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## Adsorptive Removal of Formaldehyde by Chemically Bamboo Activated Carbon with addition of Ag nanoparticle: Equilibrium and Kinetic

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**Abstract.** Carbon was prepared from dried waste bamboo (*Dendrocalamus asper*) using chemical activation with KOH. The carbon was prepared with the activating agent in a mass ratio of KOH and dried bamboo (3:1) at 800°C. Using impregnation technique, the bamboo-based activated carbon has developed with modified Ag nanoparticle (Ag-AC) to capture formaldehyde. The Ag-AC has characteristics of moderate surface area of 685 m<sup>2</sup>/g and average pore size of 2.7 nm. The adsorption equilibriums and kinetics of formaldehyde on Ag-AC measured. The influences of initial formaldehyde on adsorption performance have measured in a batch system. The equilibrium data were evaluated by isotherm models of Langmuir, Freundlich, and  $10^{\text{th}}$  order. The Langmuir model well describes the adsorptive removal of formaldehyde on Ag-AC in this study. Pseudo-first-order and pseudo-second-order kinetic equations were applied to test the experimental data. The pseudo-second-order exhibited the best fit for kinetic study.

### 1 Introduction

The quality of indoor air has become areas of concern since it related to human health, directly. Maintaining adequate indoor air quality through controlling ambient concentrations of formaldehyde is recommended, although the relationships between air quality and formaldehyde rarely determined. Formaldehyde with a high level can cause symptoms such as eye, nose, and throat irritation, chest pains, bronchitis, and prolonged formaldehyde has associated with reduced pulmonary function and asthma [1,2]. General methods are being used to purify the polluted air, including adsorption [3], membrane, absorption, plasma decomposition [4], and photocatalytic oxidation [5] techniques. Considering adsorption has a promising method for controlling low concentration of formaldehyde because of its simple technique and low energy requirement. Activated Carbon (AC) of bamboo (*Dendrocalamus asper*) is extremely porous material with a moderate surface area and the most widely used as adsorbent. The AC is used for capturing indoor air pollutants at ambient temperature is considered economically feasible. However, AC is a non-polar adsorbent that sometimes exhibits low selectivity with polar contaminants. In recent years, modified AC with metal nanoparticle was considered a potential adsorption and oxidation catalytic process due to increased capacity to formaldehyde. Chemically bamboo AC was produced by with the activating agent in a mass ratio of KOH and dried bamboo (3:1) at 800°C. Silver nanoparticles are impregnated on AC to enhance the

adsorption capacity of removal formaldehyde. An interesting fact was that the Ag nanoparticles dispersed on commercial coal-based AC to effectively remove formaldehyde from air [6]. Silver also possesses high applicability for use in controlling indoor air due to its non-toxic feature [7].

In this study, AC was modified with silver nanoparticles (Ag-AC) to improve their adsorption capacities. It is important to determine the Ag-AC by submitting it for the adsorption of formaldehyde and evaluating the equilibrium isotherms and kinetics adsorption of formaldehyde.

### 2 Materials and Methods

#### 2.1 Materials

The AC had conducted as Rengga's method [7]. The Bamboo Chars and KOH were mixed with 1:3 of weight ratio. The immersed bamboo chars in KOH solution was dehydrated until it formed the slurry. Subsequently, the slurry was activated in a tube-electronic heater at heating rate of 10°C/min in the presence of nitrogen gas with flowing rate of 200 mL/min. The activation process was continued up to temperature of 800°C and maintained for 1 h. The activated product was washed with HCl followed by rinsing with distilled water until the pH of the washing effluent was 6. The activated carbon was dried at 105 °C and grinded into 200 mesh prior to characterization and adsorption study.

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## 2.2 Synthesis Adsorbent

Silver nanoparticle synthesized was conducted with reduction process similar to Gaedhi's method [8] in the previous paper [7]. As much 4 mL of 0.1 M AgNO<sub>3</sub> added to 1 L of an aqueous solution with 0.15 wt% of the soluble starch. The mixture vigorously stirred for 1 h under atmosphere. The pH of resulting solution adjusted to 8.0 by adding NaOH solution. The mixture was maintained at 50 °C until a change of the reaction solution became yellow. The next step, silver nanoparticles in solution was mixed with 107.9 g of activated carbon under magnetic stirring, resulting in the deposition of about 4 wt% of silver nanoparticles on AC. The AC-Ag was filtered then washed with distilled water and dried at 105 °C until its constant weight.

## 2.3 Batch equilibrium studies

The adsorption tests were performed to determine the addition amounts of Ag-AC (0.5 g) to capacity adsorption of formaldehyde at various gas-phase concentrations. The test was done into 10 mL glass vial equipped with a septum, and subsequently the vials was shaken at 25 °C for 24 h. The inlet and outlet concentrations of formaldehyde in the batch tests were collected using a 0.5 mL gas-tight syringe. Furthermore, the sample was analyzed immediately by using gas chromatography equipped with a flame ionization detector (GC/FID, Shimadzu GC-2014). In the experiments, procedures of kinetic tests were identical to equilibrium tests of adsorption. The samples were drawn at suitable time intervals. Each experiment continued until equilibrium conditions were reached when no further decrease in the formaldehyde concentration was measured. Initially on the Ag-AC did not contain any formaldehyde. Furthermore, formaldehyde adsorbed was calculated by using Eq. 1.

$$q_e = \frac{V(c_0 - c_t)}{m}$$

where  $q_e$  is formaldehyde adsorbed at equilibrium [mg/g];  $V$  is a volume of adsorbate [L];  $m$  is the mass of modified adsorbent [g]. The value of  $c_0$  is formaldehyde concentration initially, and  $c_t$  is formaldehyde concentration after adsorption at equilibrium time [ppm], respectively. The best-fit model determined for a system of formaldehyde adsorption, experimental data substituted for equilibrium isotherm model equations and kinetic models. Parameter values of equations were obtained using non-linear regression analysis. The correlation coefficient ( $R^2$ ) of the fit these values, were also reported for each model.

## 3 Results and Discussion

The characteristics of pore are shown in Table 1. We calculated the mesoporosity of AC by BET and BJH

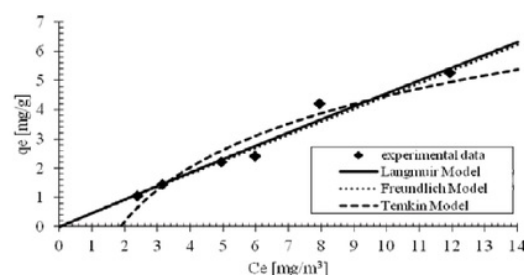
methods with surface area and average size of pore of 685 m<sup>2</sup>/g and 2.7 nm, respectively.

**Table 1.** The Characteristics of Pores

$S_{BET}$ [m <sup>2</sup> /g]	$D_{BET}$ [nm]	$V_{meso\ BJH}$ [cm <sup>3</sup> /g]	$V_{micro\ BJH}$ [cm <sup>3</sup> /g]	$V_{BJH}$ [cm <sup>3</sup> /g]
685	2.7	0.2009	0.0336	0.2355

## 3.1 Isotherm Equilibrium

Formaldehyde adsorption equilibrium, formaldehyde removal increases with an increase in the initial formaldehyde concentration. Formaldehyde at equilibrium concentration was indicated the maximum of adsorption capacity is shown in Fig. 1. The adsorption process reaches an equilibrium state, occurs when the molecules distribute between gas and solid phase. The data of formaldehyde adsorption equilibrium were analyzed using Langmuir, Freundlich, and Temkin isotherms, based on the equilibrium of the sorption between the adsorbate and adsorbent. The Langmuir model describes a monolayer sorption with a homogenous distribution both energies and sites of sorption. The Freundlich model describes the adsorption distribution energy for heterogeneous sites. The Temkin isotherm assumption is the adsorption heat of all the molecules in AC decreases linearly with coverage. In order, removal formaldehyde from indoor for optimization adsorption system is important to be done to establish the most appropriate correlation in equilibrium curve in Fig 1.



**Figure 1.** Representation of formaldehyde adsorption on g-AC for Langmuir, Freundlich, and Temkin Models

The Langmuir, Freundlich, and Temkin isotherm equations with their equation are presented in Table 2. In the Langmuir model,  $Q_m$  is the maximum monolayer coverage adsorption capacity [mg/g], and  $K_L$  is the Langmuir isotherm constant [L<sup>3</sup>/g]. In the Freundlich model, the values of  $K_F$  and  $n$  are Freundlich isotherm constant and adsorption intensity. The value of  $1/n$  is the heterogeneity of the adsorbent. The value of  $1/n$  is higher than 1 for favorable adsorption process. In Temkin model,  $B = RT/b$ ,  $R$  is the universal gas constant [8.314 J/K mol], and  $T$  is the absolute temperature [K]. The

value of  $B_T$  is constant related to the heat of adsorption [ $\text{m}^3/\text{g}$ ], and  $A_T$  is the equilibrium binding constant.

The isotherm constants of adsorption models and correlation coefficients are reported in Table 2. The adsorption of formaldehyde on Ag-AC for Langmuir isotherm is shown range studies with high correlation coefficient ( $R^2 > 0.96$ ). Therefore, the conclusion of isotherm that Langmuir model is the best isotherm to predict the adsorption of formaldehyde on Ag-AC adsorbent. It reported that Ag-AC had a relatively suitable adsorption capacity of 157.223 mg/g in this work. The Langmuir model predictions for concentration formaldehyde of 50  $\text{mg}/\text{m}^3$  using commercial coal-based carbon is 15.8 mg/g [5] that is close to this study (21 mg/g). Confirmation of the experimental data to the Langmuir isotherm equation proves that the surface of the Ag-AC for the adsorption of formaldehyde consists of homogeneous sites. The results are in agreement with previous researchers to the adsorption of pollutants using AC prepared from bone char [9].

**Table 2.** Parameter of Isotherm Constant and Correlation Coefficient for Adsorption of Formaldehyde on Ag-AC

Model	Non-linear equation	Parameter	Value
Langmuir	$q_e = \frac{Q_m \cdot K_L \cdot C_e}{1 + K_L \cdot C_e}$	$Q_m$ [mg/g]	157.2
		$K_L$ ( $\text{m}^3/\text{mg}$ )	23
		$R^2$	0.966
		$1/n$	1.005
Freundlich	$q_e = K_F \cdot C_e^{1/n}$	$K_F$ [ $\text{mg}/\text{g}$ ] $[\text{m}^3/\text{mg}]$	0.466
		$R^2$	0.962
Temkin	$q_e \left( \frac{RT}{b_T} \right) \cdot \ln(A_T \cdot C_e)$	$\left( \frac{RT}{b_T} \right)$ [J/mol]	2.668
		$A_T$ [ $\text{m}^3/\text{mg}$ ]	0.536
		$R^2$	0.925

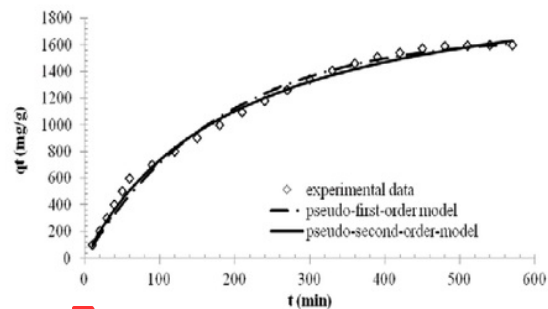
### 3.2 Adsorption kinetic

The kinetics of adsorption of formaldehyde on Ag-AC were investigated by using pseudo-first-order and pseudo-second-order models. The pseudo-first-order and pseudo second-order kinetic models of  $k_1$ ,  $k_2$ , and  $R^2$  are presented in Table 3. The value of  $k_1$  is the adsorption rate constant (1/h) and  $k_2$  is the rate constant of second-order adsorption (g/mg.h). The values of  $q_e$  and  $q_t$  are the amounts of formaldehyde adsorbed at equilibrium and time [mg/g], respectively. It can see that the values of  $R^2$  obtained did not trend consistently at low formaldehyde initial concentration for the pseudo-first-order model. The pseudo-second-order model predicts the behavior of the concentration range adsorption. Between the experimental and the calculated  $q_e$  values (Fig. 2) have a good agreement. Besides that the experimental  $q_e$  value agree with the estimated  $q_e$  values obtained from non-linear. Obtained that the correlation coefficient values  $R^2 = 0.995$  for the second-order kinetic model, it is clear that the adsorption follows this model. The value of  $R^2$  was close to unity, indicating that Ag-AC fitted well to this model for the adsorption of formaldehyde. Between these

models, the criterion for their applicability is based on judgment on the respective  $R^2$  with the high values indicates that the pseudo-second-order model. Agreed the results of the work performed by previous researchers for adsorption of formaldehyde pollutants using activated carbon prepared from bone char [9]. Wu et al. [10] reported that the pseudo second-order model were suitable for adsorption of adsorbate with low molecular weight on smaller adsorbent particles as described for this work.

**Table 3.** Kinetic Parameter for The Adsorption of Formaldehyde on Ag-AC

Kinetic Model	Non-linear equation	Parameter	Value
Pseudo-first-order	$q_t = q_e [1 - \exp(-k_1 \cdot t)]$	$k_1$ [1/h] $R^2$	0.333 0.991
pseudo-second order	$q_t = \frac{q_e^2 \cdot k_2 \cdot t}{1 + q_e \cdot k_2 \cdot t}$	$k_2$ [g/mg.h] $R^2$	0.00015 0.995



**Figure 2.** Pseudo-first-order and pseudo-second-order kinetics plot for the adsorption of formaldehyde on Ag-AC

### 4 Conclusions

The present investigation showed that low concentration formaldehyde by chemically activated carbon of bamboo modified with Ag nanoparticles is an effective adsorption. In the equilibrium condition, adsorption using modified activated carbon is appropriate with Langmuir isotherm model, as well as Freundlich and Temkin models. Isothermal equilibrium model is suitable Langmuir monolayer adsorption model, with adsorption capacity of 157.223 mg/g. For kinetics adsorption of formaldehyde using modified activated carbon, the uses of pseudo-second-order model is more appropriate rather than pseudo-first-order-model.

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