

## Supplementary information

**Table S1.** HPLC-HR-MS data of identified phenolic compounds from *Lavandula viridis* L'Hér and *Thymus lotocephalus* G. Lopez & R. Morales extracts.

Peak	Compound identity	Chemical formula	Theoretical exact mass [M-H] <sup>-</sup> (m/z)	Delta ppm (error)	RT (min)	MSI MI level*
1	O-Caffeoylquinic acid <sup>a</sup>	C <sub>16</sub> H <sub>18</sub> O <sub>9</sub>	353.0867	2.103	1.48	2
2	Caffeic acid hexoside <sup>a</sup>	C <sub>15</sub> H <sub>18</sub> O <sub>9</sub>	341.0867	-0.007	2.9	2
3	Caffeic acid hexoside <sup>a</sup>	C <sub>15</sub> H <sub>18</sub> O <sub>9</sub>	341.0867	0.143	3.5	2
4	Fertaric acid <sup>a</sup>	C <sub>14</sub> H <sub>14</sub> O <sub>9</sub>	325.0557	0.774	4.23	2
5	Caffeic acid <sup>a</sup>	C <sub>9</sub> H <sub>8</sub> O <sub>4</sub>	179.0338	-0.604	4.59	1
6	Ferulic acid <sup>a</sup>	C <sub>10</sub> H <sub>10</sub> O <sub>4</sub>	193.0495	-0.374	7.59	1
7	Luteolin-7-O-glucuronide <sup>c</sup>	C <sub>21</sub> H <sub>18</sub> O <sub>12</sub>	461.0714	0.569	8.11	2
8	Rosmarinic acid <sup>a</sup>	C <sub>18</sub> H <sub>16</sub> O <sub>8</sub>	359.0761	0.488	9.57	1
9	Salvianolic acid A isomer I <sup>a</sup>	C <sub>26</sub> H <sub>22</sub> O <sub>10</sub>	493.1129	0.508	10.11	2
10	Salvianolic acid A isomer II <sup>a</sup>	C <sub>26</sub> H <sub>22</sub> O <sub>10</sub>	493.1129	0.358	10.97	2
11	Salvianolic acid I <sup>a</sup>	C <sub>27</sub> H <sub>22</sub> O <sub>12</sub>	537.1027	1.499	11.13	2
12	Salvianolic acid B <sup>a</sup>	C <sub>36</sub> H <sub>30</sub> O <sub>16</sub>	717.1450	-0.049	11.23	2
13	Herniarin <sup>b</sup>	C <sub>10</sub> H <sub>8</sub> O <sub>3</sub>	175.0389	-0.429	12.14	2
14	Salvianolic acid A isomer III <sup>a</sup>	C <sub>26</sub> H <sub>22</sub> O <sub>10</sub>	493.1129	-0.132	12.8	2
15	Apigenin <sup>c</sup>	C <sub>15</sub> H <sub>10</sub> O <sub>5</sub>	269.0444	0.172	14.8	1

Class of compounds: Phenolic acids (<sup>a</sup>Hydroxycinnamic acids); <sup>b</sup>Coumarin derivative; Flavonoids (<sup>c</sup>Flavones).

\*Metabolite standards initiative metabolite identification (MSI MI) levels. Reference standards were available for all compounds identified at MSI MI level 1.

**Table S2.** Summary of HPLC-HR-MS parameters for quantification of phenolic compounds in *Lavandula viridis* L'Hér and *Thymus lotocephalus* G. Lopez & R. Morales extracts.

Compound	Linear range (mg/L)	Intercept	Slope	R <sup>2</sup>	LOD (mg/mL)	LOQ (mg/mL)
Rosmarinic acid	0.1-100	-103410	1598747	0.9992	0.06	0.19
Caffeic acid	0.1-10	1436398	3950274	0.9990	0.50	1.68
Ferulic acid	0.2-60	2487	9717	0.9995	0.37	1.25
Gallic acid	0.2-15	-330994	2547719	0.9998	0.00	0.01
Chlorogenic acid	0.2-15	-243627	1363025	0.9993	0.00	0.01
Luteolin	0.2-15	1464030	5454860	0.9990	0.29	0.96
Apigenin	0.2-15	185082	6649959	0.9995	0.00	0.01
Naringenin	0.2-15	-314843	2143048	0.9996	0.37	1.23
<i>p</i> -Coumaric acid	0.2-15	-2024	73034	0.9999	0.22	0.72