Supplementary Material

Cadmium and Lead remediation using magnetic and non-magnetic sustainable biosorbents derived from *Bauhinia purpurea* pods

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Models name	Equation	Parameters	References
(a) Kinetic models			
Psuedo First order	$q_t = q_e (1 - e^{-k_1 t})$	$k_1(h^{-1})$ is the first order adsorption rate constant, q_e , is the Pb ²⁺ amount adsorbed at equilibrium and q_t is the Pb ²⁺ adsorbed at time "t".	1, 2
Psuedosecond order	$\frac{t}{q_{t}} = \frac{1}{k_{2}q_{e}^{2}} + \frac{t}{q_{e}}$	$k_2(g.mg^{-1} h^{-1})$ is the second order rate constant, q_e is the amount adsorbed at equilibrium, q_t is the amount adsorbed at time "t" and $k_2q_e^2$ represents the initial sorption rate.	3
Freundlich	$q_e = K_F C_e^{1/n}$	(b) Equilibrium models K_F -constant indicative of the relative adsorption capacity of adsorbent (mg/g), 1/n- a constant indicative of the intensity of the adsorption,q _e -adsorption capacity (mg/g), , C _e -equilibrium concentration of solute (mg/L),	4
Sips or Langmuir- Freundlich	$q_{e} = \frac{K_{LF} C_{e}^{n_{LF}}}{1 + (a_{LF} C_{e})^{n_{LF}}}$	K_{LF} , a_{LF} and n_{LF} are the sips constants. q_e -solute amount adsorbed per unit weight (mg/g), C_e is equilibrium concentration (mg/L),	5
Redlich Peterson	$q_e = \frac{K_{RP}C_e}{(1 + a_{RP}C_e^{\beta_{RP}})}$	K_{RP} , a_{RP} and β_{RP} are Redlich-Peterson constants and the exponent, β , lies between 0 and 1. q_e - solute amount adsorbed per unit weight (mg/g) and C_e is equilibrium concentration (mg/L).	6
Koble– Corrigan	$q_e = \frac{AC_e^{n_{KC}}}{1 + (bC_e)^{n_{KC}}}$	A, b and n_{KC} are the ships constants. q_e - solute amount adsorbed per unit weight (mg/g) and C_e is equilibrium concentration (mg/L).	7
Radke and Prausnitz	$\mathbf{q}_{\mathbf{e}} = \frac{\mathbf{a} \mathbf{b} \mathbf{C}_{\mathbf{e}}^{\beta}}{\mathbf{a} + \mathbf{b} \mathbf{C}_{\mathbf{e}}^{\beta-1}}$	a, b, and β are Radke and Prausnitz constants. q _e - solute amount adsorbed per unit weight (mg/g) and C _e is equilibrium concentration (mg/L)	8
Toth	$q_e = \frac{K_T C_e}{(1 + B_T C_e^{\beta_T})^{\frac{1}{\beta}} T}$	K_T , B_T , and β_T are Toth constants. q_e -adsorption capacity (mg/g), C_e -equilibrium concentration (mg/L)	9

Table SM1: Adsorption isotherm and kinetic models used in the study



Figure SM1. Schematic diagram for Pb²⁺/Cd²⁺ adsorption unto KPP/MKPP (A) batch mode and (B) continuous mode, and KPP recovery



Figure SM2. BET surface area plot for (A) KPP and (B) MKPP



Figure SM3. Pore size distribution plots for (A) KPP and (B) MKPP.



Figure SM4. FT-Raman spectra of (A) KPP and (B) MKPP before and after Pb²⁺ -and Cd²⁺ adsorption





Figure SM5. Effect of adsorbent dose on Pb^{2+} adsorption by (A) KPP and (B) MKPP [pH 4.5; initial Pb^{2+} and Cd^{2+} concentration = 10 mg/L; particle size = 30-50 B.S.S mesh and T = 25 °C]





Figure SM6. Effect of adsorbent dose on Cd^{2+} adsorption by (A) KPP and (B) MKPP [pH 5.0; initial Cd^{2+} concentration = 20 mg/L; particle size = 30-50 B.S.S mesh and T = 25 °C]



Figure SM7. Effect of adsorbate concentrations on Pb^{2+} adsorption by (A) KPP and (B) MKPP [pH 4.5; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S mesh and T = 25 °C]





Figure SM8. Effect of adsorbate concentrations on Cd^{2+} adsorption by (A) KPP and (B) MKPP [pH 5.0; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S mesh and T= 25 °C]



Figure SM9. Effect of temperature on Pb²⁺ by (A) KPP and (B) MKPP [pH 4.5; lead concentration= 5-100 mg/L; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S mesh]



Figure SM10. Effect of temperature on Cd^{2+} by (A) KPP and (B) MKPP [pH 5.0; cadmium concentration= 2-100-mg/L; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S mesh]



Figure SM11. Pseudo-second order kinetic plots for Pb^{2+} removal by (A) KPP and (B) MKPP at different doses [initial lead concentration = 10 mg/L; pH = 4.5, particle size = 30-50 B.S.S. mesh]



Figure SM12. Pseudo-second-order kinetic plots for Cd^{2+} removal by (A) KPP and (B) MKPP at different doses (initial cadmium concentration = 20 mg/L; pH = 5.0, particle size = 30-50 B.S.S. mesh)





Figure SM13. Pseudo-second-order kinetic plots for Pb^{2+} removal by (A) KPP and (B) MKPP at different concentration [pH = 4.5, adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP), particle size = 30-50 B.S.S. mesh]



Figure SM14. Pseudo-second-order kinetic plots for Cd^{2+} removal by (A) KPP and (B) MKPP at different concentrations [pH = 5.0,-adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size = 30-50 B.S.S. mesh]



Figure SM15. Freundlich adsorption isotherm of Pb^{2+} by (A) KPP and (B) MKPP at different temperatures [pH= 4.5; initial lead concentration range= 2-100 mg/L; T = 25 °C; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S. mesh]



Figure SM16.Temkin adsorption isotherm of Pb^{2+} by (A) KPP and (B) MKPP at different temperatures [pH= 4.5; initial lead concentration range= 2-100 mg/L; T = 25 °C; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S. mesh]



Figure SM17. Freundlich adsorption isotherm of Cd^{2+} by (A) KPP and (B) MKPP at different temperatures [pH= 5.0; initial cadmium concentration range= 2-100 mg/L; T = 25 °C; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S. mesh]



Figure SM18.Temkin adsorption isotherm of Cd^{2+} by (A) KPP and (B) MKPP at different temperatures [pH= 5.0; initial cadmium concentration range= 2-100 mg/L; T = 25 °C; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S. mesh]



Figure SM19. Adsorption isotherms of Pb²⁺ on (A) KPP and (B) MKPP in the absence and presence of Cd²⁺ and Cu²⁺. The sample with interfering Cd²⁺ and Cu²⁺ was in 1:1:1 molar ratio [pH= 4.5; adsorbent concentration= 4.0 g/L; particle size= 30-60 B.S.S. mesh].



Figure SM20. Cd²⁺ removal on [A] KPP and [B] MKPP in Cd²⁺-Cu²⁺-Pb²⁺system. [Initial pH= 5.0; adsorbent dose= 1 g/L (KPP) and 2 g/L (MKPP); particle size= 30-50 B.S.S mesh; Cd²⁺, Cu²⁺ and Pb²⁺molar ratio= -1:1:1]

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