

**Table S1.** Pseudo-first-order and pseudo-second-order model parameters.

<b>Pseudo-first-order model</b>		
<b><math>k_1</math> (min<sup>-1</sup>)</b>	<b><math>q_e</math> (μg/mg)</b>	<b><math>R^2</math></b>
6.43	61.41	0.94
<b>Pseudo-second-order model</b>		
<b><math>k_2</math> (μg mg<sup>-1</sup>min<sup>-1</sup>)</b>	<b><math>q_e</math> (μg/mg)</b>	<b><math>R^2</math></b>
0.013	63.36	0.98

**Table S2.** Langmuir isotherm constants obtained for the cypermethrin adsorption.

<b><math>Q_0</math> (μg/g)</b>	<b><math>K_L</math> (L/μg)</b>	<b><math>R_L</math></b>	<b><math>R^2</math></b>
588.24	0.0027	0.79	0.98

**Table S3.** Freundlich isotherm constants obtained for the cypermethrin adsorption.

<b>n</b>	<b><math>K_F</math> (μg/L)</b>	<b><math>R^2</math></b>
1.53	5.56	0.88