

## Supplemental material

**Table A1.** Experimental domain and codification of independent variables in the *CCCD* factorial design with 5 range levels.

<b>CODED VALUES</b>	<b>NATURAL VALUES</b>		
	<i>t (min)</i>	<i>T (°C)</i>	<i>S (%)</i>
<b>-1.68</b>	20	25	0
<b>-1</b>	40,3	37,2	20.3
<b>0</b>	70	55	50
<b>+1</b>	99,7	72,8	79.7
<b>+1.68</b>	120	85	100

**Table A2.** Mathematical models of the extraction process derived from the second-order polynomial model with interactions of Eq. (1).

<i>Extraction yield</i>	$Y_{Yield} = 40.8 + 1.75t + 2.92T - 3.86S - 5.56S^2 + 0.63tS + 0.43TS$	Eq. (3)
<i>Hydrolysable tannins (Hta)</i>	$Y_{Elg} = 9.34 - 2.68t + 2.34T - 6.5S + 2.38t^2 - 3.04S^2 + 1.76TS$	Eq. (4)
<i>Flavonoids (Fla)</i>	$Y_{Fla} = 5.56 - 0.72t - 1.12S + 0.53t^2 + 1.28S^2 - 0.65tS + 0.89TS$	Eq. (5)
<i>Total Phenolics (Phe)</i>	$Y_{Phe} = 14.9 - 3.4t + 2.11T - 7.62S + 2.91t^2 + 4.32S^2 - 0.87tS + 2.65TS$	Eq. (6)
<i>Compound P5</i>	$Y_{P5} = 0.37 - 0.31t + 0.37t^2 + 0.37T^2 + 0.412S^2 - 0.12tT - 0.39tS + 0.15TS$	Eq. (7)
<i>Compound P7</i>	$Y_{P7} = 7.88 - 1.68t + 2.09t^2 - 6.15S + 2.81t^2 + 1.02tT + 1.25tS$	Eq. (8)
<i>Compound P8</i>	$Y_{P8} = 1.22 - 0.17T - 0.43S + 0.33t^2 + 0.28S^2 - 0.2tT - 0.37tS + 0.39TS$	Eq. (9)
<i>Compound P9</i>	$Y_{P9} = 1.55 - 0.23t - 0.13T - 0.17S + 0.15S^2 + 0.19tT + 0.14tS + 0.24TS$	Eq. (10)
<i>Compound P11</i>	$Y_{P11} = 1.06 - 0.06t - 0.29S + 0.19S^2$	Eq. (11)
<i>Compound P13</i>	$Y_{P13} = 0.64 - 0.1t - 0.05T + 0.05S + 0.15S^2 - 0.05tS$	Eq. (12)
<i>Compound P14</i>	$Y_{P14} = 0.51 - 0.08t + 0.04T + 0.02S + 0.04t^2 + 0.06S^2 + 0.02tT - 0.04tS + 0.07TS$	Eq. (13)
<i>Compound P1,2,3,4,6,10,12</i>	$Y_{Minor} = 2.61 - 0.89t - 0.53S + 1.35S^2 - 1.12tT - 1.40tS + 1.53TS$	Eq. (14)

**Table A3:** Parametric results of the dose-response model of Eq. (2) for of all the response values assessed in terms of the variation of the *S/L* ratio: *Extraction yield* (%), T. hydrolysable tannins (*Hta*), T. Flavonoids (*Fla*), T. Phenolics (*Phe*) and major (*P5, P7, P8, P9, P11, P13* and *P14*) and minor (*P1, P2, P3, P4, P6, P10* and *P12*) compounds. The underlined values were obtained with an increasing dose-response model of Eq. (2) using the  $m_{99\%}$  as response criteria.

	<b><i>K</i></b>	<b><i>m</i><sub>50%</sub> or <i>m</i><sub>99%</sub></b>	<b><i>a</i></b>	<b><i>R</i><sup>2</sup></b>
<i>Extraction yield</i>	54.96±3.85	258.40±25.24	0.22±0.02	0.9241
<i>Hydrolysable tannins (Hta)</i>	85.73±5.14	<u>165.69</u> ±7.12	0.86±0.01	0.9836
<i>Flavonoids (Fla)</i>	35.35±4.45	31.01±3.89	0.44±0.03	0.9436
<i>Total Phenolics (Phe)</i>	205.40±18.49	<u>337.80</u> ±11.56	0.33±0.01	0.9482
<i>Compound P5</i>	8.65±0.78	<u>112.81</u> ±3.39	0.84±0.08	0.9479
<i>Compound P7</i>	47.62±2.38	<u>124.35</u> ±7.39	0.79±0.08	0.9696
<i>Compound P8</i>	17.92±1.43	10.28±0.72	0.43±0.01	0.9189
<i>Compound P9</i>	4.31±0.30	379.82±25.19	1.00±0.05	0.9457
<i>Compound P11</i>	3.12±0.09	253.13±24.25	1.02±0.09	0.9249
<i>Compound P13</i>	3.58±0.29	17.62±0.18	0.44±0.01	0.9563
<i>Compound P14</i>	2.58±0.13	28.93±0.29	0.94±0.09	0.8648
<i>Compound P1,2,3,4,6,10,12</i>	116.59±8.16	<u>482.16</u> ±8.89	0.79±0.04	0.9676