



## Two-Dimensional Transient Modeling of Energy and Mass Transfer in Porous Building Components using COMSOL Multiphysics

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(Received May 1, 2016; accepted August 6, 2016)

### ABSTRACT

This paper reports on a transient heat, air and moisture transfer (HAM) model. The governing partial-differential equations are simultaneously solved for temperature and capillary pressure through multi-layered porous media, including the non-linear transfer and storage properties of materials. Using partial differential equations functions, some thermo-physical properties of porous media are converted into coefficients depending on temperature and capillary pressure. Major features of the model are multi-dimensional and transient coupling of heat, air and moisture transport. The coupled equations are solved using the COMSOL Multiphysics time-dependent solver. This solver enables HAM (Heat, Air, Moisture) modeling in porous media. Besides, the good agreements obtained with a 2D benchmark suggest that the model can be used to assess the hygrothermal performance of building envelope components. This paper concludes that the total heat flux in the insulated wall represents only the quarter of that crossing the uninsulated concrete roof. On the other hand, the concrete having the lowest water vapour permeability of all used materials allows maintaining the vapour pressure levels close to the initial value ( $10^3$  Pa). This induces a situation of interstitial condensation within the concrete of the roof. Being able to evaluate the hygrothermal behaviour, the proposed model may turn out to be a valuable tool to solve other building problems.

**Keywords:** Numerical modeling; Moisture; Heat; Porous media; Capillary pressure.

### NOMENCLATURE

$c_{p,m}$	dry specific heat of material	$\delta_p$	water vapour permeability
$c_{p,a}$	specific heat of dry air	$\lambda$	thermal conductivity
$c_{p,l}$	specific heat of liquid water	$\phi$	relative humidity
$g$	moisture flow	$\mu_a$	air dynamic viscosity
$k_a$	air permeability of material	$\Omega$	moisture storage capacity
$K_l$	liquid water permeability	$\rho_m$	dry density of the material
$L_V$	latent heat of vaporisation	$\rho_a$	dry air density
$M$	total moisture weight	$\rho_l$	water density
$p_a$	dry air pressure	$\rho_v$	water vapour density
$p_v$	partial water vapour pressure		
$P_C$	capillary pressure		
$P_{sat}$	saturated water vapour pressure		
$P_{atm}$	atmospheric pressure		
$q$	heat flux		
$R_V$	gas constant for water vapour		
$t$	time		
$T$	temperature		
$v$	air velocity		
$w$	moisture content		
$\alpha$	heat transfer coefficient		
$\beta_p$	vapour transfer coefficient		
		<b>Subscripts</b>	
		$a$	air
		$cond$	conduction
		$conv$	convection
		$dry$	dry condition
		$e$	exterior side of building envelope
		$eq$	equivalent
		$evap$	evaporation
		$g$	gaseous
		$h$	heat

<i>i</i>	interior side of building envelope	<i>surf</i>	surface
<i>l</i>	liquid	<i>v</i>	vapour
<i>m</i>	medium	<i>wdr</i>	wind-driven rain
<i>sat</i>	saturation state		

## 1. INTRODUCTION

HAM (Heat, Air, and Moisture) models are frequently being used to solve the combined transport of heat and moisture in building components. Nowadays, a wide variety of modelling tools are available. The state of the art and the different simulation programs for heat and mass transfers in buildings have been carried out by Woloszyn and Rode (2008). In addition, an overview of different approaches to modeling heat and moisture transfer was conducted by Woloszyn *et al.* (2009). Besides, coupled heat and moisture simulation in air-conditioned buildings and its impact on energy demands has been presented by Qin *et al.* (2009) and comparison between CFD and well-mixed zonal models was performed by Steeman *et al.* (2009). Currently, substantial published simulation programs and codes are available on moisture transfer in buildings, building materials and components, and allow the prediction of the indoor thermal environment (Korjenic and Bednar 2012; Langmans *et al.* 2012).

Similarly, there are valuable basic models for the simulation of coupled transport of heat and moisture providing a valid method where possible processes associated with damage in materials and construction elements can be predicted (van Schijndel 2007a, 2007b). In this way, Qinru *et al.* (2009) developed a numerical tool to predict combined heat, air and moisture transport in building envelopes, abbreviated as HAM-BE, by making use of commercial finite-element software, COMSOL Multiphysics is also used in the present work. The material properties being expressed as analytical or interpolation functions of moisture state variables. The convective air transport may play an important role on the prediction of both temperature and moisture content distribution within porous building envelopes (dos Santos and Mendes (2009). The dominant effect of the convective vapour flow has also been noticed.

More recently, Liu *et al.* (2013) evaluated the internal surface temperature of walls made of different materials. They concluded that when the moisture transfer is taken into account, the internal surface temperature has a cooling effect on the indoor air and is beneficial to improve the indoor thermal environment during the working hours in summer. Qin *et al.* (2009, 2010) proposed both one-dimensional and two-dimensional numerical solutions based on the control volume finite difference technique with fully implicit scheme in time. In their work, they adopted the vapour content and temperature as main driving potentials. Assessment of the risk of moisture related damage in valuable objects for cases with large temperature or humidity gradients in the air is described by coupling a model for heat and moisture transport in porous

materials to a commercial Computational Fluid Dynamics (CFD) package (CEN/TC89 WG10 2007).

Litavcova *et al.* (2014) have derived a mathematical model describing liquid water and vapour diffusion in a wet material as two separate processes. Their results are compared with classical moisture transfer solution representing transfer of both liquid water and vapour as a single moisture variable. Janssen *et al.* (2007) found that the common temporal discretisation of nonlinear equations may yield conservation errors, which can be solved by starting from the mixed transfer equation form. It should be noted that the three transport phenomena (HAM) were coupled and solved simultaneously for temperature, relative humidity and pressure considered as driving potentials by Tariku *et al.* (2010). They leaned on the so called HAMFIT method which requires less time of implementation and provides a high degree of transparency and flexibility of modeling the coupled PDEs. Dufourestel (1992) used the vapour pressure as a potential and shows that his model is equivalent to De Vries model (De Vries 1958).

A newly-developed coupled heat and moisture model including vapour and liquid moisture transport was presented by Van Bellegem *et al.* (2014). It has a broader application range, as it can be used to study vapour diffusion through porous materials and capillary moisture transport. The model lends itself to study the drying phenomena of porous materials and was validated against a drying experiment in which ceramic brick is used as a porous material.

Tamene *et al.* (2011) studied the heat transfer in a multilayered wall exposed to variable solar flux. The developed numerical code allowed them to determine influence of the coupling between heat and mass transfer on the evolution of temperature and moisture in presence of variable solar flux and external temperature.

A recent study showing the impact of the temperature dependency of the sorption curves on the hygrothermal behavior of a hemp concrete building envelope was carried out by Tran Le *et al.* (2016). They concluded that taking into account the influence of temperature on the sorption curves is necessary for better prediction of the hygrothermal behavior of a hemp concrete envelope.

A 3D model for heat and water vapour transport in porous materials was integrated into a commercial CFD package (Steeman *et al.* 2009). This new model makes it possible to take the effect of indoor air distributions into account when simulating the hygric response of porous objects. The coupled model was validated using climate chamber experiments.

Attempts of combining CFD and HAM can be found

in Mortensen *et al.* (2007) and Defraeye *et al.* (2012). Defraeye applied an external coupling procedure and used a separate CFD and HAM model which were then combined by transferring and setting the boundary conditions from one model to the other. These recent attempts to combine existing HAM models with CFD showed to be promising. They allow a more detailed and accurate modelling of the influence of convection on the heat and moisture transport in porous materials. However, the development of these hygrothermal models is still ongoing.

The current article presents a modelling and a simulation approach that tries to bridge the accuracy requirements with acceptable solver performance. The major aim of the method described in this paper is to present an accurate scheme with an acceptable simulation time that can be used to evaluate the hygrothermal performance of building components. A steady state air mass balance equation is solved together with a transient energy and moisture balance equation. The calculation domain consists of porous materials in which air transport is treated as Darcy flow.

The governing partial-differential equations (PDEs) of the three transport phenomena are coupled and solved simultaneously for temperature and capillary pressure. The model accommodates non-linear transfer and storage properties of materials, moisture transfer by vapour diffusion, capillary liquid water transport and convective heat and moisture transfer through multi-layered porous media. The PDEs are derived in such a way that each PDE is described with a single driving potential, which is continuous across the interfaces of adjoining materials. The three-coupled transient HAM equations were simultaneously solved using the COMSOL Multiphysics time-dependent solver. The solver is based on an explicit scheme with variable time stepping. The user can predefine the maximum time step so that it matches with the boundary conditions change periods.

## 2. DEVELOPMENT OF GOVERNING EQUATIONS

Simultaneous heat and moisture transfer in a porous material involves complex physical phenomena. The strength of this complexity depends on how the mutual effect of heat on mass transfer is dealt with. Physical models that form the basis for various software tools used to predict the heat, air and moisture response of building envelopes seem quite diverse. In the coupled transfer, the moisture transport in building materials appears under two different phases: liquid and vapour. The vapour phase is split into diffusion and convection parts. Indeed, the diffusive flow of vapour is engendered by vapour pressure gradient and the corresponding conductivity represents the permeability of the vapour. As for, the convective vapour flow, it is advected by the moving air (Steeman *et al.* 2009). The modeling of the transfer in the vapour phase by the gradient of capillary pressure as conductive potential is the most appropriate approach and the

most used one in this kind of modeling (Carmeliet *et al.* 2004). For the liquid flow, moisture content gradient has been used as the driving potential in some hygrothermal tools, and moisture diffusivity was used as the moisture transfer conductivity.

According to the principle of the preservation of the combined transport of heat and humidity of a representative elementary volume, which is defined as being large enough when compared to pore dimensions but small enough compared to the size of the sample, governing equations of the coupled transfer in building materials can be formulated. In order to simulate heat and moisture transfers in multi-layered wall, different kinds of transport equations and boundary conditions are required (Hens 2007). Hereafter, these equations are outlined according to the considered medium. These equations require dedicated boundary conditions in order to close the problem and solve the coupled equations. Note that, to characterize the humid air mixture, the assumption of ideal gas was made for dry air and water vapour.

### 2.1 Air Transfer

In this work, the transfer of air through a porous medium is implicitly included in the mass and energy conservation equations. The air flow  $g_a$  is considered constant and moisture content independent. In the current model, the air transport is individually considered through the dry-air mass balance and the conservation equation can be expressed by Eq. (1) as:

$$\partial \rho_a / \partial t + \nabla \cdot g_a = 0 \quad (1)$$

Poiseuille's law (Steeman *et al.*, 2009) linking the pressure gradient to the flow velocity ( $v$ ) of air is given in Eq. (2):

$$v = -(k_a / \mu_a) \nabla p_a \quad (2)$$

Equation (3) allows calculating air flow under the following way:

$$g_a = \rho_a v \quad (3)$$

where  $k_a$  (kg/m.s.Pa) is the air permeability of the material,  $\mu_a$  (Pa.s) is air dynamic viscosity,  $p_a$  (Pa) is the dry air pressure,  $\rho_a$  (kg/m<sup>3</sup>) is the dry air density.

Therefore, the dry air transport given in Eq. (4) below can be written by substituting Eq. (3) in Eq. (1). This leads to the following equation:

$$\partial \rho_a / \partial t = v \cdot \nabla \rho_a \quad (4)$$

### 2.2 Moisture Transfer

It should be noted that the moisture transfer through composite walls, even in a one-dimensional process is a complex phenomenon that involves coupled transfers of liquid, vapour and heat. The airflow through a structure is driven by air pressure differences. The HAM-model, used in this study, is

a finite element model based on the standard partial differential equations of coupled heat, air and moisture transfer in porous building materials (Hagentoft *et al.* 2004). Moisture transfer is split into vapour and liquid flow parts as shown in Eq. (5):

$$g = g_v + g_l \quad (5)$$

In Eq. (6), the liquid flux is described by Darcy's law:

$$g_l = -K_l \nabla P_C \quad (6)$$

where  $K_l$  (s) is the liquid water permeability and  $P_C$  is the capillary pressure.

Note that vapour flow given in Eq. (7) is split into diffusion and convection driven parts:

$$g_v = -\delta_p \nabla p_v + \rho_v v \quad (7)$$

where  $p_v$  is the partial water vapour pressure,  $v$  is the air velocity, and  $\rho_v$  is the water vapour density.

The mass balance is expressed in Eq. (8):

$$\partial w / \partial t + \nabla \cdot g = 0 \quad (8)$$

$w$  (kg/m<sup>3</sup>) being the moisture content and  $t$  (s) is the time.

The moisture content is written in Eq. (9) in terms of temperature and capillary pressure derivative in the following way

$$\partial w / \partial t = (\partial w / \partial P_C) \cdot (\partial P_C / \partial t) \quad (9)$$

Relative humidity is often chosen as a flow potential, since it is continuous at the interface of layers of materials having different moisture storage properties. This potential is linked to the capillary pressure via relationship given by Eq. (10) (Kelvin's law):

$$P_C = \rho_l R_v T \ln \phi \quad (10)$$

Where  $\rho_l$  is the water density and  $R_v$  is the gas constant for water vapour.

Combining the six last equations, one can express in Eq. (11) the governing moisture content equation in a porous material

$$\begin{aligned} \frac{\partial w}{\partial P_C} \frac{\partial P_C}{\partial t} = & \nabla \cdot (K_l \nabla P_C) + \nabla \cdot \left( \frac{\rho_v}{\rho_l} \nabla P_C \right. \\ & \left. + \left( \phi \frac{\partial p_{sat}}{\partial T} - \frac{p_v \ln \phi}{T} \right) \nabla T \right) - v \cdot \nabla \rho_v \end{aligned} \quad (11)$$

### 2.3 Heat Transfer

For transient problems, in the absence of heat generation within the medium, the heat flow written in Eq. (12) consists of two parts (Eqs. 12-13), namely the conduction and convection parts, respectively.

$$q = q_{cond} + q_{conv} \quad (12)$$

The conductive part is given by

$$q_{cond} = -\lambda \nabla T \quad (13)$$

where  $T$  (K) is the temperature and  $\lambda$  (W/m.K) is the thermal conductivity. Equation (14) expresses the convection of sensible and latent heat part, it is given by

$$q_{conv} = v \rho_a c_{p,a} \nabla T + g_v L_v \quad (14)$$

where  $L_v$  (J/kg) is the latent heat of vaporization,  $\rho_a$  (kg/m<sup>3</sup>) is the dry air density, and  $c_{p,a}$  (J/kg.K) is the specific heat of dry air. Note that the specific heat capacity refers to humid air, while the second term accounts for latent heat.

Equation (15) gives the energy balance equation

$$-\nabla \cdot q = (c_{p,m} \rho_m + c_{p,l} w) \left( \frac{\partial T}{\partial t} \right) \quad (15)$$

where  $c_{p,m}$  (J/kg.K) is the dry specific heat of material,  $\rho_m$  (kg/m<sup>3</sup>) is the dry density of the material, and  $c_{p,l}$  (J/kg.K) is the specific heat of liquid water.

Equations (12) - (15) allow us writing the recovered equation (Eq. 16) for heat transfer as:

$$\begin{aligned} (c_{p,m} \rho_m + c_{p,l} w) (\partial T / \partial t) = & \nabla \cdot (\lambda \nabla T) \\ & + L_v \nabla \cdot (\delta_p \nabla p_v) - v \cdot (L_v \nabla \rho_v + \rho_a c_{p,a} \nabla T) \end{aligned} \quad (16)$$

### 2.4 Mathematical Modelling

This section addresses the model to be implemented for solving the coupled governing Eq. (11) and (16) which can be written in term of coefficients, and by considering temperature  $T$  as independent variable for heat transfer and capillary pressure  $P_C$  as an independent variable for moisture transfer. In fact, it seems appropriate to assume that the moisture transfer is driven by gradients in capillary pressure (Tariku 2008). Thereby, this quantity can be regarded as a potential.

The capillary pressure  $P_C$  given in Eq. (17) is defined as the pressure difference between the liquid and the gaseous phase:

$$P_C = p_g - p_l \quad (17)$$

Moreover, the relationship between the partial water vapour pressure and the relative humidity can be expressed by Eq. (18) as:

$$p_v = \phi P_{sat} \quad (18)$$

where  $\phi$  is the relative humidity, and  $P_{sat}$  is the saturated water vapour pressure.

In Eqs. (19) - (20), we simply express the pressure, density gradients of water vapour, and the time derivative of the moisture content, in terms of capillary pressure and temperature according to the following formulation:

$$\nabla p_v = (\partial p_v / \partial T) \cdot \nabla T + (\partial p_v / \partial P_C) \cdot \nabla P_C \quad (19)$$

$$\nabla \rho_v = (\partial \rho_v / \partial T) \cdot \nabla T + (\partial \rho_v / \partial \phi) (\partial \phi / \partial P_C) \cdot \nabla P_C \quad (20)$$

Note that in the last equation, the time derivative of the moisture is none other than the moisture storage capacity, which is the slope of water retention curve ( $\Omega = \partial w / \partial P_C$ ).

By injecting Eqs. (10) - (18) in Eqs. (19) - (20) and using Eq. (9), the governing equations (11) and (16) can be rewritten in term of PC and T gradients only, as shown in Eqs. (21) - (22):

$$\begin{aligned} (c_{p,m} \rho_m + c_{p,l} w) \frac{\partial T}{\partial t} = \\ \nabla \cdot \left[ (\lambda + L_v \delta_p \phi P'_{sat}) \nabla T - \frac{L_v \delta_p \phi P_{sat}}{\rho_l R_v T} \nabla P_C \right] \\ + v \cdot \left[ - \left( \rho_a c_{p,a} + L_v \frac{\partial \rho_v}{\partial T} \right) \nabla T - \left( \frac{L_v \phi}{\rho_l R_v T} \cdot \frac{\partial \rho_v}{\partial \phi} \right) \nabla P_C \right] \quad (21) \end{aligned}$$

$$\begin{aligned} \xi \frac{\partial P_C}{\partial t} = \nabla \cdot \left[ -\delta_p \phi P'_{sat} \nabla T + \left( K_l + \frac{\delta_p \phi P_{sat}}{\rho_l R_v T} \right) \nabla P_C \right] \\ + v \cdot \left[ \frac{\partial \rho_v}{\partial T} \nabla T - \left( \frac{\phi}{\rho_l R_v T} \cdot \frac{\partial \rho_v}{\partial \phi} \right) \nabla P_C \right] \quad (22) \end{aligned}$$

$P'_{sat} (= \partial P_{sat} / \partial T)$  being the derivative of saturation vapour pressure.

To go further, the current model consists of converting, via MatLab, the measurable physical properties of the material such as  $K_l$ ,  $\phi$ ,  $\delta_p$  and  $\lambda$  which depend on moisture content  $w$  into partial differential equations (PDEs),  $C_{11}, C_{12}, D_{11}, D_{12}, \Omega$  and  $C_T$  which are dependent on  $P_C$  and  $T$  (Qin *et al.* 2010). Hence, the above expressions may be simplified under the form given in Eqs. (23) - (24) :

$$\begin{aligned} C_T (\partial T / \partial t) = \nabla \cdot (C_{11} \nabla T + C_{12} \nabla P_C) \\ + v \cdot (D_{11} \nabla T + D_{12} \nabla P_C) \quad (23) \end{aligned}$$

$$\begin{aligned} \Omega (\partial P_C / \partial t) = \nabla \cdot (C_{21} \nabla T + C_{22} \nabla P_C) \\ + v \cdot (D_{21} \nabla T + D_{22} \nabla P_C) \quad (24) \end{aligned}$$

where  $C_T$  is the specific heat of the medium defined as a function of dry air and liquid water heat capacities.

After simple rearrangements, Eqs. (23) - (24) can be rewritten under the simple matrix form given in Eq. (25):

$$d_a \begin{bmatrix} \partial T / \partial t \\ \partial P_C / \partial t \end{bmatrix} = \nabla \cdot \left( C \nabla \begin{bmatrix} T \\ P_C \end{bmatrix} \right) + \beta \cdot \nabla \begin{bmatrix} T \\ P_C \end{bmatrix} \quad (25)$$

where damping ( $d_a$ ), diffusion ( $C$ ) and convection

( $\beta$ ) matrices are respectively defined by Eqs. (26) - (27) - (28):

$$d_a = \begin{bmatrix} C_T & 0 \\ 0 & \Omega \end{bmatrix} \quad (26)$$

$$\begin{aligned} C = \begin{bmatrix} C_{11} & C_{12} \\ C_{21} & C_{22} \end{bmatrix} \\ = \begin{bmatrix} \lambda + L_v \delta_p \phi P'_{sat} & -\frac{L_v \delta_p \phi P_{sat}}{\rho_l R_v T} \\ -\delta_p \phi P'_{sat} & K_l + \frac{\delta_p \phi P_{sat}}{\rho_l R_v T} \end{bmatrix} \quad (27) \end{aligned}$$

$$\begin{aligned} \beta = v \cdot \begin{bmatrix} D_{11} & D_{12} \\ D_{21} & D_{22} \end{bmatrix} = \\ v \cdot \begin{bmatrix} - \left( \rho_a c_{p,a} + L_v \frac{\partial \rho_v}{\partial T} \right) & -\frac{L_v \phi}{\rho_l R_v T} \frac{\partial \rho_v}{\partial \phi} \\ \frac{\partial \rho_v}{\partial T} & - \left( \frac{\phi}{\rho_l R_v T} \frac{\partial \rho_v}{\partial \phi} \right) \end{bmatrix} \quad (28) \end{aligned}$$

### 3. BOUNDARY CONDITIONS

In building envelope simulation, the boundary conditions (BCs) and initial conditions (ICs), carefully selected and combined, greatly help to achieve good accuracy without resorting to the use of prohibitive computing times. In this framework, external BCs of building envelopes belong to three main groups (Künzel et Kiessel 1997; Hagentoft 2002) that are moisture saturation, constant heat and moisture flow, and heat/moisture flow through surface resistance film fixed on external surface. Equations. (29-30) given below provide the exterior boundary conditions based on Dirichlet assumptions. As for the internal surface of the wall, the temperature and pressure are maintained constant.

#### 3.1 Moisture Boundary Conditions

Originally, the moisture supply from the environment  $g_{n,e}$  consists of two terms, namely the moisture supply due to wind-driven rain  $g_{wdr}$  and the evaporation due to unequal vapour pressure between material and the surrounding air  $g_{evap}$ . Moreover, this form is extended as shown in with the runoff module that uses the excess moisture to calculate the liquid film flow at the surface and the amount of runoff. This occurs whenever the moisture supply, due to the wind-driven rain ( $wdr$ ), exceeds the possible amount of absorbed moisture at the material surface:

$$g_{n,e} = \underbrace{\beta_{p,e} (p_{v,e} - p_{surf,e})}_{g_{evap}} + g_{wdr} + g_{runoff} \quad (29)$$

where  $\beta_{p,e}$  is the surface vapour transfer coefficient,  $p_{v,e}$  is the water vapour pressure of the

outdoor air, and  $p_{surf,e}$  is the water vapour pressure at the surface of the building envelope part.

Noting that benchmarking of the above numerical model supposes that moisture content at surface is limited to the saturated value, and then no wind-driven rain and runoff water are considered in the calculations. Therefore, Eq. (29) reduces to the evaporation term  $g_{evap}$ .

For the internal side of the wall, the moisture flux is obtained according to the following relationship:

$$g_{n,i} = \beta_{p,i} (p_{v,i} - p_{surf,i}) \quad (30)$$

where  $\beta_{p,i}$  is the vapour transfer coefficient of the interior surface,  $p_{v,i}$  is the water vapour pressure of the indoor air and  $p_{surf,i}$  is the water vapour pressure of the interior surface.

### 3.2 Heat Boundary Conditions

The heat flow across the exterior surface expressed in Eqs. (31) - (32) given below, includes the effects of conduction, convection, latent heat flow due to vapour transfer and sensible heat flow due to rain absorption, so radiation is not considered, except for the influence on  $T_{eq}$  (longwave radiation)

$$q_{n,e} = \alpha_e (T^{eq} - T_{surf,e}) + \beta_{p,e} (L_V + C_{p,m} T) (p_{v,e} - p_{surf,e}) + g_{wdr} c_{p,l} T^{eq} \quad (31)$$

where  $\alpha_e$  is the convective heat transfer coefficient of the exterior surface,  $T^{eq}$  is the equivalent exterior temperature and  $T_{surf,e}$  is the temperature of the exterior surface.

Likewise, heat flux through internal surface of the building envelope,  $q_{n,i}$ , is given by:

$$q_{n,i} = \alpha_i (T_i - T_{surf,i}) + L_V \beta_{p,i} (p_{v,i} - p_{surf,i}) \quad (32)$$

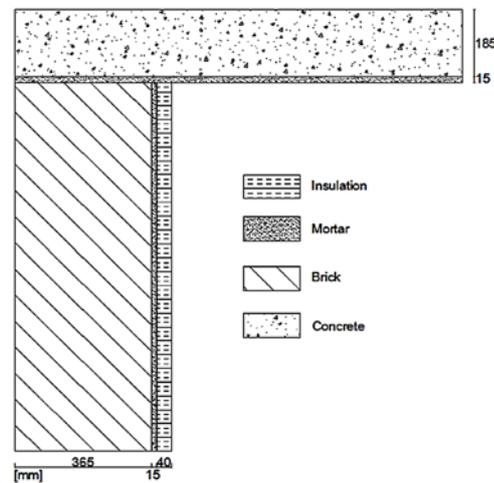
where  $\alpha_i$  is the heat transfer coefficient of the interior surface,  $T_i$  is the temperature of the indoor air, and  $T_{surf,i}$  is the temperature of the interior surface.

## 4. NUMERICAL HANDLING OF THE TWO-DIMENSIONAL CASE

The validation process has been achieved. For this, the working of the aforementioned model equations has been recently implemented by Maliki *et al.* (2014, 2015) considering the 1D benchmark exercise #1. The latter arises from a series of benchmark cases from a work outcome of the EU-initiated project for standardization of heat, air and moisture calculation methods (European project known as HAMSTAD-WP2) (Hagentoft *et al.* 2004; Hagentoft 2002).

The purpose of this study is to extend the assessment of such model to two-dimensional cases. An overview of the roof structure and walls is depicted in Fig.1. Structure details for the 2D air permeable roof-wall corner are also provided. It is represented by three layers, namely mortar, brick and concrete. This is supplemented by transfer coefficients and boundary conditions.

The HAMSTAD benchmark Exercise #5 considered here deals with interstitial condensation occurring at the contact surface between two materials. This test case is challenging (Hagentoft *et al.* 2004) as it involves severe climatic load that causes surface condensation on the exterior surface due to nighttime cooling (low equivalent temperature), and frequent occurrences of wetting and drying of the wall due to the alternating rain and solar radiation loads.



**Fig. 1. Structure details for the 2D air permeable roof-wall corner.**

To complete this overview, Table 1 gathers constituent materials' properties. First though, the simulation of the benchmark exercise #1 involving condensation occurring at the contact surface between the two materials has enabled us to appraise the model's ability to correctly predict the two-dimensional case considered here. The required results are temperature field, heat flux crossing the wall from inside and water vapour pressure distribution across the computational domain.

**Table 1 Properties of considered materials**

Thermophysical properties	Brick	Mortar	Insulation	Concrete
$\mu_{dry} [-]$	7.5	50	5.6	120
$\xi [kg/m^3]$	9	5	1.4	85
$K_j [kg/msPa]$	7e-12	4.6e-11	5.5e-5	6e-11
$\lambda_{dry} [W/mK]$	0.682	0.6	0.06	1.6
$\rho_m [kg/m^3]$	1600	230	212	2300
$C_p [J/kgK]$	1000	920	1000	850

To undertake numerical simulation, data of the roof/wall corner, including material properties, initial condition and boundary settings, are input as constants, analytical expressions or interpolation files. This is complemented by coefficients of the conservation equation.

Needing to be able to use meshes that smoothly change in size near two different layers internal interface, where the solution is expected to change abruptly, a regular 811 mesh points has been applied producing 1496 triangular elements of normal size of only few millimeters as shown in Fig. 2.

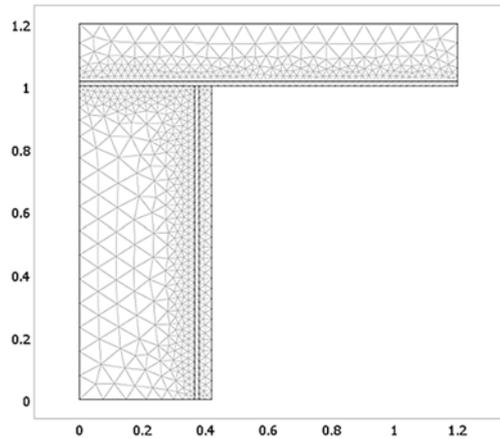


Fig. 2. Meshing generation of the roof/wall corner.

#### 4.1 Initial Conditions

For the whole construction, initial conditions are the following:

- Relative humidity:  $\phi = 60\%$
- Temperature:  $T = 298\text{ K}$
- Water vapour pressure:  $P_v = 10^3\text{ Pa}$

#### 4.2 Boundary Conditions

For heat and moisture, a data file supplies the hourly values for a period lasting over one year. The simulation period chosen for this study case is of 20 days (480h). For intermediate values of time, they are obtained by interpolation. Fig. 3 shows the variation for a period of twenty days (simulation duration) for:

- The outside equivalent temperatures  $T_e$  encompassing both the temperatures of the ambient air and that of the radiation, Note that the internal temperature  $T_i$  is set to  $20^\circ\text{C}$  in order to insure optimal interior thermal comfort.
- External and internal air pressure variations  $P_e$  and  $P_i$ .

The surface transfer coefficients are given by:

- For heat:

$$\alpha_{e,e(\text{wall \& roof})} = 25\text{ W m}^{-2}\text{ K}^{-1}$$

$$\alpha_{e,i(\text{wall})} = 8\text{ W m}^{-2}\text{ K}^{-1}; \alpha_{e,i(\text{roof})} = 25\text{ W m}^{-2}\text{ K}^{-1}$$

- For moisture:

$$\beta_{p,e} = 1.83 \cdot 10^{-7}\text{ s m}^{-1}; \beta_{p,i} = 5.88 \cdot 10^{-8}\text{ s m}^{-1}$$

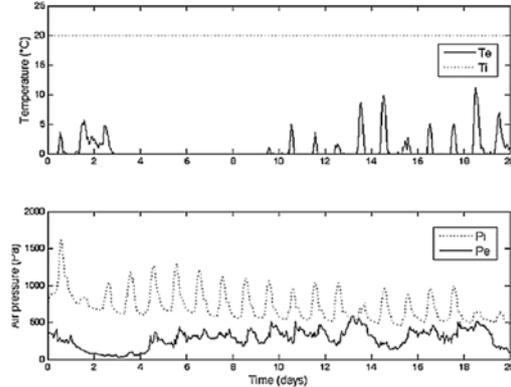


Fig. 3. Boundary conditions for the simulation period.

### 4.3 Results and Discussion

#### 4.3.1 Heat Transport

Before going further in our investigations, note first of all that we treated the heat transfer through the wall by conduction, convection and diffusion. A linear profile within the different layers is observed due to the very weak effect of the convection (no air pressure difference). In all the computations, the curves obtained are similar with those of the benchmark case as shown in Fig. 4. The highest temperature discrepancy registered is 0.5 K.

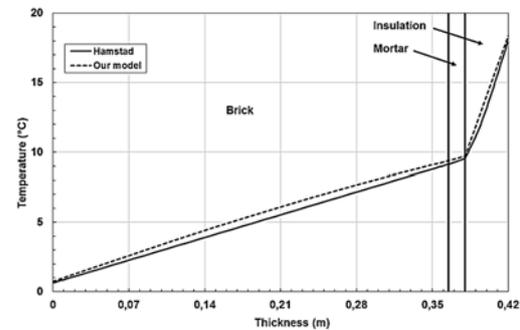
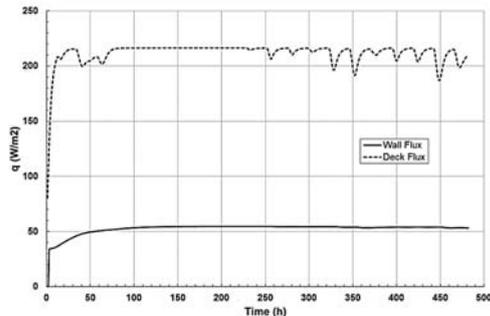


Fig. 4. Temperature profile through the wall (20 days).

Figure 5 plots the heat flux crossing the corner from the interior for the wall and roof during the first 20 days (480 hours). It should be pointed out that after 48 hours, the curve adopts a linear behavior for the heating of the wall flow, whereas the heat flux on the roof exhibits some peaks. The graph tendency of heat flux from interior to the roof depends on the period of the day; the energy amount crossing the wall is conversely proportional to outside temperature. Indeed, an ascending tendency corresponds to the period of the day included between 3 pm and 8 am when the average temperature is globally negative with a maximum outgoing flux reached at about 8 am.

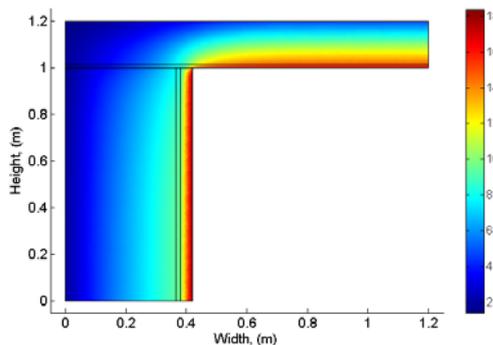
The downward parts of the graph represents the flux

behavior during the rest of the day where the average temperature is globally positive, the lowest heat flux value is registered at 3 pm. This is understandable, in so far as the insulating material's temperature serves as a regulator helping to maintain the heat flow at a lower level. The total heat flux by conduction and convection for the wall represents the quarter of that crossing the roof.



**Fig. 5. Time-dependent change of heat flux from interior to the wall and to the roof.**

The insulating material facing the inner surface of the wall regulates the temperature and increase outstandingly the thermal inertia of the wall, limits considerably the heat flow dissipating towards the outside and thus energy losses as shown in Fig. 4. The roof deprived of any insulation adopts a thermal behavior opposite to that of the wall. Indeed, the distribution of the temperature is perfectly regular along the thickness of the roof, leading to a very important waste in heat flow during the simulation period; this can be easily checked by comparing the flow crossing the wall to that crossing the roof (Fig. 4). The insulating material in the wall limits the heat flow dissipating towards the outside and limits considerably energy losses as shown in Fig. 6.



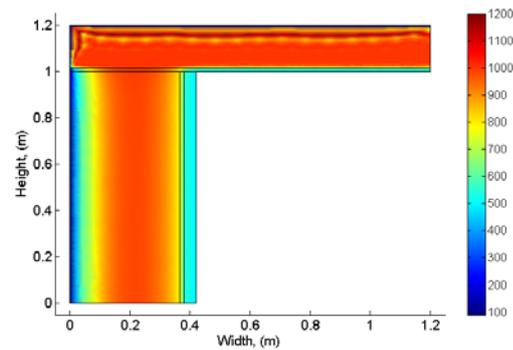
**Fig. 6. Temperature distribution after 480 hours.**

#### 4.3.2 Moisture Transport

Moisture distribution was also investigated. From Fig. 7, one can observe that the calculated vapour pressure reaches very quickly the interior vapour pressure evaluated to 480 Pa. This is due to the high air permeability of the insulating layer. Also, the vapour pressure dissipating through the porous brick after 480 hours occurs in a very slow way when compared against the two other layers. On the other hand, the concrete having the lowest water vapour permeability of all used materials allows maintaining

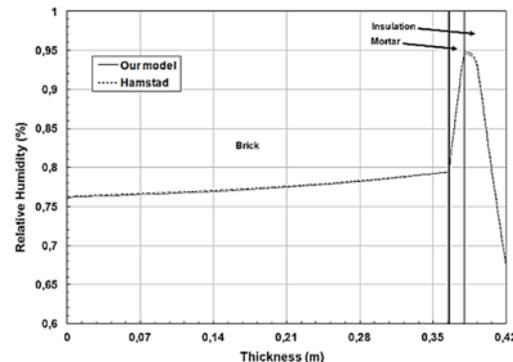
the vapour pressure levels close to the initial value ( $10^3$  Pa).

A situation of interstitial condensation is handled within the concrete of the roof. This is understandable by the very long movement of water vapour due to the low liquid diffusion coefficient of the concrete ( $6.10^{-11}$  Kg/(m.s.Pa) ). So it can be concluded that the concrete roof having a quasi airtight behaviour, contribute to a better maintain of the internal relative humidity essential for the users comfort.



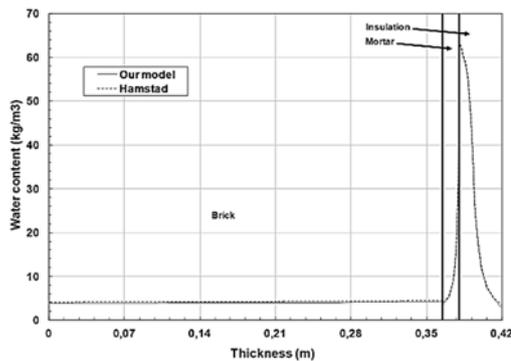
**Fig. 7. Water vapour pressure distribution after 480 hours.**

The relative humidity and moisture content profiles of the wall system after 480 hours are presented in Figs. 8 and 9, respectively. As can be seen in these figures, the simulation results of the model are in very good agreement with the HAMSTAD benchmark's solutions. In the whole building hygrothermal modeling, the coupling of building enclosure and indoor environment is through interior surfaces, and therefore, it is important to accurately predict the hygrothermal states of these surfaces to obtain useful results (Tariku *et al.* 2004).



**Fig. 8. Relative humidity evolution through the wall (after 20 days).**

As shown in the validation work, one can state that our numerical tool may be appropriate for predicting the transient HAM transport in 2D building envelopes subjected to various levels of moisture loads in the building envelope and to changing boundary conditions. It should be noted that the prediction of a problem belonging to the current context can be achieved with little time depending on the complexity of the case considered.



**Fig. 9. Moisture content profile across the wall (after 20 days).**

## 5. CONCLUSION

The present study numerically investigates the coupled heat and mass transport in porous buildings materials. The model considers the capillary pressure and the temperature gradients as driving potentials of the coupled heat and moisture transfer through porous materials of building envelope. Such a model has been successfully benchmarked against a two-dimensional numerical test case.

The governing PDEs (partial differential equations) of the three transport phenomena are coupled and solved simultaneously for potentials driving herein adopted. The model accommodates non-linear transfer and storage properties of materials, moisture transfer by vapour diffusion, capillary liquid water transport and convective heat and moisture transfer through multi-layered porous media.

The PDEs are derived in such a way that each PDE is described with a single driving potential, which is continuous across the interfaces of adjoining materials.

The COMSOL-Multiphysics solver has been chosen to solve the governing equations of HAM transport. The good agreement obtained with the considered benchmark, evidences the potential of the present model and suggests that its development and implementation are promising; thus, it can be further coupled with an indoor model to create a whole building hygrothermal model, to take into account the multi-dimensional heat, air and moisture transfer. The approach borrowed here provides a reliable and efficient model for simulating the coupled heat, air and moisture transfer through multilayer building materials.

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