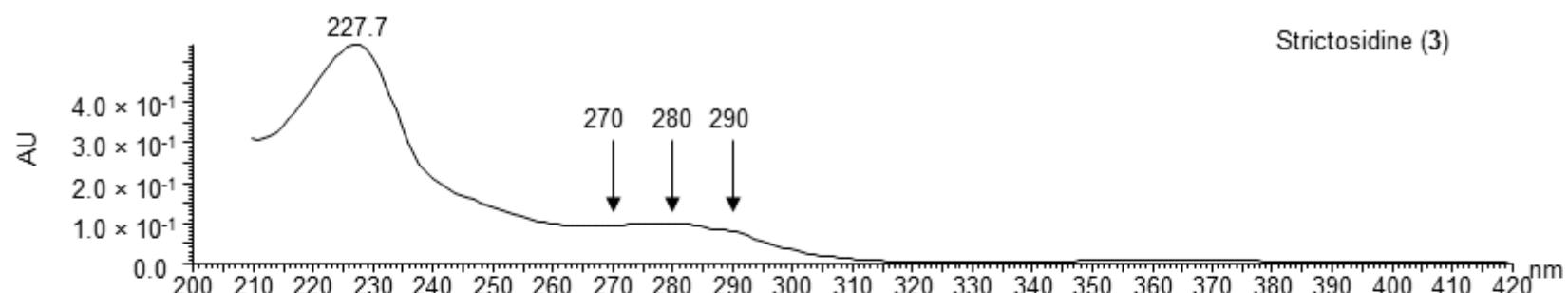


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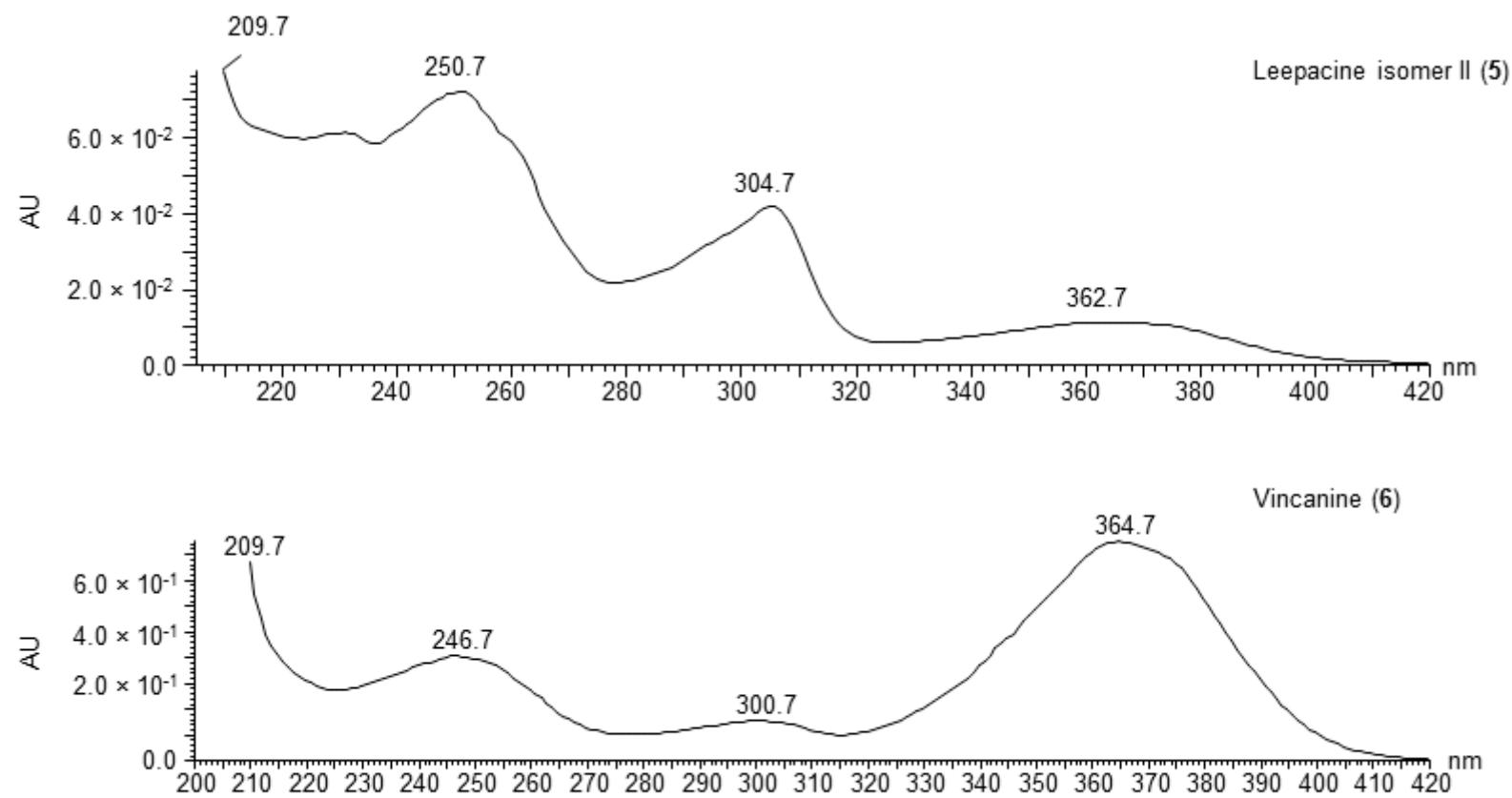


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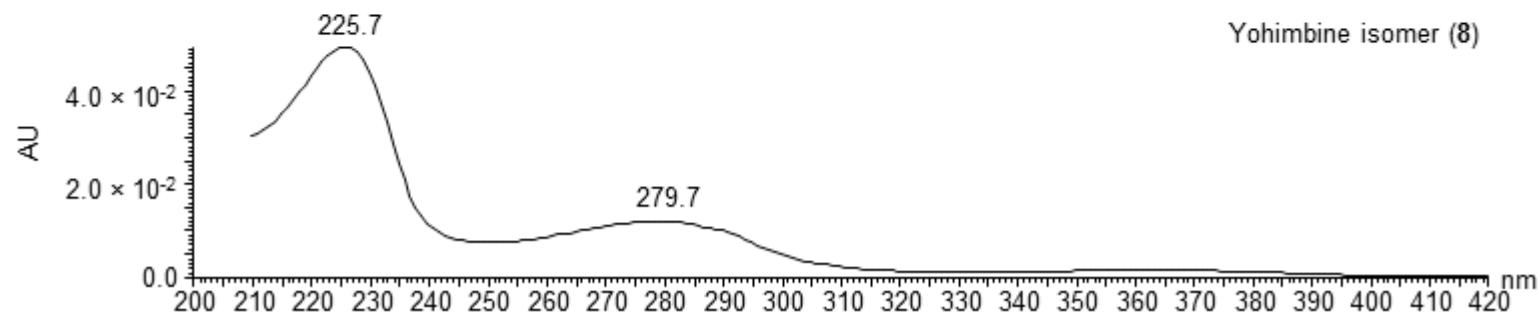
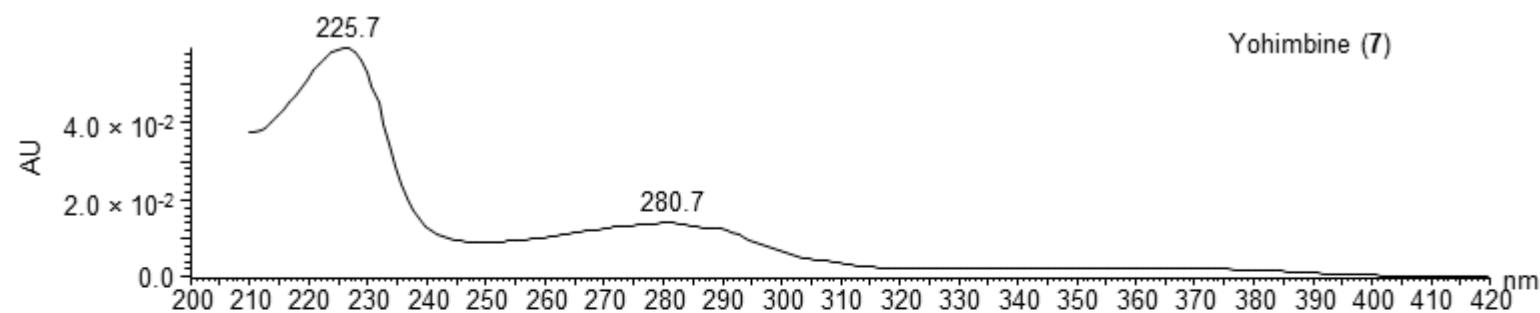


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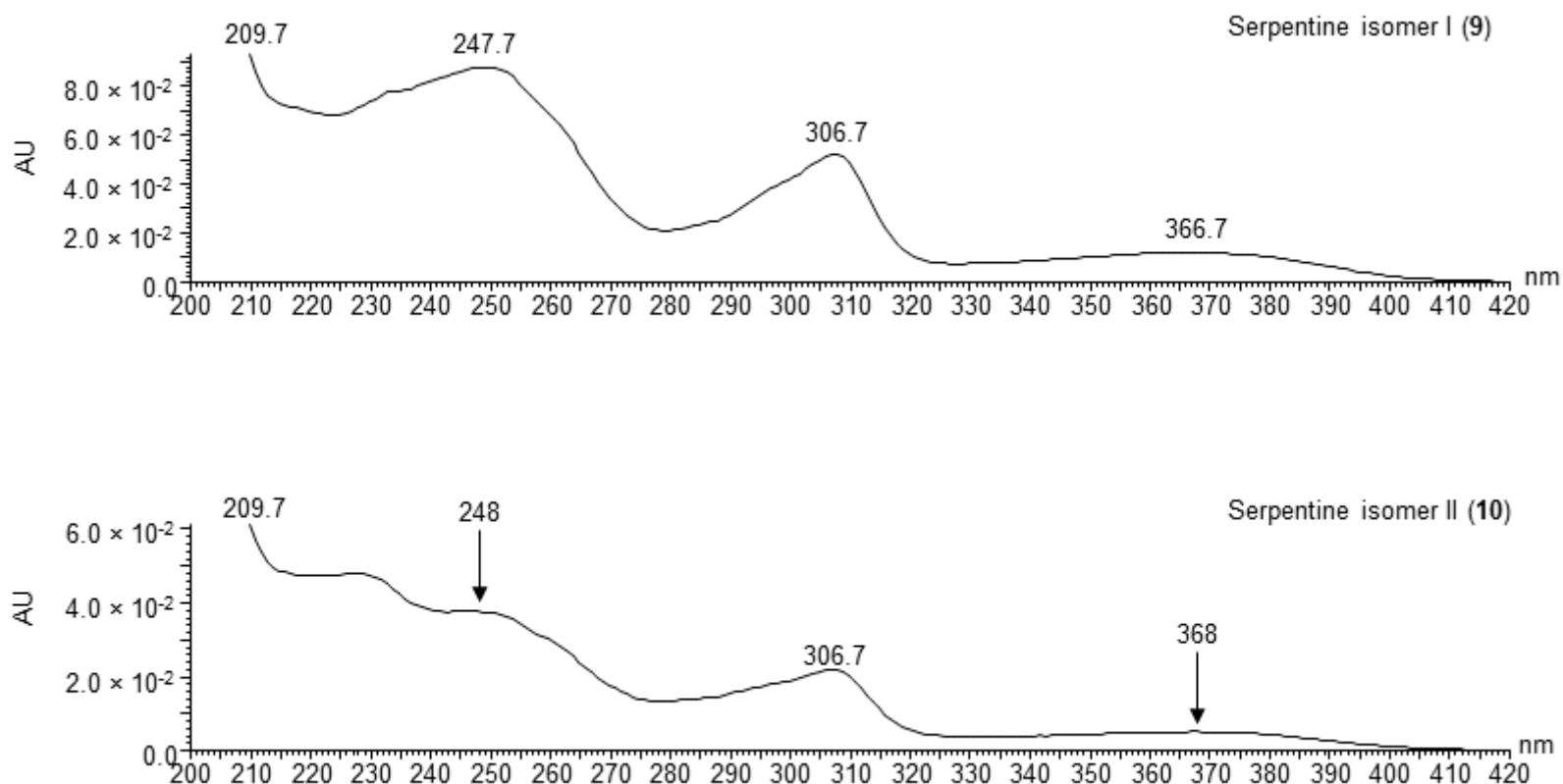


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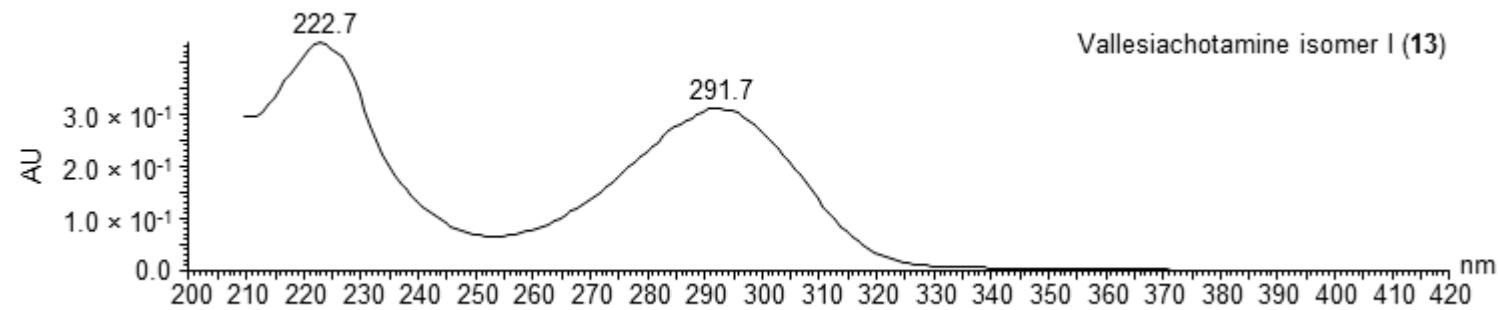
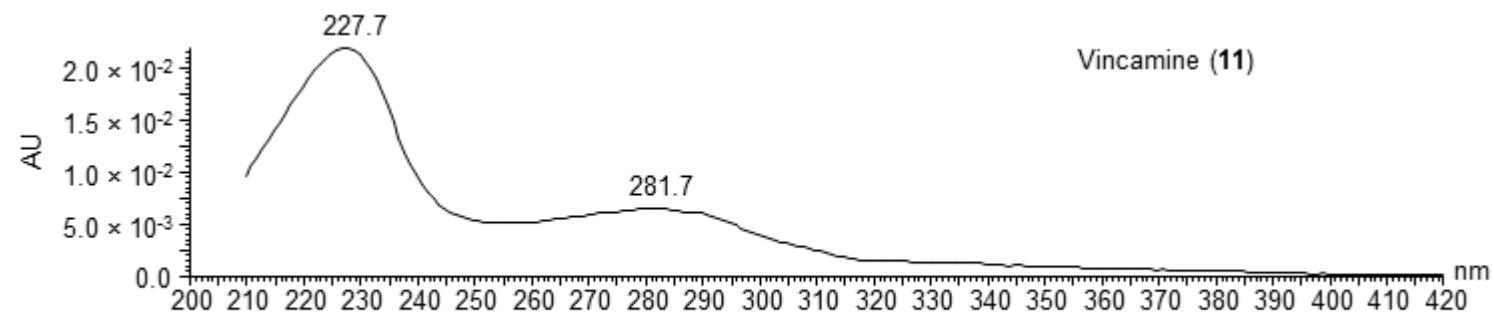


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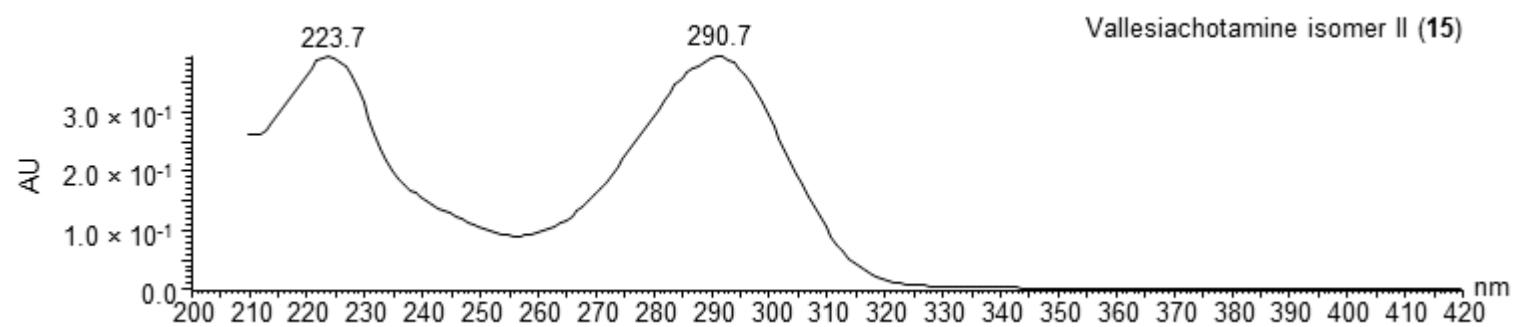
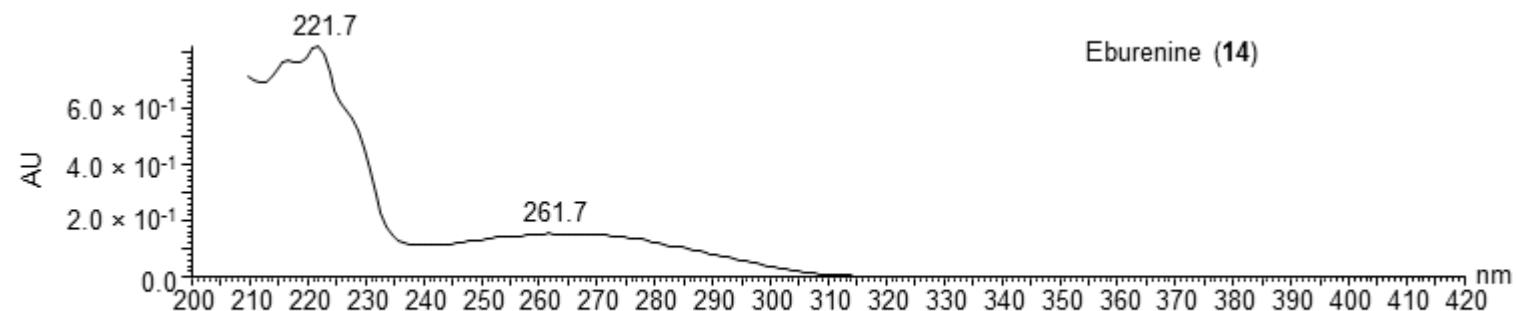


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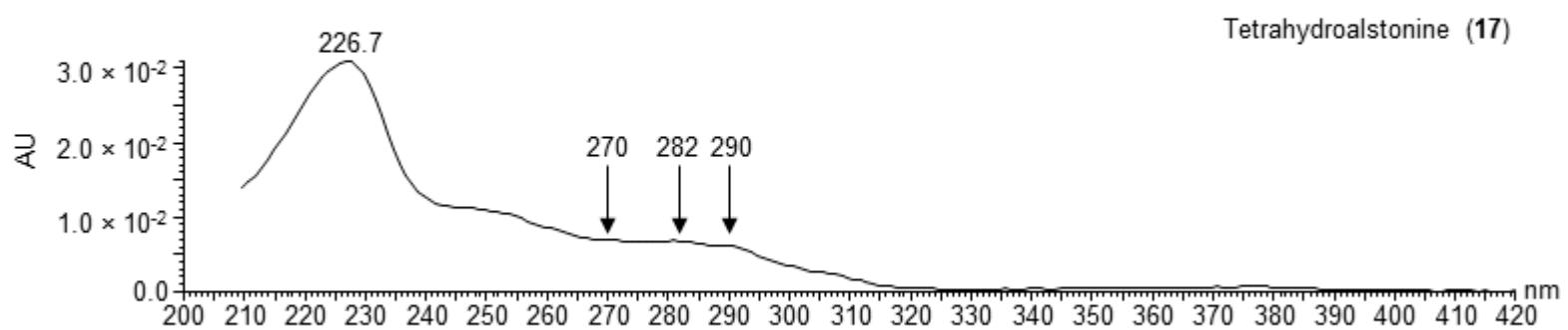
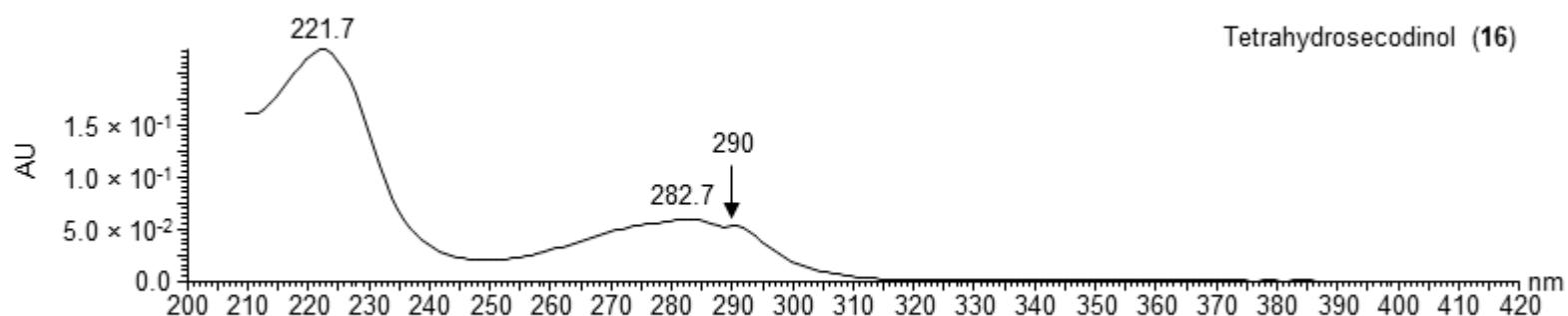


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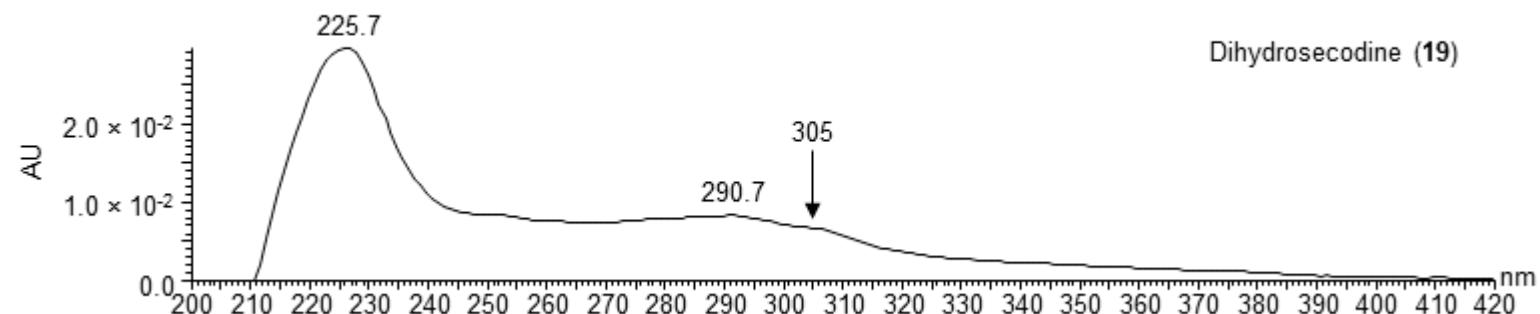
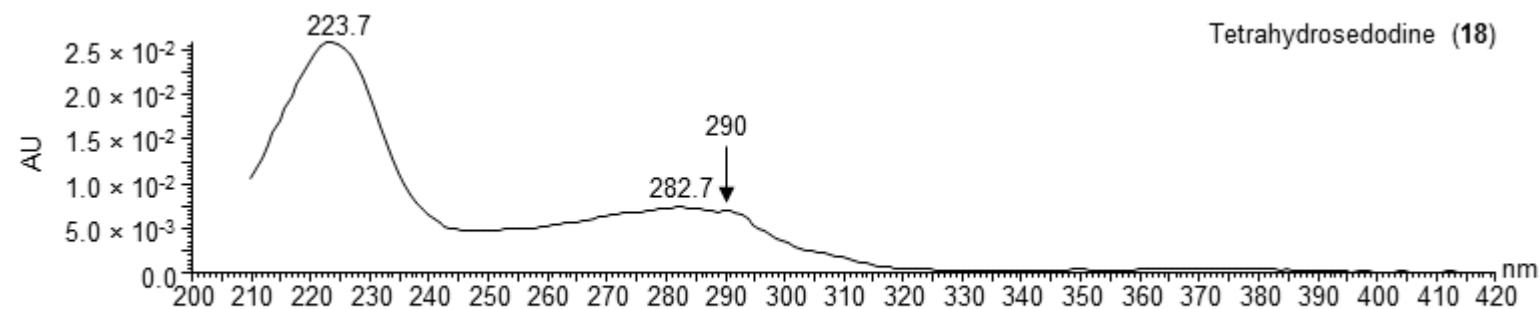
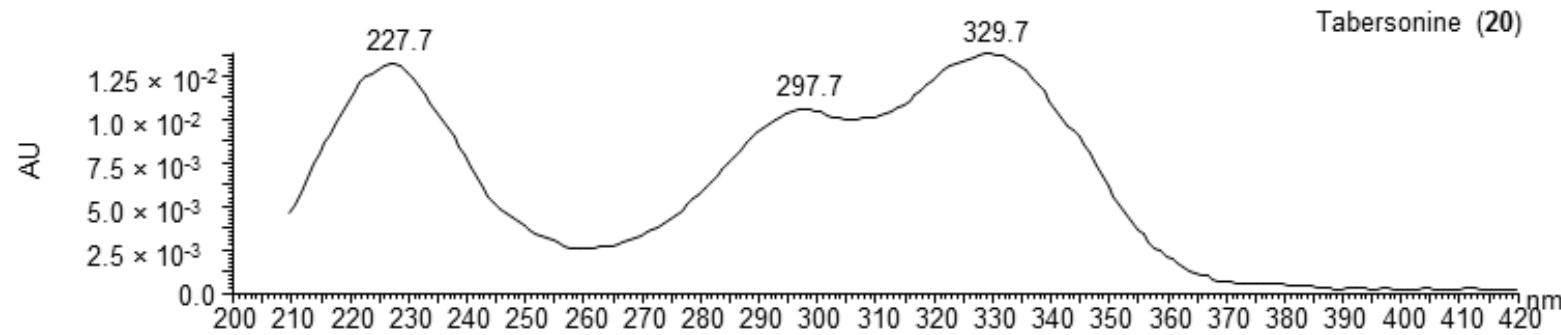
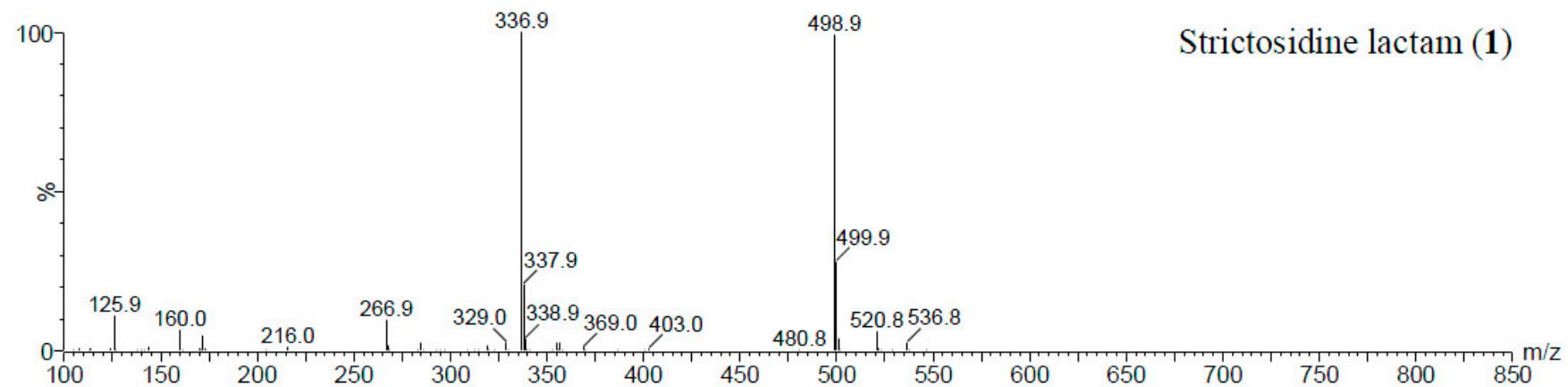


Figure S1. *Cont.*



**Figure S1.** UV spectra of *R. stricta* alkaloids in UPLC-UV analysis. Data for vincadifformine (**12**) is presented in Figure S3.



**Figure S2. Cont.**

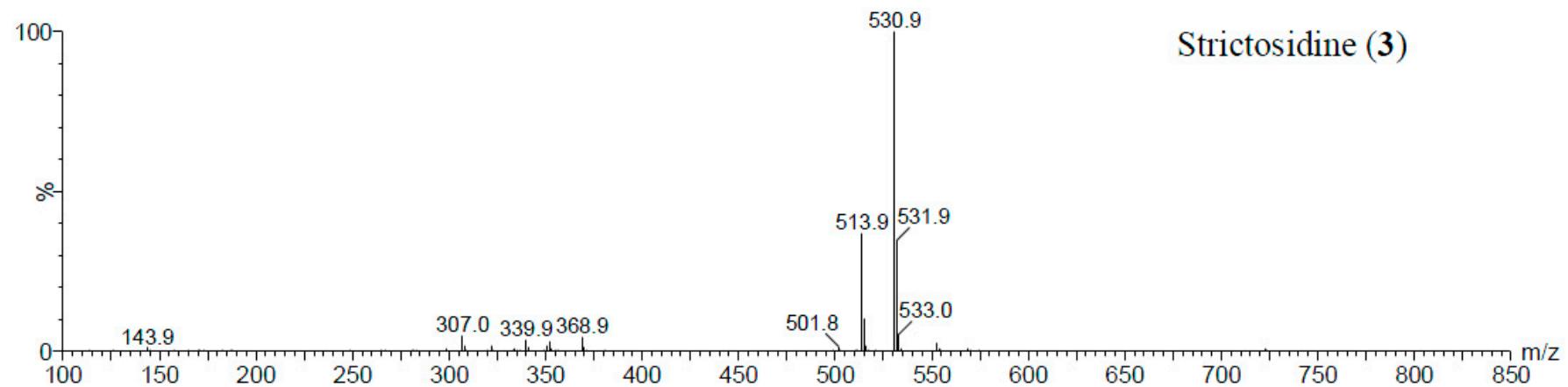
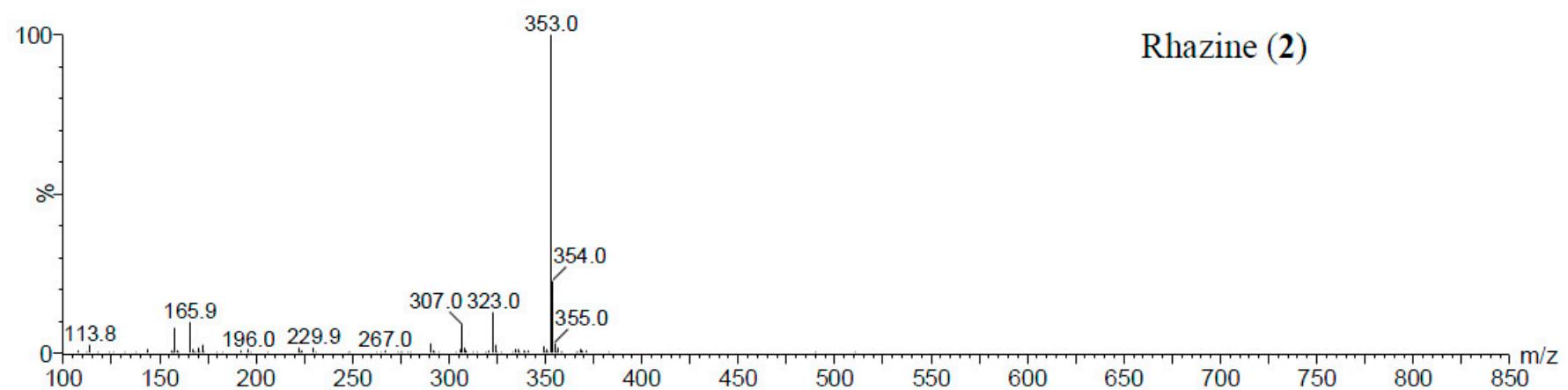


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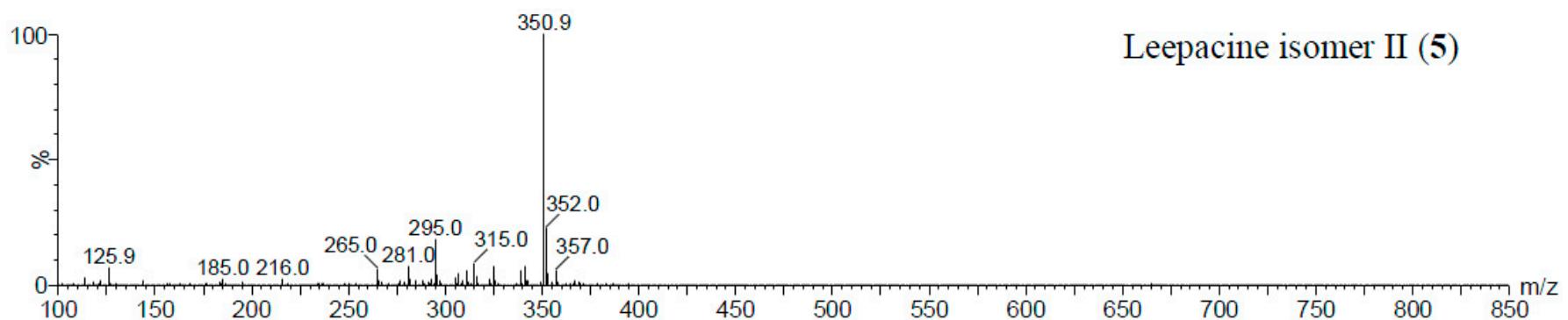
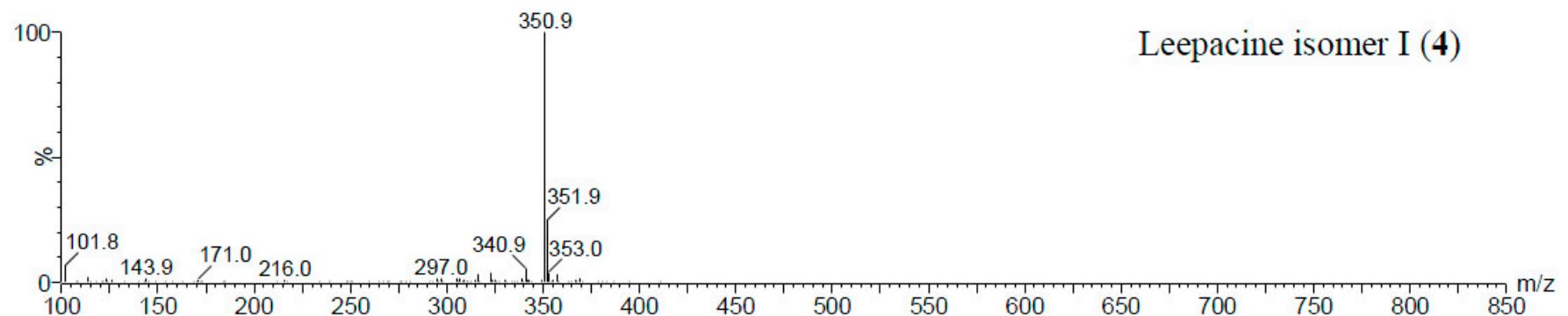


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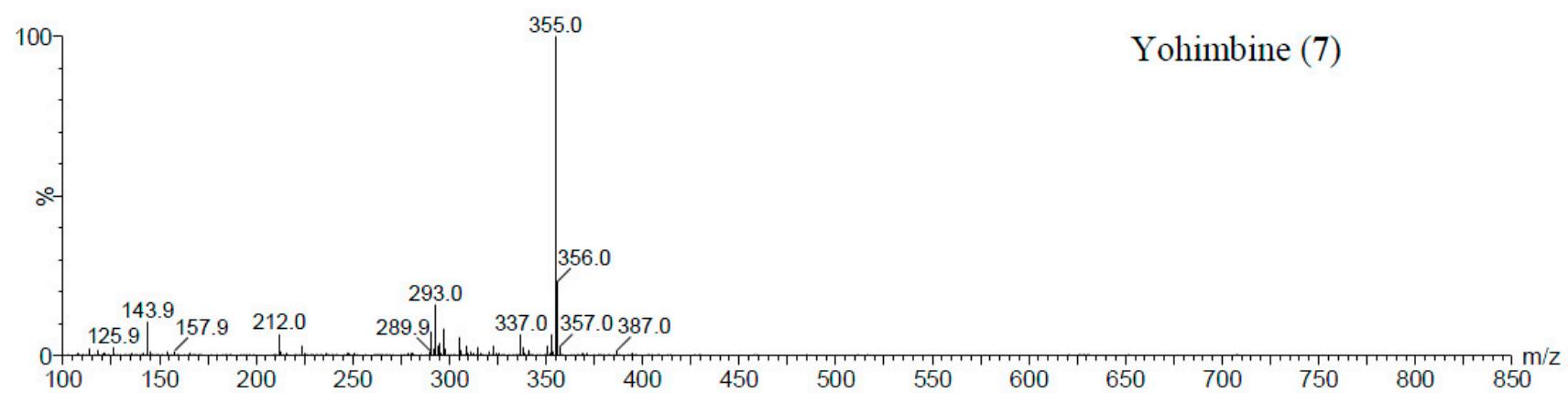
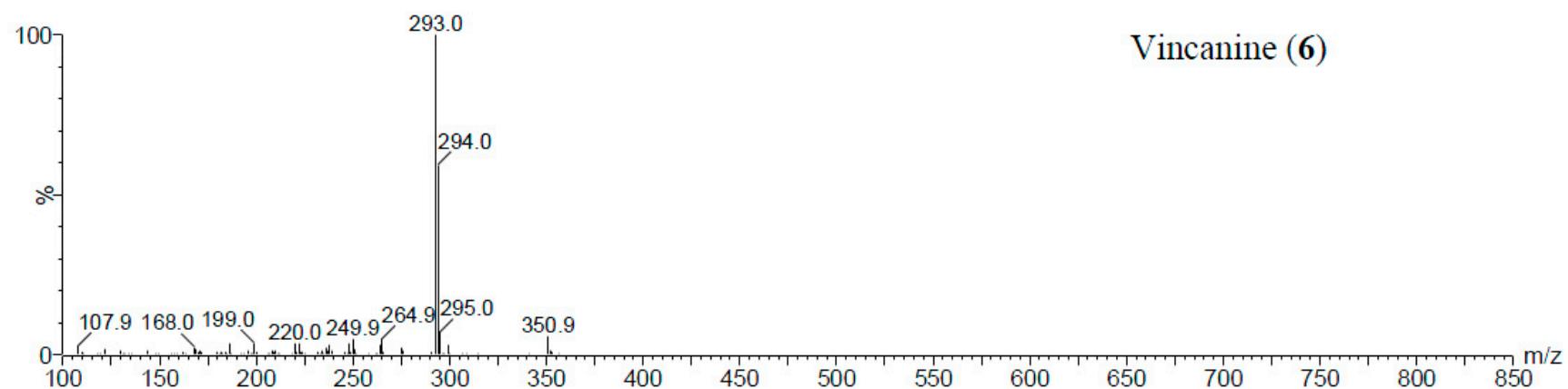


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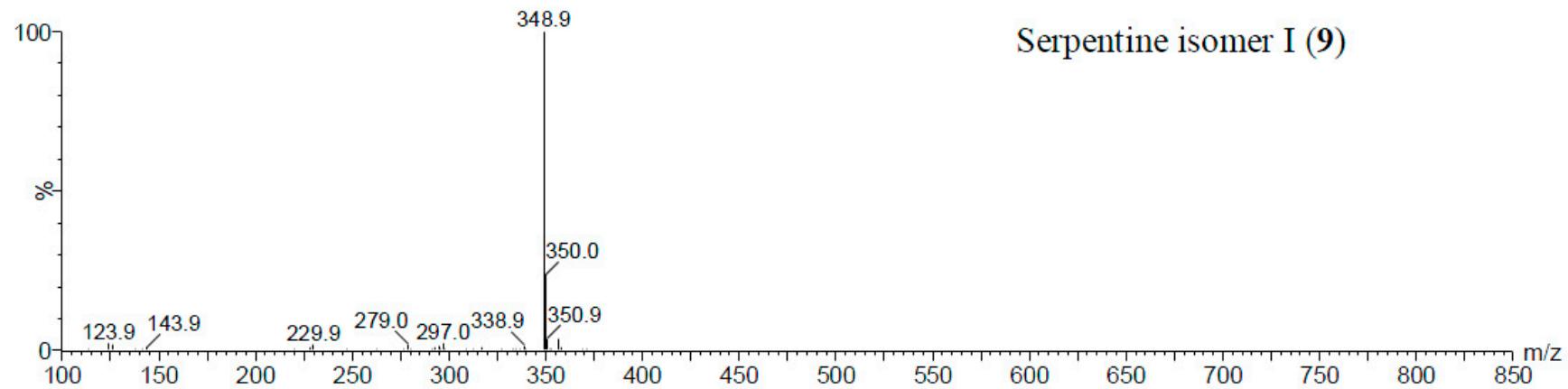
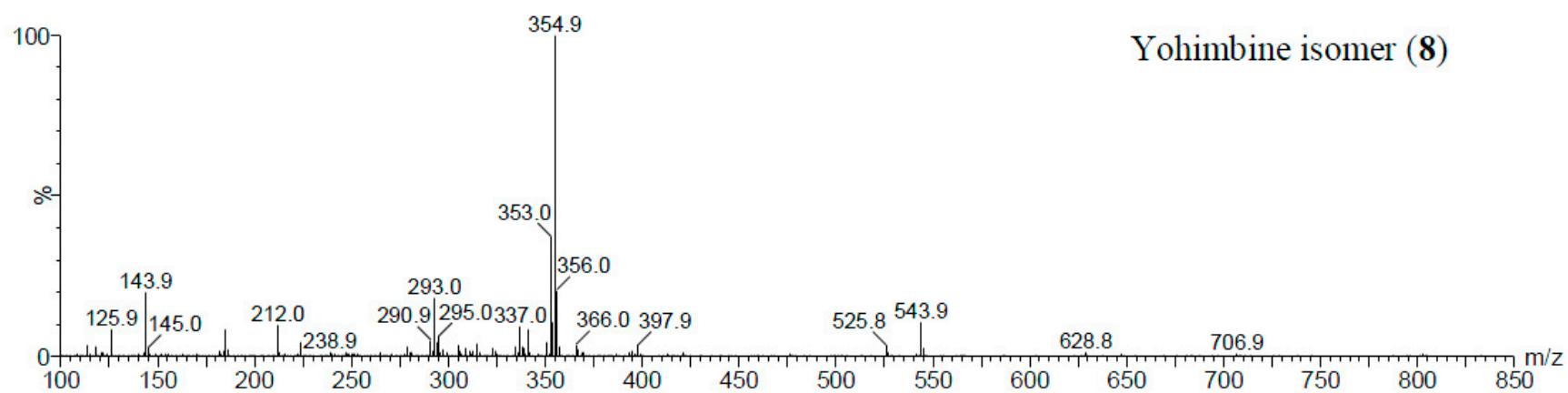


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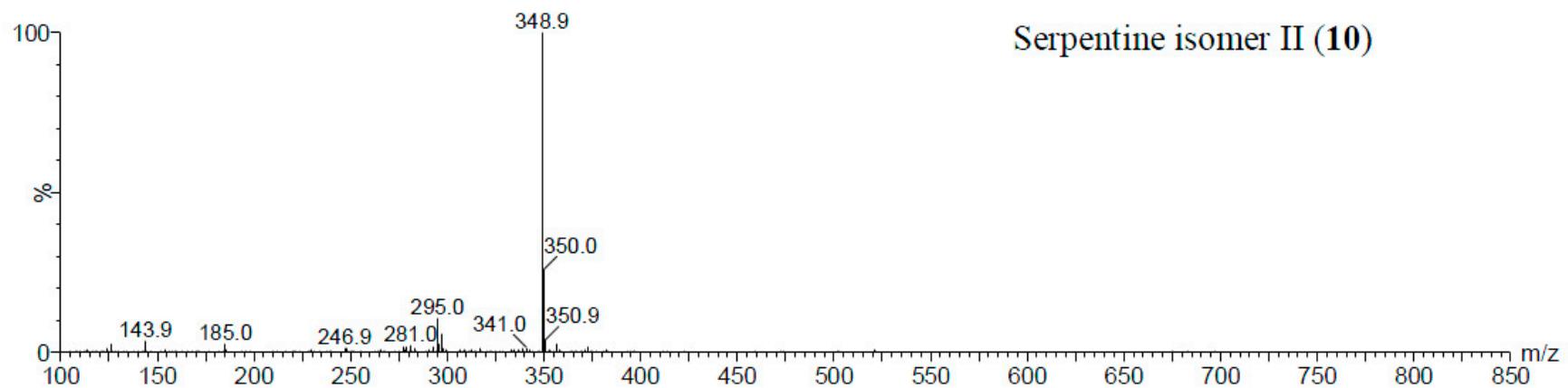
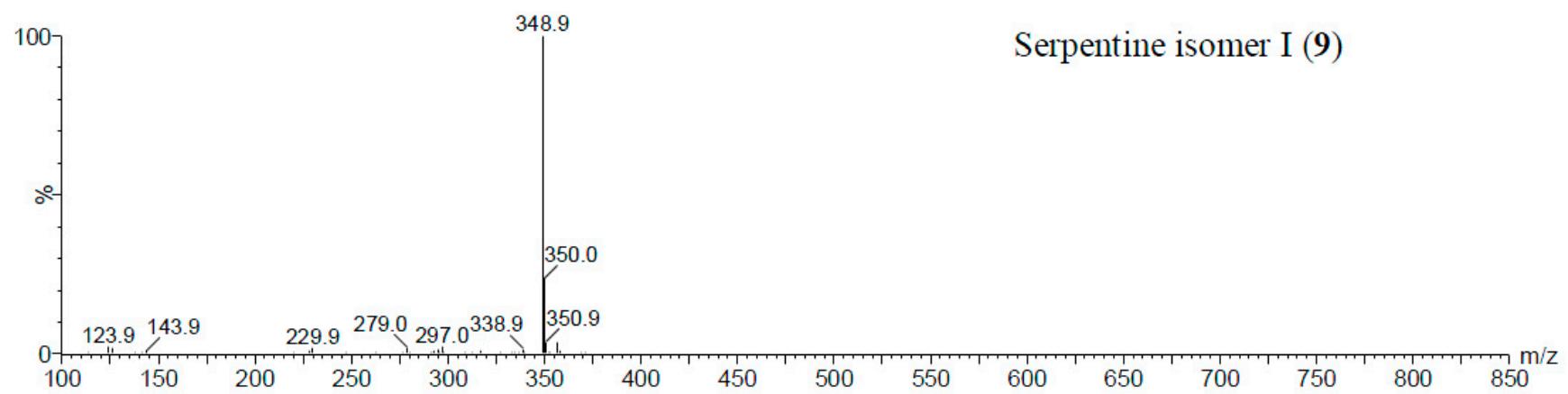


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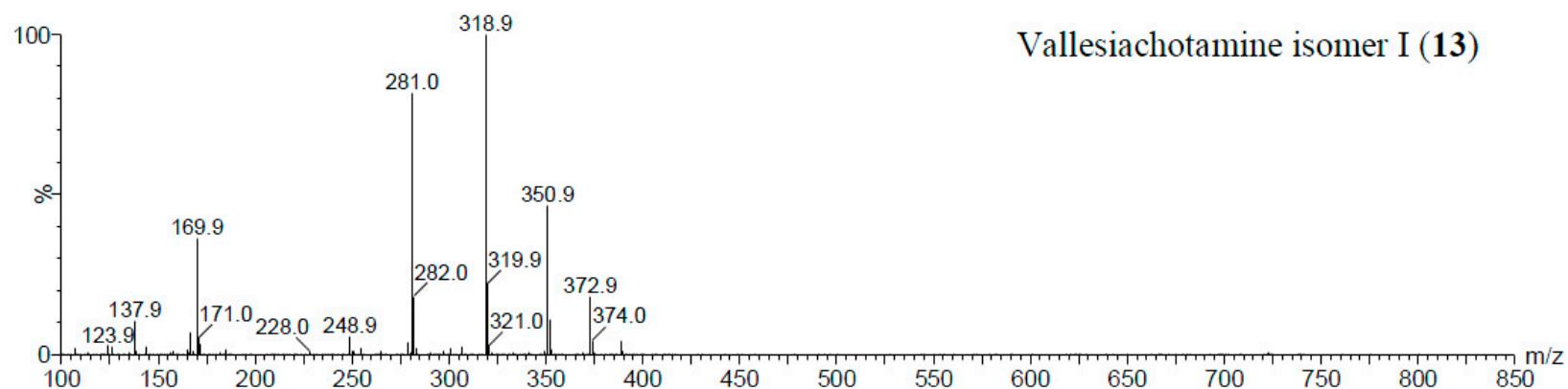
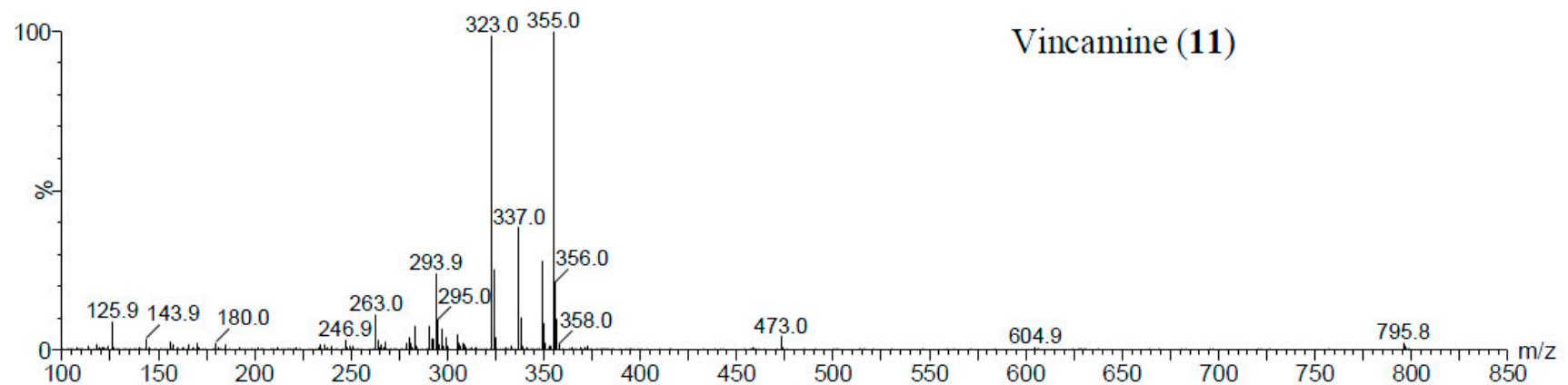


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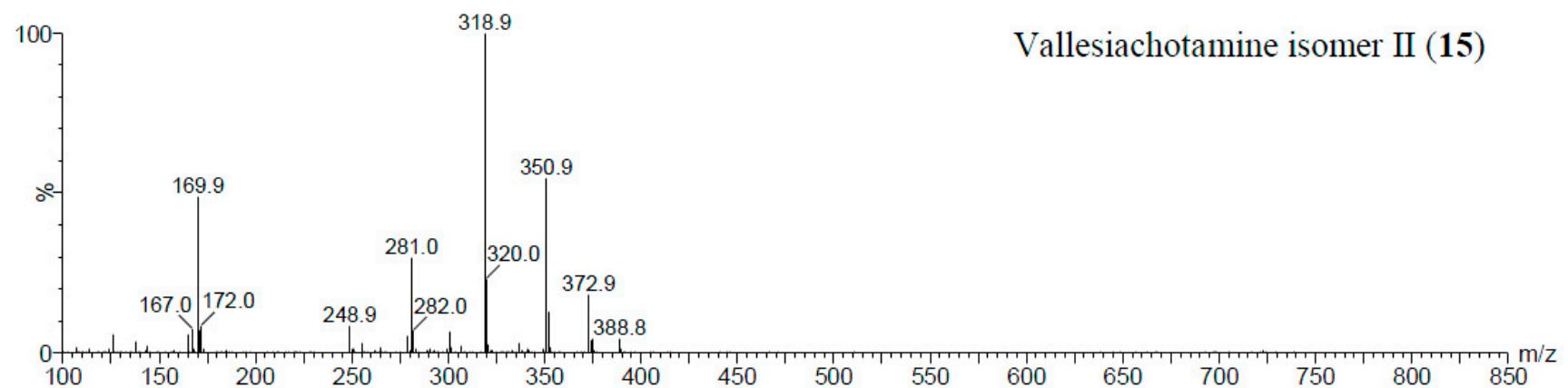
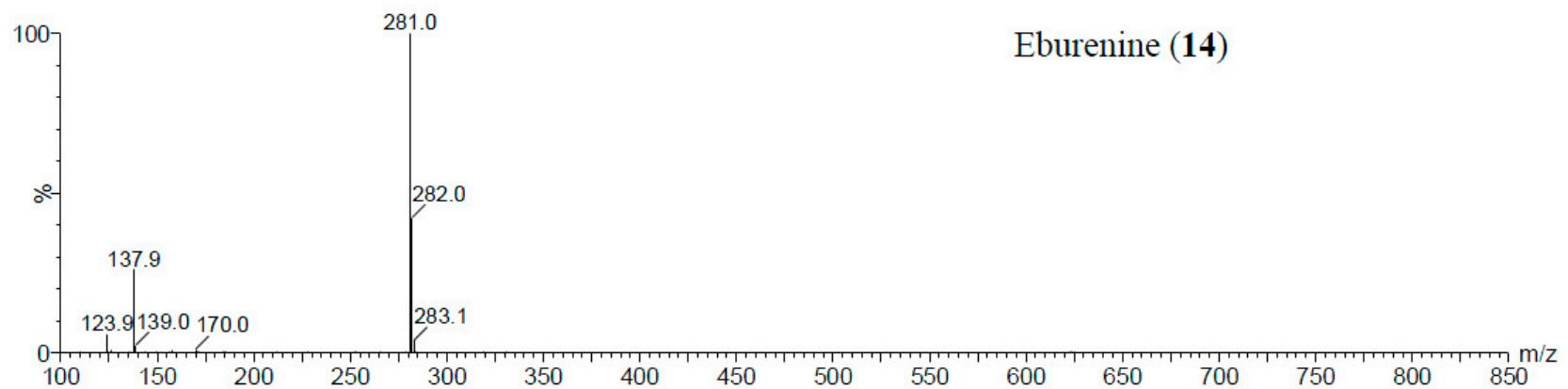


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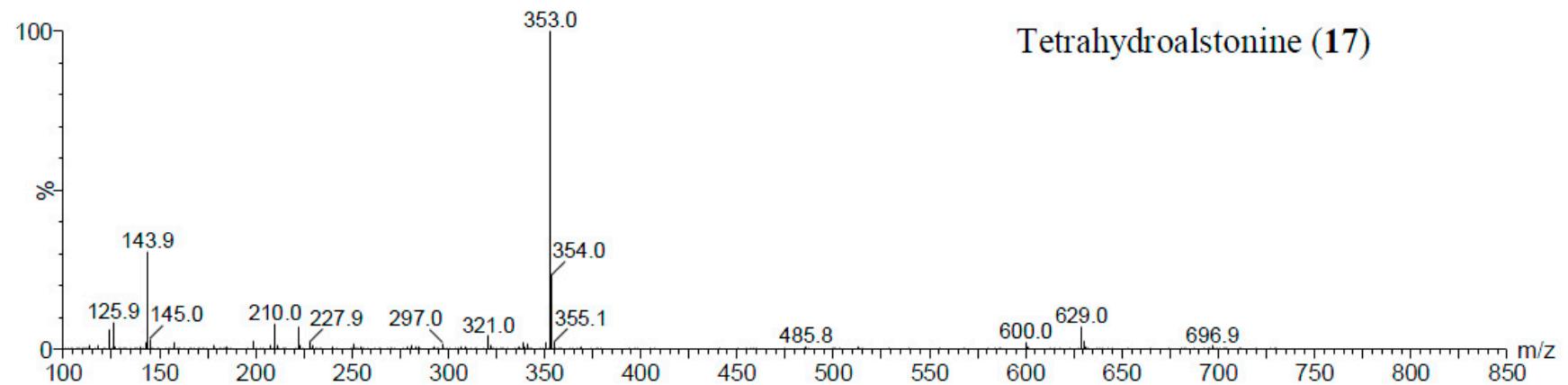
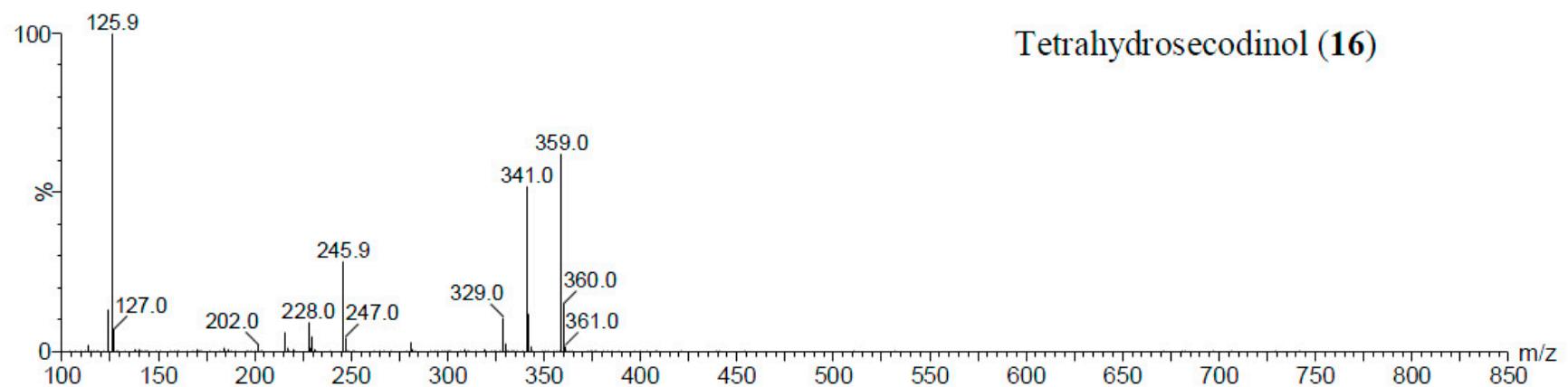


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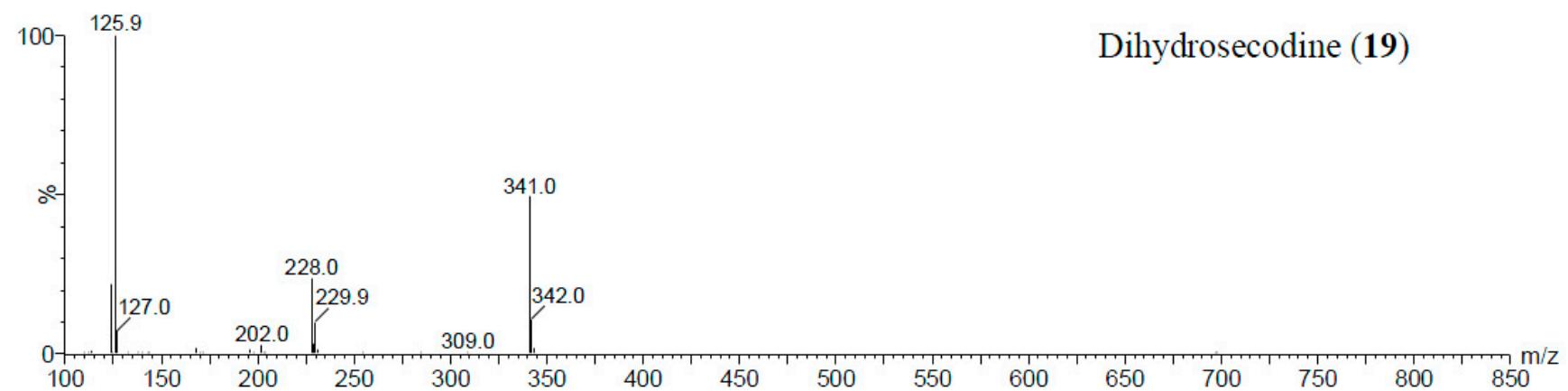
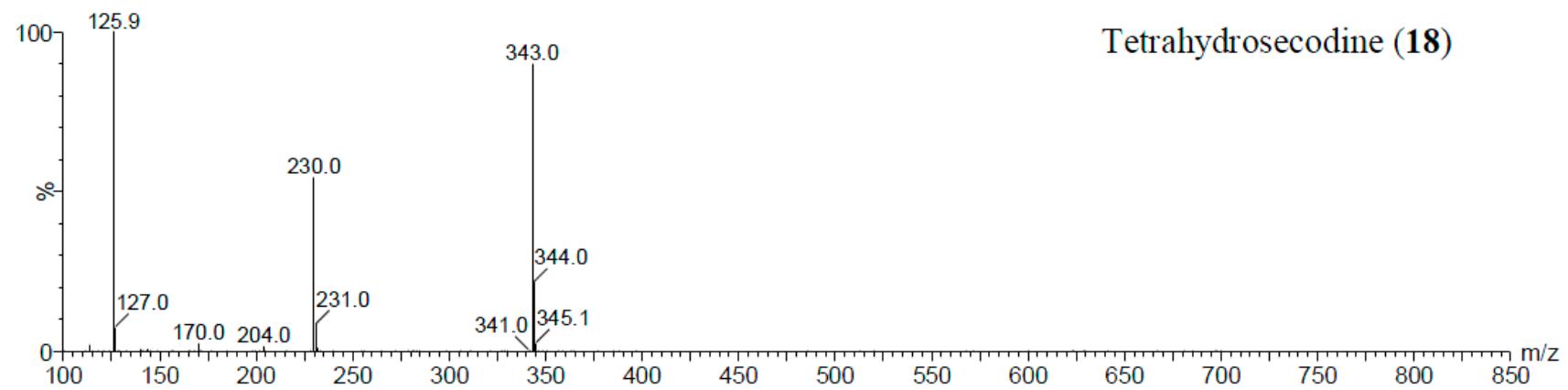
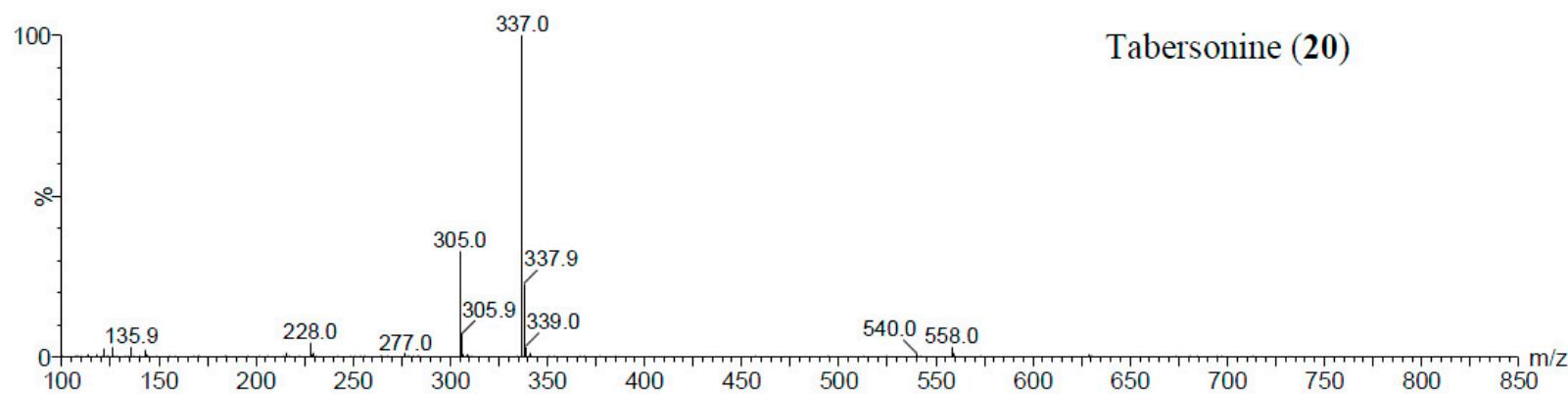
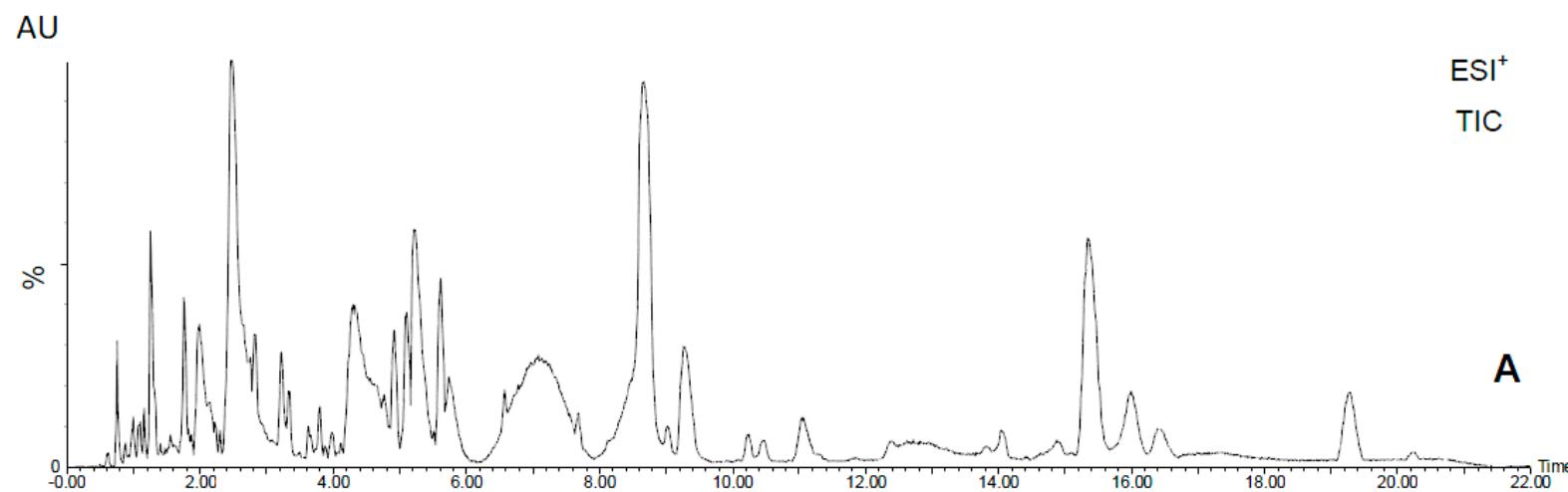


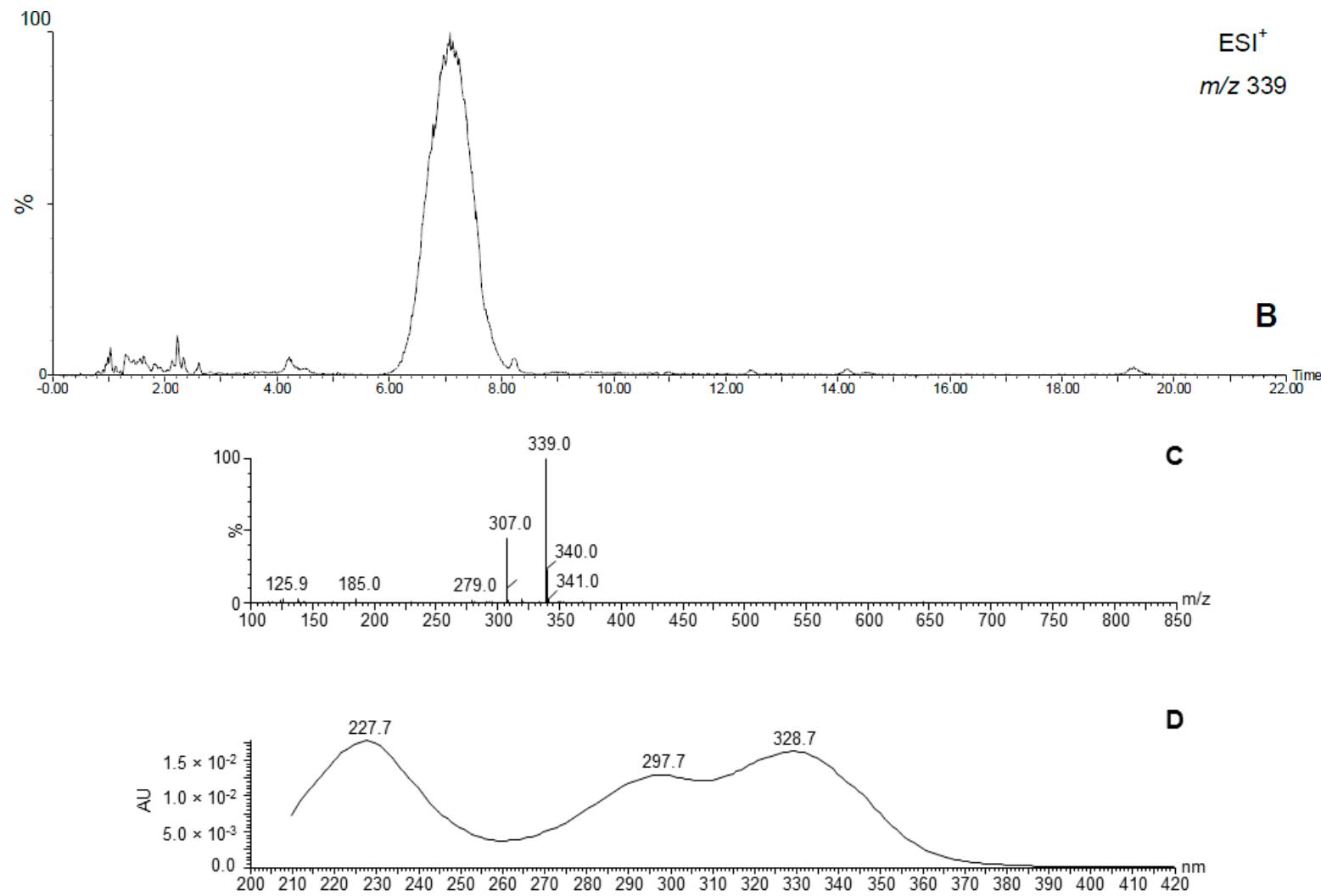
Figure S2. Cont.



**Figure S2** UPLC-MS spectra of *R. stricta* alkaloids in positive ionization mode (ESI<sup>+</sup>). Data for vincadifformine (**12**) is presented in Figure S3.



**Figure S3. Cont.**



**Figure S3** UPLC-PDA-MS-TIC analysis of *R. stricta* in positive ionization mode (**A**), extracted ionogram of vincadifformine (**12**, channel *m/z* 339) (**B**), its UPLC-MS spectrum (**C**), and UV spectrum (**D**).