

Synthesis of Thiamine-Modified Bentonite for Pretreatment of Pharmaceutical Wastewater

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Abstract

A series of sodium bentonite and thiamine were used as raw materials to prepare a new adsorbent – thiamine-modified bentonite. The adsorbent was applied in the pretreatment of pharmaceutical wastewater. The optimum preparation conditions of thiamine-modified bentonite were studied. The experiments showed that the removals of COD (chemical oxygen demand) reached 47.98 % under optimal condition. The FTIR results indicated that bentonite has been modified by thiamine. The X-ray diffraction result showed that the interplanar distance of modified bentonite was increased by intercalating thiamine. From the analyses of the SEM photographs, it can be clearly observed that the bentonite, after thiamine insertion reaction, loses its foliated structure and exhibits a rougher surface. Among several isotherm equations, Langmuir and Freundlich adsorption isotherms were investigated. They were widely used to describe equilibrium data for water and wastewater treatment applications. Adsorption isotherms correlate well with the Langmuir isotherm model. Detailed isothermal and kinetic studies show that the modified bentonite removes organic pollutants from pharmaceutical wastewater by physical adsorption processes.

Keywords

Bentonite, thiamine, pharmaceutical wastewater, pretreatment

Introduction

The pharmaceutical industry is one of the 12 industries in the State Environmental Protection Plan focusing on governance in China. The pharmaceutical industry accounts for 1.7 % of China's total industrial output, and the pharmaceutical wastewater discharge was about 2 % of total industrial discharges.¹ As the production processes and drug products are different and characteristic, the composition of pharmaceutical wastewater is complicated. According to the different production processes, pharmaceutical wastewater is divided into fermentation, chemical synthesis, mixing and formulating, biological engineering, extraction, and Chinese traditional medicine. Generally, pharmaceutical wastewater is characterized by complicated elements, intricate species of organic pollutants, high concentrations of suspended substances and ammonia, dark colour, high toxicity, and low ratio of BOD and COD.²

At present, many common conventional methods of treatment have been applied for pharmaceutical wastewater. The up-flow anaerobic stage reactor (UASR) was used for treatment of pharmaceutical wastewater containing macrolide antibiotics, and the COD reduction was 70–75 %.³ The anaerobic treatment was applied to pharmaceutical wastewater of chemical synthesis-based drug. With differ-

ent treatment processes, the COD removal efficiency was different.⁴⁻⁶ Biological treatment was also applied to improve the treatability of the pharmaceutical wastewater at the sources.

However, more than 50 pharmaceutical compounds have been detected over the last years in pharmaceutical wastewater samples.⁷⁻¹¹ Moreover, many of these pharmaceuticals are designed to be persistent and lipophilic and are not eliminated by wastewater treatment or biodegraded in the environment. The removal rate of pharmaceutical compounds increases significantly after pretreatment.^{12,13}

Physicochemical treatment is commonly used in the pre-processing stage and biochemical follow-up stage, and the common methods include coagulation, flotation, adsorption, electrolysis, advanced oxidation technology, and wet air oxidation. Coagulation and sedimentation method of physical treatment is the first choice in physical and chemical processing technology. This method can effectively remove the pollutants and improve the biodegradability of wastewater. For better dispersion, small particles and large specific surface area, bentonite is used for wastewater treatment.^{14,15} The modified bentonites with thiamine as modifier are used for pretreatment of fosfomycin sodium epoxy mother liquor wastewater. This kind of modified bentonite is a new and green adsorbent, which reduces the chemical sludge and salt content. The ion exchange process in modification is shown in Fig. 1.¹⁶

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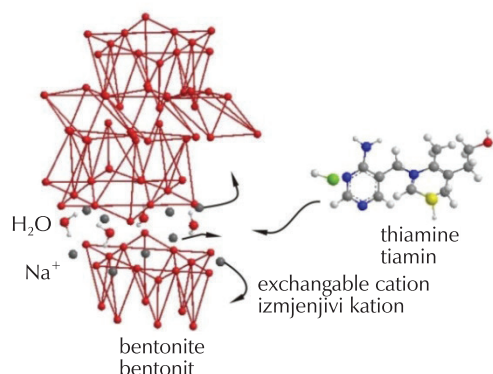


Fig. 1 – The modified reaction of bentonite and thiamine
Slika 1 – Reakcija modifikacije bentonita tiaminom

Experiments

Materials

Raw material: The sodium bentonite used in the experiments was purchased from Heishan, Liaoning Province, China. The basic properties are shown in Table 1.

The modifier: A certain quantity of thiamine (Guangfu Institute of Fine Chemical, China).

Table 1 – Properties of natural bentonite
Tablica 1 – Svojstva prirodnog bentonita

pH	8–9
Colloid value/ml (15 g) ⁻¹ Koloidnost/ml (15 g) ⁻¹	100
Cation exchange capacity/mmol 100 ⁻¹ g ⁻¹ Kationskoizmjenjivački kapacitet/mmol 100 ⁻¹ g ⁻¹	62
Swelling/ml g ⁻¹ Bubrenje/ml g ⁻¹	15
Methylene blue absorption/g g ⁻¹ Apsorpcija metilenskog modrila/g g ⁻¹	23–25

Fosfomycin sodium ethylene wastewater (FSE wastewater), a kind of pharmaceutical wastewater, was taken from a pharmaceutical factory in Shenyang, China, with characteristics of high organic concentration, a pungent odour, and pH of 9–10. The COD of FSE wastewater was up to $2.060\text{--}2.950 \cdot 10^5 \text{ mg l}^{-1}$.

Test method

The potassium chromate method was applied to determine COD.

Absorption test

The experiments were performed in a jar test apparatus consisting of six identical beakers of 250 ml, each equipped with a stirrer. The FSE wastewater was treated

under conditions of various bentonite dosages, modifier concentrations, stirring times, particle sizes, and pH values. The natural bentonite was added to the beakers with rapid agitation (400 rpm) for some time. Thereafter the particles were allowed to settle for 15 minutes, then the upper solution was taken for COD measurement by the potassium chromate method. The optimal preparing conditions of thiamine-modified bentonite were obtained through single-factor experiments.

Preparation of the modified bentonite

Thiamine 0.4 g was dissolved in 1 l hydrochloric acid solution at concentration of 5 % on a magnetic stirrer. Then, 2.4 g of the sodium bentonite ($\leq 74 \mu\text{m}$ size fraction) was added and the mixture was stirred continuously for 40 minutes. Later, the modified bentonite was separated from the mixture by centrifugation and washed several times using distilled water until a negative result for the presence of chloride was obtained with AgNO_3 . The modified bentonite was then dried at 30 °C for 24 h in a hot air oven, crushed into powder in agate mortars, and stored in moisture-free conditions for further use.

The natural bentonite was modified through the ion exchange reaction using thiamine. Thus, the functional group on the bentonite surface was added. The thiamine-modified bentonite was applied in the pretreatment of FSE wastewater.

Characterization

Fourier Transform Infrared (FTIR) spectrometer, X-ray diffraction and scanning electron microscopy (SEM) were used for characterization on natural bentonite and modified bentonite.

FTIR: The FTIR was used to identify the surface functional groups of two bentonites.

The FTIR spectra of thiamine bentonite samples were recorded in the transmission mode at room temperature by means of the 470 FT Thermo Nicolet instrument using the KBr pellet technique (1 : 20) with the resolution of 2 cm^{-1} . The KBr was dried at 200 °C for 24 h. Then 560 mg of KBr was homogenized with the bentonite sample in a ball grinder. The tablets (radius 1 cm and thickness 0.1 cm) were prepared using the hydraulic press.

X-ray diffraction: The interplanar distance of two bentonites was measured by X-ray diffraction.

The X-ray diffraction measurements were performed using a Brook D8-Advance instrument (Germany). The conditions of these measurements were as follows: $\text{Cu K}\alpha$ radiation, Ni filter for wavelength discrimination, pulse height analyser, scintillation detector, divergence slit 0.3° , receiving slit 0.15° , range of 2θ $3\text{--}30^\circ$, step size 0.02° , and count time per step 20 s.

SEM: The bentonite powders were measured by a JSM-6360LV SEM with High-Low vacuum (JEOL). The bentonites were enlarged by certain multiples. The three-dimensional surface conformation of natural bentonites and modified bentonites was observed.

Adsorption process of FSE wastewater

The 0.3 g of modified bentonite was mixed with 200 ml of wastewater sample at three different temperatures and pH = 3. Then the mixture was stirred for 50 minutes and left to set. The COD value was measured by the potassium dichromate method (Chinese national standards GB11941-89). The amount of organic pollutants adsorption was calculated by the following mass-balance equation:

$$q_e = (C_0 - C_e) \cdot V/m \quad (1)$$

q_e – equilibrium adsorption capacity, C_0 – initial COD value of FSE wastewater, C_e – equilibrium COD value of FSE wastewater, V – volume of FSE wastewater, m – mass of adsorbent.

The nonlinear and linear forms of Langmuir isotherm equation were given in the following equations:¹⁷

$$q_e = q_{\max} C_e K_L / (1 + C_e K_L) \quad (2)$$

$$1/q_e = 1/q_{\max} + 1/(K_L q_{\max} C_e) \quad (3)$$

q_{\max} – maximum adsorption capacity of the adsorbent; K_L – Langmuir equilibrium constant.

The nonlinear and linear forms of Freundlich isotherm equation were given in the following equations:¹⁸

$$q_e = K_F C_e^{1/n} \quad (4)$$

$$\log(q_e/\text{mg g}^{-1}) = \log K_F + 1/n \log(C_e/\text{mg l}^{-1}) \quad (5)$$

q_e – equilibrium adsorption capacity; C_e – equilibrium COD value of FSE wastewater; K_F – characteristic constant of Freundlich model related to the adsorption capacity; n – parameter characterizing the heterogeneity of the system.

The thermodynamic parameters of the adsorption, i.e. the standard Gibbs energy ΔG^\ominus , standard enthalpy ΔH^\ominus , and entropy ΔS^\ominus , were calculated using the Van't Hoff equation.

$$\Delta G^\ominus = -RT \ln K_{\text{ads}} \quad (6)$$

$$\ln K_{\text{ads}} = -\Delta H^\ominus/RT + \Delta S^\ominus/R \quad (7)$$

where, the slope and intercept of the plot of $\ln K_{\text{ads}}$ versus $1/T$ were used to determine the ΔH^\ominus and ΔS^\ominus . R is the gas constant and T is the absolute temperature.

Results and discussion

Optimal conditions of modified bentonite preparation

Effect of bentonite dosage on COD removal rate

The influence of different bentonite dosages on adsorption is shown in Fig. 2. The COD removal rate increased with an increase in the bentonite dosage up to 4 g l⁻¹. When the bentonite dosage was higher than 4 g l⁻¹, the COD removal rate dropped.

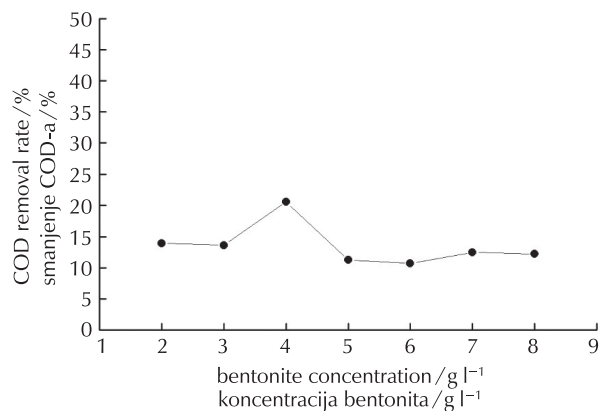


Fig. 2 – Effect of bentonite dosage on COD removal rate
Slika 2 – Učinak koncentracije bentonita na smanjenje COD-a

The bentonite was the matrix of modified bentonite. The greater the matrix, the more modified bentonite was prepared. But the bentonite could not be modified by thiamine if the bentonite surpassed the exchanged capability.

Effect of the modifier concentration on COD removal rate

The effect of the modifier concentration on COD removal rate is shown in Fig. 3. The COD removal rate increased gradually with a modifier concentration in the range of 0 – 5 g l⁻¹. The COD removal rate was close to a stable equilibrium state when the modifier concentration was in the range of 5 g l⁻¹ to 13 g l⁻¹. Then the COD removal rate decreased.

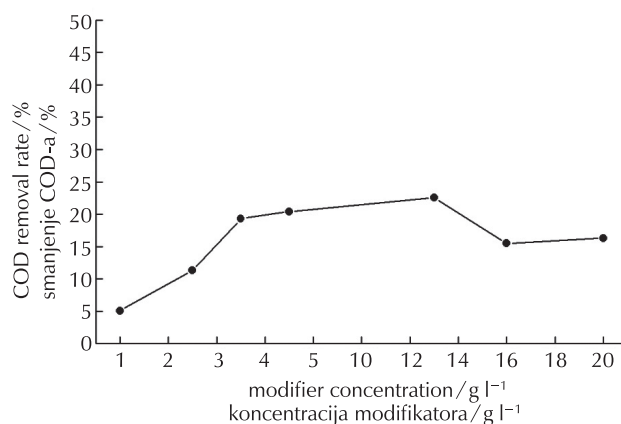


Fig. 3 – Effect of the modifier concentration on COD removal rate

Slika 3 – Učinak koncentracije modifikatora na smanjenje COD-a

The COD value of FSE wastewater was 2.060 – 2.950 · 10⁵ mg l⁻¹ which meant that a large number of the organic particles existed in the processing solution. The organic particles, bentonite and thiamine, were in a congested state. The positive charge of bentonite and negative charge of thiamine interacted as the function of irregular Brownian motion. Therefore, the bentonite was modified by thiamine on the surface or in the interlaminar

region. The organic particles were adsorbed by the modified bentonite (at modifier concentration of $0-5 \text{ g l}^{-1}$). When the flocculation particles were large enough, the collision force effect between particles was not obvious. Then Brownian motion tended to balance (at modifier concentration in the range of $5-13 \text{ g l}^{-1}$). When the modifier concentration increased continually, the excess thiamine existed with the form of organic matter in wastewater. The COD removal rate decreased (at modifier concentration higher than 13 g l^{-1}).

Effect of stirring time on COD removal rate

The dependence of COD removal rate on adsorption time is stated in Fig. 4. The dosage of the modified bentonite was 5 g l^{-1} . Firstly, the COD removal rate increased, then achieved a balance, and finally decreased when stirring time ranged from 20 to 140 minutes. The COD removal rate reached the maximum at stirring time of 40 minutes.

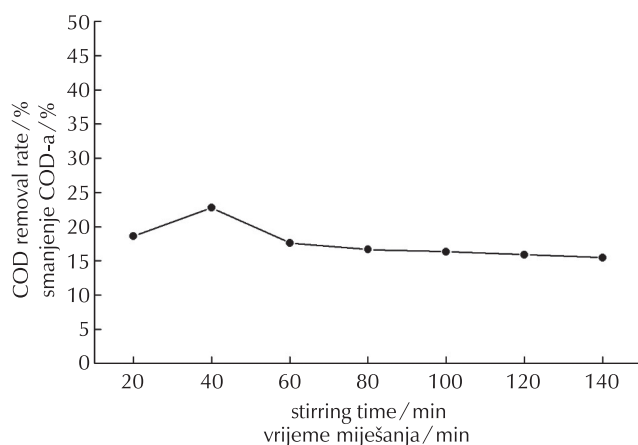


Fig. 4 – Effect of stirring time on COD removal rate
Slika 4 – Učinak vremena miješanja na smanjenje COD-a

Using mechanical rabble, the bentonite is modified by thiamine sufficiently. As displayed in Fig. 4, the removal capacity of the modified bentonite increased with the increase in stirring time and reached a maximum value at 40 minutes. With a further increase in stirring time, there was a steady decrease in the percentage of adsorption.

Effect of particle size on COD removal rate

The particle size in the range of $0.063-0.45 \text{ mm}$ presented a slight effect on COD removal (Fig. 5). The dosage of the modified bentonite was 5 g l^{-1} . A peak COD removal effect was observed at particle size of 0.076 mm .

The shape of modified bentonite in wastewater was expressed by particle size. The aggregates could not form when particle size was too small (smaller than 0.076 mm). The specific surface area decreased when particle size was too large. Therefore, the activity and adsorption ability decreased at particle size larger than 0.076 mm . The flocculation ability of modified bentonite at particle size of 0.076 mm was the highest.

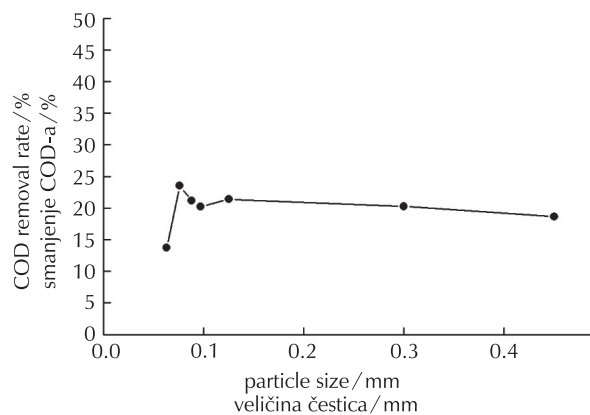


Fig. 5 – Effect of particle size on COD removal rate
Slika 5 – Učinak veličine čestica na smanjenje COD-a

The shape of modified bentonite in wastewater was expressed by particle size. The aggregates could not form when particle size was too small (smaller than 0.076 mm). The specific surface area decreased when particle size was too large. Therefore, the activity and adsorption ability decreased at particle size larger than 0.076 mm . The flocculation ability of modified bentonite at particle size of 0.076 mm was the highest.

Effect of the modifier pH on COD removal rate

The dosage of the modified bentonite was 5 g l^{-1} . From the dependency demonstrated in Fig. 6, it was clear that the modifier pH greatly influenced COD removal rate. At pH of 2–3, the COD removal rate showed an upward trend and the removal efficiency was very good. However, with the further increase in pH value, the COD removal rate decreased.

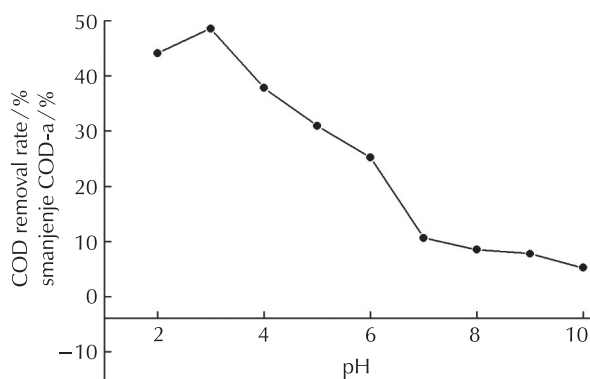


Fig. 6 – Effect of pH on COD removal rate
Slika 6 – Učinak pH na smanjenje COD-a

The adsorption capability of modified bentonite was impacted significantly by the physical and chemical properties of the thiamine. The modifier pH was decisive to physical and chemical properties of thiamine. Thiamine was stable at acidic pH. In addition, at low pH, the bentonite

surface was positively charged, which could be beneficial to the modification process due to the ion exchange. The molecular structure of thiamine was unstable in the alkaline solution, and it was susceptible to oxidation and decomposition. As illustrated in the figure, when pH was 2, the removal efficiency reached the maximum.

Through optimization of modified-bentonite preparation conditions, the COD removal rate of pharmaceutical wastewater reached 47.98 %. With a higher concentration of organic contaminant in pharmaceutical wastewater, a multi-step treatment process was needed. The adsorption process on thiamine-modified bentonite could be applied in the pretreatment process.

Characterization on sodium bentonite and modified bentonite

FTIR: Each functional group of bentonite was unique and corresponded to the fundamental vibrational bands of the infrared spectrum, respectively. The spectra of natural and modified bentonite are presented in Fig. 7.

The functional groups of two bentonites and their corresponding wave number are summarized in Table 2.

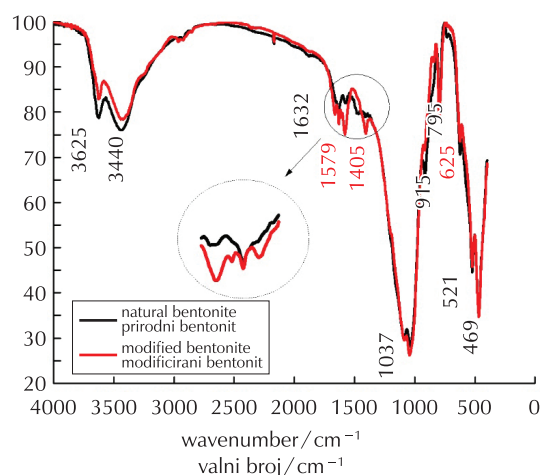


Fig. 7 – FTIR of two bentonites (A. natural bentonite; B. modified bentonite)

Slika 7 – FTIR dvaju bentonita (A – prirodni bentonit, B – modificirani bentonit)

The basic infrared spectra of natural bentonite still existed in modified bentonite. The interlayer structure of two sili-

Table 2 – Functional groups of two bentonites
Tablica 2 – Funkcijske skupine dvaju bentonita

Peak assignment	Natural bentonite Prirodni bentonit	Modified bentonite Modificirani bentonit
	Wavenumber / cm ⁻¹ Valni broj / cm ⁻¹	
Si-OH-Al stretch istezanje Si-OH-Al	3625	3627
O-H stretch istezanje O-H	3440	3424
O-H bend svijanje O-H	1632	1632
N-H (primary amines) N-H (primarni amini)	-	1579
C-N stretch istezanje C-N	-	1405
Si-O stretch istezanje Si-O	1037	1044
Al-Al-OH bend svijanje Al-Al-OH	915	917
Si-O bend svijanje Si-O	795	796
N-H plane deformation ravninska deformacija N-H	-	625
Al-O-Si bend svijanje Al-O-Si	521	520
Si-O-Si bend svijanje Si-O-Si	469	468

con-oxygen tetrahedral layers and one aluminum-oxygen octahedral layer remained unchanged. By comparing the FT-IR spectra of the natural and modified bentonite, the intensities increase apparently at 1579 cm^{-1} , 1405 cm^{-1} and 625 cm^{-1} . There were function groups of N-H, C-N and N-H. Nitrogen element was excluded from natural bentonite. Thus, the new functional groups were obtained from the ion exchange between natural bentonite and thiamine. The modified function of natural bentonite was effective.

X-ray diffraction: The X-ray diffraction was mainly used to study the interplanar distance of the modified bentonite. The X-ray diffractograms of the sodium bentonite and modified bentonite are shown in Fig. 8.

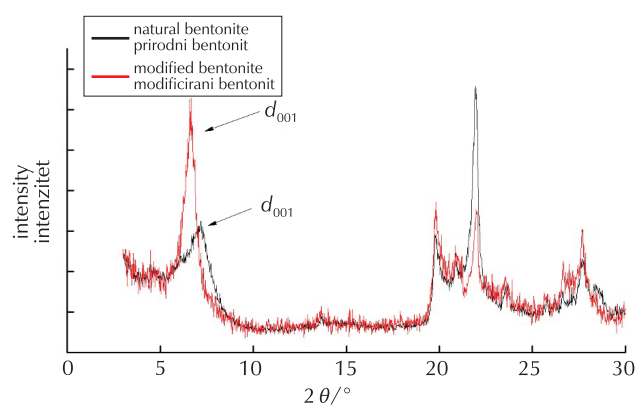


Fig. 8 – X-ray diffraction patterns of two bentonites
Slika 8 – Rendgenogrami dvaju bentonita

The small angle ($3^\circ - 10^\circ$) of XRD diffraction peaks could be used to study the structure of mesoporous materials. The angle 2θ was measured and the interplanar distance d was calculated according to X-ray diffraction spectrum and Bragg equation: $2d \sin\theta = n\lambda$. The interplanar distance of sodium bentonite was 12.2692 nm, and the interplanar distance of modified bentonite was 13.3407 nm.

The anion of thiamine has exchanged the cation on the surface of bentonite. The thiamine molecules entered the layer space to a certain extent. This indicated the presence of the thiamine molecule between the silicate layers.

SEM: The samples of sodium bentonite and modified bentonite were observed using a SEM, as shown in Fig. 9.

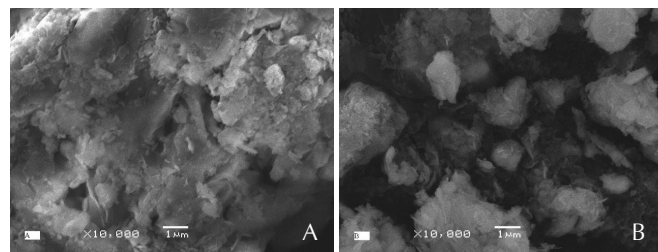


Fig. 9 – SEM images of two bentonites (A. sodium bentonite; B. modified bentonite)

Slika 9 – SEM-slike dvaju bentonita (A – natrijev bentonit, B – modificirani bentonit)

By analysing the SEM photographs, it was clearly observed that the bentonite lost its foliated structure and gained a rougher surface after the thiamine insertion reaction.

Adsorption isotherms

The interactions between adsorbate and adsorbent could be described by adsorption isotherms. The equilibrium data of FSE wastewater adsorption along with the fitting models are shown in Fig. 10.

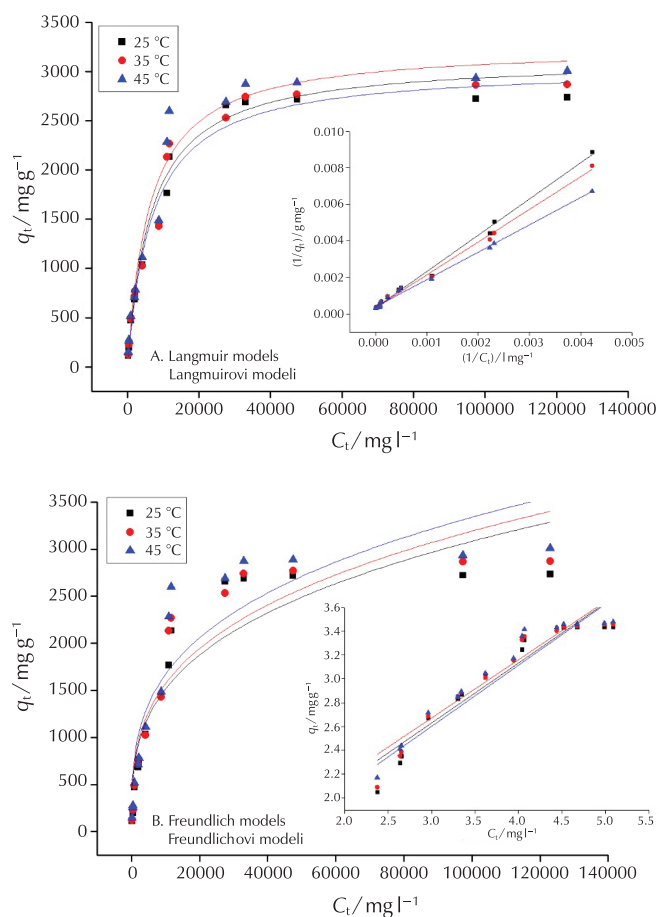


Fig. 10 – Adsorption isotherms of FSE wastewater adsorption onto modified bentonite and the fitting model (inset is the linear form of models)

Slika 10 – Adsorpcijske izoterme adsorpcije iz otpadne vode s FSE-om na modificirani bentonit i prilagođeni model (umetnut je linearni model)

The temperature-dependent parameters of Langmuir models obtained from the fitting procedure are summarized in Table 3. The data indicates that Langmuir models could represent the equilibrium data fairly well (R^2 of Langmuir $> R^2$ of Freundlich). The organic pollutant of FSE wastewater was adsorbed easily onto the surface and interlaminar of the bentonite, and monolayer adsorption was formed. Then the molecular movement was more and more regular.

Table 3 – Fitted parameters of temperature dependent form of two isotherm models
 Tablica 3 – Prilagođeni parametri dvaju izotermnih modela ovisnih o temperaturi

Langmuir model Langmuirov model				
temperature temperatura	equation jednadžba	q_0 mg g ⁻¹	K_L l mg ⁻¹	R^2
25 °C	$q_e = \frac{0.4711 C_e}{1 + 0.0001504 C_e}$	31.32	0.0001504	0.9795
35 °C	$q_e = \frac{0.5250 C_e}{1 + 0.0001609 C_e}$	32.62	0.0001609	0.9675
45 °C	$q_e = \frac{0.4424 C_e}{1 + 0.0001451 C_e}$	30.48	0.0001451	0.9810
Freundlich model Freundlichov model				
temperature temperatura	equation jednadžba	$1/n$	K_F mg g ⁻¹ /(l mg ⁻¹) ⁻ⁿ	R^2
25 °C	$q_e = 87.82 C_e^{0.3091}$	0.3091	87.82	0.8609
35 °C	$q_e = 91.85 C_e^{0.3082}$	0.3082	91.85	0.8644
45 °C	$q_e = 104 C_e^{0.3011}$	0.3011	104.27	0.8404

In order to further investigate the adsorption mechanism, the adsorption thermodynamic parameters of organic contaminants on the modified bentonite were calculated, and are listed in Table 4.

Table 4 – Adsorption thermodynamic parameters
 Tablica 4 – Termodinamički parametri adsorpcije

T/K	k_0/kg^{-1}	$\Delta G^\ominus/kJ mol^{-1}$	$\Delta H^\ominus/kJ mol^{-1}$	$\Delta S^\ominus/J K mol^{-1}$
298	0.0054	-12.9363		57.28658
308	0.0056	-13.2773	4.1351	56.53377
318	0.0060	-13.5259		55.53774

The adsorption reaction was an endothermic process as the ΔH^\ominus value was greater than zero. So the elevated temperature was beneficial to the adsorption reaction. When the enthalpy changed in the range of 0–63 kJ mol⁻¹, the adsorption reaction was a physical one. And the heat of adsorption caused by the molecular forces ranged from 4 to 10 kJ mol⁻¹, and the molecular force was a van der Waals force. Hence, it could be said that the adsorption of FSE wastewater on modified bentonite is a physical adsorption. Meanwhile, the ΔS^\ominus value decreased with the rise in temperature. This was in line with the rule of polymer entropy that decreased by adsorption reaction.

Conclusion

A new adsorption material was synthesized using sodium bentonite and thiamine, the optimal preparation conditions of modified bentonite were: bentonite dosage of 4 g l⁻¹, modifier concentration of 13 g l⁻¹, stirring for 40 minutes, particle size of 0.076 mm, and modifier pH 3. The preparation technology was non-toxic with no secondary pollution. The thiamine-modified bentonite was applied in the pretreatment process of pharmaceutical wastewater and the COD removal rate reached 47.98 %.

FTIR analysis results of modified bentonite showed that all characteristic peaks of natural bentonite had been retained, and the interlayer structures of two silicon-oxygen tetrahedral layers and one aluminum-oxide octahedral layer were constant. Three new characteristic peaks were observed at 1579 cm⁻¹, 1425 cm⁻¹, and 625 cm⁻¹. These peaks were all amino functional groups and originated from thiamine. Therefore, the natural bentonite was modified by ionization of thiamine. Analysis results of X-ray diffraction of modified bentonite showed that the interplanar distance increased at varying degrees, the thiamine had entered the interplanar space of bentonite, which improved the adsorption performance. The SEM characterization results showed that the modified bentonite clumps were smaller with higher dispersion and clear structure hole, the adsorption performance increased.

Adsorption thermodynamic processes of organic pollutants in FSE wastewater on modified bentonite were employed to correlate Langmuir models. The adsorption reaction was a physical adsorption process. Temperature increase was conducive to the reaction.

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List of symbols and abbreviations

Popis simbola i kratica

C_0	– initial COD, mg l^{-1} – početni COD, mg l^{-1}
C_e	– equilibrium COD, mg l^{-1} – ravnotežni COD, mg l^{-1}
d	– spacing between the planes in the lattice – razmak između susjednih ravnina u rešetki kristala
ΔG^\ominus	– standard Gibbs energy, kJ mol^{-1} – standardna Gibbsova energija, kJ mol^{-1}
ΔH^\ominus	– standard enthalpy, kJ mol^{-1} – standardna entalpija, kJ mol^{-1}
k_0	– adsorption constant, kg^{-1} – adsorpcijska konstanta, kg^{-1}
K_{ads}	– adsorption equilibrium constant – adsorpcijska ravnotežna konstanta
K_F	– Freundlich constant, $(\text{mg g}^{-1})/(\text{l mg}^{-1})^{-n}$ – Freundlichova konstanta, $(\text{mg g}^{-1})/(\text{l mg}^{-1})^{-n}$
K_L	– Langmuir equilibrium constant, l mg^{-1} – Langmuirova konstanta ravnoteže, l mg^{-1}
m	– mass of adsorbent, g – masa adsorbenta, g
n	– heterogeneity parameter, reflection order – parametar heterogenosti, red refleksije
q_e	– equilibrium adsorption capacity, mg g^{-1} – ravnotežni adsorpcijski kapacitet, mg g^{-1}
q_{max}	– maximum adsorption capacity, mg g^{-1} – maksimalni adsorpcijski kapacitet, mg g^{-1}
R	– gas constant, $8.314 \text{ J K}^{-1} \text{ mol}^{-1}$ – plinska konstanta, $8,314 \text{ J K}^{-1} \text{ mol}^{-1}$
ΔS^\ominus	– standard entropy, J K mol^{-1} – standardna entropija, J K mol^{-1}
T	– temperature, K – temperatura, K
V	– volume, l – obujam, l
θ	– Bragg angle, $^\circ$ – Braggov kut, $^\circ$
λ	– wavelength, nm – valna duljina, nm
BOD	– biochemical oxygen demand, mg l^{-1} – biokemijska potreba za kisikom, mg l^{-1}
COD	– chemical oxygen demand, mg l^{-1} – kemijska potreba za kisikom, mg l^{-1}
UASR	– up-flow anaerobic stage reactor – anaerobni stupnjeviti reaktor s uzlaznim tokom
FSE	– fosfomicin sodium ethylene

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SAŽETAK

Modificiranje bentonita tiaminom za predobradu otpadne vode iz farmaceutske industrije

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Bentonit je modificiran tiaminom i primijenjen u predobradi vode iz farmaceutske industrije. Pročavani su optimalni uvjeti pripreme bentonita. Eksperimenti su pokazali da se pod optimalnim uvjetima kemijska potreba za kisikom (COD) smanjuje za 47,98 %. FTIR ukazuje na modifikaciju bentonita tiaminom, a rendgenska difrakcija na povećanje udaljenosti među slojevima bentonita interkaliranjem tiamina. Bentonit nakon interkalacije tiamina gubi lisnatu strukturu i pokazuje grublju površinu. Od nekoliko izotermnih jednadžbi, istraživane su Langmuirove i Freundlichove adsorpcijske izoterme. One se široko primjenjuju za opisivanje ravnotežnih podataka u obradi i pročišćavanju voda. Adsorpcijske izoterme dobro koreliraju s Langmuirovim izotermnim modelom. Detaljne izotermne i kinetičke studije pokazuju da modificirani bentonit fizičkim adsorpcijskim procesima uklanja organske onečišćujuće tvari iz farmaceutske otpadne vode.

Ključne riječi

Bentonit, tiamin, farmaceutske otpadne vode, predobrada

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