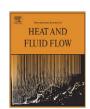
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Development and assessment of a coupled strategy for conjugate heat transfer with Large Eddy Simulation: Application to a cooled turbine blade

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ABSTRACT

Although Large Eddy Simulation (LES) is identified today as the most promising method for turbulent flow problems, few applications of LES coupled to heat transfer solvers in solids have been published. This paper describes a coupling strategy of a LES solver and a heat transfer code within solids on parallel architectures. The numerical methods used in both solvers are briefly recalled before discussing the coupling strategy in terms of physical quantities to exchange (fluxes and temperatures), stability and parallel efficiency. The stability study is performed using an amplification matrix analysis on a one-dimensional case and allows the determination and optimization of coupling parameters. The coupled tool is then applied to a cooled turbine blade model where results demonstrate both the efficiency of the parallel implementation and the quality of the results. Coupled and non-coupled simulations are compared to experimental results and discussed in terms of cooling efficiency and flow structures.

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1. Introduction

Determination of heat loads, as wall temperatures and heat fluxes, is a key issue in combustion (Lakshminarayana, 1996; Lefebvre, 1999; Schiele and Wittig, 2000; Dunn, 2001; Bunker, 2007): the interaction of hot gases and reacting flows with colder walls is an important phenomenon in combustion chambers and a main design constraint in gas turbines. After combustion, the interaction of the hot burnt gases with the high pressure stator and the first turbine blades conditions the temperature and pressure levels reached in the combustor, and therefore the engine efficiency. Numerical simulations of the thermal interaction between fluid flows and solids offer new design paths to diminish development costs through important reductions of the number of experimental tests.

When a Computational Fluid Dynamics (CFD) solver is coupled to a heat transfer solver, the accuracy of the coupled tool is generally controlled by the fluid code. Conventional CFD techniques use Reynolds Averaged Navier–Stokes (RANS) methods. Such approaches cannot capture all complex effects in turbomachinery flows and looking for more precise flow solvers is a usual objective

in this field. Direct Numerical Simulation (DNS) methods cannot offer such an alternative solution because of their computational cost. Recent progresses in Large Eddy Simulation (LES) (Sagaut, 2000; Mahesh et al., 2004; Poinsot et al., 2005) and the continuously increasing computer power offered by the newly developed parallel computer architectures, allow to accurately predict turbulent flows in complex geometries (Acharya et al., 2001; Azzi et al., 2002; Rozati, 2007; Boudier et al., 2007). As LES is still a computationally expensive method, the aim of this paper is to develop and assess a suitable strategy based on LES to efficiently converge to steady thermal states.

There are two basic approaches to solve Conjugate Heat Transfer (CHT) problems. The first one is a direct coupling where the different fields are solved simultaneously in a large system of equations by a monolithic solver (Kao and Liou, 1997; Han et al., 2001; Rahman et al., 2005; Luo and Razinsky, 2007; Ganesan et al., 2007). The second approach consists in solving each set of field equations separately with dedicated solvers that exchange boundary conditions (Heselhaus et al., 1995; Sondak and Dorney, 2000; Papanicolaou et al., 2001; Garg, 2002; Bohn et al., 2005). This solution has the advantage of using existing state-of-the-art codes to solve fluid and solid equations and of being able to exchange one solver with another easily (Alonso et al., 2006). The main drawback of this coupling methodology is that an adapted CHT framework is requested for the simulations especially on parallel machines. The performances of such a coupling framework are linked to (1) the

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strategy to couple the solvers in an accurate and stable fashion as well as (2) the exchange of information between the solvers in an efficient and scalable fashion when using a large number of processors Alonso et al., 2006.

This paper (Alonso et al., 2006) investigates important issues for fully parallel and coupled CHT based on LES for flow models:

- The boundary conditions applied to the fluid and solid codes, including the variables shared by the codes, are critical for the accuracy and stability of the computations,
- In gas turbine applications, the time scales of the flow and of the solid are generally very different. As a consequence, the frequency of exchanges between the codes and the synchronization of the solvers influence the stability as well as the restitution time of the computations.

The document is organized as follows. The fluid and solid solvers are presented in Section 2. Section 3 is dedicated to the CHT framework. The coupling methodology (boundary conditions and synchronization of the solvers) is described in Section 3.1. The stability study (Section 3.2) is performed on a one-dimensional problem using an amplification matrix analysis. The strategy is validated (Section 4) through comparisons of pressure and temperature fields on the experimental film-cooled blade. After a short description of the configuration (Section 4.1) and of solid and fluid discretizations (Section 4.2), an adiabatic LES is discussed (Section 4.3). Thermal results given by the CHT methodology are presented in Section 4.4 which also describes the main flow structures controlling heat transfer. Finally, Section 4.5 proposes efficiency analyses of the coupled strategy.

2. Numerical approach

2.1. Governing equations for flow LES models

LES of reacting flows involves the spatial Favre filtering operation that reduces for spatially, temporally invariant and localised filter functions (Vreman et al., 1994) to:

$$\widetilde{f(\mathbf{x},t)} = \frac{1}{\rho(\mathbf{x},t)} \int_{-\infty}^{+\infty} \rho(\mathbf{x}',t) f(\mathbf{x}',t) G(\mathbf{x}'-\mathbf{x}) d\mathbf{x}'$$
 (1)

where *G* denotes the filter function.

In the mathematical description of compressible turbulent flows the primary variables are the volumic mass fraction $\rho(\mathbf{x},t)$, the velocity vector $u_i(\mathbf{x},t)$ and the total energy $E(\mathbf{x},t) \equiv e_s + 1/2u_iu_i$.

The fluid follows the ideal gas law, $p=\rho rT$ and $e_s=\int_0^T C_p dT-p/\rho$, where e_s is the sensible energy, T the temperature, C_p the fluid heat capacity at constant pressure and r is the mixture gas constant. The LES solver takes into account changes of heat capacity with temperature using tabulated values of heat capacities. The viscous stress tensor and the heat diffusion vector use classical gradient approaches. The fluid viscosity follows Sutherland's law and the heat diffusion coefficient follows Fourier's law. The application of the filtering operation to the instantaneous set of compressible Navier–Stokes transport equations yields the LES transport equations (Poinsot et al., 2005) which contain so-called Sub–Grid Scale (SGS) quantities that need modelling (Sagaut, 2000; Ferziger, 1977). The unresolved SGS stress tensor $\overline{\tau_{ij}}^t$ is modelled using the Boussinesq assumption (Smagorinsky, 1963; Pope, 2000; Chassaing et al., 2000):

$$\overline{\tau_{ij}}^t - \frac{1}{3}\overline{\tau_{kk}}^t \delta_{ij} = -2\bar{\rho}\nu_t \widetilde{S}_{ij} \quad \text{with } \widetilde{S}_{ij} = \frac{1}{2} \left(\frac{\partial \tilde{u}_i}{\partial x_j} + \frac{\partial \tilde{u}_j}{\partial x_i} \right) - \frac{1}{3} \, \frac{\partial \tilde{u}_k}{\partial x_k} \delta_{ij}$$

In Eq. (2), \widetilde{S}_{ij} is the resolved strain rate tensor and v_t is the SGS turbulent viscosity. The Wall Adapting Linear Eddy (WALE) model (Nicoud and Ducros, 1999) is chosen to model the SGS viscosity:

$$v_{t} = (C_{w} \Delta)^{2} \frac{(s_{ij}^{d} s_{ij}^{d})^{3/2}}{(\widetilde{S}_{ij} \widetilde{S}_{ij})^{5/2} + (s_{ij}^{d} s_{ij}^{d})^{5/4}} \quad \text{with } s_{ij}^{d} = \frac{1}{2} (\widetilde{g}_{ij}^{2} + \widetilde{g}_{ji}^{2}) + \frac{1}{3} \widetilde{g}_{kk}^{2} \delta_{ij}$$
(3

In Eq. (3), Δ denotes the filter characteristic length (approximated by the cubic-root of the cell volume), C_w is a model constant equal to 0.5 and $\widetilde{g_{ii}}$ is the resolved velocity gradient.

The SGS energy flux \overline{q}_i^t is modelled using a SGS turbulent heat conductivity obtained from v_t by $\lambda_t = \bar{\rho} v_t C_p / P r_t$, where $P r_t = 0.7$ is a constant turbulent Prandtl number:

$$\overline{q_i}^t = -\lambda_t \frac{\partial \widetilde{T}}{\partial x_i} \tag{4}$$

In Eq. (4), \tilde{T} is the Favre filtered temperature which satisfies the modified filtered state equation $\bar{p}=\bar{p}r\tilde{T}$ (Moin et al., 1991; Erlebacher et al., 1992; Ducros et al., 1996; Comte, 1996). Although the performances of the closures could be improved through the use of a dynamic formulation (Moin et al., 1991; Lilly, 1992; Germano et al., 1992; Ghosal and Moin, 1995; Meneveau et al., 1996) Eqs. (2)–(4) are considered sufficient to address the present preliminary investigation.

The WALE subgrid model is used in conjunction with no-slip wall conditions. This model is designed to provide correct levels of turbulent viscosity down to the wall and no wall model is required. Total pressure and total temperature with velocity angle are imposed using the Navier–Stokes Characteristic Boundary Condition (NSCBC) formalism (Poinsot and Lele, 1992) at the inlets of the fluid domain. Static pressures are enforced at outlet boundaries in characteristic NSCBC form.

2.2. Governing equations for solid heat transfer models

Heat transfer in solid domains is described by the energy conservation:

$$\rho_{s}C_{s}\frac{\partial T(\mathbf{x},t)}{\partial t} = -\frac{\partial q_{i}}{\partial x_{i}}$$

$$\tag{5}$$

where T is the temperature, ρ_s the density, C_s is the heat capacity and q the conduction heat flux. The heat diffusion follows Fourier's law:

$$q_i = -\lambda_s \frac{\partial T}{\partial \mathbf{x}_i} \tag{6}$$

where λ_s is the heat conductivity of the medium. The solid solver takes into account local changes of heat capacity and conductivity with temperature.

2.3. Numerical schemes

(2)

The parallel LES code (Schönfeld et al., 1999; Moureau et al., 2005; Mendez and Nicoud, 2008; Roux et al., 2008) solves the full compressible Navier–Stokes equations using a cell-vertex/finite element approximation and Taylor-Galerkin weighted residual central distribution scheme (Donea and Huerta, 2003). This explicit scheme, which provides third-order accuracy on hybrid meshes, is particularly adequate for low-dissipation requirements of LES applications (Colin and Rudgyard, 2000). Boundary conditions are handled with the NSCBC formulation (Poinsot et al., 2005; Moureau et al., 2005).

The parallel conduction solver is based on the same data structure and uses an explicit scheme for time advancement.

3. Conjugate heat transfer

During this work, only a thermal steady state solution within the blade structure was sought for. Since the two solvers are explicit and time dependent, the coupling strategy between the solvers requires a specific approach discussed in Section 3.1. The stability of the approach is studied in Section 3.2 by determining stability limits in a simplified 1D case.

3.1. Coupling strategies

The resolution of CHT problems on massively parallel machines involves two main issues:

- The boundary conditions applied to the fluid and solid codes control the precision and stability of the computations (Section 3.1.1).
- The frequency of exchanges between the codes and the synchronization of the solvers influence the stability as well as the restitution time of the computations (Section 3.1.2).

3.1.1. Boundary conditions

At a fluid/solid interface Γ_{fs} , heat flux and temperature are continuous:

$$\begin{cases} \phi_f^l = \phi_s^l \\ T_f^l = T_s^l \end{cases}, \quad l \in \Gamma_{fs}$$
 (7)

with (ϕ_f^I, ϕ_s^I) the fluid and solid heat fluxes at the interface point Iand $(T_f^{i,j}, T_s^{i,j})$ the fluid and solid temperatures at the same location. Thus, natural variables to share between the codes are ϕ and T at the interfaces. However, it is usually not efficient to directly impose ϕ or T on either one of the solver. Typical strategies to treat CHT problems (Sondak and Dorney, 2000; Bohn et al., 2005) consist in choosing an adequate composition of variables ϕ_s^I and T_s^I coming from the solid computation as boundary conditions for the fluid domain. The flow solution is computed with this set of boundary conditions during n_f iterations of the fluid solver. Similarly, an adequate composition of variables ϕ_f^l and T_f^l coming from the fluid computation is imposed as boundary conditions for the solid domain during $n_{\rm s}$ iterations of the thermal code. To reach a thermal steady state, these steps are repeated until convergence of boundary variables ϕ^{I} and T^{I} . The solution retained here is to write the continuity of heat flux and temperature across the interface Γ_{fs} (Eq. (7)) in the following form (Chemin et al., 2006):

$$\begin{cases} T_f^I = T_s^I \\ \phi_f^I + hT_f^I = \phi_s^I + hT_s^I \end{cases}$$
 (8)

where *h* is a positive numerical coupling relaxation parameter (CRP) which has the dimensions of a convection coefficient. As recommended by previous studies to improve the stability of the coupled scheme (Giles, 1997; Chemin et al., 2006; Radenac et al., 2006; Roe et al., 2008), the temperature from the solid is imposed to the fluid domain with a Dirichlet condition:

$$T_f^{l,n} = T_s^l, \quad \text{with } n = 1, n_f \tag{9}$$

The resulting mixed boundary condition for the structure is:

$$\phi_s^{l,n} = \phi_f^l + h(T_f^l - T_s^{l,n}), \text{ with } n = 1, n_s$$
 (10)

Both $\phi_s^{l,n}$ and $T_s^{l,n}$ of the mixed formulation (Eq. (10)) converge to steady state between two successive updates of ϕ_f^l and T_f^l . The stability of the mixed condition for the solid domain used in conjunction with Dirichlet boundary temperature for the fluid depends on the value of the CRP h and on update frequency. Section 3.1.2 gives

arguments concerning update frequencies adopted within this study and Section 3.2 proposes a stability analysis of the coupled strategy based on a simplified configuration.

3.1.2. Synchronization of the solvers

In a gas turbine, a blade submitted to the hot stream exiting from a combustion chamber has a characteristic time scale τ_s , based on a reference length and thermal diffusivity, of the order of a few seconds, while the flow-through time along the blade τ_f is less than 1 ms. Hence, the convergence to a thermal steady state of the conjugate system is controlled by the solid.

Let us consider that between two updates of the boundary values, the flow is advanced in time of a quantity $\alpha_f \tau_f$ and the solid is advanced of a time $\alpha_s \tau_s$ where α_f and α_s are two non-dimensional constants. Then, two limit cases are of interest: (1) $\alpha_s = \alpha_f$ ensures that both solid and fluid converge to steady state at the same rate and (2) $\alpha_f \tau_f = \alpha_s \tau_s$ ensures that the two solvers are synchronized in physical time. This last case is needed for fully coupled unsteady computation. As for the present application only the steady state temperature field in the solid is requested, case (1) $\alpha = \alpha_s = \alpha_f$ is used. In the following, α is called the coupling synchronization time parameter (CSTP). The numbers of iterations of the fluid and solid solver, respectively, n_f and n_s , are linked to fluid and solid time steps Δt_f and Δt_s by:

$$n_f = \alpha \tau_f / \Delta t_f$$

$$n_s = \alpha \tau_s / \Delta t_s$$
(11)

On a parallel machine, codes for the fluid and for the structure may be run together or sequentially. Fig. 1 shows how heat fluxes and temperature are exchanged in a mode called Sequential Coupling Strategy (SCS) (Duchaine et al., 2008) or staggered solution procedure (Felippa et al., 2001): at synchronization n_{cvl} of the solvers (the subscript cpl refer to CouPLing quantities), after a physical time $\alpha \tau_f$ the fluid solver provides fluxes and temperatures to the solid solver which then starts and gives back temperatures (physical duration $\alpha \tau_s$). In SCS, the codes are loaded into the parallel machine sequentially and each solver uses all available processors (P). Most existing applications that use SCS are based on steady state codes (i.e. RANS solvers for the fluid) (Divo et al., 2002; Heidmann et al., 2003; Mercier and Tesse, 2006; Verstraete et al., 2007). Another solution is Parallel Coupling Strategy (PCS) (Duchaine et al., 2008), or parallel staggered procedure (Felippa et al., 2001), where both solvers run together using informations obtained from the other solver at the previous coupling iteration (Fig. 2). PCS is employed to study transient or unsteady phenomena (Montenay

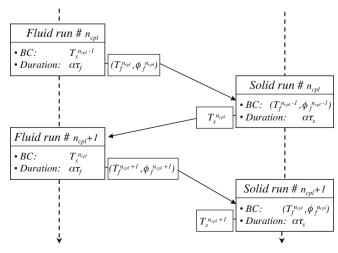


Fig. 1. Sequential coupling strategy SCS.

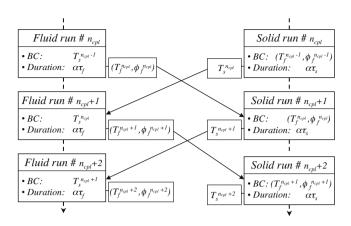


Fig. 2. Parallel coupling strategy PCS.

et al., 2000; Hegab et al., 2001; dos Santos et al., 2008; Duchaine et al., 2008). In the case of PCS, the two solvers must share the $P = P_s + P_f$ processors. To first order, ensuring load balancing requires that the P_s and P_f processors dedicated to the solid and the fluid, respectively, must be such that:

$$\frac{P_f}{P} = \frac{1}{1 + \mathcal{F}_s/\mathcal{F}_f} \tag{12}$$

where \mathcal{T}_s and \mathcal{T}_f are the execution times of the solid and fluid solvers on one processor to compute physical times $\alpha \tau_s$ and $\alpha \tau_f$, respectively. Perfect scaling for both solvers is assumed in Eq. (12). Efficient implementation of PCS requires a software to manage the parallel execution of the solvers as well as the data exchanges during their execution. In order to insure the performance of the PCS, a code coupler, initially developed for ocean–atmosphere coupling (Lagarde et al., 2001; Buis et al., 2005), has been used here.

Both SCS and PCS have been tested: to deal with a fully non stationary coupled problem with synchronization in physical time $(\alpha_f \tau_f = \alpha_s \tau_s)$, PCS is recommended (Hegab et al., 2001). However, when studying the convergence to a steady thermal state where $\alpha = \alpha_f = \alpha_s$, the choice of the method is open and will be addressed in Section 4.5.

3.2. Stability of the coupling methodology

The coupling methodology used to reach a steady thermal state is composed of the following set of boundary conditions:

$$\begin{cases} \text{Fluid (Dirichlet)}: \ T_f^{l,i_f} = T_s^l, & i_f = 1, n_f \\ \text{Solid (mixed)}: \ \phi_s^{l,i_s} = \phi_f^l + h(T_f^l - T_s^{l,i_s}), & i_s = 1, n_s \end{cases}$$
 (13)

Boundary conditions of Eq. (13) are applied to fluid and solid domains during the n_f and n_s iterations of the solvers between two consecutive updates. In Eq. (13), the coupling relaxation parameter h and the coupling synchronization time parameter α are adjustable. The other free parameter is the mode of synchronization: SCS or PCS. Before applying the CHT methodology to a full configuration, it is interesting to study its stability in terms of h and α . This is done here in a simplified 1D case with the PCS.

3.2.1. Description of the simplified conjugate heat transfer problem

In coupled simulations, modes which are normal to the interface between the fluid and the solid can become unstable. Hence, the stability of the PCS with Dirichlet/mixed set of boundary conditions is analyzed here with a 1D purely thermal diffusion problem $\mathcal{P}_{\mathcal{FMF}}$ (Simplified conjugate heat transfer problem). As mentioned by Giles (Giles, 1997), since there is no velocity compo-

nent normal to the solid boundaries, it is appropriate in a 1D model to omit convection terms. Stability for problem $\mathcal{P}_{\mathcal{FCAF}}$ is a necessary condition for real applications but may not be sufficient. Nevertheless, understanding 1D instabilities clearly gives insight into the potential instabilities in 2D and 3D complex computations.

The problem $\mathscr{P}_{\mathscr{FGMF}}$ is a system composed of two domains with constant properties. Domain #1 has the thermal properties (thermal conductivity, heat capacity and density) of plexiglass (Table 1) while domain #2 has the thermal properties of air at 328 K (Table 2). The interface between the domains at x=0 is noted Γ . Dirichlet conditions on temperature are imposed at $x=b_1(b_1<0)$ and $x=b_2(b_2>0)$. To solve the problem $\mathscr{P}_{\mathscr{FGMF}}$, the heat diffusion equation is considered in both domains:

$$\rho_{i}C_{i}\frac{\partial T_{i}(x)}{\partial t} = \lambda_{i}\frac{\partial^{2}T_{i}(x)}{\partial x^{2}}, \quad i = 1, 2$$
(14)

with the boundary and continuity conditions:

$$\begin{cases}
T_{1}(x = b_{1}) = T_{b_{1}} \\
T_{2}(x = b_{2}) = T_{b_{2}} \\
\phi_{1}(x = 0) = \phi_{2}(x = 0) \\
T_{1}(x = 0) = T_{2}(x = 0)
\end{cases}$$
(15)

For the sake of simplicity, Eq. (14) is discretized using a forward Euler time differencing and a classical second-order central difference scheme on uniform grids in each domain:

$$T_{i,i}^{n+1} = T_{i,i}^n + \mathcal{F}_i(T_{i,i+1}^n - 2T_{i,i}^n + T_{i,i-1}^n), \quad i = 1, 2$$
(16)

where $\mathscr{F}_i = \frac{\lambda_i \Delta t_i}{\rho_i C_i \Delta x_i^2}$, Δt_i and Δx_i are the Fourier number, the time step and the mesh size of domain i, respectively. For a periodic domain with uniform thermo-physical properties, the stability of the discretization given by Eq. (16) is ensured when $\mathscr{F}_i \leqslant 0.5$ (Burnett, 1981). In Eq. (16), $T_{i,j}^n$ is given by:

$$T_{i,j}^{n} = \begin{cases} T_{1}(b_{1} + j\Delta x_{1}, n\Delta t_{1}) & j = 1, N_{1} \ n = 1, n_{1} \ (\text{solid}) \\ T_{2}(j\Delta x_{2}, n\Delta t_{2}) & j = 1, N_{2} \ n = 1, n_{2} \ (\text{fluid}) \end{cases}$$
(17)

where N_1 and N_2 are the number of mesh points of domain #1 and #2, respectively. Previous works dealing with numerical stability of coupled thermal computations (Giles, 1997; Chemin et al., 2006; Radenac et al., 2006; Roe et al., 2008) were done with the very restrictive conditions $n_1 = n_2 = 1$ and $\Delta t_1 = \Delta t_2$. In this study, the coupled solvers conserve their own time step based on stand-alone numerical stability. Moreover, the solvers can integrate Eq. (16) with different number iterations n_1 and n_2 between two updates. To mimic the conditions of the blade heat transfer problem, the spatial and time discretizations Δx_i and Δt_i are similar to those of the

Table 1Thermal characteristics of plexiglass.

Thermal conductivity	λ_s	$0.184~W~m^{-1}~K^{-1}$
Heat capacity	C_s	$1450 \mathrm{J kg^{-1} K^{-1}}$
Density	$ ho_s$	$1190 \ kg \ m^{-3}$
Thermal diffusivity	D_s	$1.07\times 10^{-7}\ m^2\ s^{-1}$

Table 2Thermal characteristics of air at 328 K.

Thermal conductivity	λ	$2.6\times 10^{-2}\;W\;m^{-1}\;K^{-1}$
Heat capacity	C_p	$1015 \mathrm{J kg^{-1} K^{-1}}$
Density	ho	$0.266 \ kg \ m^{-3}$
Thermal diffusivity	D_f	$9.63\times 10^{-5}\ m^2\ s^{-1}$

computation with heat conduction solver for domain #1 and the LES code for domain #2.

The discretized heat flux used for boundary conditions are approximated with first-order differences. After the synchronization number n_{cpl} in PCS (Fig. 2), the boundary conditions applied to domain #1 and #2 during n_1 and n_2 iterations, respectively are:

$$\begin{cases}
T_{1,1}^{i_{1},n_{cpl}} = T_{b_{1}} \\
\phi_{1,N_{1}}^{i_{1},n_{cpl}} = \bar{\phi}_{2,1}^{n_{cpl}-1} + h(T_{2,1}^{n_{2},n_{cpl}-1} - T_{1,N_{1}}^{i_{1},n_{cpl}}) \\
T_{2,1}^{i_{2},n_{cpl}} = T_{1,N_{1}}^{n_{1},n_{cpl}-1} \\
T_{2,1}^{i_{2},n_{cpl}} = T_{b_{2}} \\
\text{with } i_{1} = 1, n_{1} \quad \text{and } i_{2} = 1, n_{2}
\end{cases} \tag{18}$$

where $\bar{\phi}_{2,1}^{n_{cpl}-1}$ is the mean value of the heat flux $\phi_{2,1}^{i_2,n_{cpl}-1}$ at the interface in domain #2 over the n_2 iterations. This mean heat flux is introduced in the mixed condition due to the use of a LES solver, intrinsically non stationary, for target applications.

3.2.2. Stability of the solid discretization

Before studying the coupled problem $\mathcal{P}_{\mathcal{FGMF}}$, let us consider only domain #1 to analyze the effect of the mixed boundary condition (Eq. (13)) on stability. Domain #1 is discretized with Eq. (16). The mixed boundary condition is applied at x=0 and a Dirichlet condition on temperature at $x=b_1$:

$$\begin{cases} T_{1,1}^{i_1} = T_{b_1} \\ \phi_{1,N_1}^{i_1} = \phi_b + h(T_b - T_{1,N_1}^{i_1}) \\ \text{with } i_1 = 1, n_1 \end{cases}$$
 (19)

In Eq. (19) ϕ_b , T_b and T_{b_1} are constant heat flux and temperatures. The vector of unknowns U_1 is introduced as:

$$U_1 = [T_{1,1} \quad \cdots \quad T_{1,N_1} \quad T_b \quad \phi_b]^t \tag{20}$$

The constants T_b and ϕ_b are introduced in the vector U_1 at this level of the study in order to prepare the notations used during the analysis of the coupled problem $\mathscr{P}_{\mathscr{SCMF}}$. The solutions at time $i_1\Delta t_1$ and $(i_1+1)\Delta t_1$ are linked by:

$$U_1^{i_1+1} = M_1(\mathcal{F}_1, h)U_1^{i_1} \tag{21}$$

In Eq. (21), $M_1(\mathcal{F}_1, h)$ is the discretization operator of Eq. (16) augmented by boundary conditions Eq. (19):

$$M_{1}(\mathcal{F}_{1},h) = \begin{bmatrix} 1 \\ \mathcal{F}_{1} & 1 - 2\mathcal{F}_{1} & \mathcal{F}_{1} \\ & \ddots & \\ & \mathcal{F}_{1} & 1 - \mathcal{F}_{1}(1+\mathcal{D}) & \mathcal{F}_{1}\mathcal{D} & \frac{\mathcal{F}_{1}\mathcal{D}}{h} \\ & & 1 \end{bmatrix}$$

$$(22)$$

where \mathscr{D} is a Biot like number based on the mesh size in domain Δx_1 , the conductivity λ_1 and the coupling relaxation parameter h used in Eq. (8):

$$\mathscr{D} = \frac{h\Delta x_1}{\lambda_1} \tag{23}$$

The value of the spectral radius $\rho(M_1(\mathscr{F}_1,h))$ of the operator $M_1(\mathscr{F}_1,h)$ determines the stability of the discretization (Hirsch, 1988; Allaire and Craig, 2007). Indeed, a requirement for a stable numerical scheme is:

$$\rho(M_1(\mathscr{F}_1, h)) \leqslant 1 \tag{24}$$

When the dimension of matrix $M_1(\mathcal{F}_1,h)$ is sufficiently small (i.e. when the number of discretization nodes N_1 is small), the eigen values of M_1 can be expressed analytically. For $N_1 = 4$, the condition of stability (Eq. (24)) gives a bounded region of stability for \mathcal{D} :

$$\mathscr{D} \leqslant \frac{\mathscr{F}_{1}^{3} - 12\mathscr{F}_{1}^{2} + 20\mathscr{F}_{1} - 8}{-3\mathscr{F}_{1}^{3} + 8\mathscr{F}_{1}^{2} - 4\mathscr{F}_{1}} \tag{25}$$

The stability limit of \mathcal{D} is a function of the Fourier number \mathcal{F}_1 : above $\mathcal{F}_1 = 0.5$, the scheme is always unstable as expected for an explicit formulation. When \mathcal{F}_1 goes to 0 (Δt_1 tends to 0), the upper limit \mathcal{D}_u of \mathcal{D} is ∞ : the scheme becomes unconditionally stable. On the interval $\mathscr{F}_1 \in]0,0.5]$, the derivative $\partial \mathscr{D}_u/\partial \mathscr{F}_1$ is negative. The use of larger time steps (increasing \mathcal{F}_1) reduce the range of \mathcal{D} leading a stable scheme. Decreasing @ also has a positive effect on convergence (see Section 4.5) so that the optimum point for CHT in this case is $\mathcal{F}_1 = 0.5$ and $\mathcal{D} = 2.33$. Fig. 3 compares the analytical evolution of \mathcal{D}_u given by Eq. (25) for $\mathcal{F}_1 \in]0,0.5]$ and numerical solutions obtained with a discrete procedure. The discrete procedure is validated with the case $N_1 = 4$ (Fig. 3) and used to measure the sensibility of the stability domain to N_1 . Results show that the curve $\mathcal{D}_u(\mathcal{F}_1)$ weakly depends on the number of nodes N_1 : the case $N_1 = 51$ leads to the same stability limit as $N_1 = 4$. When the number of nodes N_1 increases, the optimal point for convergence speed goes to $\mathcal{F}_1 = 0.5$ and $\mathcal{D} \approx 2$.

3.2.3. Stability analysis of the coupled discretizations

Now that the stability of domain #1 with a mixed boundary condition has been studied, the stability of the coupled problem $\mathscr{P}_{\mathscr{T}\mathscr{K}\mathscr{T}}$ is investigated. Between the synchronizations n_{cpl} and $n_{cpl}+1$, solutions U_1 and U_2 of domains #1 and #2 are advanced in time using discretization operators $M_1(\mathscr{F}_1,h)$ and $M_2(\mathscr{F}_2)$ similar to the definition given in Eq. (22):

$$\begin{cases}
U_1^{n_{cpl}+1} = \left(\prod_{n=1}^{n_1} M_1\right) U_1^{n_{cpl}} = M_1^{n_1} U_1^{n_{cpl}} \\
U_2^{n_{cpl}+1} = \left(\prod_{n=1}^{n_2} M_2\right) U_2^{n_{cpl}} = M_2^{n_2} U_2^{n_{cpl}}
\end{cases}$$
(26)

where the numbers of iterations n_1 and n_2 between two synchronization events are linked to time-steps of domains #1 and #2:

$$n_1 = \alpha \tau_1 / \Delta t_1
 n_2 = \alpha \tau_2 / \Delta t_2$$
(27)

The characteristic time scales τ_1 and τ_2 in Eq. (27) are chosen in order to mimic the behavior of CHT of the turbine vane given in Table 5. Thus, $\tau_1 = \frac{\rho_1 c_1 b_1^2}{\lambda_1}$ is a diffusion time in domain #1 while τ_2 is a convection time taken in the condition of the blade simulations. For the same reason, the time-step Δt_1 is obtained with a Fourier

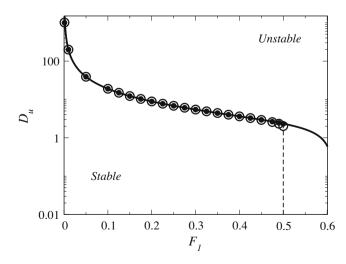


Fig. 3. Upper limit of stability \mathscr{D}_u for the discretisation Eq. (22) as a function of Fourier Number \mathscr{F}_1 : —, analytical solution for $N_1=4$; \bigcirc , computed numerically with $N_1=4$; \bigcirc , computed numerically with $N_1=51$.

number \mathscr{F}_1 equal to 0.5 while Δt_2 is of the order of the time-step in LES computations, viz. $\mathscr{F}_2 << 0.5$ (the CFL stability limit associated to acoustic waves is usually much more restrictive than diffusion in compressible LES).

Based on Eq. (26), the discretization of the whole system can be expressed as:

$$U^{n_{cpl}+1} = \mathcal{M}(\alpha, h)U^{n_{cpl}}$$
(28)

where U is a composition of the nodal temperatures in the domains and heat fluxes at the interface. The value of the spectral radius $\rho(\mathcal{M}(\alpha,h))$ of the operator $\mathcal{M}(\alpha,h)$ determines the stability of the discretization (Hirsch, 1988; Allaire and Craig, 2007). Fig. 4 gives the stability map of the operator $\mathcal{M}(\alpha,h)$ for α in [0.001,3] and $\mathcal{D}=h\Delta x_1/\lambda_1$ from 0 (Neumann boundary condition) to 2.3. Above a threshold $\mathcal{D}=\mathcal{D}_m$, the mixed condition is stable. As shown on Fig. 4, \mathcal{D}_m is a function of α : if the CSTP is sufficiently small $(\alpha<0.02)$, the Neumann condition is stable. Then, until $\alpha=0.5$, the value of \mathcal{D} that ensures numerical stability increases with α . Beyond this point, \mathcal{D}_m remains approximatively constant when the CSTP increases. The stability domain is also limited by an upper bound \mathcal{D}_M (Fig. 4). The limit of \mathcal{D}_M when α increases is $\mathcal{D}_u\approx 2$ obtained during the stability analysis of domain #1 alone.

As a summary, the stability analyses performed for a simple 1D problem have pointed out a stability domain in terms of (α, \mathcal{D}) where α stands for the frequency between two successive updates of the coupled boundary conditions and $\mathcal{D} = h\Delta x_1/\lambda_1$ where h is a coupling relaxation factor defined in the mixed boundary condition (Eq. (13)). For a given value of the coupling synchronization time parameter α , a range of h allows numerical stability of the coupled scheme. This range of CRP is bounded by an upper limit determined by the solid discretization whereas the lower bound is linked to the coupling scheme.

An important information not accessible from these analyses still misses: since the aim of the film-cooled turbine vane simulations is to find a stationary thermal state of the blade in a reduced computational time, it is crucial to understand the effect of the CSTP α and of the CRP h on convergence. Section 4.5 presents sensitivity analyses obtained on the blade configuration when varying α and h.

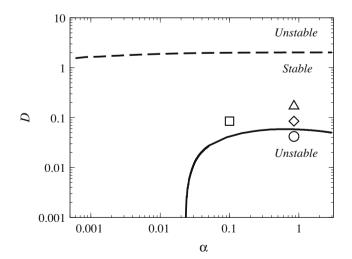


Fig. 4. Stability map in the (α, \mathscr{D}) space of the simplified conjugate heat transfer problem $\mathscr{P}_{\mathscr{P}\mathscr{C}\mathscr{X}\mathscr{F}}$ (Eq. (16) and (18)): —, lower stability limit \mathscr{D}_m : — upper stability limit \mathscr{D}_m . α is the coupling synchronization time parameter (Eq. (27)) and \mathscr{D} is a function of the coupling relaxation parameter h (Eq. (23)). Positions of some computations done on T120D blade (Section 4): point (α_1, h_2) \square ; point (α_2, h_1) \bigcirc ; point (α_2, h_2) \Diamond and point (α_2, h_3) \triangle . The parameters (α_1, h_1) are given in Table 6.

4. Application of the coupled strategy to the T120D blade

This section presents the application of the conjugate heat transfer (CHT) strategy to the T120D blade. After describing the configuration, an adiabatic computation is analyzed. It show the potential of the LES to predict accurately complex turbomachinery flows. Then, CHT results are presented. Finally the performance of the CHT methodology is discussed.

4.1. Description of the T120D blade

The T120 cascades were designed by Rolls Royce Deutschland for the European project AITEB (Haselbach and Schiffer, 2007) (Aerothermal Investigations on Turbine Endwalls and Blades). The experiments were conducted in the High-speed Cascade Wind Tunnel of the Institute of Jet Propulsion (Sturm et al., 1985; Homeier et al., 2005; Gomes et al., 2009).

The highly-loaded high-pressure turbine airfoil of the T120 cascades was designed in order to have a large separation on the pressure side. Therefore the profile presents a strong concave curvature on the pressure side starting shortly after the leading edge. Together with the high pitch to chord ratio relatively strong adverse pressure gradients are present on the pressure side resulting in flow separation. Typically for high-pressure turbine blades, a high change of direction is imposed on the flow. The blade is designed for high exit Mach numbers. The aerodynamical and geometrical data at design conditions are listed in Table 3.

The film cooling device of the T120D blade is composed of three jet rows located on the pressure side in the region of the main flow separation (Fig. 5). The first row of jets is placed near the stagnation point and has cylindrical holes with compound angle against the main stream. This first row is located to protect the leading edge region which has the highest heat transfer rate over the entire airfoil (Han et al., 2000). The second jet row contains fan-shaped holes with zero compound angle located at approximately 20% of the axial chord length. The second row is supposed to re-energize the boundary layer of the main stream and therefore reduce the flow separation. A third row of cylindrical holes is placed at approximately 35% of the axial chord. This last row slightly diminishes the separation bubble on the pressure side. The temperature difference between the mainstream $(T_1^t = 333.15 \text{ K})$ and cooling $(T_c^t = 303.15 \text{ K})$ flows is limited to 30 K to facilitate measurements. The blade is made of plexiglass with a low conductivity of $0.184\,W\,m^{-1}\,K^{-1}$ that make the CHT problem difficult to treat. Experimental results include pressure data on the blade suction and pressure sides as well as temperature measurement on the pressure side (Gomes et al., 2009).

4.2. Computational domains

The computational domains for the fluid and the structure contain one spanwise pitch of the film cooling hole pattern corre-

Table 3 Experimental settings for T120D blade.

Blade height to chord ratio		1.5
Pitch to chord ratio		1.007
Turning	$\Delta \beta$	120°
Inlet angle	β_1	138°
Inlet Mach number	Ma_1	0.295
Outlet Mach number	Ma_2	0.87
Outlet Reynolds number	Re_2	390,000
Inlet total pressure	P_1^t	26737.0 Pa
Inlet total temperature	T_1^{t}	333.15 K
Cooling total pressure	P_c^t	29143.3 Pa
Cooling total temperature	T_c^t	303.15 K
Outlet static pressure	P_2	16325 Pa

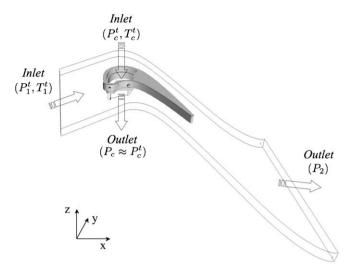


Fig. 5. Fluid computational domain.

sponding to 8 mm along the z-axis (Fig. 5), with periodicity enforced on each side. This simplification neglects end-wall effects but retains the three-dimensionality of the flow and greatly reduces the number of tetrahedral cells required to model the blade: about 6.5 million cells for the fluid and 600,000 for the solid. For the fluid region, periodicity condition is also assumed in the ydirection. Specific care is devoted to the tetrahedral cell isotropy in the wall regions; the maximum values of grid spacings on the blade surface expressed in wall units are about $dx^+ \approx dy^+ \approx$ $dz^+ \approx 40$. In the dilution zone, grid spacings are smaller than 5 wall units. Such resolutions are fine enough to capture major aerodynamic events at the wall reasonably well. The size of one spanwise pitch of hole pattern allows these grid spacings with a moderate number of cells. As shown in Fig. 5, the three film-cooling holes and the plenum used to inject the cooling air are also included in the fluid domain. Hence, the distribution of mass flow rate within the holes is a result of the simulations. The skin meshes are the same for the fluid and the solid so that no interpolation error is introduced at this level when CHT is simulated. The computational domains for both fluid and solid parts are considered fine enough to assess the coupling methodology in a reasonable time.

4.3. Adiabatic computation of the T120D blade

The description of the LES results with adiabatic wall conditions justifies the use of LES for CHT. The main flow structures are first emphasized before comparing computational results and experimental data.

Fig. 6 depicts an instantaneous snapshot of vorticity (left) and a field of mixture fraction showing the path of cooling air in the main stream (right). The LES predicts an intense turbulence intensity and mixing in the region of the three jets. Downstream from the jets, the strong acceleration on the pressure side relaminarizes the flow and forces the cooling air against the blade surface. At the beginning of the suction side, the boundary layer is rather laminar. Then, the flow accelerates up to supersonic velocities. A weak shock appears at a reduced abscissa of about 0.75 and destabilizes the boundary layer. Vortex shedding develops behind the blade.

Fig. 7 shows that jet 1, directly exposed to the main stream, mixes rapidly with hot gases. Protected by the concave shape of the blade and by the first jet, the cooling air of the second hole penetrates more into the flow until it mixes with the third jet. Jet 3 is aligned with the main stream and remains coherent until it impacts the blade in a region between reduced abscissa of 0.5–0.6.

The adiabatic LES results are compared to the experiment in terms of pressure profiles on the blade (Fig. 8) and of temperature profiles on the pressure side (Fig. 9). Pressure fields are displayed in terms of isentropic Mach numbers $M_{is}(\mathbf{x})$ computed by:

$$Ma_{is}(\mathbf{x}) = \sqrt{\frac{2}{\gamma - 1} \left[\left(\frac{P_1^t}{P_w^t(\mathbf{x})} \right)^{\frac{\gamma - 1}{\gamma}} - 1 \right]}$$
 (29)

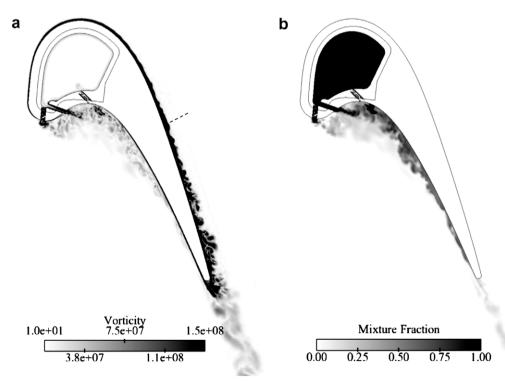


Fig. 6. Instantaneous snapshot of (a) vorticity and (b) distribution of cooling air within a cutting plane at constant *z* passing thought jet 2. The dashed line on (a) represents the approximate position of the shock at 75% of the axial chord.

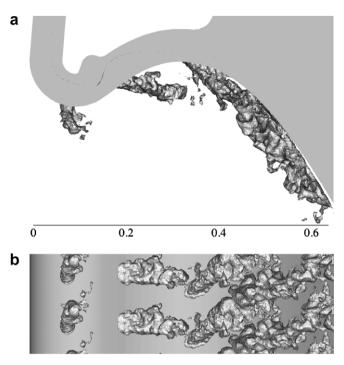


Fig. 7. Instantaneous isosurface of temperature T = 318 K: (a) z view and (b) y view

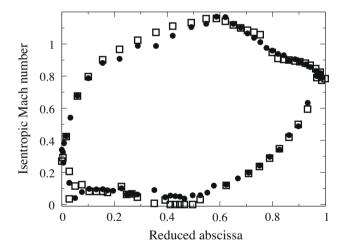


Fig. 8. Time-averaged isentropic Mach number (Eq. (29)) along the blade at pressure probe locations as a function of reduced abscissa: \bullet , adiabatic LES; \Box , experiment.

where γ is the isentropic coefficient, P_1^t and $P_w^t(\mathbf{x})$ are the total pressure of the mainstream and at the wall at location \mathbf{x} . Fig. 8 displays a time-averaged distribution of isentropic Mach number obtained by LES and experimentally at static pressure tap locations. The flow strongly accelerates on the suction side up to supersonic conditions $(Ma_{is}(\mathbf{x}) > 1)$ and features a weak shock at about 75% of the axial chord. On the pressure side, the flow is exposed to strong adverse pressure gradients shortly downstream of the stagnation point. The flow separates at a reduced abscissa of about 0.2 (position of the second jet) reattaching at about 0.45, after the third jet. After reattachment, the flow is strongly accelerated leading to values for the acceleration parameter that are beyond common relaminarization limits (Mayle, 1991). Although the shock position on the

suction side is not perfectly captured, the overall agreement between LES and experimental results is fair.

Table 4 compares numerical and experimental values of the total mass flow rate of air $Q=\sum_j Q_j$ passes through the three holes as well as of the blowing ratios M_i of the jets. Proper descriptions of the mass flow rate and blowing ratio of each jet are crucial to predict heat transfer with a good accuracy. LES gives a good approximation of the total mass flow rate across the holes. The numerical simulation also gives the distribution of air among the three jets, not available experimentally. Since the total coolant pressure is the same for all three coolant rows, the mass flow rate is almost equally distributed between the holes. Blowing ratios are expressed in terms of velocity U and density ρ as:

$$M_j = \frac{\rho_j U_j}{\rho_{\infty} U_{\infty}}, \quad j = 1, 2, 3$$
 (30)

where j refers to jet quantities and ∞ to local main stream values. Equal coolant mass flow $Q_j = \rho_j U_j$ per hole is assumed for the determination of experimental blowing ratios of Table 4 and the LES validates this hypothesis. The agreement between experimental and numerical blowing ratios is rather good (Table 4).

Wall temperatures $T_w^t(\mathbf{x})$ are presented in Fig. 9 in terms of cooling efficiency defined by:

$$\Theta(\mathbf{x}) = \frac{T_1^t - T_w^t(\mathbf{x})}{T_1^t - T_c^t} \tag{31}$$

where T_1^t and T_c^t are the total temperatures at the inlets of the mainstream and plenum, respectively. Time and spanwise averaged cooling efficiencies for adiabatic LES and RANS computations as well as experimental results are plotted as function of relative axial chord on Fig. 9. The RANS computation is done with Fluent using a $k - \omega/SST$ turbulence model (Menter et al., 1993). As expected, in the region of the jets (reduced abscissa up to 0.45) the cooling efficiencies obtained with the adiabatic simulations are lower than the experimental values: adiabatic temperature fields over-predict the real one. Downstream of the impact of the jets on the blade, the adiabatic LES fits the experimental level of Θ whereas the RANS computation over-estimates it. In the experience, the film of colder air that forms after the interaction between the jets and the surface of the blade maintains the wall temperature close to adiabatic one. Hence, the LES captures fairly well the air mass flow through the jets as well as the mixing of the cooling air with the main stream. That is not the case for the RANS simulation. Indeed, even if the RANS computation reproduces the real air mass flow rates ejected by the holes, the simulation does not describe mixing correctly. As a result, the jets remain coherent on a too long distance without mixing with the hot stream and too much cold air impacts the blade surface.

Both LES and RANS simulations exhibit a non-physical peak of Θ near the trailing edge. This peak is due to an over-expansion near the trailing edge which does not appear in the experiment. The round trailing edge of the T120D blade profile and a lack of resolution in this region where there is a very strong acceleration of the flow cause this difficulty in the computation (Denton et al., 1999; Mei et al., 2005).

From these results, LES appears to be a good candidate to treat CHT problems in complex turbomachinery flows.

4.4. Coupled computation of the T120D blade

This sub-section presents a fully coupled simulation of the T120D blade obtained with a two-step methodology:

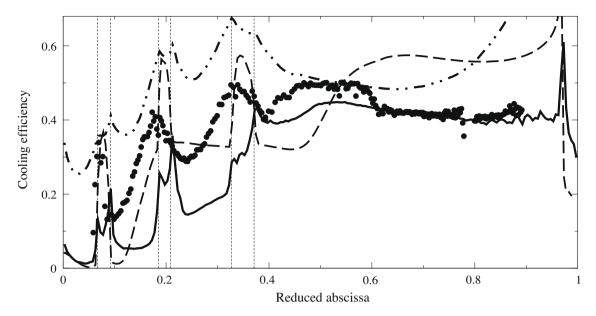


Fig. 9. Time and spanwise averaged cooling efficiency (Eq. (31)) versus abscissa on the pressure side as a function reduced abscissa: ●, experiment from UNIBW; — adiabatic LES: — adiabatic RANS: — · · — coupled LES.

Table 4 Mass flow rate of air through the three jets Q_i and blowing ratios (Eq. (30)) of each jet M_i of the T120D blade.

	$Q_1 \ (g \ s^{-1})$	$Q_2 \ (g \ s^{-1})$	$Q_3 \ (g \ s^{-1})$	$Q = \sum_j Q_j \left(g s^{-1} ight)$	M_1	M_2	M_3
Experiment	-	-	-	0.250	2.03	2.33	3.78
LES	0.083	0.090	0.075	0.248	2.1	2.3	3.0

- 1. Initialization of the coupled calculation that includes:
 - a converged adiabatic fluid simulation (presented in Section 4.3).
 - a converged solid computation with imposed boundary temperatures given by the adiabatic fluid solution,
- Coupled simulation with a parallel coupling strategy (PCS on Fig. 2).

The maximum value of the Biot mesh number $\mathcal{D}=h\Delta x_s/\lambda_s$ that gives a stable mixed condition for the solid discretization of the blade (Eq. (13)) with a Fourier number of 0.5 is $\mathcal{D}_u\approx 1.8$. The solid mesh size Δx_s used to compute \mathcal{D}_u corresponds to the discretization size at the fluid/solid interface (i.e. $\min\Delta x$ for the solid in Table 5). The conductivity of the solid media is given in Table 1. It is interesting to note that the value $\mathcal{D}_u\approx 2$ highlighted in Section 3.2 for 1D simple problem is a good approximation for the complex 3D problem. As a result, the values of the coupling synchronization time parameter α and of the coupling relaxation parameter h used for the CHT simulation are chosen so that the point (α, h) belongs to the stability domain of the simple 1D problem $\mathcal{P}_{\mathcal{PCMT}}$ (Fig. 4) presented in Section 3. These coupling parameters correspond to (α_1, h_2) in Table 6.

Table 5Minimum length, time-steps and characteristic times of fluid and solid discretization for the T120D blade.

	Fluid	Solid
$\min \Delta x$	$1.23 \times 10^{-4} \text{ m}$	$1.41 \times 10^{-4} \text{ m}$
Δt	$9.65\times10^{-8}\;\text{s}$	$1.74\times10^{-2}\;\text{s}$
τ	$1.2\times10^{-1}\;\text{s}$	233.6 s

The coupled simulation was performed on 32 processors of an IBM JS21. The synchronization in CPU time (Eq. (12)) was reached with 30 processors allocated to the fluid solver and two for the solid one. The converged state is obtained in 10 characteristic solid time scales $\tau_{\rm s}$ (Fig. 11) and requires about 4800 CPU h. As a previous study on this configuration has shown that the pressure distribution over the blade is not affected by the coupling (Duchaine et al., 2008), the analysis focuses on thermal aspects.

At the converged state, the net heat flux through the blade reaches zero (i.e. the mean temperature of the blade is stabilized, Fig. 11). Fig. 9 shows measurements, adiabatic and coupled results of the cooling efficiency $\Theta(x)$ spanwise and time averaged along pressure side. As mentioned previously, the adiabatic temperature field (solid line) over-predicts the real one. The main contribution of conduction throughout the blade is to reduce the wall temperature on the pressure side and thus to increase $\Theta(x)$ (dot dashed line). The global form of the reduced temperature from the coupled simulation matches the experimental data correctly. Differences in the levels are explained by the insufficient wall resolution used for the LES computation as well as to the experimental difficulties and uncertainties for temperature measurements and processing (i.e. spanwise average). The strong acceleration caused by the blade induces large thermal gradients not-well resolved by the simulation that leads to an underestimation of the thermal fluxes as well as to non-physical values of cooling efficiency at the trailing edge.

Table 6Values of the coupling synchronization time parameter and coupling relaxation parameter used in Section 4.

$\alpha_1 = 0.1$	$\alpha_2 = 0.85$	$\alpha_3 = 5.00$
$h_1 = 50$	$h_2 = 100$	$h_3 = 200$

Fig. 10 compares experimental and numerical cooling efficiency fields on a 2D plot of the pressure side. The computation matches the experimental visualization. Fig. 10 evidences the thermal effects of the cooling jets on the vane. Jet 1 is folded back against the surface by the main stream but detaches rapidly and mixes with the hot gases due to the curvature of the blade. Downstream of the second hole, a streak with higher efficiency and a spot with enhanced cooling close to the ejection location indicate a partially attached jet. A streak with a lower surface temperature is also visible downstream of the third jet. Jet 3 seems to be the most active in the cooling process: it protects the blade from the hot stream until a reduced abscissa of 0.5 and it impacts the vane at reduced abscissa between 0.5 and 0.6. The curvature of the pressure side induces a slack film coverage: the concave shape of the surface spreads the cooling air laterally along the spanwise direction, as explained by Schwarz and Goldstein (Schwarz and Goldstein. 1989).

4.5. Efficiency analyses of the coupled strategy

This last sub-section analyzes the efficiency of the coupling methodology with regard to the adjustable parameters used in the boundary conditions, the coupling synchronization time parameter α and the coupling relaxation parameter h:

$$\begin{cases} \text{Fluid (Dirichlet)}: \ T_f^{l,i_f} = T_s^l, & i_f = 1, n_f \\ \text{Solid (mixed)}: \ \phi_s^{l,i_s} = \phi_f^l + h(T_f^l - T_s^{l,i_s}), & i_s = 1, n_s \end{cases}$$
 (32)

The number of iterations of fluid (n_f) and solid (n_s) solvers between two consecutive updates of boundary conditions are:

$$n_f = \alpha \tau_f / \Delta t_f$$

$$n_s = \alpha \tau_s / \Delta t_s$$
(33)

In Eq. (33), Δt_f and Δt_s are the time-steps of the fluid and solid solvers, respectively. τ_f and τ_s are characteristic times of the flow and solid phenomena. The different values used for the CSTP α_k and CRP h_k in this section are summarized in Table 6. As the implementation of SCS is simpler than PCS, SCS is preferred for efficiency analyses. As the formulations for SCS and PCS are different, it is useful to check whether the implementations of these strategies give the same results: Fig. 11 compares $SCS(\alpha_1, h_2)$ and $PCS(\alpha_1, h_2)$ and shows that both strategies converge to the same thermal state at the same rate. As a reference for computational

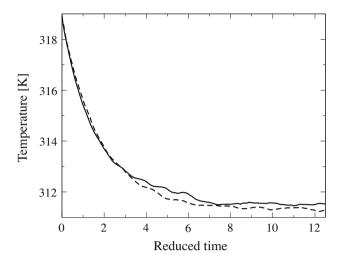


Fig. 11. Evolutions of the mean temperature in the blade as a function of reduced solid time t_s/τ_s for $SCS(\alpha_1, h_2)$ — and $PCS(\alpha_1, h_2)$ — . The parameters (α_1, h_2) are given in Table 6. As a reference for restitution times, one solid characteristic time τ_s requires almost a total of 295 CPU h for the LES and the heat conduction solvers on a SGI Altix ICF8200

costs and restitution times, one solid characteristic time τ_s requires almost a total of 295 CPU h for the LES and the heat conduction solvers on a SGI Altix ICE8200.

To evaluate the impact of h in Eq. (32) on restitution times for the T120D CHT problem, three computations with SCS are compared: $SCS(\alpha_2, h_1)$, $SCS(\alpha_2, h_2)$ and $SCS(\alpha_2, h_3)$ (Table 6). Fig. 12a depicts the evolutions of the mean temperature in the blade as a function of reduced solid time t_s/τ_s for these simulations. The retained values h_1 , h_2 and h_3 with α_2 give stable coupled schemes that converge to the same thermal steady state. Fig. 4 shows that the predictions from the 1D problem $\mathcal{P}_{\mathcal{PCAP}}$ are not directly applicable to the complex 3D CHT case: following the 1D analysis, the points (α_2, h_2) and (α_2, h_3) are stable while the point (α_2, h_1) is not stable. The convergence rate is influenced by the choice of the CRP: simulations converge faster for slower values of h (Fig. 12a). Indeed from h_3 to h_1 the CPU time needed to reach a steady state is divided by approximatively 2.5.

As done for the CRP h, three computations with SCS are then compared to evaluate the effect of the CSTP α in Eq. (33) on convergence: $SCS(\alpha_1, h_2)$, $SCS(\alpha_2, h_2)$ and $SCS(\alpha_3, h_2)$ (Table 6). Fig. 12b shows the evolutions of the mean temperature in the blade as a function of reduced solid time t_s/τ_s for these simulations. The

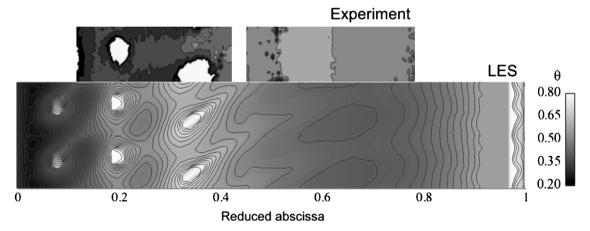


Fig. 10. 2D plot of time averaged cooling efficiency on the pressure side: comparison of experimental results and coupled simulation. The scale of Θ corresponds only to the LES field.

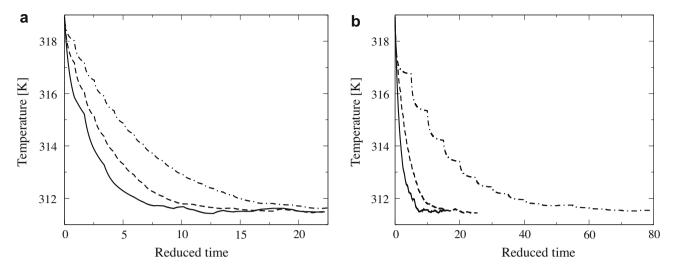


Fig. 12. Evolutions of the mean temperature in the blade as a function of reduced solid time t_s/τ_s for (a) $SCS(\alpha_2, h_1)$ —, $SCS(\alpha_2, h_2)$ — and $SCS(\alpha_2, h_3)$ — · · ; and for (b) $SCS(\alpha_1, h_2)$ —, $SCS(\alpha_2, h_2)$ — and $SCS(\alpha_3, h_2)$ — · · . The parameters (α_i, h_j) are given in Table 6. As a reference for restitution times, one solid characteristic time τ_s requires almost a total of 295 CPU h for the LES and the heat conduction solvers on a SGI Altix ICE8200.

numerical schemes obtained with α_1 , α_2 and α_3 combined with h_2 are stable and converge to the same thermal steady state. Fig. 12b shows that the choice of the CSTP directly affects the convergence rate of the coupled simulations. Small values of α enhances the restitution times of the computations: high update frequencies of interface variables allow the computations to follow the steepest descent to converged state.

To conclude, good convergence rates are obtained for small values of both the CSTP α and CRP h within the stability domain. When α is sufficiently small, SCS becomes inefficient compared to PCS. Indeed, below a certain value of the CSTP, the computational time due to the resolution of fluid and solid equations is smaller than the management times due to launching of the codes as well as initialization and finalization of the computations (reading of meshes and initial solutions, writing of final solutions...). PCS allows to avoid these restart steps and is therefore recommended. It is also important to note that when the CSTP α becomes too small (i.e. decreasing the number of iterations of the solvers), the execution time of the fluid and solid solvers between two successive updates of boundary conditions can become lower than the time requested to data exchanges. Such a configuration is not acceptable for the efficiency of the method and a balance is needed to fix the CSTP.

5. Conclusions

The thermal loads on gas turbine components controls life-time of the system. Numerical predictions of conjugate heat transfer raise problems which make the simulations difficult to compute:

- the flows are very complex: three-dimensional unsteady turbulent effects, mixing of flows with different properties, high speeds with compressibility effects (with large variations of Mach number up to supersonic conditions).
- distinct physics must be coupled with very different length and time scales,
- the size of the computations requires parallel computations to reduce restitution times.

In the present study, a coupling strategy for conjugate heat transfer with Large Eddy Simulation (LES) is proposed and applied to a film-cooled turbine vane experimentally tested at the Institute of Jet Propulsion. The methodology is based on a partitioned approach: a LES solver exchanges information with a heat transfer code within solids via a supervisor. It is shown that the LES solver is well adapted to capture the flow that develops around the film-cooled turbine blade.

The coupling methodology is first assessed in term of stability and efficiency. At the fluid/solid interface, a mixed condition on wall temperature and heat fluxes is written based on a coupling relaxation parameter h. A stability analysis on a simplified onedimensional problem allows the determination and optimization of coupling parameters. Two parameters control both the stability and cost of the coupled simulations: the mesh Biot number of the solid domain $\mathcal{D} = h\Delta x_s/\lambda_s$ (where Δx_s and λ_s are, respectively, the solid discretization and the solid conductivity) and the coupling synchronization time parameter α which defined the time between two coupling events. The coupled tool is ported on massively parallel architectures and applied to an experimental film-cooled turbine vane. Results demonstrate both the efficiency of the parallel implementation and the quality of the results. Thermal convergence of the blade is accelerated by high frequencies of exchanges between the codes (small values of the coupling synchronization time parameter α). Coupled and non-coupled simulations are compared to experimental results and discussed in terms of cooling efficiency and flow structures.

The comparison between adiabatic LES and measurements is good for the isentropic Mach number along the blade. Reasonable aerodynamic predictions are a necessary condition to reproduce the blade skin temperature field. As the LES computation gives a correct air flow distribution within the cooling system and reproduces the mixing between the hot stream and the cooling flow, the trends of computed cooling efficiency agree reasonably well with experimental data. Discrepancies in the levels are mainly due to a lack of resolution in the fluid flow near the wall. Finally, the coupled simulation is used to describe the role of the three rows of cooling jets in heat transfer.

Further studies will concern the analyze of mesh resolution as well as the use of thermal wall model to enhance thermal predictions. The effect of the number of spanwise pitch of the film cooling hole pattern including jet to jet interactions will also be investigated.

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