

Hydrothermal synthesis of Mg/Al-layered double hydroxide modified water hyacinth hydrochar for remediation of wastewater containing mordant brown dye

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Supporting Information

S1: Instruments

The WH, MgAl LDHs, MgAl@WH were characterized using a diversity of analytical techniques. Fourier transform infrared (FTIR) spectra were recorded using a Nicolet-Nexus 670 spectrometer equipped with a diamond ATR accessory. The measurement employed a resolution of 4 cm⁻¹ and 32 scans. The crystal structure of the sample was analyzed using powder X-ray diffraction (XRD) on a PANalytical MPD X'Pert PRO diffractometer. The measurements employed 45 kV, 40 mA Ni-filtered Cu K α 1 radiation at room temperature. X-ray photoelectron spectra (XPS) were obtained on a Thermo-Fisher ESCALAB 250 spectrometer equipped with a micro-focused, monochromated Al K α X-ray source operating at 15 kV and a double-focusing, full 180° spherical sector electron analyzer. Scanning electron microscopy (SEM): A Hitachi SU-70 field-SEM with an energy of 5.0 kV was used to take the SEM images. Transmission electron microscopy (TEM): A JEOL JEM1400 TEM was used to obtain the TEM images at 100 kV.

S2: Determination of pH_{PZC}

Solutions with pH between 2 and 12 were obtained by adding NaOH (0.1 M) or HCl (0.1 M) to Erlenmeyer flasks containing 50 mL of 0.01 M NaCl solution. Then, 0.15 g of the adsorbent was added to each flask; final pH measurements of the bottles were made after shaking for 48 h. The pH_{PZC} of the sample is calculated from the point where pH_{final} = pH_{initial} crosses the curve between pH_{final} and pH_{initial}.

Table S1: Remediation of various anionic dyes using MgAl LDHs and/or biochars prepared through different reported methods.

Adsorbent	Target dye	Q _e (mg/g)	Ref.
Date palm derived biochar	Eriochrome black T	309.59	1
Date palm derived biochar	Methylene orange	163.12	1
MgAl-LDH	Methylene orange	197.62	2
MgAl-Charcoal activated	direct yellow	133.33	3
hydrochar from olive waste	Congo red	99.0	4
Mg/Al/ pine sawdust biochar	Methyl Orange (MO)	21.8	5
Mg/Al/ pine sawdust biochar	Sunset Yellow FCF (SYF)	23.6	5
Biochar derived from <i>Caulerpa scalpelliformis</i>	Reactive Yellow 81	151.5	6
new biochar derived from pecan nutshell	Reactive red 141	130.0	7
Fe ₂ O ₃ -Biochar	Methyl Orange (MO)	46.6	8
functionalized magnetic bamboo biochar adsorbent	Methyl Orange (MO)	305.4	9
MgAl@WH	Mordant brown 15 dye	311.0	Present work

S3.

The Langmuir and Freundlich isotherm models are two extensively used mathematical models. The Langmuir model assumes a monolayer coverage and that all the adsorbent sorption sites are the same while the Freundlich isotherm model assumes that the coverage is multilayer and that all the adsorption sites are heterogenous. The Langmuir and Freundlich models are presented as Eq.(1) and Eq. [2], respectively, as follows:¹⁰⁻¹¹

$$\ln q_e = \ln K_f + \frac{1}{n} \ln C_e \quad (1)$$

$$\frac{C_e}{q_e} = \frac{1}{b Q_0} + \frac{C_e}{Q_0} \quad (2)$$

Where q_e , C_e , b , and Q_0 are the equilibrium adsorption capacity, the equilibrium concentration of the metal ions, the Langmuir constant, and the Langmuir monolayer adsorption capacity, respectively. $1/n$ and K_f are the adsorption intensity and the Freundlich constant, respectively.

Moreover, the essential feature of the Langmuir isotherm can be defined as R_L parameter given by Eq. (5).¹²

$$R_L = \frac{1}{1 + b c_0} \quad (5)$$

The separation factor (R_L) can be used to indicate the shape of the adsorption behavior to be either irreversible ($R_L = 0$), linear ($R_L = 1$), unfavorable ($R_L > 1$), or favorable ($0 < R_L < 1$).¹³

S4.

The PFO model usually predicts the behavior at the initial stage of the adsorption process, while PSO model predicts the behavior at all stages of the adsorption process¹⁴.

Pseudo-first-order kinetic model:

$$\log(q_e - q_t) = \log q_e - \frac{K_1 t}{2.303} \quad (3)$$

Pseudo-second-order kinetic model:

$$\frac{t}{q_t} = \frac{1}{K_2 q_e^2} + \frac{t}{q_e} \quad (4)$$

Where k_1 (min^{-1}) and k_2 ($\text{g mol}^{-1} \text{min}^{-1}$) are the rate constants. q_t and q_e are the adsorption uptake of heavy metal at time t (min) and at equilibrium. Where k_1 (min^{-1}) and k_2 ($\text{g mol}^{-1} \text{min}^{-1}$) are the rate constants. q_t and q_e are the adsorption uptake of heavy metal at time t (min) and at equilibrium.

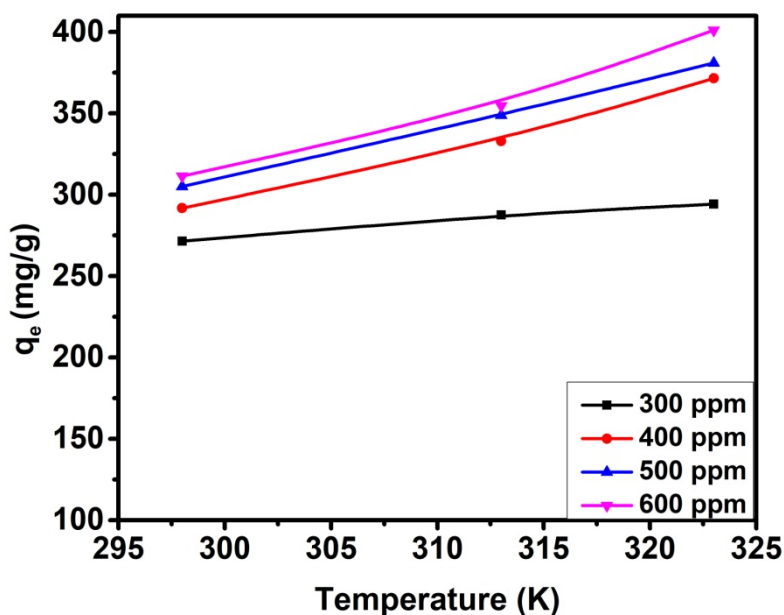


Figure S1: Effect of temperature on mordant brown dye adsorption onto MgAl@WH.

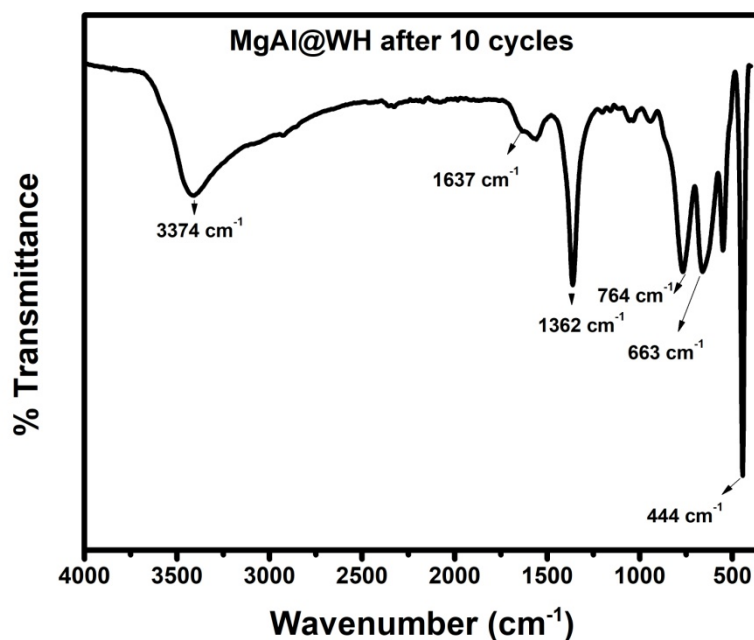


Figure S2: FTIR spectra of MgAl@WH biochar after 10 cycles.

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