

Table S1. Chemical composition of the essential oil from four locations from aerial parts of *Veronica austriaca* ssp. *jacquinii* calculated based on the masses of EOs extracted from dry plant material.

Component	Mr	St	Br	GJ	
	RI ¹	RI ²	EO ± SD (mg/g dry plant material)	EO ± SD (mg/g dry plant material)	
Sesquiterpene hydrocarbons		0.056	0.358	0.143	
E-Caryophyllene*	1424	1585	0.012 ± 0.00038	0.109 ± 0.00047	0.078 ± 0.00051
δ-Cadinene	1517	1745	0.044 ± 0.00038	0.086 ± 0.00047	0.051 ± 0.00103
<i>allo</i> -Aromadendrene	1465	1662	-	0.041 ± 0.00047	-
β-Chamigrene	1476	1724	-	-	0.014 ± 0.00103
Germacrene D	1482	1692	-	0.122 ± 0.00233	-
Oxygenated sesquiterpenes		2.014	1.443	1.505	
Spathulenol	1577	2101	-	0.086 ± 0.00047	0.023 ± 0.00051
β-Caryophyllene oxide*	1581	1955	0.017 ± 0.00038	0.029 ± 0.00047	0.025 ± 0.00051
γ-Eudesmol	1632	2175	-	0.012 ± 0.0014	-
α-Bisabolol oxide	1748	2511	-	0.017 ± 0.00047	-
Hexahydrofarnesyl acetone	1839	2113	1.997 ± 0.00038	1.299 ± 0.00047	1.457 ± 0.00051
Phenolic compounds		0.024	0.125	0.043	
Methyl eugenol	1403	2005	-	0.0588 ± 0.00047	-
(Z)-Methyl isoeugenol	1451	2070	0.024 ± 0.00114	0.066 ± 0.00047	0.043 ± 0.00051
Acids, alcohols and esters		1.341	2.226	2.969	
1-Hexadecanol	1874	2371	-	0.027 ± 0.0014	-
Hexadecanoic acid	1959	2912	1.015 ± 0.00076	2.199 ± 0.000467	2.804 ± 0.00257
Oleic acid	2133	2998	0.089 ± 0.00038	-	0.026 ± 0.00154
Octadecanol acetate	2209	2211	0.237 ± 0.00038	-	0.116 ± 0.00051
1-Heptatriacotanol	2309	2309	-	0.023 ± 0.00051	-
Hydrocarbons		0.06114	0.095	0.129	
Eicosane*	2000	2000	-	0.054 ± 0.00047	0.023 ± 0.00051
Heneicosane*	2100	2100	0.02014 ± 0.00076	0.016 ± 0.00047	-
Docosane*	2200	2200	0.014 ± 0.00038	0.014 ± 0.00047	0.042 ± 0.00051
Tricosane*	2300	2300	-	-	0.032 ± 0.00051
Tetracosane*	2400	2400	-	0.011 ± 0.00047	0.032 ± 0.00051
Pentacosane*	2500	2500	0.027 ± 0.00152	-	-
Total identification (mg/g)		3.496	4.247	4.789	
				5.878	

Retention indices (RI) were determined relative to a series of *n*-alkanes (C8–C40) on capillary columns VF5-ms (RI¹) [24] and CPWax 52 (RI²) [25]; Identification method: RI, comparison of RIs with those listed in a homemade library, reported in the literature [24], and/or authentic samples; comparison of mass spectra with those in mass spectral libraries NIST02 and Wiley 9; * co-injection with reference compounds; SD, standard deviation.

Table S2. Chemical composition of the hydrosols from four locations from aerial parts of *Veronica austriaca* ssp. *jacquinii* calculated based on the masses of volatile compounds extracted from hydrosols from dry plant material

Mr	St	Br	GJ
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Component	RI ¹	RI ²	Hy ± SD (mg/g dry plant material)			
Monoterpene hydrocarbons			0.047	-	-	-
α-Thujene	924	1032	0.023 ± 0.00028	-	-	-
β -Phellandrene	1002	1194	0.024 ± 0.00083	-	-	-
Oxygenated monoterpenes			0.585	0.708	0.238	0.679
<i>trans</i> -Linalool oxide*	1088	1434	0.010 ± 0.00110	-	-	-
<i>n</i> -Nonanal	1100	1389	0.120 ± 0.00028	0.118 ± 0.00042	-	-
Borneol	1176	1719	0.043 ± 0.00028	0.021 ± 0.00042	-	-
Camphor	1151	1499	0.060 ± 0.00028	0.039 ± 0.00042	-	0.160 ± 0.00045
Pinocarvone	1160	1565	0.055 ± 0.00028	-	-	-
<i>trans</i> - <i>p</i> -Mentha-1(7),8-dien-2-ol	1187	1803	0.212 ± 0.00028	0.220 ± 0.00042	0.220 ± 0.00030	0.288 ± 0.00090
Hexyl 2-methyl butanoate	1233	1425	0.035 ± 0.00028	0.131 ± 0.00126	-	0.197 ± 0.00045
Menthyl acetate	1294	1550	0.050 ± 0.00083	0.179 ± 0.00042	0.018 ± 0.00030	0.034 ± 0.00271
Sesquiterpene hydrocarbons			0.189	0.155	0.047	0.300
<i>E</i> -Caryophyllene*	1424	1585	0.073 ± 0.00028	0.056 ± 0.00042	0.019 ± 0.00059	0.033 ± 0.00045
δ-Cadinene	1517	1745	0.065 ± 0.00028	-	0.028 ± 0.00177	0.108 ± 0.00362
<i>allo</i> -Aromadendrene	1465	1662	0.042 ± 0.00028	-	-	0.056 ± 0.00045
β-Chamigrene	1478	1724	0.009 ± 0.00028	-	-	-
Germacrene D	1482	1692	-	0.099 ± 0.00042	-	0.103 ± 0.00045
Oxygenated sesquiterpenes			0.088	0.398	0.285	0.194
Spathulenol	1577	2101	-	-	-	0.056 ± 0.00045
<i>β</i> -Caryophyllene oxide*	1581	1955	0.060 ± 0.00028	0.053 ± 0.00042	0.033 ± 0.00030	0.023 ± 0.00045
γ-Eudesmol	1632	2175	-	-	-	-
α-Muurolol	1645	2181	-	0.052 ± 0.00042	-	-
α-Cadinol	1655	2208	-	0.103 ± 0.00042	-	-
α-Bisabolol	1685	2210	0.015 ± 0.00083	-	0.015 ± 0.00030	0.060 ± 0.00045
α-Bisabolol oxide	1748	2511	-	-	0.009 ± 0.00030	0.023 ± 0.00045
Hexahydrofarnesyl acetone	1839	2113	0.013 ± 0.00028	0.190 ± 0.00168	0.228 ± 0.00059	0.032 ± 0.00045
Phenolic compounds			1.197	1.582	1.854	2.243
Thymol*	1289	2154	0.230 ± 0.00138	0.397 ± 0.00084	0.103 ± 0.00030	0.189 ± 0.00045
Thymol acetate	1349	-	0.101 ± 0.00028	0.095 ± 0.00042	-	0.110 ± 0.00136
Methyl eugenol	1403	2005	0.834 ± 0.00055	0.980 ± 0.00042	1.713 ± 0.00030	1.891 ± 0.00045
(Z)-Methyl isoeugenol	1451	2070	0.032 ± 0.00028	0.110 ± 0.00252	0.038 ± 0.00030	0.053 ± 0.00045
Acids, alcohols and esters			0.191	0.557	0.206	0.281
1-Hexadecanol	1874	2371	-	-	0.072 ± 0.00030	-
Hexadecanoic acid	1959	2912	0.126 ± 0.00028	0.264 ± 0.00084	0.067 ± 0.00030	0.085 ± 0.00045
Oleic acid	2133	2998	0.008 ± 0.00028	0.204 ± 0.00042	0.014 ± 0.00030	0.171 ± 0.00045
Octadecanol acetate	2209	-	0.042 ± 0.00028	0.050 ± 0.00042	0.017 ± 0.00059	0.025 ± 0.00045
1-Heptatriacotanol	2309	2309	0.015 ± 0.00028	0.039 ± 0.00042	0.036 ± 0.00030	-
Hydrocarbons			0.279	0.380	0.114	0.461
Eicosane*	2000	2000	0.042 ± 0.00110	-	0.013 ± 0.00030	0.062 ± 0.00045

Heneicosane*	2100	2100	0.020 ± 0.00028	-	0.009 ± 0.00030	0.025 ± 0.00271
Docosane*	2200	2200	0.032 ± 0.00028	-	0.011 ± 0.00030	0.054 ± 0.00045
Tricosane*	2300	2300	0.017 ± 0.00028	0.036 ± 0.00084	-	-
Tetracosane*	2400	2400	-	0.020 ± 0.00042	0.026 ± 0.00030	0.031 ± 0.00045
Pentacosane*	2500	2500	0.018 ± 0.00028	0.010 ± 0.00042	-	-
Hexacosane*	2600	2600	0.070 ± 0.00028	0.129 ± 0.00042	0.029 ± 0.00059	0.038 ± 0.00136
Heptacosane*	2700	2700	0.087 ± 0.00028	0.135 ± 0.00042	0.009 ± 0.00030	0.047 ± 0.00045
Octacosane*	2800	2800	0.025 ± 0.00028	0.171 ± 0.00042	0.017 ± 0.00059	0.179 ± 0.00045
Total identification			2.576	3.670	2.744	4.158

(mg/g)

Retention indices (RI) were determined relative to a series of *n*-alkanes (C8–C40) on capillary columns VF5-ms (RI¹) [24] and CPWax 52 (RI²) [25]; Identification method: RI, comparison of RIs with those listed in a homemade library, reported in the literature [24], and/or authentic samples; comparison of mass spectra with those in mass spectral libraries NIST02 and Wiley 9; * co-injection with reference compounds; SD, standard deviation.

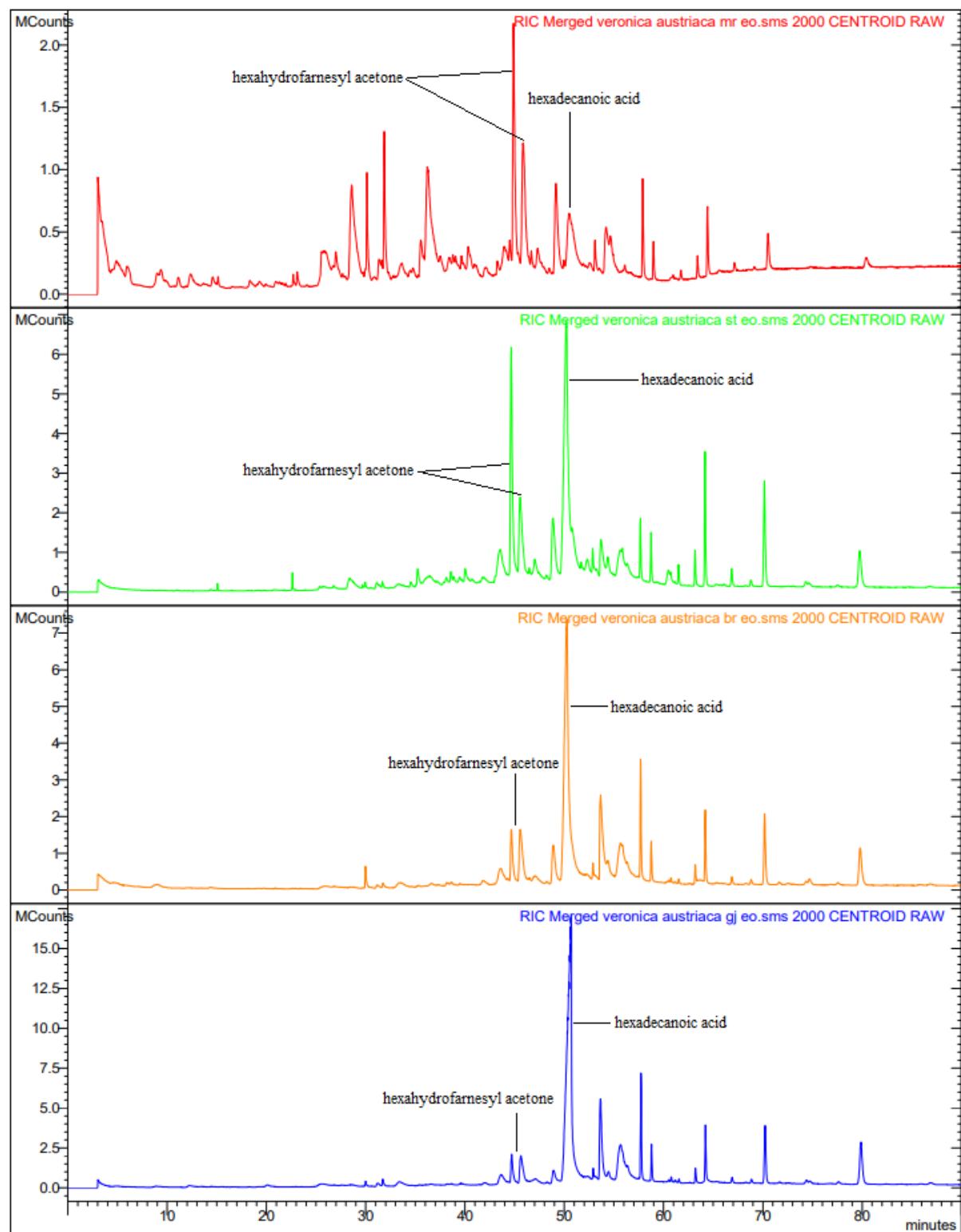


Figure S1. RIC chromatograms of the four samples of EOs of *V. jacquinii* with marked major compounds

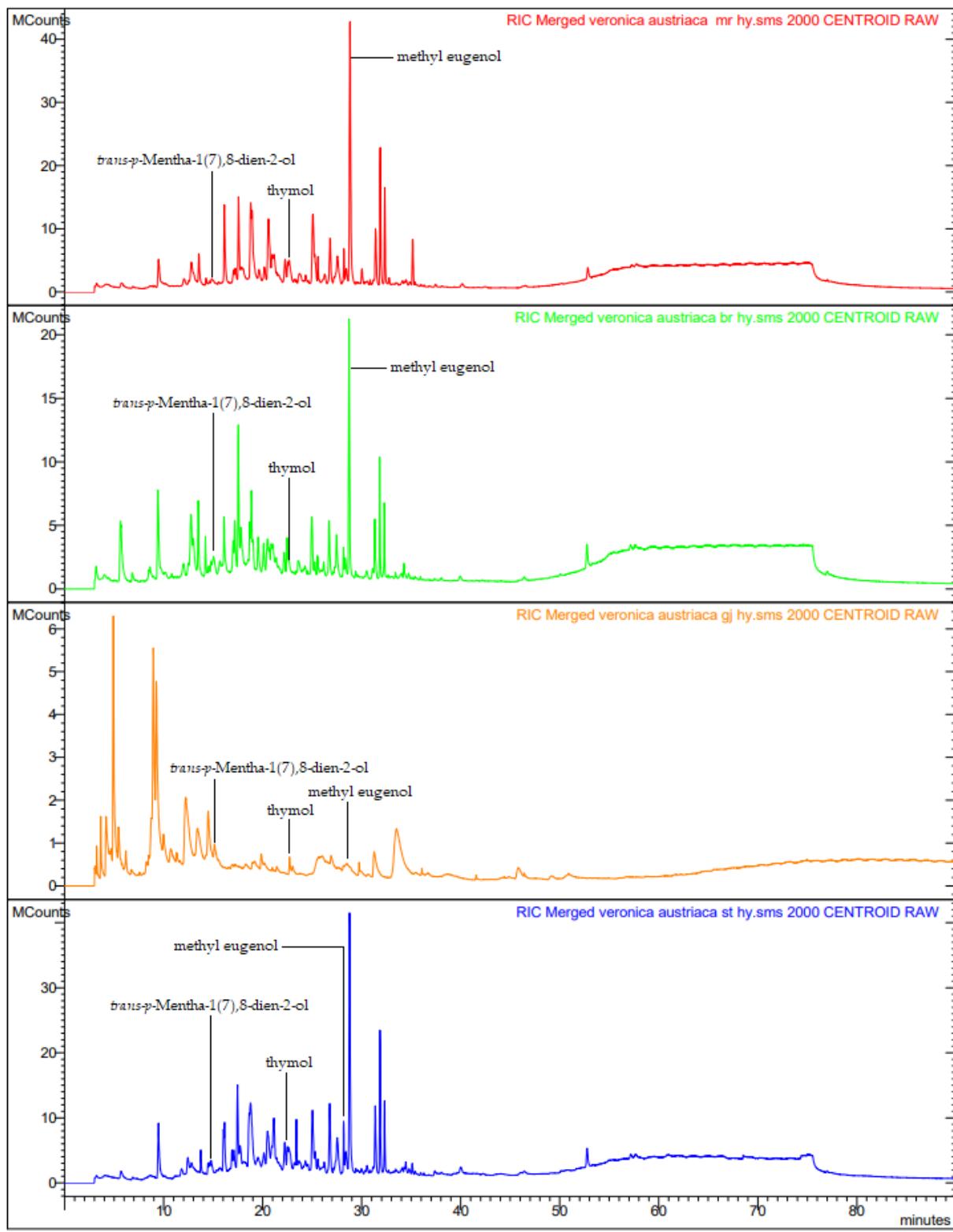


Figure S2. RIC chromatograms of the four samples of hydrosols of *V. jacquinii* with marked major compounds