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# Modal reduction for a problem of heat transfer with radiation in an enclosure

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## Abstract

We propose an original method to reduce a radiation dominated heat transfer problem. To that end, the Amalgam Reduced Order Modal Model (AROMM) method, which preserves the integrity of the geometry, is coupled to the classical radiosity method. The treated application is a complex titanium piece heated by two radiant panels placed in a furnace. The AROMM method reduces by a factor 100 the computation time needed to obtain the solution, while keeping an average error below 0.1%. These very good performances open the way for control or identification procedure in a reasonable time for these kind of applications.

*Keywords:* Reduced model, modal reduction, finite elements, radiosity, thermal radiation, industrial furnace

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## 1. Introduction

Industrial applications involving thermal phenomena in enclosures are frequently encountered : climatic chambers, honeycomb structures, industrial furnaces, ... For realistic configurations, as none of the three modes of heat transfer can be neglected, the problem is complex and a numerical approach is mandatory.

Numerical resolution of heat conduction problems in solids is carried out using classical numerical methods, such as finite elements. However, for complex geometries needing a fine meshing, the time required to solve the problem can become unbearably long. It is especially true if the model is implied in a control or identification loop. To that end, reduced models have been developed [1] [2]. Reduced modal models correspond to a generalization of the Fourier time space separation method :

$$T(M, t) \approx \sum_{i=1}^{\tilde{N}} \tilde{x}_i(t) \tilde{V}_i(M) . \quad (1)$$

From the knowledge of the reduced base  $\tilde{V}_i$ , the unknowns become the excitation states  $\tilde{x}_i$ , and the dimension of the problem might be greatly reduced. Modal methods do not degrade the geometry, and give access to the entire temperature field.

Different methods are based on this principle : the most popular is the proper orthogonal decomposition method (POD) which requires the knowledge of thermal fields from experimental or numerical data [3], [4], [5], [6]. The Modal Identification Method (MIM) directly identifies the modal model without determining modes. It is particularly adapted to inverse problems [7] [8]. Finally, in the Amalgam Reduced Order Modal Model (AROMM) method, a modal base is calculated by solving an eigenvalue problem adapted to the considered physical problem. The reduced model is obtained by reducing the initial base by the amalgam method. Initially developed for linear problems [9] [10], this method has since been extended to nonlinear problems [11] [12].

Numerical modeling for natural convection phenomena in enclosures is more delicate, since the entire volume of the cavity is meshed. The grid is refined near the walls to properly capture the dynamics in the boundary layers, leading to a significant increase of the number of degrees of freedom. To limit this number, the simplest solution is to only simulate the solid parts. The internal fluid is then simply modeled by a single temperature and it exchanges a convective heat flux with the walls via a convective coefficient, which is either known from correlations or experimentally. A more sophisticated solution is to employ reduced modal models [13] to simulate the fluid flow. They indeed capture the main characteristics of the flow while highly decreasing the computational complexity.

The addition of radiative phenomena exacerbates the difficulty. The heat radiation is indeed characterized by its spectral dependence, its integration into space, as well as its strong nonlinearity in temperature. When the temperature differences are small, the radiation is modeled by a linearized global coefficient [14]. For simple configurations, analytic relations that express the radiative fluxes

30 exchanged between surfaces might be used [15]. Otherwise, numerical techniques adapted to heat radiation have been developed. The radiosity method is the reference for gray and diffuse surfaces. First studies focused on very simple geometries : the view factors remain easy to calculate and the number of degrees of freedom is low. For simple cavities, both convection in the fluid and heat radiation between the surfaces were considered [16]. In other works the phenomenon of conduction  
35 in the walls has been added [17] [18]. In the 1980s, finite element formulations specific to radiation problems have been developed [19] [20] [21]. In particular, high order finite elements were used for the calculation of radiative fluxes [22] [23]. Other methods exist : Recently the Control Volume - Lattice Boltzmann Method has been introduced by [24], and ANSYS<sup>®</sup> proposes a specific radiation module using a Hybrid Finite Element Boundary Integral Method [25].

40 **The development of these powerful tools has paved the way to design issues involving radiative transfer. The objective of design is to find the value of a physical quantity that optimizes a process. This is mostly done by iterative procedures which integrate the above mentioned methods. As an example, [26] presents a technique to find the best distribution of heating sources in an oven, in order to obtain the temperature of the heated object as homogeneous as possible. Other functions can be minimized, such as finding the lowest possible wall temperature [27].** As in conduction or  
45 convection problems, reduced models could be employed to decrease the computation time needed to solve these optimization processes. As the access to the whole temperature field is imperative for the resolution of radiative dominated problems, modal reduction techniques are appropriate. However, to date, we have not found any study using modal reduction for radiation problems. This  
50 article presents an adaptation of the Amalgam Reduced Order Modal Model (AROMM) method for heat radiation and conduction problems in an enclosure.

The paper is structured as follows. In Section 2, the physical description of the considered problem is presented, with a classical finite element formulation. Section 3 investigates the reduced method allowing to solve this problem. Section 4 is dedicated to results and discussion about  
55 reduced model. Section 5 offers a conclusion to the paper.

## 2. Resolution by the finite element method of a thermal problem in an enclosure

### 2.1. Physical problem

A titanium object with a complex shape is placed on a stand in an industrial furnace. This object is heated by two radiant tubes surrounded by parabolic reflectors located at the top of the  
60 furnace (see Fig. 1). The different thermophysical properties of these elements are given in Table 1.

The purpose of this study is to compute the temperature distribution in the titanium object for several exterior boundary conditions and different radiant tubes temperatures. For each scenario, we assume that :

- 65 — the furnace exchanges heat with the outside. This is modeled by an equivalent global exchange coefficient  $h_{ext}$ . The outside temperature is supposed to be constant at  $T_{ext} =$

	Heat capacity $c [J.m^{-3}.K^{-1}]$	Thermal conductivity $k [W.m^{-1}.K^{-1}]$	Emissivity $\varepsilon$	Thickness $e [m]$
Piece	$2.35 \times 10^6$	21.9	0.8	0.001
Stand	$3.95 \times 10^6$	16.3	0.95	0.005
Wall	$0.18 \times 10^6$	45	0.95	0.01
Tube	$3.4 \times 10^6$	45	0.95	0.01
Reflector	$3.95 \times 10^6$	16.3	0.3	0.001

TABLE 1: Thermophysical characteristics of the different components

293.15K ;

- the convective exchanges between the interior surfaces of the furnace are represented by a constant coefficient  $h_{int} = 5 W.m^{-2}.K^{-1}$  ;
- the radiant tubes are considered to be at constant temperature  $T_{tube}$ .

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Table 2 presents the parameter values for the different scenarios.

	$h_{ext} [W.m^{-2}.K^{-1}]$	$T_{tube} [K]$
Scenario 1	5	673.15
Scenario 2	1	973.15
Scenario 3	0.5	1273.15

TABLE 2: Variable boundary conditions

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The initial condition is the ambient temperature  $T_0 = 293.15K$ , and the simulation duration to reach the steady state is  $5 \times 10^4 s$ . Considering the high temperature of the tubes, the heating of the titanium piece is mainly driven by radiative transfer. The complex geometry creates masks, and a fine detailed resolution of the radiative flux is mandatory to correctly represent the temperature evolution.

## 2.2. Finite element formulation of the problem

### 2.2.1. Adapting the heat equation for thin shells in an enclosure

Considering the above characteristics, the maximum Biot Number is  $Bi = \frac{he}{2k} \leq 0.072$  (with  $e = 0.5 cm$ ,  $k = 16.3 W.m^{-1}.K^{-1}$  and  $h_{max} = 4\varepsilon\sigma T_{max}^3 \simeq 500 W.m^{-2}.K^{-1}$ ), validating a shell model hypothesis [28] [29]. In a shell model, the temperature field depends only on local coordinates  $(\eta, \zeta)$  of a plane  $\Omega$ , which corresponds to the main surface of the original 3D domain (see Fig. 2). Both faces  $\Omega_{(+)}$  and  $\Omega_{(-)}$ , characterized by normal vectors of opposite signs (Fig. 3.a), receive a heat flux

80

$\varphi$ . The weak formulation of the heat equation yields :

$$\int_{\Omega} e c \frac{\partial T}{\partial t} f d\Omega = - \int_{\Omega} e k \vec{\nabla} T \cdot \vec{\nabla} f d\Omega + \int_{\Omega_{(+)}} \varphi f d\Omega + \int_{\Omega_{(-)}} \varphi f d\Omega, \quad (2)$$

85 where  $f \in H_1(\Omega)$  ( $H_1(\Omega)$  being the Hilbert space such that  $f$  and its derivatives are square-integrable functions) is the test function of the variational formulation on  $\Omega$ .

Faces  $\Omega_{(+)}$  and  $\Omega_{(-)}$  can be rearranged based on their location, outside or inside the enclosure (Fig. 3.b).

For external surfaces  $\Omega_{ext}$ , a simple global exchange coefficient models furnace insulation :

$$\int_{\Omega_{ext}} \varphi f d\Omega = \int_{\Omega_{ext}} h_{ext} (T_{ext} - T) f d\Omega. \quad (3)$$

90 For interior surfaces  $\Omega_{int}$ , the heat flux is divided into distinct convective and radiative phenomena :

$$\int_{\Omega_{int}} \varphi f d\Omega = \int_{\Omega_{int}} h_{int} (T_{int}(T) - T) f d\Omega + \int_{\Omega_{int}} \varphi_{rad} f d\Omega, \quad (4)$$

where  $T_{int}(T)$  is the **unique** air temperature inside the furnace. Assuming that the heat capacity of the internal fluid is negligible compared to that of the walls, a simple heat balance provides the expression of indoor temperature :

$$T_{int}(T) = \frac{\int_{\Omega_{int}} h_{int} T d\Omega}{\int_{\Omega_{int}} h_{int} d\Omega}. \quad (5)$$

Equation (2) finally becomes :

$$\begin{aligned} \int_{\Omega} e c \frac{\partial T}{\partial t} f d\Omega &= - \int_{\Omega} e k \vec{\nabla} T \cdot \vec{\nabla} f d\Omega \\ &\quad - \int_{\Omega_{int}} h_{int} T f d\Omega + \int_{\Omega_{int}} h_{int} T_{int}(T) f d\Omega \\ &\quad - \int_{\Omega_{ext}} h_{ext} T f d\Omega + \int_{\Omega_{ext}} h_{ext} T_{ext} f d\Omega \\ &\quad + \int_{\Omega_{int}} \varphi_{rad} f d\Omega, \end{aligned} \quad (6)$$

95 where  $T_{int}$  is defined by relation (5).

### 2.2.2. Modeling the radiative heat flux with the radiosity method

In this problem, diffuse-gray bodies exchanging radiative heat flux are placed in a nonparticipating media. The radiative heat flux exchanged by a surface depends, in a complex way, on the temperature distribution of the surrounding surfaces. This flux can be obtained by the net-radiation method [17] [30]. The considered surface is first discretized in  $N_p$  elementary surfaces  $\Omega_j^e$ , on which the radiative flux is considered constant,  $\varphi_{rad} = \bar{\varphi}_j$ . **These elementary surfaces are named ‘patches’.**

The  $N_p^2$  view factors  $F_{ij}$ , which are the proportion of flux leaving a **patch**  $\Omega_i^e$  towards another elementary surface  $\Omega_j^e$ , are then computed with the polar-plane-based method [31] [32], which is an adaptation of the hemicube method [33]. An hemisphere centered on  $\Omega_i^e$  (Fig. 4) is discretized in  $N^q$  quadrangles. Their size is defined so that the value of the elementary view factors towards  $\Omega_j^e$  is constant. This avoids to compute and keep track of the elementary view factors. The hemisphere is then projected on a plan parallel to  $\Omega_i^e$ . Surface  $\Omega_j^e$  is finally projected on this plan. The view factor is retrieved from the sum of the quadrangles covered by this projection.

The radiosity method relates the mean flux  $\bar{\varphi}_i$  exchanged by **patch**  $\Omega_i^e$  to the set of mean temperatures  $\bar{T}_j$ , with  $j \in [1, N_p]$  :

$$\forall j \in [1, N_p] \quad \sum_{i=1}^{N_p} \left[ \frac{\delta_{ji}}{\varepsilon_i} - \left( \frac{1}{\varepsilon_i} - 1 \right) F_{ji} \right] \bar{\varphi}_i = - \sum_{i=1}^{N_p} (\delta_{ji} - F_{ji}) \sigma \bar{T}_i^4, \quad (7)$$

where  $\delta_{ji}$  is the Kronecker delta. **This sign convention ensures that exchanged flux  $\bar{\varphi}_i$  is negative if the surface  $S_i$  emits more flux than it absorbs.** Relation (7) can be written in matrix form :

$$\mathbf{A} \bar{\varphi} = \mathbf{B} \bar{\mathbf{T}}^4. \quad (8)$$

The mean flux exchanged by a **patch**  $\bar{\varphi}_j$  expresses as :

$$\bar{\varphi}_j = \sum_{i=1}^{N_p} r_{ji} \bar{T}_i^4, \quad (9)$$

where  $r_{ji}$  are the elements of  $\mathbf{R}_{rad} [N_p, N_p] = \mathbf{A}^{-1} \mathbf{B}$ .

### 2.2.3. Variational formulation with radiative heat flux

The variational expression of the exchanged radiative flux is derived from Eqs. (4) and (9) :

$$\begin{aligned}
\int_{\Omega_{int}} \varphi_{rad} f d\Omega &= \sum_{j=1}^{N_p} \int_{\Omega_j^e} \varphi_{rad} f d\Omega \\
&= \sum_{j=1}^{N_p} \bar{\varphi}_j \int_{\Omega_j^e} f d\Omega \\
&= \sum_{j=1}^{N_p} \sum_{i=1}^{N_p} \left( r_{ji} \bar{T}_i^4 \right) \int_{\Omega_j^e} f d\Omega.
\end{aligned} \tag{10}$$

115 The addition of this term to the variational formulation Eq. (6) leads to :

$$\begin{aligned}
\int_{\Omega} e c \frac{\partial T}{\partial t} f d\Omega &= - \int_{\Omega} e k \vec{\nabla} T \cdot \vec{\nabla} f d\Omega \\
&\quad - \int_{\Omega_{int}} h_{int} T f d\Omega - \int_{\Omega_{ext}} h_{ext} T f d\Omega \\
&\quad + \int_{\Omega_{int}} h_{int} T_{int} f d\Omega + \int_{\Omega_{ext}} h_{ext} T_{ext} f d\Omega \\
&\quad + \sum_{j=1}^{N_p} \sum_{i=1}^{N_p} \left( r_{ji} \bar{T}_i^4 \right) \int_{\Omega_j^e} f d\Omega.
\end{aligned} \tag{11}$$

### 2.3. Numerical resolution

The spatial discretization of the problem defined by Eqs. (5) and (11) through Lagrange finite elements brings out the following matrix problem (hereafter called detailed model) :

$$\mathbf{C} \frac{d\mathbf{T}}{dt} = [\mathbf{K} + \mathbf{H}_{int} + \mathbf{H}_{ext}] \mathbf{T} + \mathbf{U}_{cpl} T_{int}(\mathbf{T}) + \mathbf{U}_0 + \bar{\mathbf{R}}_{rad} \bar{\mathbf{T}}^4. \tag{12}$$

Vector  $\mathbf{T}$  contains the temperature value at the  $N$  discretization points, and  $\mathbf{U}_0$  is a vector representing the external known solicitations (**the received external heat flux**  $\int_{\Omega} h_{ext} T_{ext}$ ).  $\mathbf{C}$ ,  $\mathbf{K}$ ,  $\mathbf{H}_{int}$ ,  $\mathbf{H}_{ext}$  are  $[N \times N]$  **symmetric sparse** matrices :  $\mathbf{C}$  is the thermal inertia matrix,  $\mathbf{K}$  the conductivity matrix and  $\mathbf{H}_{int}$  and  $\mathbf{H}_{ext}$  are accommodation matrices. Vector  $\mathbf{U}_{cpl}$  represents the convective



	$N_t$	$N_p$	$(\langle S_{\Omega_i^e} \rangle) (m^2)$
Piece	2 891	5 782	$5.02 \times 10^{-4}$
Stand	1 562	3 124	$1.15 \times 10^{-3}$
Wall	3 566	3 566	$2.80 \times 10^{-3}$
Tubes	11 144	22 288	$3.90 \times 10^{-5}$
Reflectors	5 039	10 078	$2.60 \times 10^{-3}$
Total	24 202	44 838	

TABLE 3: Number of triangular elements ( $N_t$ ) and number of patches ( $N_p$ ) of the different components

exchange with the air inside the furnace, at temperature  $T_{int}(\mathbf{T})$ . This latter term is computed thanks to the discretization of Eq. (5) :

$$T_{int}(\mathbf{T}) = \mathbf{D} \mathbf{T}. \quad (13)$$

Vector  $\bar{\mathbf{T}}$  of dimension  $[N_p]$  contains mean temperatures of every patch  $\Omega_i^e$  and is easily estimated by :

$$\bar{\mathbf{T}} = \mathbf{U}_R \mathbf{T}. \quad (14)$$

Matrix  $\mathbf{U}_R$  of dimension  $[N_p \times N]$  is the average operator which allows to compute the mean temperature on the  $N_p$  patches from the temperature at the  $N$  nodes of the mesh.

Finally, radiation matrix  $\bar{\mathbf{R}}_{rad} [N \times N_p]$  allots the mean heat flux density from the  $N_p$  patches to the  $N$  nodes. This matrix is computed with the following relation

$$\bar{\mathbf{R}}_{rad} = (\mathbf{U}_R)^T \mathbf{S} \mathbf{R}_{rad},$$

where  $\mathbf{S}$  is a diagonal matrix of dimension  $[N_p \times N_p]$  such that  $S_{ii}$  contains the surface of the patch  $i$ . The components of vector  $\bar{\mathbf{T}}^4$  contains the components of the vector  $\bar{\mathbf{T}}$  to the power 4.

#### 2.4. Results

All calculations were performed on a laptop with a 6-core Intel<sup>®</sup> Xeon<sup>®</sup> E-2176M @ 2.7GHz and 64GB of RAM. The discretization used for all the calculations is a meshing of  $N = 12\,167$  nodes which define  $N_t = 24\,202$  triangular elements. Table 3 details the meshing component by component. The mesh has only been refined on the radiant tubes, the source of luminance, so that the direction of the emitted flux is well defined, and also on the platinum object, which is the critical parameter. The determination of the patches is an open question. In this study, the patches coincide with the triangles of the mesh. As inner surfaces are doubled (they radiate via their two faces  $\Omega_{(+)}$  and  $\Omega_{(-)}$ , see Fig. 3),  $N_p = 44\,838$  patches  $\Omega_j^e$  emerge. First step is to perform the preliminary calculation needed by radiation. The computation time needed to get the  $N_p^2$  view factors is  $t_{CPU}^{F_{ij}} \approx 1h$ . The computation of  $\mathbf{A}^{-1}\mathbf{B}$  lasts  $t_{CPU}^{rad} \approx 8h$ . These two calculations are performed only once and are therefore reused for each simulation.

The temporal discretization of equation (12) is done with a semi-implicit second-order scheme with variable time steps [34].

135 The numerical simulation of  $5 \times 10^4$  s requires a CPU time that is dependent on nonlinearities, as it varies between 4 and 6 hours.

Figure 5 presents temperature fields at  $t = 3000$  s for the three scenarios, characterized by very different heat source temperature. The first remark concerns the temperature range. As the maximum temperature goes from  $673.15$  K to  $973.15$  K (**scenario 1 and 2**), the minimum temperature increases only of  $55$  K. But when another increase of  $300$  K is applied on the radiant tubes temperature (**scenario 3**), minimum temperature gains more than  $150$  K, underlining the non-linear effect of radiation. The second remark concerns the shape of the temperature fields. As the heat transfer by radiation is predominant, mask effect plays a significant role. This is particularly apparent on the titanium object. For scenario 3, its top surface is at almost  $800$  K, while the bottom one, which is not directly exposed to the radiant tubes, is around  $600$  K. Finally, even if the temperature contour on the titanium item barely changes from one simulation to another, its temperature amplitude is obviously dependent on the applied scenario. This preliminary study shows that the radiation has to be modeled in all its complexity to give significant results. This fine modeling results in an important CPU time, which forbids parametric study, hence the need for reduced models considering radiation.

### 3. Modal method

In modal methods, the temperature field  $T(M, t)$  is searched as a weighted sum of elementary functions  $V_i$  called modes :

$$T(M, t) = \sum_{i=1}^{\infty} x_i(t) V_i(M). \quad (15)$$

Weighting coefficients  $x_i(t)$  are now the unknowns, called excitation states. Virtually, the sum does not reach infinity. When the modes are obtained after spatial discretization, the number of modes numerically accessible corresponds to the number of mesh points  $N$ .

155 Different modal approaches exist. In this study, modes are eigenfunctions of the Laplace operator, with Neumann Boundary conditions :

$$- \int_{\Omega} e k \vec{\nabla} V_i \cdot \vec{\nabla} f d\Omega = z_i \int_{\Omega} e c V_i f d\Omega, \quad (16)$$

where  $z_i$  is the associated eigenvalue, homogeneous to the inverse of a time.

Orthogonality relationships play a fundamental role in modal methods. They ensure that the decomposition (15) is unique. Modes defined by Eq. (16) verify the following relations :

$$\forall i, j \in [1, N], \quad \int_{\Omega} e c V_i V_j d\Omega = \delta_{ij}, \quad (17)$$

$$\forall i, j \in [1, N], \quad \int_{\Omega} e k \vec{\nabla} V_i \cdot \vec{\nabla} V_j d\Omega = z_i \delta_{ij}. \quad (18)$$

160 *3.1. Reduction step*

The purpose of the modal reduction is to approach the thermal field with a reduced base  $(\tilde{z}_i, \tilde{V}_i)$ , of dimension  $\tilde{N} \ll N$  :

$$\tilde{T}(M, t) = \sum_{i=1}^{\tilde{N}} \tilde{x}_i(t) \tilde{V}_i(M). \quad (19)$$

This is performed by the amalgam reduction method. In this method, dominant modes are selected. The remaining modes, called minors, are aggregated to the dominant ones, instead of being discarded. The resulting reduced modes, named amalgamated modes, are then a linear combination of the original modes :

$$\forall i \in \{1, \dots, \tilde{N}\} \quad \tilde{V}_i(M) = V_{i,0} + \sum_{p=1}^{N_i} \alpha_{i_p} V_{i_p}. \quad (20)$$

The specificity of the amalgam method is to keep every mode of the original base. To ensure the orthogonality of the reduced basis, each original eigenvector  $V_i$  is used only once. The distribution of the original modes  $V_i$  in the reduced basis  $\tilde{V}_i$ , and the determination of the weighting coefficient  $\alpha_{i_p}$  are obtained by minimizing in the modal space an energy difference on a reference simulation (more details can be found in [11]). The reference simulation shall trigger the relevant modes, and should be readily computed. In the case of a radiative problem, all surfaces are considered as black-bodies in the reference simulation. The computation is greatly simplified, as  $\mathbf{A}$  becomes the identity matrix in Eq. (8).

*3.2. Modal formulation of the problem*

*3.2.1. Reduced formulation*

The modal formulation is obtained by replacing temperature by its decomposition on the Neumann modes in Eq. (11). Neumann modes are also used for test functions  $f$ . The following equation

175 is obtained :

$$\begin{aligned}
\sum_{p=1}^{\tilde{N}} \int_{\Omega} e c \tilde{V}_p \tilde{V}_q d\Omega \frac{\partial \tilde{x}_p}{\partial t} &= - \sum_{p=1}^{\tilde{N}} \int_{\Omega} e k \vec{\nabla} \tilde{V}_p \cdot \vec{\nabla} \tilde{V}_q d\Omega \tilde{x}_p \\
&- \sum_{p=1}^{\tilde{N}} \left( \int_{\Omega_{int}} h_{int} \tilde{V}_p \tilde{V}_q d\Omega + \int_{\Omega_{ext}} h_{ext} \tilde{V}_p \tilde{V}_q d\Omega \right) \tilde{x}_p \\
&+ \sum_{p=1}^{\tilde{N}} \left( \int_{\Omega_{int}} h_{int} T_{int} \tilde{V}_q d\Omega + \int_{\Omega_{ext}} h_{ext} T_{ext} \tilde{V}_q d\Omega \right) \\
&+ \sum_{p=1}^{\tilde{N}} \sum_{j=1}^{N_p} \sum_{i=1}^{N_p} \left( r_{ji} \bar{T}_i^4 \right) \int_{\Omega_j^e} \tilde{V}_q d\Omega.
\end{aligned} \tag{21}$$

The temperature of the air inside the enclosure, as well as  $\bar{T}_i$ , the mean temperature on the patch  $\Omega_i^e$ , can also be expressed under a modal formulation :

$$T_{int} = \frac{\sum_{i=1}^{\tilde{N}} x_i \int_{\Omega_{int}} h_{int} \tilde{V}_i d\Omega}{\int_{\Omega_{int}} h_{int} d\Omega}, \quad \bar{T}_i = \frac{\sum_{j=1}^{\tilde{N}} x_j \int_{\Omega_i^e} \tilde{V}_j d\Omega}{\int_{\Omega_i^e} d\Omega}. \tag{22}$$

The modal formulation of  $T_{int}$  in Eq. (22) is injected in Eq. (21). However, as  $\bar{T}_i^4 \neq \bar{T}_i^4$ , it can not be done for  $\bar{T}_i$ . The repeated change of basis between modal and temperature space will deteriorate the expected gain in computation time.

### 3.3. Numerical resolution

180 The set of equations (21) and (22) can be written under matrix form :

$$\tilde{\mathbf{V}}^T \mathbf{C} \tilde{\mathbf{V}} \frac{d\tilde{\mathbf{X}}}{dt} = \tilde{\mathbf{V}}^T [\mathbf{K} + \mathbf{H}_{int} + \mathbf{H}_{ext} + \mathbf{H}_{cpl}] \tilde{\mathbf{V}} \tilde{\mathbf{X}} + \tilde{\mathbf{V}}^T \mathbf{U}_0 + \tilde{\mathbf{V}}^T \mathbf{R}_{rad} \bar{\mathbf{T}}^4, \tag{23}$$

with

$$\bar{\mathbf{T}} = \mathbf{U}_R \tilde{\mathbf{V}} \tilde{\mathbf{X}}, \tag{24}$$

where  $\tilde{\mathbf{V}} \left[ \tilde{N} \times \tilde{N} \right]$  is the matrix containing the  $\tilde{N}$  discrete reduced modes  $\tilde{V}_p$ , and  $\mathbf{C}$ ,  $\mathbf{K}$ , etc. are finite element matrices defined in Eq. (12). Compact matrices like  $\tilde{\mathbf{V}}^T \mathbf{C} \tilde{\mathbf{V}}$  are dense and of dimension  $\left[ \tilde{N} \times \tilde{N} \right]$ .

## 4. Results and discussion

### 185 4.1. Reduced model creation

The creation of the reduced model demands preliminary computations. As for the finite element model, view factors have to be computed, and the inversion of matrix  $\mathbf{A}$  in Eq. (8) has to be performed. This is done in 9 hours. The second step consists in the computation of the modal base (Eq. (16)), done in 4 minutes. Figure 6 depicts typical examples of Neumann eigenmodes. They can be seen as temperature fields, solution of the steady-state heat equation where the volume solicitation is proportional to the temperature. For homogeneous domain, these eigenmodes are basically a combination of trigonometric functions. As the order of the eigenmodes increases, its time constant decreases, yielding modes that vanish quickly. This occurs with an increase of the spatial frequency, as, for instance, shown by mode 599. Typically, with a simple truncation reduction procedure, this mode will be discarded, together with the spatial information it contains. With the amalgam procedure, this mode is kept, but its time behaviour will be imposed by another mode.

The amalgam procedure requires a reference field. As stated above, this reference field is obtained by a finite element simulation with all surfaces considered as black bodies. By doing so, the numerical problem is greatly simplified. Moreover, the reference simulation is performed over a very short period of 1 800s instead of 50 000s. With these two simplifications, the reference simulation is done in about 2 minutes. The amalgam procedure in itself is performed in ten seconds. Figure 7 shows amalgamated modes. From the Neumann eigenmodes, the amalgam procedure built new modes related to the solicitations. This is evidenced by mode 3, where the temperature contour on the titanium object is similar to those shown in Fig. 5, while mode 12 provides a contribution to the parabolic reflectors.

### 4.2. Reduced Model utilization

Figure 8 illustrates the performance of AROMM method in terms of accuracy. The modal temperature field computed with a reduced model of order 200 (compared to the 12 167 nodes of the original model) is represented for scenario 3 at  $t = 3000$  s, and should be confronted to Fig. 5 (c) (same scale and contours are used). Visually the agreement between these two fields is excellent. In particular, contours produced by the complex radiative heat transfer are remarkably well recovered. This is confirmed by the difference field  $|T - \tilde{T}|$  represented in the same figure. This field is erratic, which is characteristic of modal reduction. The difference does not exceed 26 K, which has to be related to the 1 200 K of radiant tubes. The average difference of 3.3 K (which represents 0.2% of 1200 K) is very satisfactory. Note also that the maximum temperature reached by the titanium piece is particularly well predicted by the modal model with a difference of prediction with the finite element model of less than 1.5 K.

A sensitivity analysis on the reduction order has been conducted. To quantify its results the

following quantities are defined :

$$\langle \varepsilon \rangle = \frac{1}{V} \frac{1}{\tau} \frac{\int_{\tau} \int_{\Omega} |T - \tilde{T}| d\Omega}{\max_{\tau, \Omega}(T) - \min_{\tau, \Omega}(T)} \quad , \quad \varepsilon_{max} = \frac{\max_{\tau, \Omega} |T - \tilde{T}|}{\max_{\tau, \Omega}(T) - \min_{\tau, \Omega}(T)} ,$$

where  $\tau$  is the simulation duration and  $V$  the volume of domain  $\Omega$ .

220 The main conclusions are summarized in Fig. 9. Figure 9 (a) represents  $\langle \varepsilon \rangle$  and  $\varepsilon_{max}$  as a function of the order of the modal model for the three scenarios. The CPU time required to solve the problem is also represented in Fig. 9 (b). For the three scenarios, even for highly reduced model (20 modes), results are decent, with  $\varepsilon_{max} \approx 10\%$  and  $\langle \varepsilon \rangle$  below 1%. However, despite those very good relative errors, the absolute errors are less satisfying : for scenario 3, the maximum error with a reduced model of order 20 reaches 100 K. Still, about an order of magnitude is found between the maximum and mean errors : the maximum error is localized in both time and space. Obviously, errors decrease monotonously as the model order increases. With 200 modes,  $\langle \varepsilon \rangle$  is below 0.1% ( $\approx 1 K$ ) while  $\varepsilon_{max}$  is under 3%. The choice of the order is a question of compromise between accuracy and computation time. The latter logically increases with the order of the model. It still remains very small compared to the computation time required for the resolution by classical finite element model (several hours). With a reduced model of order 200, CPU time for scenario 3 is 4 minutes, which is almost a hundred time faster than the FE model which requires 6 hours. As the reduced model has proven its efficiency on three scenarios characterized by different thermal solicitations (different global exchange coefficients and different radiant tube temperature), it is suitable for parametric study, or control loops.

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## 5. Conclusion

The purpose of this study was to demonstrate the feasibility and the relevance of reducing a thermal radiation problem for a realistic configuration, while keeping the complexity engendered by the radiation and the geometry. The employed method combines a modal reduction technique and the radiosity method. Indeed, modal methods give access to the whole temperature field, which is mandatory in radiative dominated problems. In this study, we used a Neumann base reduced by the amalgam method. The computation of the modal model in itself is effortless (less than 7 minutes). However, some computations required by the radiosity method could not be avoided, such as the view factors determination or the inversion of the radiosity matrix. The overall entrance ticket for using a modal model with an accurately modeled radiation might seem expensive, as it requires nine hours of computation. However, this time has to be compared to a finite elements simulation which can demand up to six hours of CPU time.

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To be relevant to other methods, the resulting reduced model has to be used several times and not computed for a one-shot study. The modal reduced model has proven to be robust, as three different configurations were tested, each having different boundary conditions and different radiative solicitations. With a modal model of order 200, the relative maximal error was found to be less

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than 3% while the average error in time and space is below 0.1%. At that order, the computation effort is very low, as only 4 minutes are needed to solve a duration of 50 000 s, compared to the 6 hours required by the finite elements problem. The gain in computation time is of the order of 100 or more, whatever the configuration tested. These very satisfying performances open access to real time control or identification procedure with radiation.

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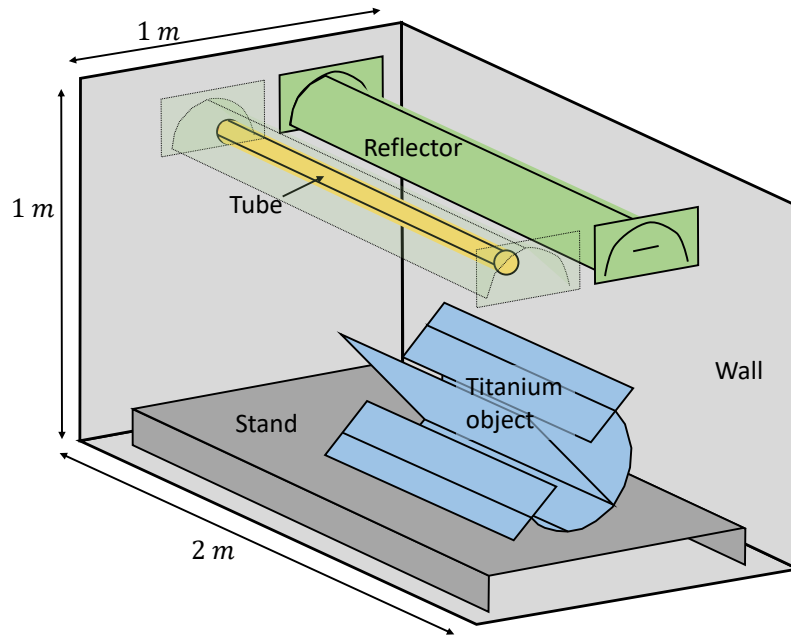


FIGURE 1: The considered geometry

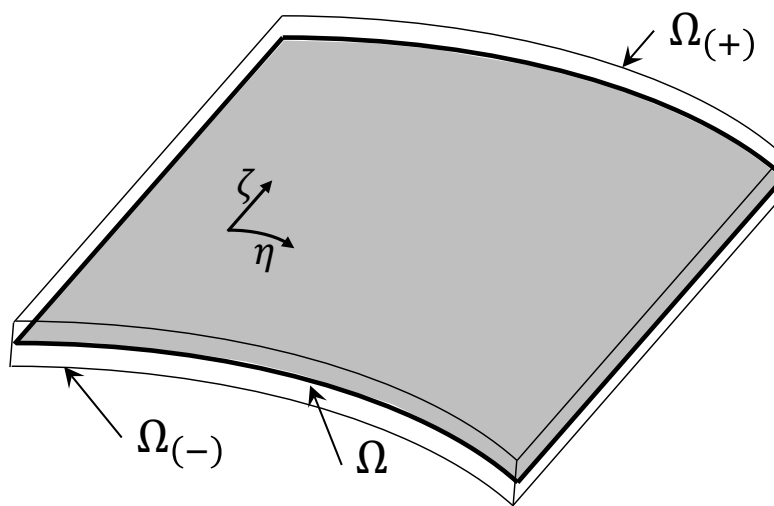


FIGURE 2: Transformation of a 3D domain to a shell configuration

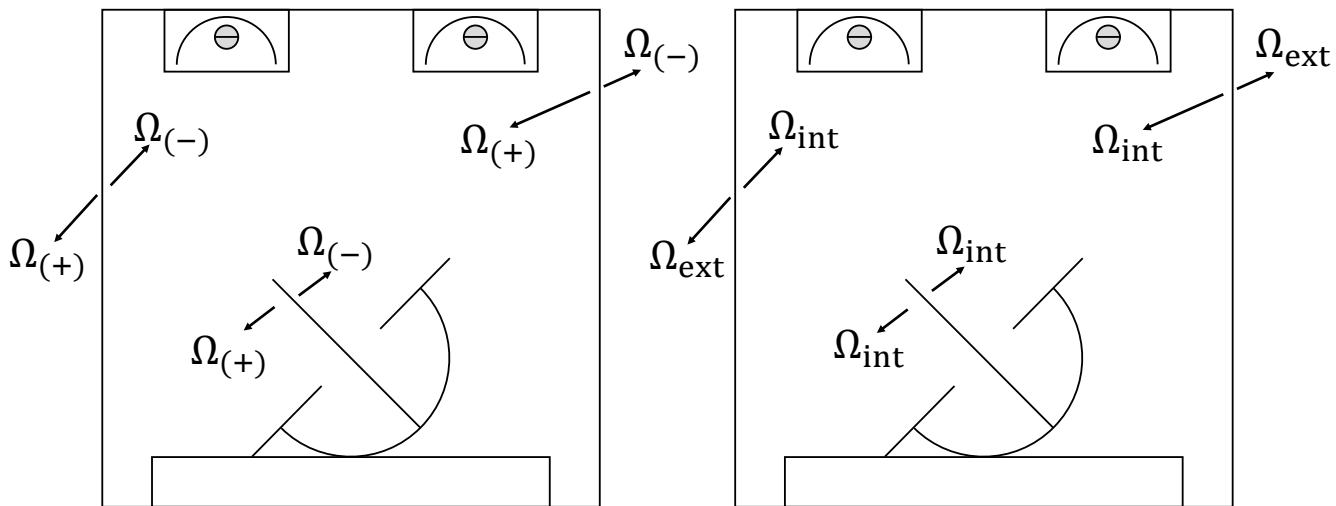


FIGURE 3: Decompositions of boundaries in a shell configuration

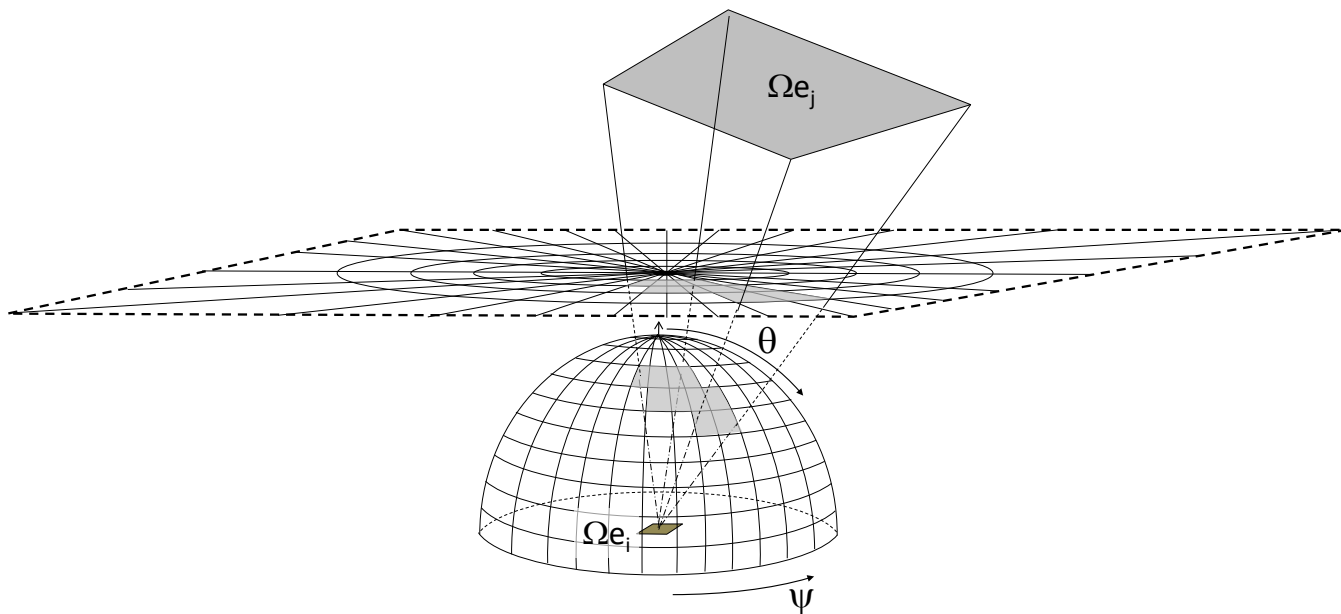


FIGURE 4: Projection using polar-plane-based method

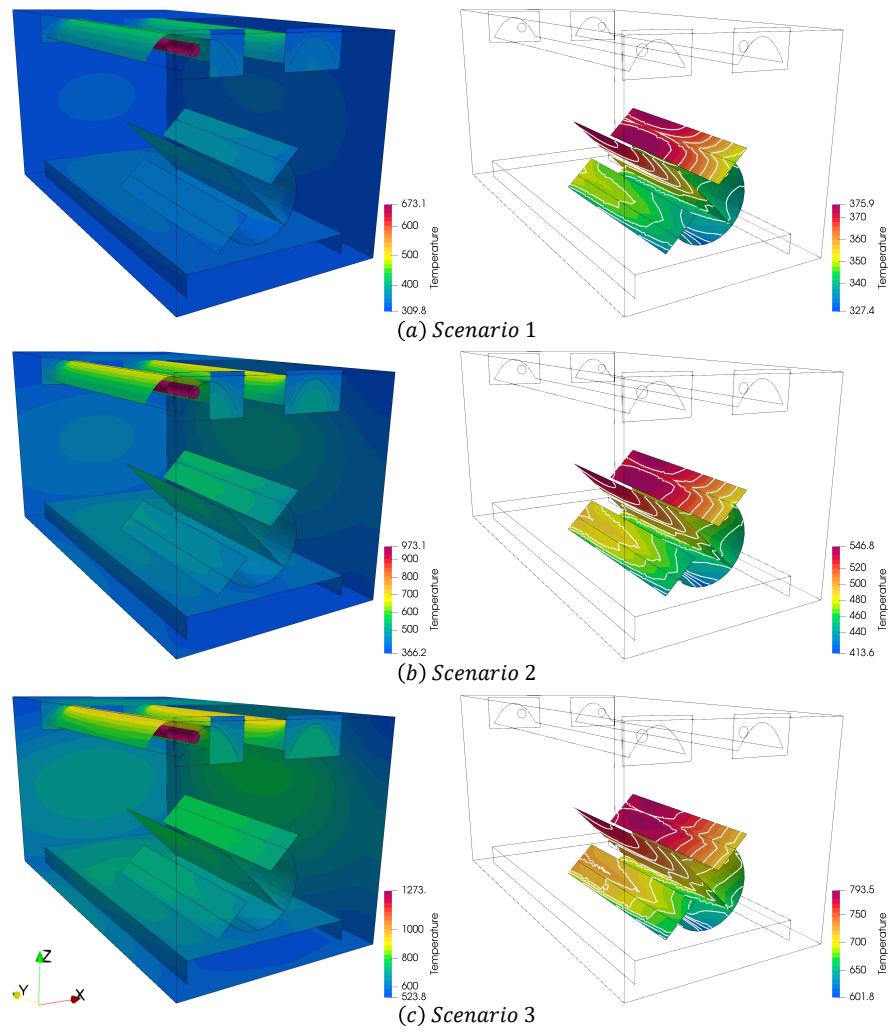


FIGURE 5: Temperature fields for the three scenarios at time 3 000 s. The left column represents the complete thermal scene, whereas the right column focuses on titanium item

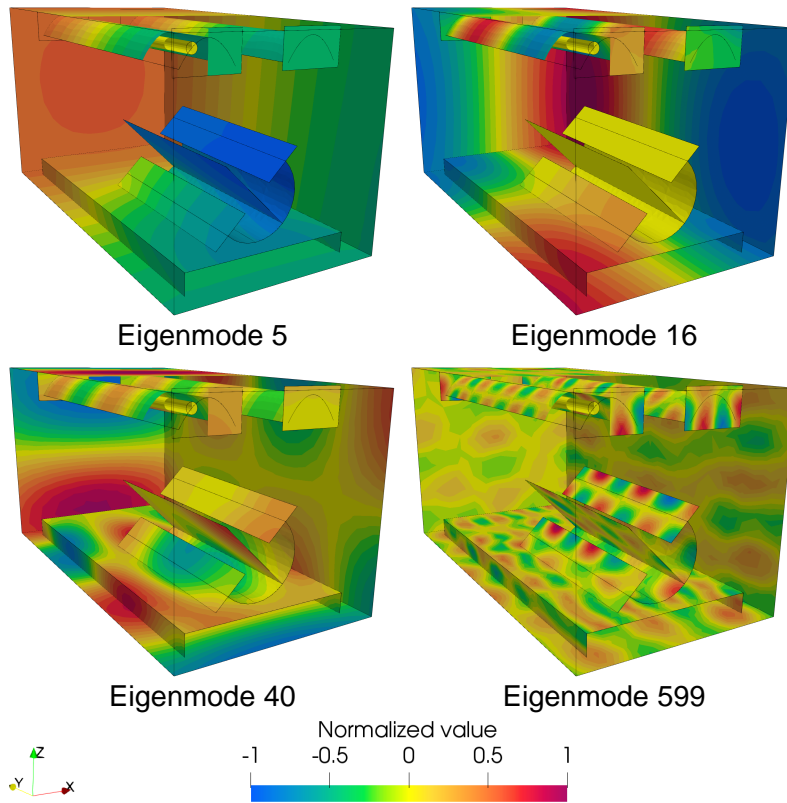


FIGURE 6: Examples of Neumann eigenmodes

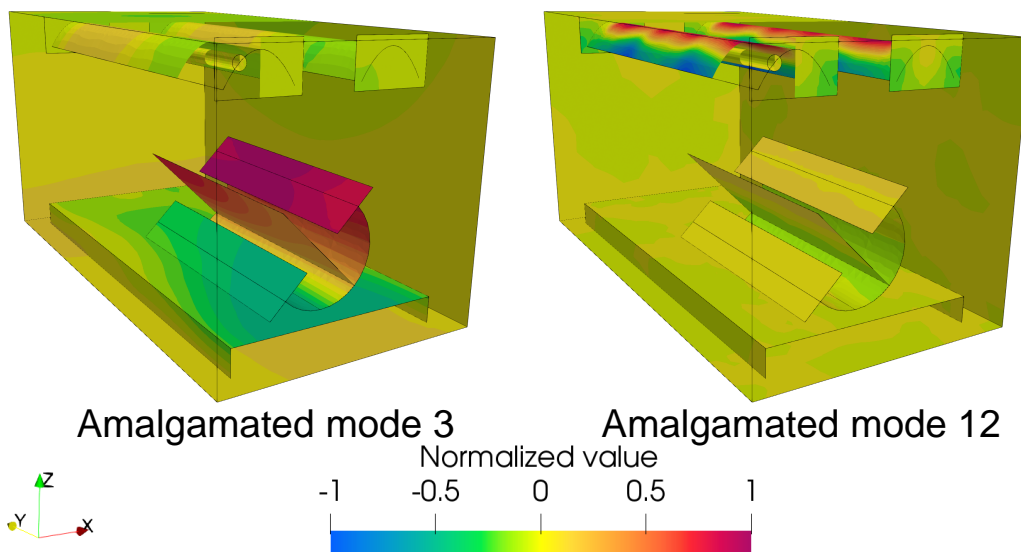


FIGURE 7: Examples of amalgamated modes

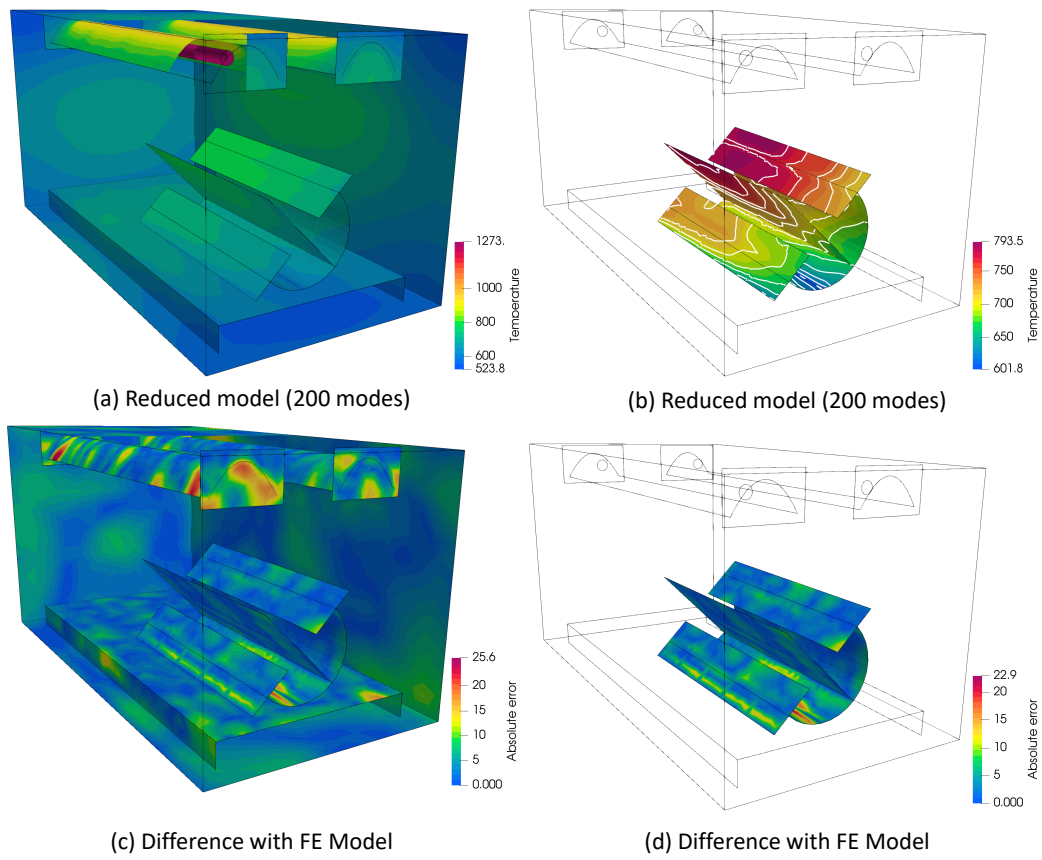
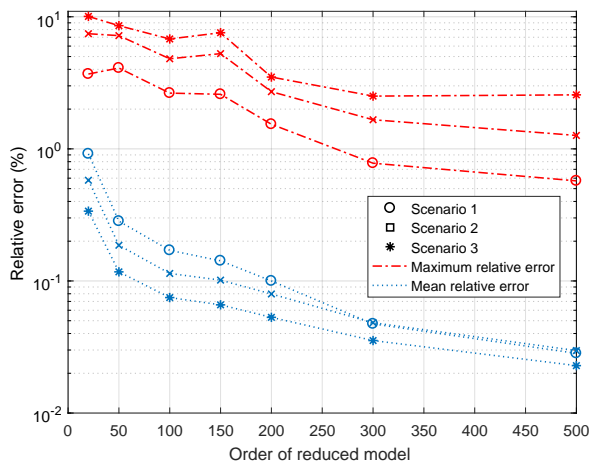
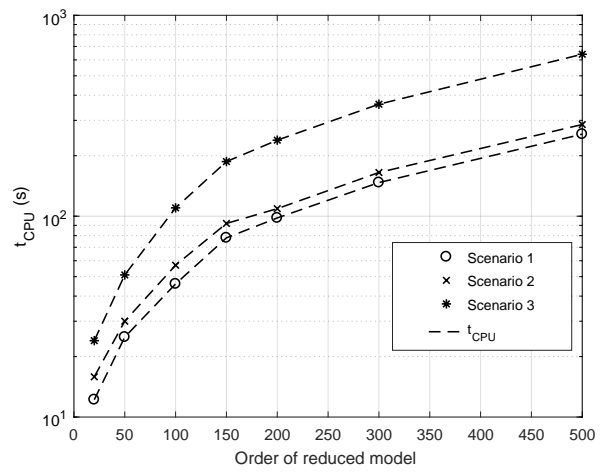


FIGURE 8: Temperature and error fields for scenario 3 at  $t = 3000$  s



(a)



(b)

FIGURE 9: Relative error and CPU time in function of the reduction order