

Table S1 Regression equation, LODs and LOQs of quantitative method for four alkaloid

compound	linear range(ng/mL)	regression equation	R	LODs(ng/mL)	LOQs(ng/mL)
Sipeimine- 3 β -D-glucoside	7.867~4028	$Y = 181.8X + 4129.5$	0.9994	0.1229	0.2458
Sipeimine	1.982~1015	$Y = 375.08X + 4374.8$	0.9991	0.2478	0.4956
Peimisine	4.320~2212	$Y = 329.69X + 6301.7$	0.9993	0.1350	0.5400
Yibeinoside A	7.969~4080	$Y = 147.32X + 304.87$	0.9998	0.06226	0.2490

Table S2 The results of precision, stability and repetition experiment

Compound	Precision degree (n=6)		Repeatability (n=6)		Stability (n=5)	
	content(%)	RSD(%)	content (%)	RSD(%)	content (%)	RSD(%)
Sipeimine- 3 β -D-glucoside	0.0338	4.89	0.0316	6.53	0.0328	6.28
Sipeimine	0.0509	1.26	0.0475	5.26	0.0501	3.05
Peimisine	0.0098	1.37	0.0095	5.14	0.0100	2.09
Yibeinoside A	0.0176	2.386	0.0169	6.95	0.0181	2.63

Table S3 The results of recovery experiments

Component	Orginal (μ g)	Added (μ g)	Measured (μ g)	Recovery (%)	Average recovery (%)	RSD (%)
Sipeimine- 3 β -D-glucoside		25.16	67.04	103.70		
	40.95	40.28	81.99	101.89	101.8	1.95
		55.38	96.18	99.73		
Sipeimine		35.52	95.60	97.35		
	61.02	60.90	125.58	106.01	103.5	5.15
		86.28	153.39	107.06		
Peimisine		8.30	19.90	103.37		
	11.32	11.06	22.48	100.90	100.7	2.82
		16.59	27.53	97.71		
Yibeinoside A		15.30	37.80	100.59		
	22.41	20.40	42.78	99.85	99.80	0.82
		30.60	52.69	98.95		