

*Supporting Information for*

Exploring the Potential of Biochar Derived from Chinese Herbal  
Medicine Residue for Efficient Removal of Norfloxacin

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## **Characterization**

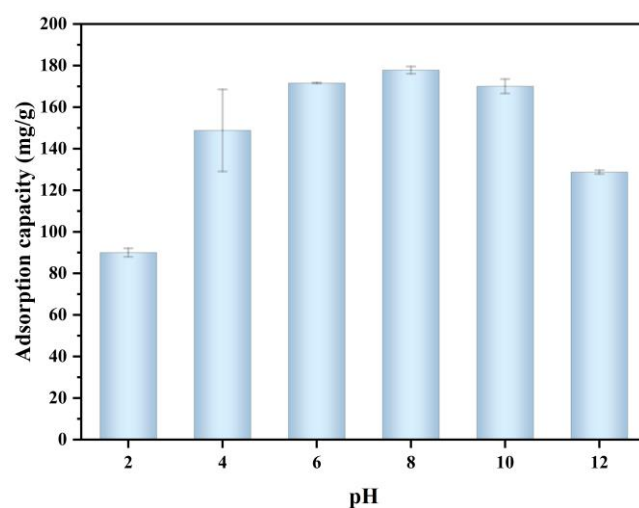
A physisorption analyzer (Autosorb IQ<sub>2</sub>, Quantachrome) was used to determine the pore structures of ABL biochars. An elemental analyzer (EA, Elementar UNICUBE) was utilized to analyze the element composition. The scanning electronic microscopy (SEM, Zeiss Gemini 300) coupled with energy dispersive spectroscopy (EDS, Zeiss Gemini 300) was used to examine the surface morphology and chemical composition of biochar samples. A Fourier transform infrared spectrometer (FTIR, Nicolet iS5, Thermo Scientific) was used to determine the surface functional groups, while an X-ray photoelectron spectroscopy (XPS, Thermo-Scientific K-Alpha) was utilized to investigate their chemical compositions. By using an X-ray diffractometer (XRD, Bruker D8), the crystal structure was investigated. Finally, a thermogravimetric analyzer (TGA, TA550) was used to determine the thermal stability.

## FIGURE CAPTIONS

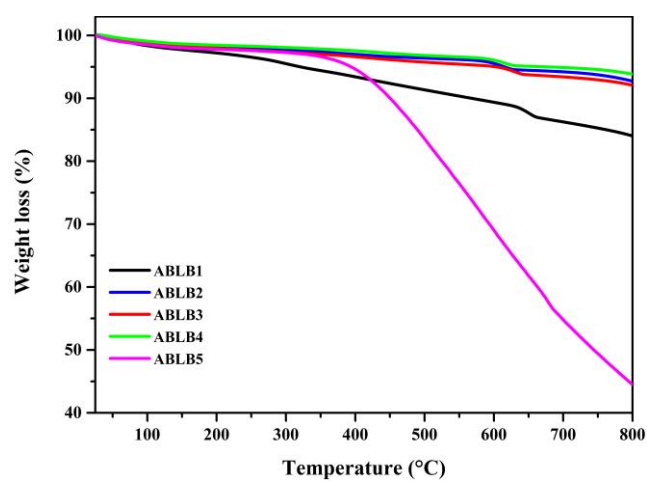
- Figure S1      Effect of pH on the adsorption capacity of NOR.
- Figure S2      TGA curve of ABL biochar samples.
- Figure S3      NOR adsorption effect comparison of ABLB4 and commercial biochar.
- Figure S4      Response surface method (RSM) regression analysis.
- Figure S5      Full-wavelength scan of NOR.
- Figure S6      SEM-EDX microphotographs of ABLB4.
- Figure S7      Temkin isotherm of ABLB4.
- Figure S8      UV full-wavelength scanning spectra of NOR with time. (b) UV full-wavelength scanning spectra of NOR at different concentration.

## TABLE CAPTIONS

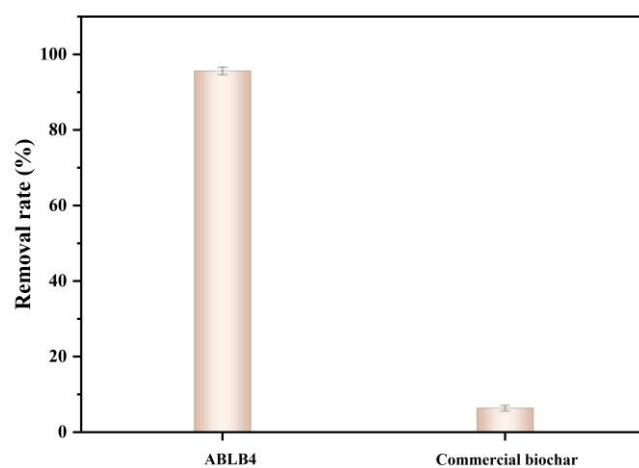
Table S1	Detailed information of involved formulas and models.
Table S2	Kinetic constants of pseudo-first-order and pseudo-second-order models.
Table S3	Kinetic constants for the IPD model.
Table S4	Lists of maximum NOR adsorption capacity on various biochar and other adsorption materials.
Table S5	Thermodynamic parameters for NOR adsorption.
Table S6	Energy consumption of biochar used in industrial wastewater production (\$/t).
Table S7	Elemental composition of raw ABL residues.
Table S8	ICP-OES analysis of raw ABL residues.



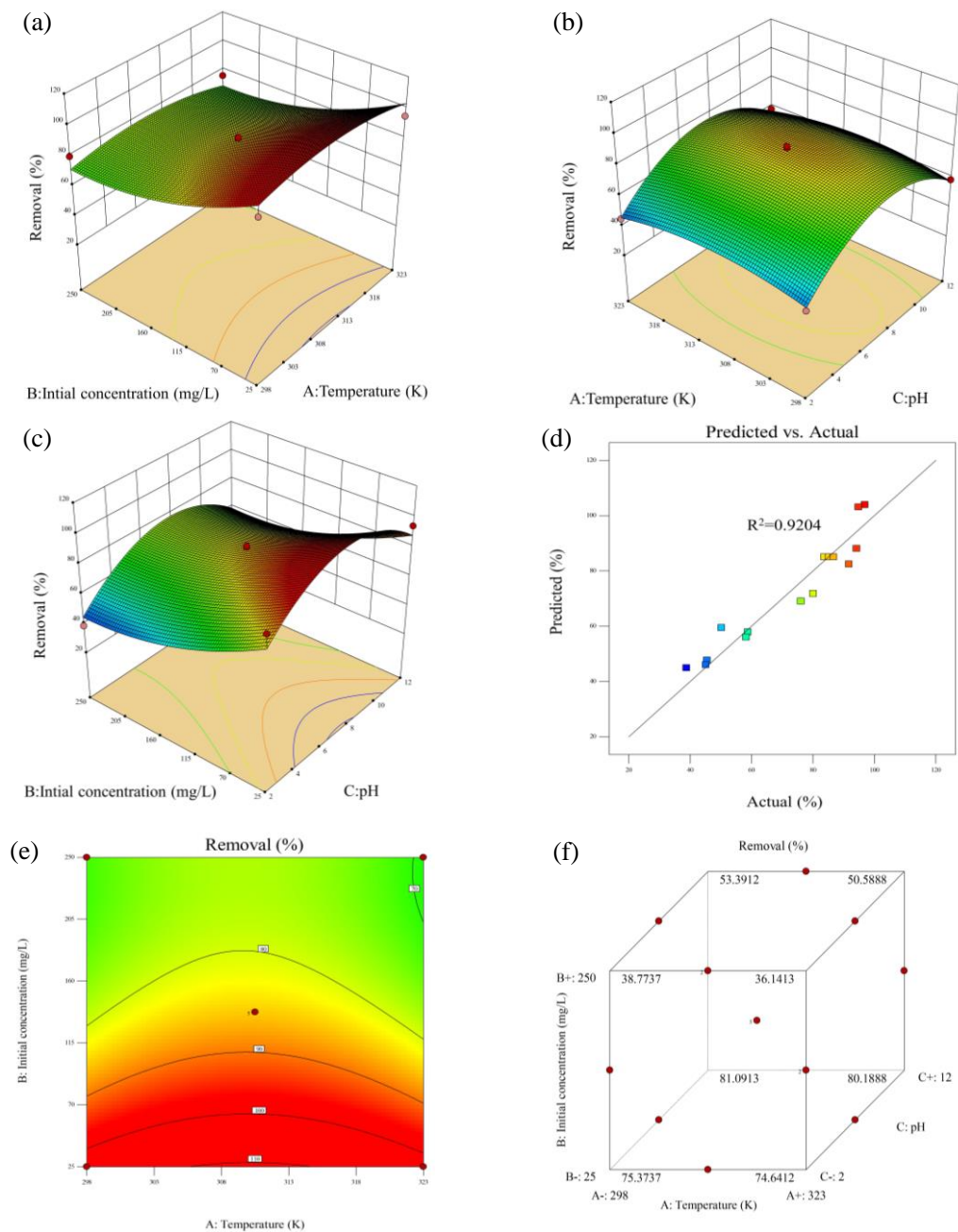
**Figure S1.** Effect of pH on the adsorption capacity of NOR (initial concentration 200 mg/L, adsorbent dosage 10 mg, contact time 1440 min).



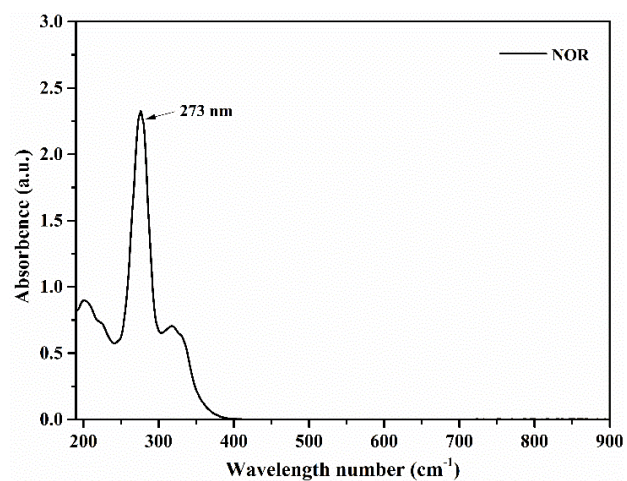
**Figure S2.** TGA curve of ABL biochar.



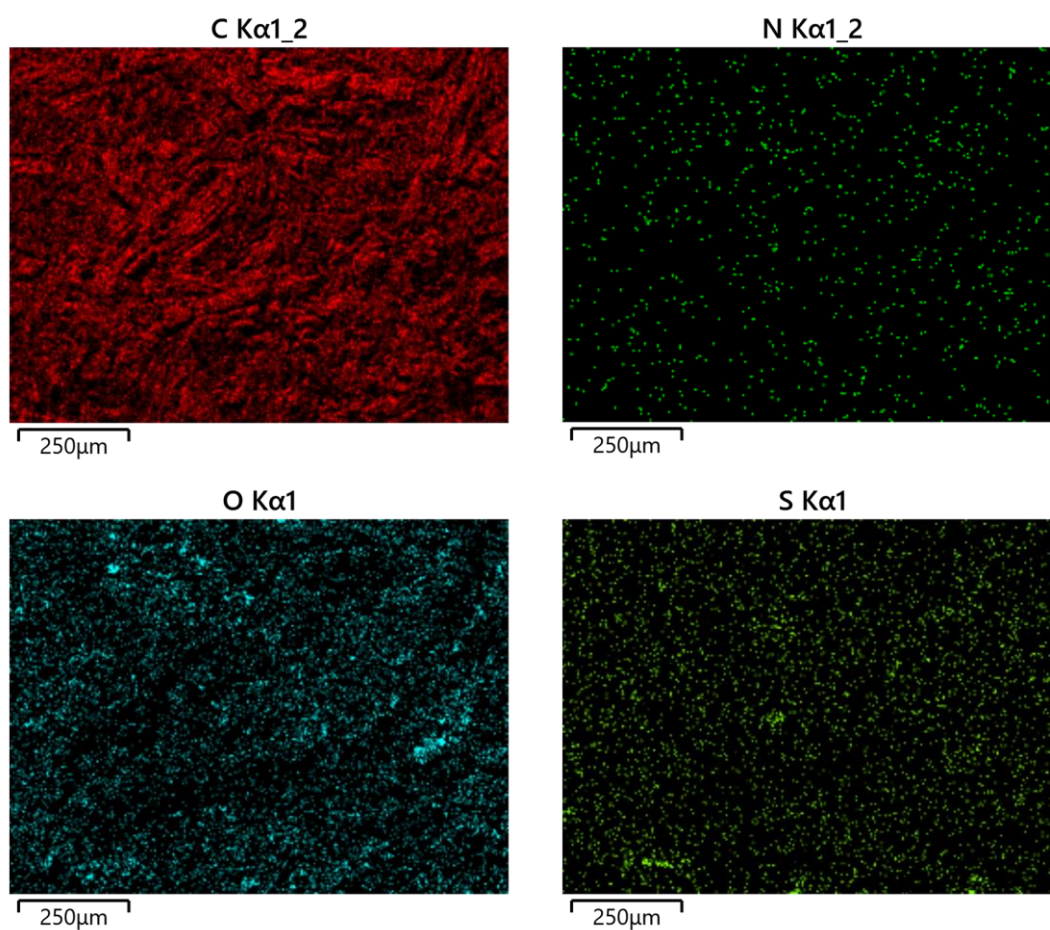
**Figure S3.** NOR adsorption effect comparison of ABLB4 and commercial biochar.



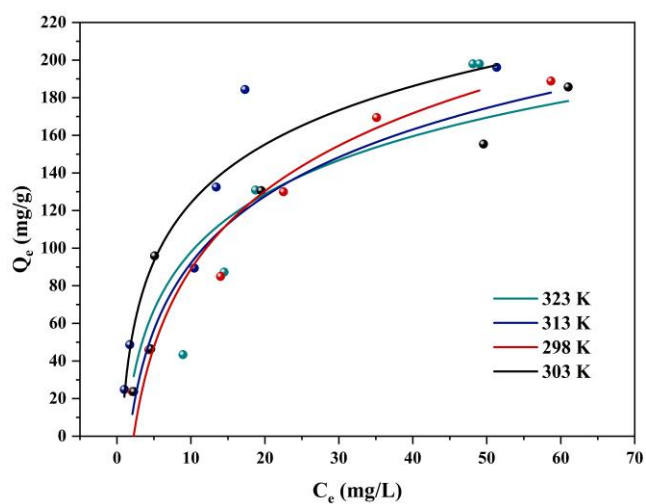
**Figure S4.** (a)-(c) Optimization of the NOR removal rate using response surface methodology. (d) Comparison between the predicted value with the actual removal rate. (e)-(f) contour plot and cube optimization for NOR removal rate.



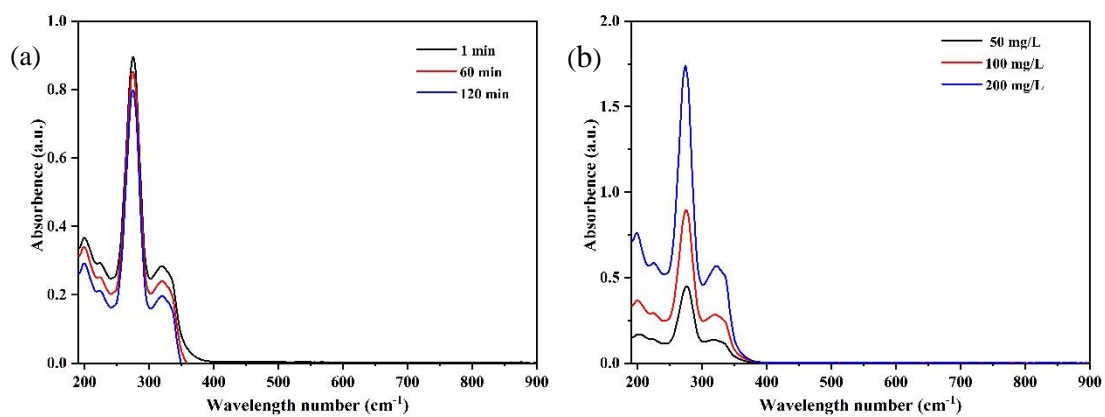
**Figure S5.** Full-wavelength scan of NOR.



**Figure S6.** SEM-EDX microphotographs of ABLB4.



**Figure S7.** Temkin isotherm of ABLB4.



**Figure S8.** (a) UV full-wavelength scanning spectra of NOR with time. (b) UV full wavelength scanning spectra of NOR at different concentration.



**Table S1**

Detailed information of involved formulas and models.

Formulas and models	Equation	Parameters
Adsorption capacity	$q_e = \frac{(C_0 - C_e)V}{m}$	$q_e$ (mg/g) are the adsorbed amount at an equilibrium concentration ( $C_e$ , mg/g) $C_0$ (mg/L): initial NOR concentration $C_e$ (mg/L): equilibrium NOR concentration $V$ (L): reaction solution volume of NOR $m$ (g): the sorbent mass
Removal rate	$E = \frac{(C_0 - C_e)}{C_0}$	$C_0$ (mg/L): initial NOR concentration $C_e$ (mg/L): equilibrium NOR concentration
Pseudo-first order	$q_t = q_e(1 - e^{-k_1 t})$	$q_e$ and $q_t$ (mg/g) are the adsorbed amount at an equilibrium concentration ( $C_e$ , mg/g) and a predetermined time ( $t$ , min) $k_1$ (min <sup>-1</sup> ) is the rate constants for the PFO model
Pseudo-second order	$q_t = \frac{k_2 q_e^2 t}{1 + k_2 q_e t}$	$k_2$ (g/mg min) is the rate constants for the PSO model
Weber-Morris intraparticle diffusion	$q_t = K_i t^{0.5} + C_i$	$K_i$ (mg/(g·min <sup>0.5</sup> )) is the adsorption rate constants of intra-particle diffusion model and $C_i$ is the constant for film thickness of the intra-particle diffusion model.
Langmuir	$q_e = \frac{q_{\max} K_L C_e}{1 + K_L C_e}$	$q_{\max}$ (mg/g) is the maximum adsorption capacity of the adsorbent $q_e$ (mg/g) is the adsorption capacity at equilibrium $C_e$ (mg/L) is the adsorbate concentration at equilibrium $K_L$ (L/mg) is the constant for the affinity between the adsorbate and the adsorbent
Freundlich	$q_e = K_F C_e^{1/n}$	$K_F$ (mg/g)/(mg/L) <sup>n</sup> is the Freundlich constant $n$ is a dimensionless Freundlich intensity parameter
Van der Hoff equation	$\Delta G = -RT \ln K_d$  $\ln K_d = \frac{\Delta S}{R} - \frac{\Delta H}{RT}$	$\Delta G$ , $\Delta H$ and $\Delta S$ are the Gibbs energy (KJ/mol), activated enthalpy (KJ/mol) and the change in entropy (J/mol K), respectively; $K_d$ is the distribution coefficient; $R$ is the ideal gas constant, whose value is 8.314 J/mol·K; $T$ is absolute temperature (K).

**Table S2**

Kinetic constants of pseudo-first-order and pseudo-second-order models.

Temperature	$q_{e,exp}$ (mg/g)	Pseudo-first-order			Pseudo-second-order		
		$K_1$ (min <sup>-1</sup> )	$q_e$ (mg/g)	R <sup>2</sup>	$K_2$ (min <sup>-1</sup> )	$q_e$ (mg/g)	R <sup>2</sup>
298 K	188.54	0.2005	189.80	0.9984	0.00094	228.73	0.9967
303 K	181.36	0.1908	183.23	0.9923	0.00119	213.93	0.9858
313 K	169.58	0.2366	169.96	0.9830	0.00219	188.42	0.9826
323 K	155.39	0.3924	153.45	0.9888	0.00333	170.41	0.9844

**Table S3**

Weber-Morris intraparticle diffusion model of NOR adsorption by ABLB4.

Temperature	First linear segment			Second linear segment			Third linear segment		
	$K_{id}$ (mg/g/ min <sup>0.5</sup> )	$C_i$ (mg/g)	R <sup>2</sup>	$K_{id}$ (mg/g/ min <sup>0.5</sup> )	$C_i$ (mg/g)	R <sup>2</sup>	$K_{id}$ (mg/g/ min <sup>0.5</sup> )	$C_i$ (mg/g)	R <sup>2</sup>
298 K	72.00	35.80	0.9979	16.45	112.30	0.9839	5.14	163.80	0.7750
303 K	56.57	13.18	0.9904	19.41	95.55	0.9310	4.07	161.68	0.8311
313 K	66.52	18.72	0.9693	27.73	62.60	0.9786	5.76	142.14	0.7311
323 K	57.32	5.38	0.9907	14.44	100.63	0.9977	3.26	140.24	0.7790

**Table S4**

Lists of maximum NOR adsorption capacity on various biochar and other adsorption materials.

Raw material	BET (m <sup>2</sup> /g)	Total pore volume (cm <sup>3</sup> /g)	Adsorbent dosage (g/L)	NOR concentration (mg/L)	$q_e$ (mg/g)	Ref.
Luffa sponge	822	0.59	0.5	50-200	250	[1]
Salix mongolica	1348	-	2	25-266	210	[2]
Porous resins and carbon nanotube	160	-	5	20-100	154	[3]
Spent coffee grounds	46	-	1	10-50	70	[4]
Cauliflower roots biochar	232	-	2	10-300	31	[5]
Pristine potato biochar	90	0.12	4	4-20	5	[6]
<i>Atropa belladonna</i> L. residue	812	0.49	1	25-250	251	This study

**Table S5**

Thermodynamic parameters for NOR adsorption.

Material	$\Delta G$ (kJ/mol)				$\Delta H$ (kJ/mol)	$\Delta S$ (J/mol/K)
	298 K	303 K	313 K	323 K		
ABLB4	-6.44	-6.20	-5.29	-3.27	-43.79	-124.45

**Table S6**

Energy consumption of biochar used in industrial wastewater production (\$/t).

Sample	Raw material	Disposal cost	Energy consumption	Wages of workers	Total cost	Pay back
ABLB4	1.5	15.3	178.2	140.5	335.5	852.5
Commercial biochar	311.6	28.3	518.1	330	1188	

**Table S7**

Elemental composition of raw ABL residues.

Sample	C (%)	H (%)	O (%)	N (%)	S (%)
ABL residue	46.13	4.96	42.26	1.78	0.16

**Table S8**

ICP-OES analysis of raw ABL residues.

Sample	Si	Ca	Mg	K	Fe	Al	Na
ABL residues	0.016	0.35	0.19	0.47	0.066	0.037	0.027

## References

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