

Numerical modelling of a hemp concrete wall

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Abstract. In a global warming context associated to the abuse of energy consumption, actual researches focus more and more on reducing energy costs in the building sector. This target could be achieved by using innovative building materials, such as hemp concrete, due to its positive impacts on thermal and environmental levels. The aim of this work is to carry out a numerical study of a hemp concrete wall subjected to several indoor and outdoor conditions of temperature and relative humidity using the program object oriented SPARK. The hygrothermal behaviour of the wall is investigated taking into account heat and moisture transfer within the wall as well as hysteresis phenomenon between the sorption and desorption curves and their temperature dependency.

1 Introduction

Actually, as mentioned in the French Thermal Regulations of 2012 [1] and in the United Nations Environmental Program of 2016 UNEP [2], buildings are the most energy consuming sector. In France, they are responsible of almost 43 % of the total energy consumption. In a context of promoting energy efficiency, building sector has undergone a notable evolution towards innovative construction insulation materials [3] [4] such as the hemp concrete.

Hemp concrete is obtained by mixing hemp shives to a mineral binder (lime), water and some additives. Plentiful works are carried out on hemp concrete at the material and wall scales and its experimental characterization is widely available [5][6][7]. Recent works have shown also the importance of including hysteresis effect [8] [9][10] and temperature dependency [9][11][12] in modelling its hygrothermal behaviour to assess relative humidity profiles within the walls.

In this context, the aim of this work is to investigate numerically the hygrothermal behaviour of a hemp concrete wall subjected to real weather conditions [8] and to assess the impact of different modelling approaches on the prediction of temperature and humidity conditions within the wall. For this purpose, four models are compared: a standard model for heat and moisture transfer, a model with hysteresis effect, a model with sorption temperature dependency and finally a global model taking into account both aspects. The oriented object simulation tool SPARK [13] is used to solve the different models using the finite difference method.

2 Numerical study

2.1 Review on moisture transfer in a single building material

Coupled heat and moisture transfer models for a simple layer building materials are studied in several researches: Künzel [14] developed a model in which moisture transfer is due to relative humidity and temperature gradients. In 1997, Mendes and al. [15] showed a model based on the theory of Philip and De Vries [16], with similar assumptions as Künzel, but this time moisture transfer is governed by volumetric moisture content gradient. Thus, Umidus model [15] is created : both diffusion and capillary regimes are taken into account and the transfer of water in the vapor and liquid phases through the material can be analysed for any climate. Moreover, in 1989, Kerestecioglu and Gu [17] investigated the phenomenon using the evaporation-condensation theory in unsaturated liquid flow stage. In addition, Burch and Thomas [18] developed a computational model using finite-difference method able to estimate heat and mass transfers through composite walls under non-isothermal conditions. Moreover, the model of Ozaki and al. [19] takes into account the moisture transfers in both liquid and vapor phases coupled to the heat generated by the phase change due to the moisture sorption process.

2.2 Mathematical model

In this study, Umidus model, where moisture is transported under liquid and vapor phases, is used. The mass conservation equation is given by:

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$$\frac{\partial \theta}{\partial t} = \frac{\partial}{\partial x} \left(D_T \frac{\partial T}{\partial x} \right) + \frac{\partial}{\partial x} \left(D_\theta \frac{\partial \theta}{\partial x} \right) \quad (1)$$

With the following boundary conditions ($x = 0$ and $x = L$):

$$-\rho_l \left(D_T \frac{\partial T}{\partial x} + D_\theta \frac{\partial \theta}{\partial x} \right) \Big|_{x=0,e} = h_{M,e} (\rho_{v,a,e} - \rho_{v,s,e}) \quad (2)$$

$$-\rho_l \left(D_T \frac{\partial T}{\partial x} + D_\theta \frac{\partial \theta}{\partial x} \right) \Big|_{x=L,i} = h_{M,i} (\rho_{v,s,i} - \rho_{v,a,i}) \quad (3)$$

“e” and “i” represent respectively the outside and inside, “s” refers to the wall surface, and “a” to the ambient air. $D_T = D_{Tl} + D_{Tv}$ and $D_\theta = D_{\theta l} + D_{\theta v}$, where D_{Tl} is the liquid phase transport coefficient associated to a temperature gradient, D_{Tv} , the vapor phase transport coefficient associated to a temperature gradient, $D_{\theta l}$, the liquid phase transport coefficient associated to a moisture content gradient, $D_{\theta v}$, the vapor phase transport coefficient associated to a moisture content gradient, D_T , the mass transport coefficient associated to a temperature gradient and D_θ , the mass transport coefficient associated to a moisture gradient [20].

$$D_\theta = \frac{\delta_a P_{vs}}{\mu \rho_0 \xi} \quad (4)$$

Vapor transport coefficient under a temperature gradient is given by the relation:

$$D_{T,v} = \phi \frac{\delta_a dP_{vs}}{\rho_l \mu dT} \quad (5)$$

$D_{\theta v}$ is expressed as follows:

$$D_{\theta v} = \frac{D_{Tv} \times P_{sat}}{\xi \left(\phi \times \frac{dP_{vs}}{dT} + P_{sat} \times \frac{1}{\xi} \times \xi_T \right)} \quad (6)$$

One dimensional model of the energy conservation equation with coupled temperature and moisture for a porous media is considered, and the effect of the absorption or desorption heat is added. This equation is written as:

$$\begin{aligned} \rho_0 C_{pm} \frac{\partial T}{\partial t} = & \frac{\partial}{\partial x} \left(\lambda(T, \theta) \frac{\partial T}{\partial x} \right) \\ & + L_v \rho_l \left(\frac{\partial}{\partial x} \left(D_{T,v} \frac{\partial T}{\partial x} \right) \right. \\ & \left. + \frac{\partial}{\partial x} \left(D_{\theta,v} \frac{\partial \theta}{\partial x} \right) \right) \end{aligned} \quad (7)$$

$$C_{pm} = C_{p0} + C_{pl} \frac{\rho_l}{\rho_\theta}$$

where C_{pm} is the average specific capacity, which takes into account the dry material specific heat and the contribution of the specific heat of the liquid phase. λ is the thermal conductivity depending on moisture content and the temperature in Kelvin.

Boundary conditions take into account radiation, heat and phase change.

$$\begin{aligned} \left(-\lambda \frac{\partial T}{\partial x} \right)_{x=0,e} - L_v \rho_l \left(D_{T,v} \frac{\partial T}{\partial x} \right. \\ \left. + D_{\theta,v} \frac{\partial \theta}{\partial x} \right)_{x=0,e} \\ = h_{T,e} (T_{a,e} - T_{s,e}) \\ + L_v h_{M,e} (\rho_{ve,a,e} - \rho_{ve,s,e}) + \Phi_{ray,e} \end{aligned} \quad (9)$$

2.3 Hysteresis modelling

Hysteresis is observed between the sorption curves of hygroscopic building materials. Physically, sorption curves describe the equilibrium between the water content of the material and the relative humidity of the surrounding air. The hysteresis phenomenon is due to the “ink bottle” effect [21] which can be explained by the fact that water isn’t retained in the materials’ pores at the same manner it fulfils them. Literature provides many models describing the hysteresis phenomenon in porous materials: in fact, three global models can be distinguished: physical models developed by Mualem [22], mathematical models of Cool and Parker [23], and empirical models proposed by Pedersen [24]. New researches done by Steeman [25] and Van Belleghem [26] showed that hysteresis influences the moisture transfer in porous materials. Moreover and particularly in hemp concrete, Samri [27], and Lelievre et al. [28] have proved the importance of considering the hysteresis phenomenon in global transfer models (heat, air and moisture).

In this article, the empirical model proposed by Pedersen [24] to describe the hysteresis that occurs in hygroscopic materials like wood is used and adapted for the hemp concrete case. Main adsorption and desorption curves are given in equations (14) and (15):

$$\frac{U(\varphi)}{U_{max}} = \left(1 - \frac{\ln(\varphi)}{A_a} \right)^{\frac{-1}{n_a}} \quad (10)$$

$$\frac{U(\varphi)}{U_{max}} = \left(1 - \frac{\ln(\varphi)}{A_d} \right)^{\frac{-1}{n_d}} \quad (11)$$

where U is the actual water content (kg.kg^{-1}) for a fixed temperature, U_{max} the maximum water content, φ the relative humidity of the air and a and d represent adsorption and desorption phenomenon respectively. The coefficients A_a , A_d , n_a and n_d are calculated using the least squares method based on the experimental results of sorption curves of Lelievre et al. [8].

Water content U (kg.kg^{-1}) could be transformed into volumetric water content θ via relation (16):

$$\theta = U \times \frac{\rho_0}{\rho_t} \quad (12)$$

Pedersen's approach for hysteresis is based on the weighted values of the moisture capacity or on the slope of adsorption and desorption curves described by:

$$\xi_a = \frac{\partial U_a}{\partial \varphi} \quad (13)$$

$$\xi_d = \frac{\partial U_d}{\partial \varphi} \quad (14)$$

After a series of alternating processes of adsorption and desorption, intermediate scanning curves are modelled by [30]:

$$\xi_{hyst,a} = \frac{\gamma_a(U - U_a)^2 \xi_d + (U - U_d)^2 \xi_a}{(U_d - U_a)^2} \quad (15)$$

$$\xi_{hyst,d} = \frac{(U - U_a)^2 \xi_d + \gamma_d(U - U_d)^2 \xi_a}{(U_d - U_a)^2} \quad (16)$$

γ_a and γ_d depend on the type of the material studied. For the case of hemp concrete, they can be fitted based on the experiments of the moisture buffering value of the hemp concrete done by Lelievre et al. according to the NordTest protocol [8] as shown in Fig.1.

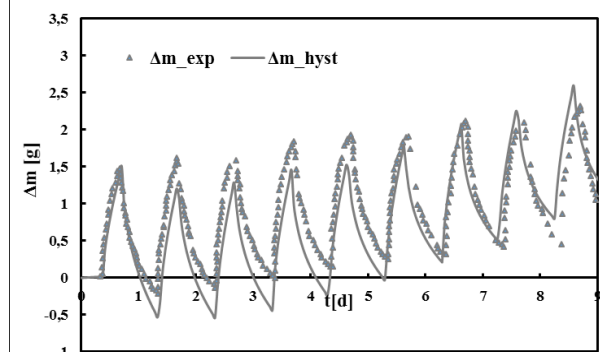


Fig. 1. Mass variations of hemp concrete specimen during the NordTest experience.

Fig.1 shows a convergence of the specimen mass during the repeating cycles (adsorption – desorption). γ_a and γ_d are set to 0.85 and 0.97 respectively in the numerical model including hysteresis.

2.4 Sorption curves temperature dependency

Temperature dependency of the sorption curves of hemp concrete is considered by allowing several sorption curves to describe the sorption characteristics at different temperature levels. So far, the study of the temperature effect on the sorption characteristics for bio-based materials is a new task. Researches have shown that the sorption capacity of materials depends on the temperature: increasing temperature entails that the

isosteric moisture content can be reached in equilibrium with a higher relative humidity [12]. Three main reasons can explain this phenomenon [30]: firstly, the microstructure alteration due to the temperature (for example, the enlargements of pores can be attributed to a rise of temperature). Secondly, the modification of water thermophysical properties with temperature and finally, the thermodynamic evolution of sorption mechanism (the exothermic process of adsorption). In the case of hemp concrete, Oumeziane [9] proved the necessity of considering the temperature effect in modelling its hygrothermal behavior. Moreover, Rode and Clorius [31] worked on the coupling between temperature and hysteresis effects and proved that an increase of temperature results in a reduction of the hysteresis loop. Poyet and Charles [32] proposed a model describing the relation between sorption characteristics at different temperatures and based on the differential heat of sorption which can be written as follows:

$$\begin{aligned} \varphi_2(T_2, \theta) &= \varphi_1(T_1, \theta) \frac{P_{sat}(T_1)}{P_{sat}(T_2)} e^{q_{st}(\theta) \frac{M_1(T_2 - T_1)}{R T_1 T_2}} \end{aligned} \quad (17)$$

where M_1 is water molar mass ($\text{kg} \cdot \text{mol}^{-1}$), R the ideal gas constant ($\text{J} \cdot \text{mol}^{-1} \cdot \text{K}^{-1}$), and q_{st} the isosteric heat ($\text{J} \cdot \text{kg}^{-1}$) calculated from two sorption isotherms at two different temperatures (T_1 and T_2) as shown in Fig.2.

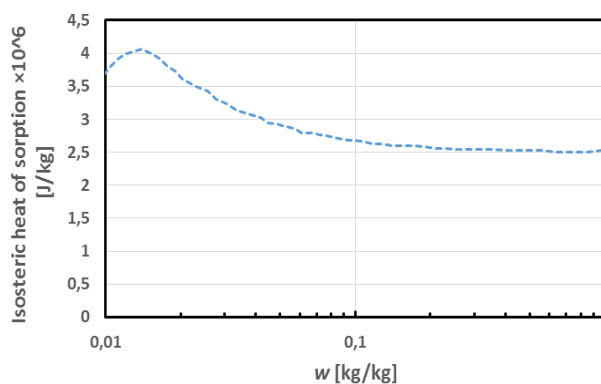


Fig. 2. Isosteric heat of sorption (Poyet and Charles Model).

2.5 Simulation environment SPARK

In order to solve the previous equation system, the Simulation Problem Analysis and Research Kernel (SPARK) tool is used [33]. Numerical solution is based on finite difference iterative method. SPARK allows solving efficiently differential equation systems. Each component is represented by an object that contains its appropriate mathematical model. Equations are defined in a generic matter called classes. Model is completed by linking objects together.

Using relation (21), main sorption curves are computed and implemented in SPARK respecting the form of equations (14) and (15), where n is a constant calibrated on the average sorption curve between main adsorption and desorption curves, U_{max} (Or θ_{max}) a

linear function of the temperature, and $A_a = A_d = A$ with A a linear function of the temperature as well. Figure 3 shows a 3D representation (φ, θ , and T) of the sorption curves.

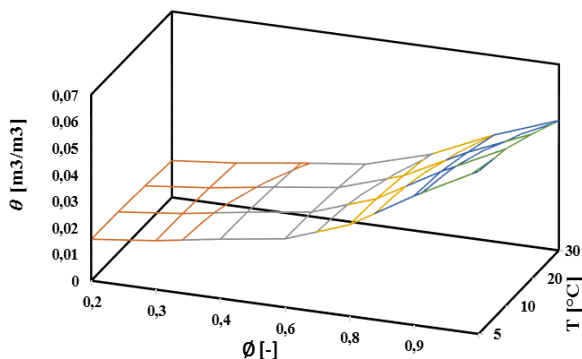


Fig. 3. 3D representation of sorption curves.

3 Numerical validation

Numerical models are investigated at the wall scale through a comparison between simulation results and experimental data provided by Lelievre et al [8]. The sample studied is a single layer wall composed of 36 cm of hemp concrete, subjected during 87 days to indoor and outdoor conditions for temperature and relative humidity as shown in Fig.4. Temperature and relative humidity through the wall are monitored at three different locations using specific sensors. Properties of hemp concrete are listed in Table 1.

Table 1. Hemp concrete properties.

Properties	Hemp concrete
Thermal conductivity [W.m ⁻¹ .K ⁻¹]	$\lambda = 0.00818 + 0.000276T + 0.0024w$
Density [Kg.m ⁻³]	$\rho = 450$
Thermal capacity [J.kg ⁻¹ .K ⁻¹]	$C_p = 1000$
Coefficient of resistance to vapor [-]	$\mu = 5$

Four models are tested and compared, labelled respectively Mod_1, Mod_2, Mod_3 and Mod_4. Their characteristics are presented in Table 2.

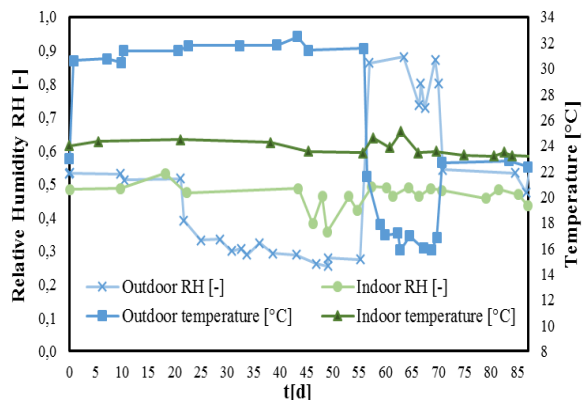


Fig. 4. Experimental test conditions.

Table 2. Numerical models characteristics.

Models	Description
Mod_1	<ul style="list-style-type: none"> Standard model for heat and moisture transfer with average sorption curve Initial conditions at 23°C and 50% relative humidity
Mod_2	<ul style="list-style-type: none"> Model for heat and moisture transfer including hysteresis Initial conditions at 23°C and 50% relative humidity from adsorption curve
Mod_3	<ul style="list-style-type: none"> Model for heat and moisture transfer including temperature dependency Initial conditions at 23°C and 50% relative humidity
Mod_4	<ul style="list-style-type: none"> Global model for heat and moisture transfer including both hysteresis and temperature dependency Initial conditions at 23°C and 50% relative humidity from desorption curve

The results for relative humidity and temperatures are plotted at the three different locations monitored in the sample: $x = 5$ cm, $x = 18$ cm and $x = 29$ cm.

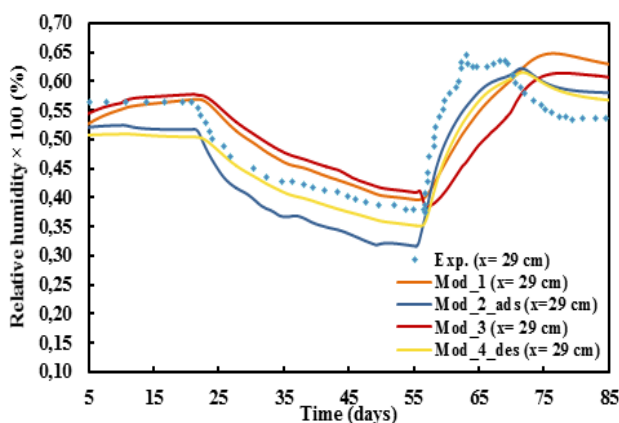


Fig. 5. Comparison between relative humidity profiles at $x = 29$ cm.

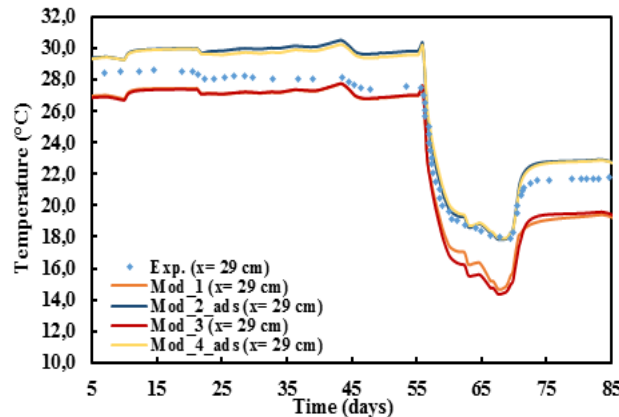


Fig. 6. Comparison between temperature profiles at $x = 29$ cm.

At $x = 29$ cm, temperature profiles obtained (Fig. 5) are acceptable for the four models up to the 57th day with

an underestimation (between 0.5 ° C and 2°C) for the Mod_1 and Mod_3 and an overestimation (from 1.5 to 2 ° C) of the temperature for Mod_2 and Mod_4. Beyond that, the temperature is advantageous for Mod_2 and Mod_4 and deteriorated for Mod_1 and Mod_3 with differences reaching 4 ° C. Fig.6 shows that when the temperature is overestimated after the first 57 days, the relative humidity is overestimated and vice versa, with remarkable advantages for Mod_4 whose deviations from the experimental do not exceed 5%. From 57 days, Mod_2 and Mod_4 represent well the dynamic appearance of the variations while Mod_1 and Mod_3 show reduced slope and delayed relative humidity peaks.

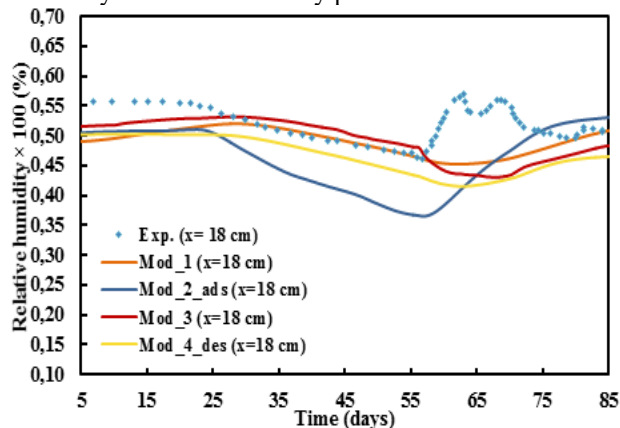


Fig. 7. Comparison between relative humidity profiles at x = 18 cm.

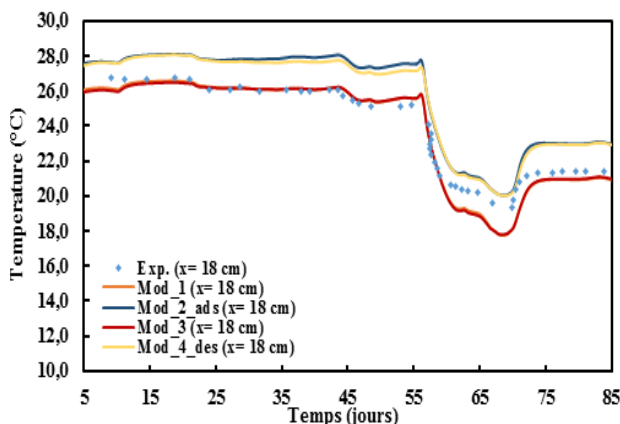


Fig. 8. Comparison between temperature profiles at x = 18 cm.

In the middle of the wall (Fig. 7), the prediction of the dynamic behaviour of relative humidity is acceptable until day 57 especially for Mod_1 and Mod_4. Beyond this, the profiles do not fit well with the experimental data for the four models. For the temperature profiles (Fig. 8), the curves obtained stick well with the experimental values for Mod_1 and Mod_3 during the 57 days, while they are overestimated by 2 ° C for Mod_2 and Mod_4. In the last twenty days, the results remain advantageous for Mod_1 and Mod_3 (maximum deviations of 1.5 ° C).

Figure 10 shows that the models are reliable for the prediction of temperature changes on the indoor side of the wall since the deviations with the experimental measurements do not exceed 1 ° C in x = 5 cm. Fig.9 gives the advantage to Mod_1 and Mod_3, and especially Mod_1, with maximum differences of about 2% with the experimental when Mod_2 succeeds in predicting the

dynamics of variations and Mod_4 presents acceptable results but a little more spread at the end (7% difference with the experimental).

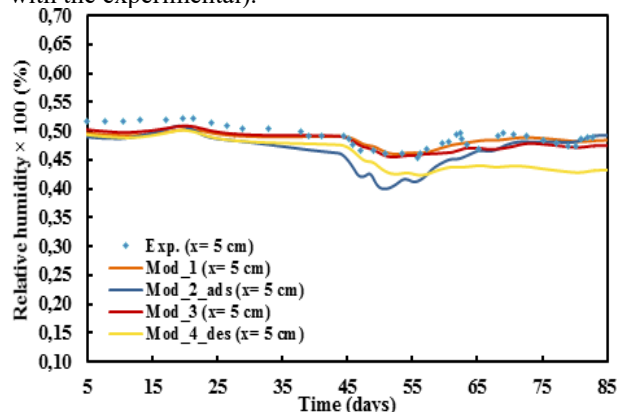


Fig. 9. Comparison between relative humidity profiles at x = 5 cm.

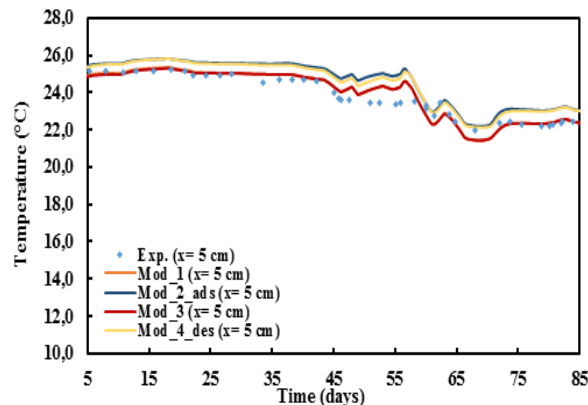


Fig. 10. Comparison between temperature profiles at x = 5 cm.

Thus, considering hysteresis improves the prediction of dynamic variations in temperature and relative humidity when the wall is subjected to these variable conditions. In addition, additional improvements are made by coupling hysteresis and temperature dependence following the results obtained by Mod_4.

4 Conclusion

This paper deals with hygrothermal behaviour study of hemp concrete via four numerical models. The hysteresis and temperature dependency phenomena are studied separately and then coupled together. Results are compared with the standard model for heat and moisture transfer. Comparison shows that predicting relative humidity profiles from the outer wall highly depends on the chosen model and presents improvements when hysteresis is considered. In parallel, temperature profiles are also improved with hysteresis at this wall side. Concerning the inner side, the four models show acceptable results for temperature and relative humidity. Thus, studying phenomena occurring within the wall requires particular attention to the selected model whereas results about investigations on indoor ambient conditions could be reliable based on one of these four models.

This work is supported by the Champagne Ardennes region and the FEDER (Fonds Européen de Développement Régional).

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