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Bundit Tengjaroenkul

Urai Tengjaroenkul

Natapol Pumipuntu3

Komkrich Pimpukdee

Sawitree Wongtangtintan

*See next page for additional authors*

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## **An In Vitro Comparative Study of Aflatoxin B1 Adsorption by Thai Clay and Commercial Toxin Binders**

### **Authors**

Bundit Tengjaroenkul, Urai Tengjaroenkul, Natapol Pumipuntu<sup>3</sup>, Komkrich Pimpukdee, Sawitree Wongtangtintan, and Piyawat Saipan

# An *In Vitro* Comparative Study of Aflatoxin B1 Adsorption by Thai Clay and Commercial Toxin Binders

Bundit Tengjaroenkul<sup>1\*</sup> Urai Tengjaroenkul<sup>2</sup> Natapol Pumipuntu<sup>3</sup> Komkrich Pimpukdee<sup>1</sup>  
Sawitree Wongtangtintan<sup>4</sup> Piyawat Saipan<sup>1</sup>

## Abstract

Twenty samples of Thai clay from fourteen provinces and seven commercial toxin binders were investigated for their adsorption capacity of aflatoxin B1 (AFB<sub>1</sub>) *in vitro*. Each sample of 5 mg/l AFB<sub>1</sub> solution was shaken at 25°C for 24 hours and supernatants of centrifuged samples were analyzed for concentrations of AFB<sub>1</sub> using a UV spectrophotometer. Adsorption capacity was calculated and applied to isothermic equations. The results indicated that Thai clays was capable of sequestering AFB<sub>1</sub> from aqueous solution differently and Thai clay from Lopburi and Lamphun provinces had the highest adsorption capacity, similar to commercial binders. S-shaped isothermic curves were observed for all samples having adsorption capacity greater than  $4 \times 10^{-3}$  mol/kg. These isothermic data were fitted using a modified Freundlich model that suggested that the samples possessed multilayered or multiple adsorption sites for the toxin. The clay from Lopburi and Lamphun had maximum adsorption capacity ( $Q_{\max}$ ) of  $4.76 \times 10^{-3}$  mol/kg and  $4.68 \times 10^{-3}$  mol/kg, respectively, whereas the commercial binders had  $Q_{\max}$  that ranged from  $4.38 \times 10^{-3}$  mol/kg to  $5.07 \times 10^{-3}$  mol/kg. Inductive couple plasma spectrometry and X-ray diffraction spectrometry of the clay samples demonstrated that the clay from Lopburi and Lamphun contained montmorillonite as a major component, similar to the bentonites. It was concluded that Thai clay from Lopburi and Lamphun provinces could absorb AFB<sub>1</sub> *in vitro* efficiently, similar to commercial toxin binders.

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**Keywords:** adsorption, isotherm, montmorillonite, mycotoxin, Thai clay

<sup>1</sup>Faculty of Veterinary Medicine, Khon Kaen University, Khon Kaen 40002, Thailand

<sup>2</sup>Faculty of Science, Chiang Mai University, Chiang Mai 50200, Thailand

<sup>3</sup>Faculty of Veterinary Sciences, Mahasarakham University, Mahasarakham 44000, Thailand

<sup>4</sup>Faculty of Agriculture, Khon Kaen University, Khon Kaen 40002, Thailand

\*Corresponding author: E-mail: btengjar@kku.ac.th

## บทคัดย่อ

### การศึกษาเปรียบเทียบประสิทธิภาพดินเหนียวในประเทศไทยและสารดูดซับในเชิงพาณิชย์ในการดูดซับสารพิษอะฟลาทอกซินปี 1 ในหลอดทดลอง

บัณฑิตย์ เต็งเจริญกุล<sup>1\*</sup> อุไร เต็งเจริญกุล<sup>2</sup> ณัฐพล ภูมิพันธ์<sup>3</sup> คมกริช พิมพ์ภักดี<sup>4</sup> สาวิตรี วงศ์ตั้งถิ่นฐาน<sup>4</sup> ปิยะวัฒน์ สายพันธ์<sup>4</sup>

การศึกษานี้ได้นำตัวอย่างดินเหนียวจาก 20 แหล่งใน 14 จังหวัดของประเทศไทยและสารดูดซับสารพิษอะฟลาทอกซินเชิงพาณิชย์ 7 ตัวอย่าง มาศึกษาเปรียบเทียบความสามารถในการดูดซับสารพิษอะฟลาทอกซินปี 1 ในหลอดทดลอง โดยการนำตัวอย่างมาแยกเขย่าในหลอดแก้วที่มีสารละลายสารพิษความเข้มข้น 5 มก./ลิตร ที่อุณหภูมิ 25 องศาเซลเซียส นาน 24 ชั่วโมง หลังจากนั้นทำการปั่นเหวี่ยง และนำสารละลายส่วนบนมาตรวจวัดปริมาณอะฟลาทอกซินด้วยเครื่องยูวีสเปกโตรโฟโตมิเตอร์ ความยาวคลื่น 362 นาโนเมตร ผลการศึกษาแสดงให้เห็นว่าดินเหนียวของไทยจากแต่ละแหล่งสามารถดูดซับสารพิษได้แตกต่างกัน โดยดินเหนียวจากจังหวัดลพบุรีและลำพูนมีประสิทธิภาพในการดูดซับสารพิษสูงกว่าดินจากแหล่งอื่น แต่ไม่แตกต่างจากสารดูดซับในเชิงพาณิชย์ และเมื่อนำข้อมูลการดูดซับสารพิษของสารตัวอย่างที่มีค่าความสามารถในการดูดซับสารพิษมากกว่า  $4 \times 10^{-3}$  โมล/กก. มาศึกษาไอโซเทอร์มพบว่า สร้างกราฟได้เป็นรูปตัวเอส ที่ประยุกต์เข้ากับสมการโมดิฟายฟรุนดลิชได้ดี ซึ่งแสดงว่าสารตัวอย่างสามารถดูดซับสารพิษได้แบบหลายตำแหน่งและหลายชั้นในโครงสร้าง นอกจากนี้ยังพบว่าดินเหนียวจากจังหวัดลพบุรีและลำพูนมีความสามารถสูงสุดในการดูดซับสารพิษที่  $4.76 \times 10^{-3}$  และ  $4.68 \times 10^{-3}$  โมล/กก. ตามลำดับ ส่วนสารดูดซับในเชิงพาณิชย์สามารถดูดซับสารพิษได้ระหว่าง  $4.38 \times 10^{-3}$  ถึง  $5.07 \times 10^{-3}$  โมล/กก. และจากการศึกษาของคัพประกอบและโครงสร้างของตัวอย่างสารดูดซับด้วยอินดิคทีฟแคลิฟลูอิดสโคป สเปกโตรเมทรีและเอ็กซ์เรย์ ดิฟเฟรคชั่น สเปกโตรเมทรีพบว่า ตัวอย่างดินจากจังหวัดลพบุรีและลำพูนมีแรมอนต์มอริลโลไนท์เป็นองค์ประกอบหลัก เช่นเดียวกับสารดูดซับเบนโทไนท์ในเชิงพาณิชย์ จากการศึกษาสรุปได้ว่าดินเหนียวจากจังหวัดลพบุรีและลำพูนมีประสิทธิภาพสูงในการดูดซับสารพิษอะฟลาทอกซินในหลอดทดลอง และไม่แตกต่างจากสารดูดซับในเชิงพาณิชย์

**คำสำคัญ:** ดินเหนียว สารดูดซับ มอนต์มอริลโลไนท์ สารพิษ ไอโซเทอร์ม

<sup>1</sup> คณะสัตวแพทยศาสตร์ มหาวิทยาลัยขอนแก่น จังหวัดขอนแก่น 40002

<sup>2</sup> คณะวิทยาศาสตร์ มหาวิทยาลัยเชียงใหม่ จังหวัดเชียงใหม่ 40002

<sup>3</sup> คณะสัตวแพทยศาสตร์ มหาวิทยาลัยมหาสารคาม จังหวัดมหาสารคาม 44000

<sup>4</sup> คณะเกษตรศาสตร์ มหาวิทยาลัยขอนแก่น จังหวัดขอนแก่น 40002

\*ผู้รับผิดชอบบทความ E-mail: btengjar@kku.ac.th

## Introduction

Aflatoxins (AF) are toxic metabolites produced by *Aspergillus flavus* and *A. parasiticus*. Aflatoxin B<sub>1</sub> (AFB<sub>1</sub>) is widely known as carcinogenic and the most hepatotoxic of natural occurring AF (Hueber et al., 2004; Godfrey et al., 2013). One strategy to detoxify AF is to add toxin binder to animal feed reducing the bioavailability of the toxin absorbed through digestive tract (Basalan et al., 2006). This approach is considered as cost-effective and practical, particularly when using in contaminated feed on an industrial scale (Hueber et al., 2004; Pimpukdee et al., 2004; Kossolova et al., 2009). To date, several *in vitro* studies have shown that binders such as aluminosilicate, bentonite, zeolite, activated charcoal, and chitin-chitosan effectively absorb AFB<sub>1</sub> *in vivo* (Khajareen et al., 2003; Pasha et al., 2007; Manafi, 2011; Khadem et al., 2012; Rao and Chopra, 2012; Sadeghi et al., 2012; Neeff et al., 2013).

Previous studies have applied methods of equilibrium isothermal analysis to characterize the adsorption of AF onto the surfaces of toxin binders

(Grant and Phillips, 1998; Pimpukdee et al., 2000). These methods provide evidence of the molecular mechanisms involved with different binders and allow comparisons of similarities and differences. Binders have distinctive molecular structures and bind AFB<sub>1</sub> differently (Phillips, 1999). A comparison of Thai clay with commercial binders for their adsorption capacity of AFB<sub>1</sub> has not been reported. Thus, the main objective of this study was to investigate the adsorption capacity and affinities of Thai clay and compare them to commercial toxin binders using a Langmuir and modified Freundlich isotherm modeling approach.

## Materials and Methods

**Chemicals:** Standard AFB<sub>1</sub> was purchased from Sigma Chemical Co. (St. Louis, USA) and seven commercial binders, three with bentonites (BN) and four without bentonites (NB), were obtained from Thai suppliers. Twenty samples of Thai clay (S) were collected from fourteen provinces as follows: Lopburi (S1), Lamphun (S2), Nakorn Ratchasima (S3-S4), Buriram (S5-S6), Phetchabun (S7), Phitsanulok (S8-S9); Lampang (S10-S11), Nakorn Srithammarat (S12), Kanchanaburi (S13),

Ratchburi (S14-S15), Suphanburi (S16), Phetburi (S17-S18), Prachuab Khirikhan (S19) and Chonburi (S20)(Table 1). All clay was sieved to achieve particle sizes less than 60  $\mu\text{m}$ . Highly purified water (18  $\text{M}\Omega\cdot\text{cm}$ ) was prepared by processing deionized water through a Milli-Q<sup>uf+</sup> system.

**Adsorption Capacity:** After mixing each clay and binder sample into 5 mg/l AFB<sub>1</sub> solutions at 0.25% (w/v) for binder to solution ratio, the mixtures were shaken at 200 x g for 24 hours at 25°C (Innova 4060 Shaker, New Brunswick Scientific, USA). After shaking, the samples were centrifuged (Beckman Coulter, USA) at 12,000 x g for 30 min, and the supernatants were analyzed to determine the concentrations of AFB<sub>1</sub> using a UV spectrophotometer (Perkin Elmer, USA) at a wavelength of 362 nm (Pimpukdee et al., 2000).

**Isothermal Adsorption:** Aflatoxin B1 solutions having the concentrations from 0.5 to 8.0  $\mu\text{g}/\text{ml}$  were prepared. Isothermic studies utilized 4.0 ml solution of AFB<sub>1</sub> at each concentration, diluted in purified water containing 10.0 mg of each sample. The equilibrating condition and the separation procedure for samples and control tubes were described as above. Supernatants were analyzed further using UV spectrophotometer at wavelength 362 nm (Grant and Phillips, 1998; Pimpukdee et al., 2000). Adsorption capacity was calculated based on the amount of AFB<sub>1</sub> left in the solution ( $C_e$ ), and the amount of AFB<sub>1</sub> adsorbed ( $q$ ) for each sample. Data were later transferred to a table curve 2D 3V (Jandel Scientific, USA) to fit a Langmuir isothermic model (LIM) or a modified Freundlich isothermic adsorption model (MFM) (Mayura et al., 1998; Pimpukdee et al., 2000).

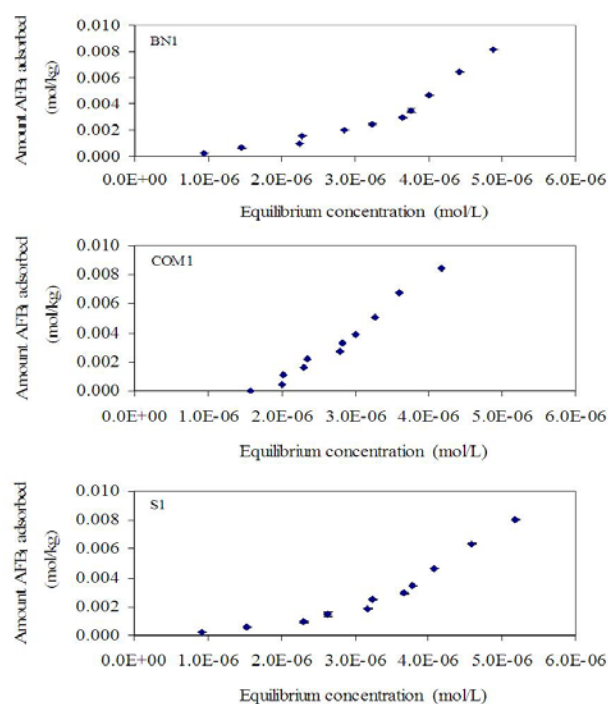
**Analyses of composition and structure:** Inductive couple plasma optical emission spectrometry (ICP-OES; Optima 3000, Perkin-Elmer, USA) was used to analyze the composition of the Thai clays and commercial binders having a high absorptive capacity for AFB<sub>1</sub>. The samples were further analyzed for their molecular structures using X-ray diffraction spectrometry (XRD; X'Pert MRD, PANalytical, B.V., The Netherlands).

## Results

Twenty samples of Thai clay and seven commercial binders were evaluated for their ability to adsorb AFB<sub>1</sub>. Table 1 shows the mean binding capacity of each commercial binder, which ranged from  $4.62 \times 10^{-4}$  to  $5.07 \times 10^{-3}$  mol/kg. The NB1 binder provided the highest adsorption capacity of AFB<sub>1</sub> ( $5.07 \times 10^{-3} \pm 3.67 \times 10^{-5}$  mol/kg). The BN1-3 had moderate adsorption capacity;  $4.74 \times 10^{-3} \pm 1.50 \times 10^{-5}$ ,  $4.65 \times 10^{-3} \pm 2.78 \times 10^{-5}$  and  $4.38 \times 10^{-3} \pm 1.51 \times 10^{-5}$  mol/kg, respectively. For the Thai clay, the highest mean adsorption capacity was found in Lopburi clay (S1) ( $4.76 \times 10^{-3} \pm 8.20 \times 10^{-5}$  mol/kg), followed by the clay from Lamphun (S2) ( $4.68 \times 10^{-3} \pm 7.41 \times 10^{-5}$  mol/kg)(Table 1). The clay from Nakorn Srithammarat (S12) had the highest distribution constant ( $6.70 \times 10^7$ ) (Table 1).

From the isothermic study, S-shaped isotherms were obtained for samples with the adsorption capacity greater than  $4 \times 10^{-3}$  mol/kg. The amount of adsorbed toxin increased linearly as the concentration of the AFB<sub>1</sub> increased. Figure 1 shows an example of isothermic plots of three representative groups: bentonite 1 (BN1), non-bentonite binder 1 (NB1), and Thai clay from Lopburi (S1). The equation for LIM ( $q = Q_{\text{max}} [K_d C_e / (1 + K_d C_e)]$ ) and the MFM ( $q = Q_{\text{max}} [K_d C_e]^n$ ), where  $q$  is the amount of AFB<sub>1</sub> adsorbed,  $Q_{\text{max}}$  is the maximum amount of AFB<sub>1</sub> adsorbed,  $C_e$  is the equilibrium concentration of AFB<sub>1</sub> in solution, and  $K_d$  is distribution constant (affinity parameter) were selected to fit the data. The LIM is generally applied to the monolayer adsorption, whereas the MFM is appropriate for the multilayer or multiple-site adsorption. With user defined functions, the isothermic data indicated that all adsorbents having the adsorption capacity greater than  $4 \times 10^{-3}$  mol/kg fitted to the MFM better than the LIM. The MFM had higher correlation coefficients ( $r^2$ ) for the adsorbents in the range of 0.80178-0.98963 and allowed a quantitative comparison of  $Q_{\text{max}}$  and  $K_d$  for adsorbents as expressed in Table 1.

The results of the composition and structure analyses demonstrated that BN1 and the Thai clay from Lopburi and Lamphun contained montmorillonite as a major component, having a deviated peak from the beam at angle  $2\theta$  approximately 6.4 (Fig 2). It was also noted that Thai clays from these two provinces contained greater percentages of montmorillonite than the other Thai clay.



**Figure 1** The S-shaped isothermic plots for AFB<sub>1</sub> adsorption to bentonite (BN1), non-bentonite clay (NB1) and Thai clay (S1), respectively. The standard deviations on each data point of adsorbed AFB<sub>1</sub> were small.

**Table 1** Comparison of isotherm fitted parameters of AFB<sub>1</sub> adsorption onto BN, NB and S samples at 25°C.

| Adsorbent | Langmuir Model(LM) <sup>a</sup> |                  |                | Modified Freundlich Model (MFM) <sup>b</sup> |                  |                |
|-----------|---------------------------------|------------------|----------------|--|------------------|----------------|
|           | r <sup>2</sup>                  | Q <sub>max</sub> | K <sub>d</sub> | r <sup>2</sup>                               | Q <sub>max</sub> | K <sub>d</sub> |
| BN1       | 0.692                           | 0.993            | 1.12E+03       | 0.97903                                      | 4.74E-03         | 2.52E+05       |
| BN2       | 0.602                           | 0.873            | 1.24E+03       | 0.93621                                      | 4.65E-03         | 2.84E+05       |
| BN3       | 0.848                           | 0.49             | 1.44E+03       | 0.88581                                      | 4.38E-03         | 3.60E+05       |
| NB1       | 0.589                           | 0.817            | 1.64E+03       | 0.96648                                      | 5.07E-03         | 4.62E+05       |
| NB2       | 0.581                           | 0.705            | 1.71E+03       | 0.93575                                      | 4.85E-03         | 4.26E+05       |
| NB3       | 0.614                           | 0.72             | 1.10E+03       | 0.98803                                      | 4.40E-03         | 3.04E+05       |
| NB4       | 0.699                           | 0.673            | 1.45E+03       | 0.8976                                       | 4.62E-04         | 4.71E+05       |
| S1        | 0.69                            | 1.026            | 1.03E+03       | 0.98905                                      | 4.76E-03         | 2.37E+05       |
| S2        | 0.64                            | 1.051            | 9.42E+02       | 0.98864                                      | 4.68E-03         | 2.31E+05       |
| S3        | 0.656                           | 0.892            | 8.73E+03       | 0.90105                                      | 3.54E-03         | 3.30E+05       |
| S4        | 0.963                           | 0.172            | 3.35E+03       | 0.96626                                      | 1.96E-03         | 2.76E+05       |
| S5        | 0.879                           | 0.55             | 8.56E+02       | 0.97249                                      | 1.66E-03         | 2.15E+05       |
| S6        | 0.899                           | 0.731            | 1.24E+03       | 0.97004                                      | 3.18E-03         | 2.54E+05       |
| S7        | 0.917                           | 0.5              | 1.53E+03       | 0.96719                                      | 1.87E-03         | 3.32E+05       |
| S8        | 0.672                           | 1.023            | 9.82E+02       | 0.98963                                      | 2.95E-03         | 2.72E+05       |
| S9        | 0.555                           | 1.092            | 1.53E+03       | 0.8645                                       | 3.12E-03         | 4.41E+05       |
| S10       | 0.746                           | 0.569            | 1.96E+03       | 0.84637                                      | 1.34E-03         | 4.79E+05       |
| S11       | 0.78                            | 0.65             | 8.94E+02       | 0.97383                                      | 9.61E-04         | 2.98E+05       |
| S12       | 0.844                           | 0.666            | 2.06E+03       | 0.95424                                      | 2.15E-03         | 6.70E+07       |
| S13       | 0.943                           | 0.148            | 2.11E+03       | 0.94789                                      | 7.89E-04         | 3.54E+05       |
| S14       | 0.667                           | 0.521            | 6.99E+02       | 0.80178                                      | 7.10E-04         | 2.38E+05       |
| S15       | 0.749                           | 0.62             | 7.77E+02       | 0.95252                                      | 8.91E-04         | 3.04E+05       |
| S16       | 0.743                           | 0.591            | 2.45E+03       | 0.88537                                      | 2.64E-03         | 4.21E+05       |
| S17       | 0.687                           | 0.544            | 7.12E+02       | 0.83265                                      | 7.5780-04        | 2.55E+05       |
| S18       | 0.816                           | 0.657            | 6.42E+02       | 0.92454                                      | 8.62E-04         | 4.76E+05       |
| S19       | 0.938                           | 0.436            | 1.35E+03       | 0.98369                                      | 1.64E-03         | 2.90E+05       |
| S20       | 0.91                            | 0.419            | 1.87E+03       | 0.9642                                       | 4.55E-04         | 3.19E+05       |

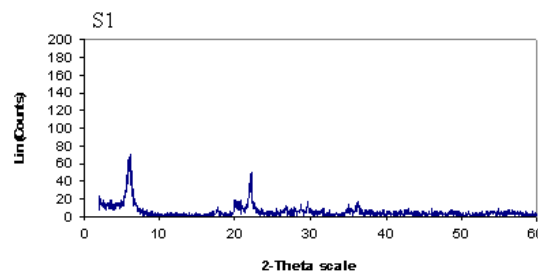
<sup>a</sup>q = Q<sub>max</sub>[K<sub>d</sub>C<sub>e</sub> / (1 + K<sub>d</sub>C<sub>e</sub>)], <sup>b</sup>q = Q<sub>max</sub>[K<sub>d</sub>C<sub>e</sub>]<sup>n</sup> where q: adsorbed amount (mol/kg), K<sub>d</sub>: distribution coefficient, C<sub>e</sub>: equilibrium concentration (mol/l), n: heterogeneity factor, BN: bentonite, NB: non-bentonite toxin binder, S: Thai clay

## Discussion

This isothermic adsorption study of AFB<sub>1</sub> Thai clay from different provinces and commercial binders was performed at 25°C to assess potential use in the agricultural industry. S-shaped isotherms were observed for samples with adsorption capacity greater than 4 × 10<sup>-3</sup> mol/kg. (Fig 1). Generally, the S isotherm is observed when a molecule does not have a strong affinity for the surface, until there is a significant amount absorbed, and the slope increases as the affinity for the surface increases (Grant and Phillips, 1998; Hinz, 2001). This occurs because the solute molecule has modified the surface or has begun to bind to the previously adsorbed molecules. The S-type curve usually appears when three conditions are fulfilled: the solute molecule (1) is mono functional, (2) has moderate intermolecular attraction, causing it to pack vertically on the absorbing layer, and (3) meets strong competition for substrate sites from molecules of the solvent or of another absorbing species.

All samples having the adsorption capacity greater than 4 × 10<sup>-3</sup> mol/kg fitted to the MFM better than the LIM. This implies that the binders possess multiple-sites or have multilayer adsorption (Hinz, 2001; Pimpukdee et al., 2004). The fitting result for

MFM indicated that the clay from Lopburi and Lamphun had relatively high average maximum adsorption capacity and were not different from other commercial toxin binders. Therefore, it was concluded that the clay from Lopburi and Lamphun was capable of absorbing AFB<sub>1</sub> efficiently. Adsorption of AFB<sub>1</sub> can occur on external surfaces, interlayer surfaces, original



**Figure 2** The X-ray diffraction spectrogram of Thai clay from Lopburi (S1) for the adsorption of AFB<sub>1</sub>.

edge sites, interlayer exchangeable cations, or on previously adsorbed molecules (Hinz, 2001; Diaz et al., 2002; Hueber et al., 2004). Previous studies have shown that montmorillonite has a high Q<sub>max</sub> for toxin, because of the extremely large and highly selectivity specific surface area, high cation exchange capacity (CEC), high cohesion and adhesion, and high equilibrium capacity (Pimpukdee et al., 2000, 2004). In

contrast, some other clay (pyrophyllite, kaolinite, illite, chlorite, talc, mica) has relatively low  $Q_{max}$  due to low CEC and relatively small internal surface areas (Pimpukdee et al., 2000).

In addition, the results of ICP-OES and XRD analyses demonstrated that the Thai clay collected from Lopburi and Lamphun contained montmorillonite as a major component and its binding capacity was similar to bentonite binders. Therefore, the composition and structure of the Thai clay from these two provinces were comparable in efficiency to commercial bentonites for binding the toxin.

In conclusion, Thai clay from Lopburi and Lamphun can absorb AFB<sub>1</sub> *in vitro* similar to other commercial toxin binders.

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