

Supporting Information

Instantaneous adsorption of synthetic dyes from the aqueous environment using kaolinite nanotubes; equilibrium and thermodynamic studies

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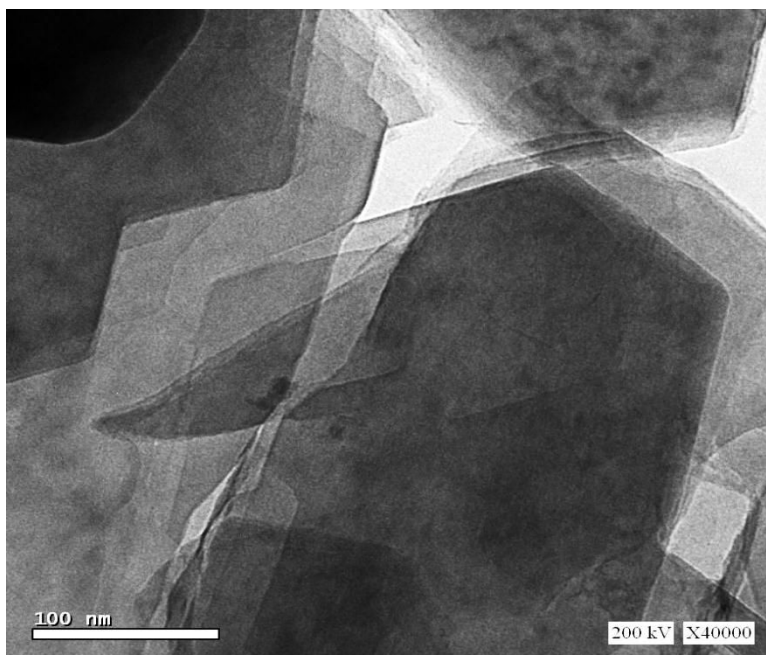


Figure S1. TEM image of raw kaolinite mineral

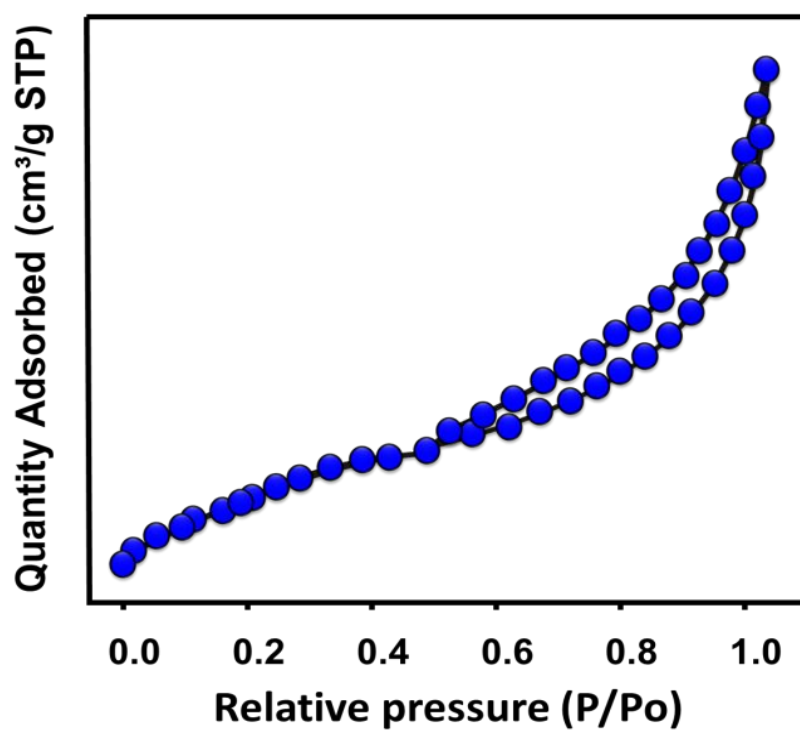


Figure S2. the nitrogen adsorption/desorption isotherm curves of KNTs

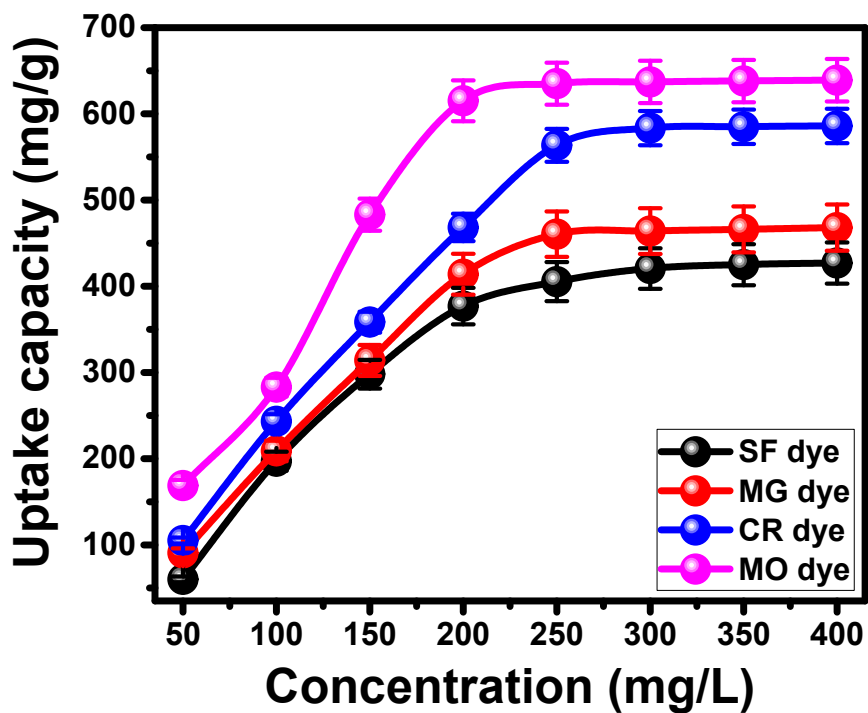


Figure S3. the uptake capacity of KNTs at different concentrations of SF, MG, CR, and MO dyes

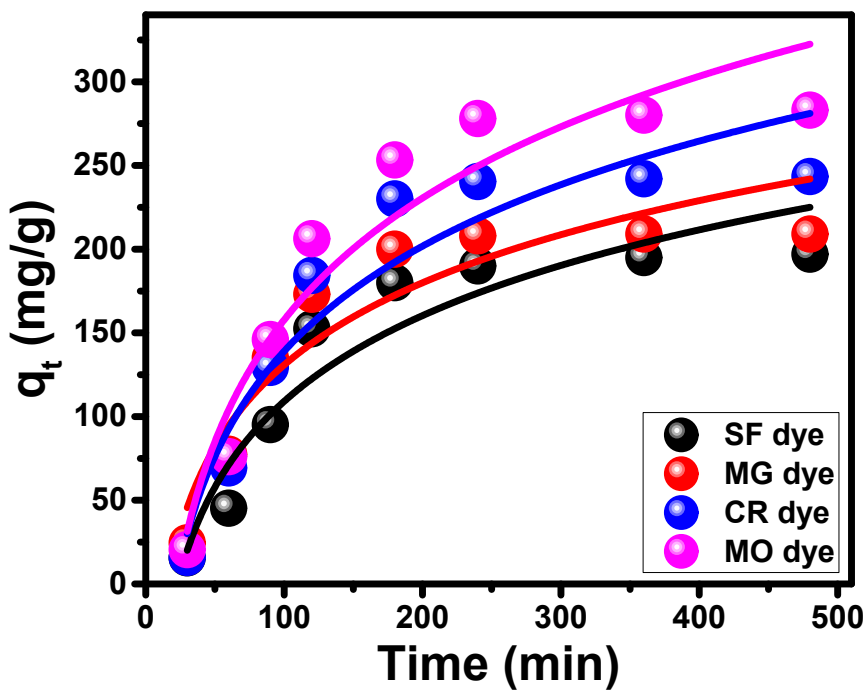


Figure S4. Fitting of the dyes adsorption results with Elovich kinetic model

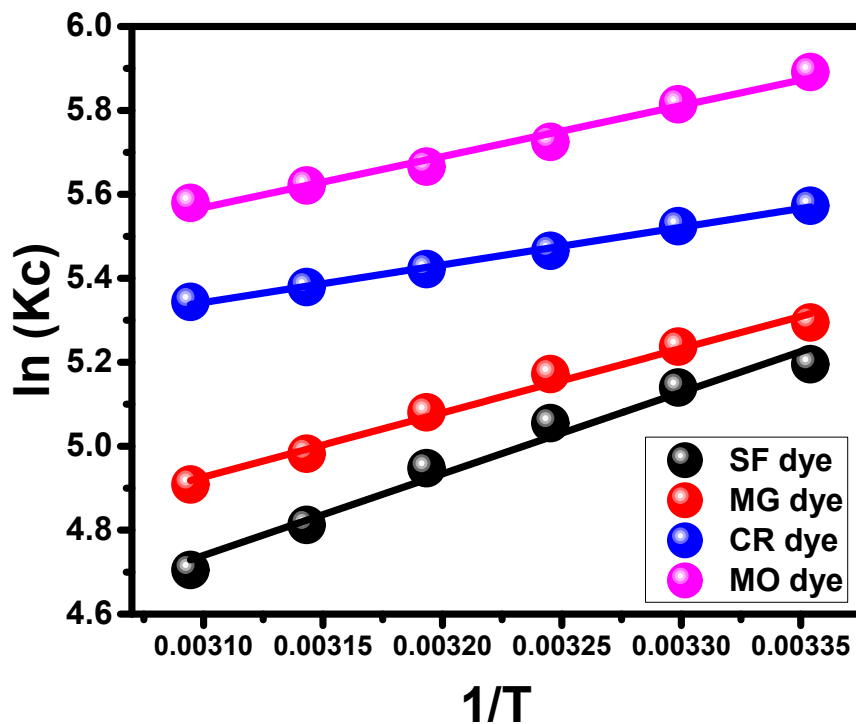


Figure S5. fitting of the dyes adsorption results with Van't Hoff equation

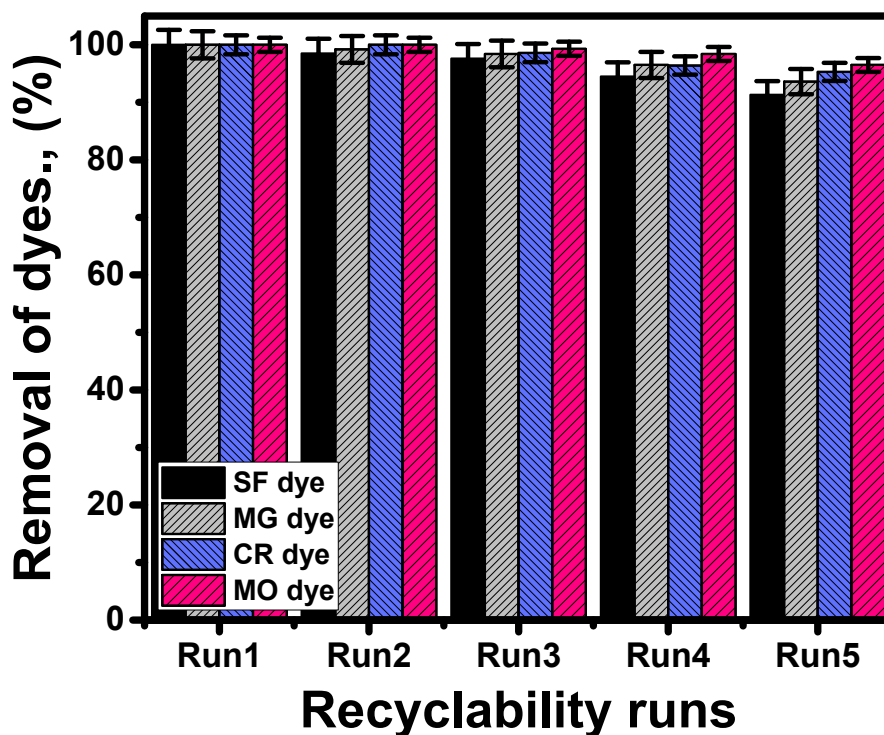


Figure S6. the removal percentages of SF, MG, CR, and MO dyes at the recyclability tests of KNTs

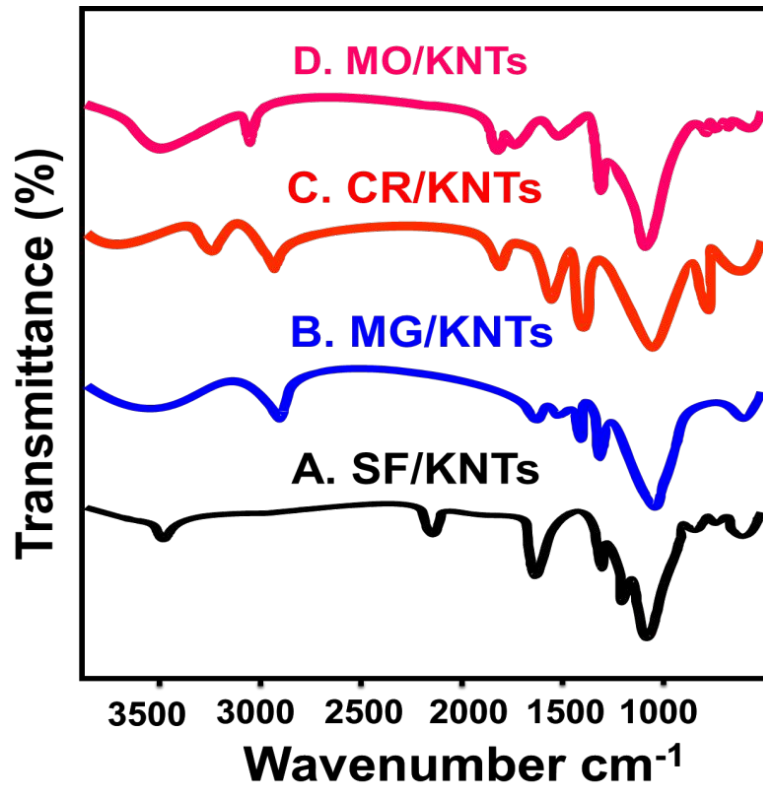


Figure S7. the FT-IR spectra of KNTs after the adsorption of SF (A), MG (B), CR (C), and MO dyes (D)

Table S1. the representative equations of the studied kinetic and isotherm model and their parameters

Kinetic models		
Model	Linear equation	Parameters
Pseudo-first-order	$\ln (q_e - q_t) = \ln q_e - k_1 t$ (Linear) $q_t = q_e (1 - e^{-k_1 t})$ (Nonlinear)	q_t (mg/g) is the adsorbed ions at time (t), and K_1 is the rate constant of the first-order adsorption (min^{-1})
Pseudo-second-order	$\frac{t}{q_t} = \frac{1}{K_2 q_e^2} + \frac{t}{q_e}$ (Linear) $q_t = \frac{q_e^2 k_2 t}{1 + q_e k_2 t}$ (Nonlinear)	q_e is the quantity of adsorbed ions after equilibration (mg/g), and K_2 is Lagergren model rate constant (g/mg min).
Elovich model	$q_t = \frac{1}{\beta} \ln (\alpha \beta) + \frac{1}{\beta} \ln (t)$ (Linear) $q_t = \frac{1}{\beta} \ln (1 + \alpha \beta t)$ (Nonlinear)	α is the initial adsorption rate (mg/min) and β is the surface saturation (g/mg)
Intra-particle diffusion	$q_t = K_p t^{1/2} + C$ (Linear) $q_t = kt^{0.5} + C$ (Nonlinear)	K_p ($\text{mg g}^{-1} \text{min}^{-0.5}$) is the intraparticle diffusion rate constant and C is the intercept of the line
Isotherm models		
Model	Equation	Parameters
Langmuir	$\frac{C_e}{q_e} = \frac{1}{b q_{\max}} + \frac{C_e}{q_{\max}}$ (Linear) $q_e = \frac{q_{\max} b C_e}{(1 + b C_e)}$ (Nonlinear)	C_e is the rest ions concentrations (mg/L), q_{\max} is the theoretical maximum adsorption capacity (mg/g), and b is the Langmuir constant (L/mg)
Freundlich	$\log q_e = (1/n) \log C_e + \log K_f$ (Linear) $q_e = K_f C_e^{1/n}$ (Nonlinear)	K_f is the constant of Freundlich model related to the adsorption capacity and n is the constant of Freundlich model related to the adsorption intensities
Dubinin–Radushkevich	$\ln (q_e) = \ln (q_m) - \beta \varepsilon^2$ (Linear) $q_e = q_m e^{-\beta \varepsilon^2}$ (Nonlinear)	β (mol^2/KJ^2) is the D-R constant, ε (KJ^2/mol^2) is the polanyi potential, and q_m is the adsorption capacity