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Deep reinforcement learning for the control of conjugate heat transfer

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Abstract

This research gauges the ability of deep reinforcement learning (DRL) techniques to assist the control of conjugate heat transfer systems governed by the coupled Navier–Stokes and heat equations. It uses a novel, "degenerate" version of the proximal policy optimization (PPO) algorithm, intended for situations where the optimal policy to be learnt by a neural network does not depend on state, as is notably the case in optimization and open-loop control problems. The numerical reward fed to the neural network is computed with an in-house stabilized finite elements environment combining variational multi-scale (VMS) modeling of the governing equations, immerse volume method, and multi-component anisotropic mesh adaptation. Several test cases of natural and forced convection in two and three dimensions are used as testbed for developing the methodology. The approach successfully alleviates the natural convection induced enhancement of heat transfer in a two-dimensional, differentially heated square cavity controlled by piece-wise constant fluctuations of the sidewall temperature. It also proves capable of improving the homogeneity of temperature across the surface of two and three-dimensional hot workpieces under impingement cooling. Various cases are tackled, in which the position of multiple cold air injectors is optimized relative to a fixed workpiece position. The flexibility of the numerical framework makes it tractable to solve also the inverse problem, i.e., to optimize the workpiece position relative to a fixed injector distribution. The obtained results showcase the potential of the method for black-box optimization of practically meaningful computational fluid dynamics (CFD) conjugate heat transfer systems. More significantly, they stress how DRL can reveal unanticipated solutions or parameter relations (as the optimal workpiece position under symmetrical actuation turns to be offset from the symmetry axis), in addition to being a tool for optimizing searches in large parameter spaces.

Keywords: Deep Reinforcement Learning; Artificial Neural Networks; Conjugate heat transfer; Computational fluid dynamics; Thermal control.

1 1. Introduction

Thermal control, defined as the ability to finesse the thermal properties of a volume of fluid 2 (and of the solid objects inside) into a certain desired state, is a field of tremendous societal and 3 economical importance. For instance, heat/cool exchangers are used in a broad range of industrial 4 applications to regulate process temperatures by heat or cool transfer between fluid media, which in turn ensures that machinery, chemicals, water, gas, and other substances remain within safe operating conditions. Green building engineering is another field whose focus is on regulating indoor 7 thermal conditions (temperature, humidity) under substantial variations of the ambient condi-8 tions to provide high-quality living and working environments. In many manufacturing processes, q thermal conditioning is also intended to improve the final mechanical (e.g., hardness, toughness, 10 resistance), electrical, or optical properties of the product, the general picture being that high 11 temperature gradients are useful to speed up the process but generally harm the quality of the 12 outcome because of heat transfer inhomogeneities caused by the increased convection by the fluid 13 14 particles. All such problems fall under the purview of this line of study.

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Numerous strategies have been implemented to control fluid mechanical systems (including con-15 jugate heat transfer systems combining thermal conduction in the solid and convective transfer in 16 the fluid), either open-loop with passive appendices (e.g., end plate, splitter plate, small secondary 17 cylinder, or flexible tail), or open-loop with actuating devices (e.g., plasma actuation, boundary 18 temperatures, steady or unsteady base bleeding, rotation) or closed-loop (e.g. via transverse mo-19 tion, perturbations of the thermal boundary layer, blowing/suction, rotation, all relying on an 20 appropriate sensing of flow variables). Nonetheless, many of the proposed strategies are trial and 21 error, and therefore require extensive and costly experimental or numerical campaigns. This has 22 motivated the development of analytical methods and numerical algorithms for the optimal control 23 of Navier–Stokes systems [1-3], and the maturing of mathematical methods in flow control and 24 discrete concepts for PDE constrained optimization. Applications to the heat equation [4] and the 25 coupled Navier–Stokes and heat equations [5–8] have also been considered, including fresh devel-26 opments meant to alter the linear amplification of flow disturbances [9], but the general picture 27 remains that the optimal control of conducting-convecting (possibly radiating) fluids has not been 28 extensively studied. 29

The premise of this research is that the related task of selecting an optimal subset of control 30 parameters can alternatively be assisted by machine learning algorithms. Indeed, the introduc-31 tion of the back-propagation algorithm [10] has progressively turned Artificial Neural Networks 32 (ANN) into a family of versatile non-parametric tools that can be trained to hierarchically extract 33 informative features from data and to provide qualitative and quantitative modeling predictions. 34 Together with the increased affordability of high-performance computational hardware, this has 35 allowed leveraging the ever-increasing volume of data generated for research and engineering pur-36 37 poses into novel insight and actionable information, which in turn has entirely transformed scientific disciplines, such as robotics [11, 12] or image analysis [13]. Owing to the ability of neural networks 38 to handle stiff, large-scale nonlinear problems [14], machine learning algorithms have also been 39 making rapid inroads in fluid mechanics, as a mean to solve the Navier–Stokes equations [15] or 40 to predict closure terms in turbulence models [16]; see also Ref. [17] for an overview of the current 41 developments and opportunities. 42

Neural networks can also be used to solve decision-making problems, which is the purpose of 43 Deep Reinforcement Learning (DRL, where the *deep* terminology generally weighs on the sizable 44 depth of the network), an advanced branch of machine learning. Simply put, a neural network trains 45 in finding out which actions or succession of actions maximize a numerical reward signal, with the 46 possibility for a given action to affect not only the immediate but also the future rewards. Successful 47 applications of DRL range from AlphaGo, the well-known ANN that defeated the top-level human 48 player at the game of Go [18] to the real-word deployment of legged robots [19], to breakthroughs in 49 computer vision (e.g., filtering, or extracting image features) [20] and optimal control problems [21, 50 22]. There is also great potential for applying DRL to fluid mechanics, for which efforts are ongoing 51 but still at an early stage. Sustained commitment from the machine learning community has 52 allowed expanding the scope from computationally inexpensive, low-dimensional reductions of the 53 underlying fluid dynamics [23–25] to complex Navier–Stokes systems [26, 27], with a handful of 54 pioneering studies providing insight into the performance improvements to be delivered in shape 55 optimization [28–30] and flow control [31–37], including recent advances assessing experimentally 56 the effectiveness of reinforcement learning control strategies [38]. The literature on thermal control 57 is even more scarce, as our literature review did not reveal any other study considering DRL-based 58 control of conjugate heat transfer aside from [39], another research effort conducted in the same 59 time frame as the present work that will be discussed further on, plus a few other publications 60 relying on dealing with energy efficiency in civil engineering from low-dimensional thermodynamic 61 models basically unrelated to the equations of fluid dynamics [40, 41]. 62

This research assesses the feasibility of using proximal policy optimization (PPO [22]) for control 63 and optimization purposes of conjugate heat transfer systems, as governed by the coupled Navier-64 Stokes and heat equations. The objective here is to keep shaping the capabilities of the method 65 (PPO is still a relatively newcomer that has quickly emerged as the go-to DRL algorithm due to 66 its data efficiency, simplicity of implementation and reliable performance) and to narrow the gap 67 between DRL and advanced numerical methods for multi-scale, multi-physics computational fluid 68 dynamics (CFD). We investigate more specifically the "degenerate" single-step PPO algorithm 69 introduced in [30] for optimization and open-loop control problems, as the optimal policy to be 70

learnt is then state-independent, and it may be enough for the neural network to get only one 71 attempt per episode at finding the optimal. Several problems of conjugate heat transfer in two and 72 three dimensions are used as testbed to push forward the development of this novel approach, whose 73 potential for reliable black-box optimization of computational fluid dynamics (CFD) systems has 74 been recently assessed for open-loop drag reduction in cylinder flows at Reynolds numbers ranging 75 from a few hundreds to a few ten thousands [42]. To the best of the authors knowledge, this 76 constitutes the first attempt to achieve DRL-based control of conjugate forced convection heat 77 transfer, while [39] is the first attempt to achieve DRL control of conjugate *natural* convection 78 heat transfer. 79

The organization is as follows: section 2 outlines the main features of the finite element CFD 80 environment used to compute the numerical reward fed to the neural network, that combines 81 variational multi-scale (VMS) modeling of the governing equations, immerse volume method, and 82 multi-component anisotropic mesh adaptation. The baseline principles and assumptions of DRL 83 and PPO are presented in section 3, together with the specifics of the single-step PPO algorithm. 84 Section 4 revisits the natural convection case of [39] for the purpose of validation and assessment 85 part of the method capabilities. In section 5, DRL is used to control conjugate heat transfer in 86 a model setup of two-dimensional workpiece cooling by impingement of a fluid. An extension to 87 three-dimensional workpieces is proposed in section 6. 88

⁸⁹ 2. Computational fluid dynamics

The focus of this research is on conjugate heat transfer and laminar, incompressible fluid flow problems in two and three-dimensions, for which the conservation of mass, momentum and energy is described by the nonlinear, coupled Navier–Stokes and heat equations

$$\nabla \cdot \boldsymbol{u} = 0, \qquad (1)$$

$$\rho(\partial_t \boldsymbol{u} + \boldsymbol{u} \cdot \nabla \boldsymbol{u}) = \nabla \cdot (-p\mathbf{I} + 2\mu\boldsymbol{\varepsilon}(\boldsymbol{u})) + \boldsymbol{\psi}, \qquad (2)$$

$$\rho c_p (\partial_t T + \boldsymbol{u} \cdot \nabla T) = \nabla \cdot (\lambda \nabla T) + \chi, \qquad (3)$$

⁹³ where \boldsymbol{u} is the velocity field, p is the pressure, T is the temperature, $\boldsymbol{\varepsilon}(\boldsymbol{u}) = (\nabla \boldsymbol{u} + \nabla \boldsymbol{u}^T)/2$ is the ⁹⁴ rate of deformation tensor, $\boldsymbol{\psi}$ and $\boldsymbol{\chi}$ are source terms (modeling, e.g., buoyancy or radiative heat ⁹⁵ transfer), and we assume here constant fluid density ρ , dynamic viscosity μ , thermal conductivity ⁹⁶ λ , and specific heat c_p .

97 2.1. The immersed volume method

The numerical modeling of conjugate heat transfer mostly depends upon a heat transfer co-98 efficient to ensure that the proper amount of heat is exchanged at the fluid/solid interface via 99 thermal boundary conditions. Computing said coefficient is no small task (as it requires solving 100 an inverse problem to assimilate relevant experimental data, which in turn requires such data to 101 be available), and is generally acknowledged to be a limiting issue for practical applications where 102 one must vary, e.g., the shape, number and position of the solid, or the fluid and/or solid material 103 properties. We thus rather use here the immerse volume method (IVM) to combine both the fluid 104 and solid phases into a single fluid with variable material properties. Simply put, we solve equa-105 tions formally identical to (1)-(3) on a unique computational domain Ω , but with variable density, 106 dynamic viscosity, conductivity, and specific heat, which removes the need for a heat transfer coef-107 ficient since the amount of heat exchanged at the interface then proceeds solely from the individual 108 material properties on either side of it. In order to ensure numerical accuracy, such an approach 109 must combine three key ingredients, that are briefly reviewed in the next paragraphs: an interface 110 capturing method, anisotropic mesh adaptation to achieve a high-fidelity description of said inter-111 face, and relevant mixing laws to describe the properties of the composite fluid. One point worth 112 mentioning is that the interface here is static, although the same numerical framework can be used 113 to dynamically track moving interfaces, and thus to encompass the effect of solid displacements. 114 This is because the solid is fixed once an action has been taken by the PPO agent, although not 115 fixed over the course of optimization, as the solid position can very well be the quantity subjected 116 to optimization, as illustrated in section 5.3.4. 117 118

3

¹¹⁹ - Level set method. The level set approach is used to localize the fluid/solid interface by the zero ¹²⁰ iso-value of a smooth function. In practice, a signed distance function ϕ is used to localize the ¹²¹ interface and initialize the material properties on both either side of it, with the convention that ¹²² $\phi > 0$ (resp. $\phi < 0$) in the fluid (resp. the solid).

- Anisotropic mesh adaptation. The interface may intersect arbitrarily the mesh elements if it 124 is not aligned with the element edges, in which case discontinuous material properties across 125 the interface can yield oscillations of the numerical solutions. We thus use the anisotropic mesh 126 adaptation technique presented in [43] to ensure that the material properties are distributed as 127 accurately and smoothly as possible over the smallest possible thickness around the interface. This 128 is done computing modified distances from a symmetric positive defined tensor (the metric) whose 129 eigenvectors define preferential directions along which mesh sizes can be prescribed from the related 130 eigenvalues. The metric used here is isotropic far from the interface, with mesh size set equal to 131 h_{∞} in all directions, but anisotropic near the interface, with mesh size equal to h_{\perp} in the direction 132 normal to the interface, and to h_{∞} in the other directions. This is written for an intended thickness 133 δ as 134

$$\boldsymbol{M} = K(\phi)\boldsymbol{n} \otimes \boldsymbol{n} + \frac{1}{h_{\infty}^2}\boldsymbol{I} \qquad \text{with} \qquad K(\phi) = \begin{cases} 0 & \text{if } |\phi| \ge \delta/2, \\ \frac{1}{h_{\perp}^2} - \frac{1}{h_{\infty}^2} & \text{if } |\phi| < \delta/2, \end{cases}$$
(4)

where $n = \nabla \phi / ||\nabla \phi||$ is the unit normal to the fluid/solid interface computed from the level set gradient. A posteriori anisotropic error estimator is then used to minimize the interpolation error under the constraint of a fixed number of edges in the mesh. A unique metric can be built from multi-component error vectors [43–46], which is especially relevant for conjugate heat transfer optimization, as it allows each learning episode to use an equally accurate mesh adapted from the velocity vector and magnitude, the temperature field, and the level set.

- Mixing laws. The composite density, dynamic viscosity and specific heat featured in equations (1)(3) are computed as the arithmetic means of the fluid and solid values, for instance the composite
density is

$$\rho = \rho_f H_\epsilon(\phi) + \rho_s (1 - H_\epsilon(\phi)), \qquad (5)$$

145 where H_{ϵ} is the smoothed Heaviside function defined as

123

$$H_{\epsilon}(\phi) = \begin{cases} 0 & \text{if } \phi < -\epsilon, \\ \frac{1}{2}(1 + \frac{\phi}{\epsilon} + \frac{1}{\pi}\sin(\pi\frac{\phi}{\epsilon})) & \text{if } |\phi| \le \epsilon, \\ 1 & \text{if } \phi > \epsilon, \end{cases}$$
(6)

and ϵ is a regularization parameter proportional to the mesh size in the normal direction to the interface, set here to $\epsilon = 2h_{\perp}$. In order to ensure continuity of the heat flux across the interface, the thermal conductivity is computed as the harmonic mean

$$\frac{1}{\lambda} = \frac{1}{\lambda_f} H_{\epsilon}(\phi) + \frac{1}{\lambda_s} (1 - H_{\epsilon}(\phi)), \qquad (7)$$

as obtained from a steady, no source, one dimensional analysis of the heat flux when the conductivity varies stepwise from one medium to the next; see [47] for detailed derivation and analysis,
and [48] for proof of the gain in numerical accuracy (with respect to the arithmetic mean model)
by comparison with analytical solutions.

153 2.2. Variational multi-scale approach (VMS)

In the context of finite element methods (that remain widely used to simulate engineering CFD systems due to their ability to handle complex geometries), direct numerical simulation (DNS) solves the weak form of (1)-(3), obtained by integrating by parts the pressure, viscous and conductive terms, to give

$$(\rho(\partial_t \boldsymbol{u} + \boldsymbol{u} \cdot \nabla \boldsymbol{u}), \boldsymbol{w}) + (2\mu\varepsilon(\boldsymbol{u}), \varepsilon(\boldsymbol{w})) - (p, \nabla \cdot \boldsymbol{w}) + (\nabla \cdot \boldsymbol{u}, q) = (\boldsymbol{\psi}, \boldsymbol{w}), \quad (8)$$

$$(\rho c_p(o_t I + \boldsymbol{u} \cdot \nabla I), s) + (\lambda \nabla I, \nabla s) = (\chi, s),$$
(9)

where (,) is the L^2 inner product on the computational domain, w, q and s are relevant test functions for the velocity, pressure and temperature variables, and all fluid properties are those mixed with the smoothed Heaviside function (6).

We use here the variational multi-scale (VMS) approach [49–51] to solve a stabilized formulation 161 of (8)-(9), which allows circumventing the Babuska—Brezzi condition (that otherwise imposes that 162 different interpolation orders be used to discretize the velocity and pressure variables, while we 163 use here simple continuous piecewise linear P_1 elements for all variables) and prevents numerical 164 instabilities in convection regimes at high Reynolds numbers. We shall not go into the extensive 165 details about the derivation of the stabilized formulations, for which the reader is referred to [52, 53]. 166 Suffice it to say here that the flow quantities are split into coarse and fine scale components, that 167 correspond to different levels of resolution. The fine scales are solved in an approximate manner 168 to allow modeling their effect into the large-scale equations. This gives rise to additional terms in 169 the right-hand side of (8)-(9), and yields the following weak forms for the large scale 170

$$(\rho(\partial_t \boldsymbol{u} + \boldsymbol{u} \cdot \nabla \boldsymbol{u}), \boldsymbol{w}) + (2\mu\boldsymbol{\varepsilon}(\boldsymbol{u}), \boldsymbol{\varepsilon}(\boldsymbol{w})) - (p, \nabla \cdot \boldsymbol{w}) + (\nabla \cdot \boldsymbol{u}, q) = (\boldsymbol{\psi}, \boldsymbol{w}) + \sum_{K \in \mathcal{T}_h} \left[(\tau_1 \mathcal{R}_M, \boldsymbol{u} \cdot \nabla \boldsymbol{w})_K + (\tau_1 \mathcal{R}_M, \nabla q)_K + (\tau_2 \mathcal{R}_C, \nabla \cdot \boldsymbol{w})_K \right],$$
(10)

$$(\rho c_p(\partial_t T + \boldsymbol{u} \cdot \nabla T), s) + (\lambda \nabla T, \nabla s) = (\chi, s) + \sum_{K \in \mathcal{T}_h} [(\tau_3 \mathcal{R}_T, \boldsymbol{u} \cdot \nabla s)_K + (\tau_4 \mathcal{R}_T, \zeta \nabla T \cdot \nabla s)_K],$$
(11)

where $(,)_K$ is the inner product on element K, we denote by $\zeta = \boldsymbol{u} \cdot \nabla T/||\nabla T||^2$ the (normalized) velocity projected along the direction of the temperature gradient, and the \mathcal{R} terms are the governing equations residuals

$$-\mathcal{R}_C = \nabla \cdot \boldsymbol{u} \,, \quad -\mathcal{R}_M = \rho(\partial_t \boldsymbol{u} + \boldsymbol{u} \cdot \nabla \boldsymbol{u}) + \nabla p - \boldsymbol{\psi} \quad -\mathcal{R}_T = \rho c_p(\partial_t T + \boldsymbol{u} \cdot \nabla T) - \chi \,, \quad (12)$$

whose second derivatives vanish since we use linear interpolation functions. In (10), $\tau_{1,2}$ are ad-hoc mesh-dependent stabilization parameters defined in [54, 55]. Conversely, in (11), $\tau_{3,4}$ are meshindependent stabilization parameters acting both in the direction of the solution and of its gradient, that proceed from the stabilization of the ubiquitous convection-diffusion-reaction equation [56, 57], whose definition is given in [58, 59].

The governing equations are solved sequentially, i.e., we solve first (10), then use the resulting 179 fluid velocity to solve (11). All linear systems are preconditioned with a block Jacobi method 180 supplemented by an incomplete LU factorization, and solved with the GMRES algorithm, with 181 tolerance threshold set to 10^{-6} for the Navier–Stokes equations, and 10^{-5} for the heat equation. 182 The time derivatives and convection terms of the Navier–Stokes equations and related VMS source 183 terms are integrated semi-implicitly using the first-order backward differentiation formula and 184 Newton–Gregory backward polynomial. The viscous, pressure and divergence terms are treated 185 implicitly with the backward Euler scheme. Finally, the VMS stabilization terms $\tau_{1,2}$ are treated 186 explicitly with the forward Euler scheme, which yields 187

$$\left(\rho\left(\frac{\boldsymbol{u}^{i+1}-\boldsymbol{u}^{i}}{\Delta t}+\boldsymbol{u}^{i}\cdot\nabla\boldsymbol{u}^{i+1}\right),\,\boldsymbol{w}\right)+\left(2\mu\varepsilon(\boldsymbol{u}^{i+1}),\,\varepsilon(\boldsymbol{w})\right)-\left(p^{i+1},\,\nabla\cdot\boldsymbol{w}\right)+\left(\nabla\cdot\boldsymbol{u}^{i+1},\,q\right)=\left(\psi^{i},\,\boldsymbol{w}\right)\\+\sum_{K\in\mathcal{T}_{h}}\left[\left(\tau_{1}^{i}\mathcal{R}_{M}^{i+1},\,\boldsymbol{u}^{i}\cdot\nabla\boldsymbol{w}\right)_{K}+\left(\tau_{1}^{i}\mathcal{R}_{M}^{i+1},\,\nabla q\right)_{K}+\left(\tau_{2}^{i}\mathcal{R}_{C}^{i+1},\,\nabla\cdot\boldsymbol{w}\right)_{K}\right],$$
(13)

188 with residuals

$$-\mathcal{R}_{C}^{i+1} = \nabla \cdot \boldsymbol{u}^{i+1}, \quad -\mathcal{R}_{M}^{i+1} = \rho(\frac{\boldsymbol{u}^{i+1} - \boldsymbol{u}^{i}}{\Delta t} + \boldsymbol{u}^{i} \cdot \nabla \boldsymbol{u}^{i+1}) + \nabla p^{i+1} - \boldsymbol{\psi}^{i}, \quad (14)$$

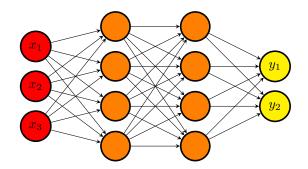


Figure 1: Fully connected neural network with two hidden layers, modeling a mapping from \mathbb{R}^3 to \mathbb{R}^2 .

where the superscript *i* refers to the solution at time $t_i = i\Delta t$. The time derivatives, convection and conduction terms of the heat equation and related VMS source terms are integrated implicitly with the backward Euler scheme (modeling the velocity after u^{i+1} wherever needed on behalf of the sequential resolution process, although we drop the dependence in the notation to ease the reading). The VMS stabilization terms $\tau_{3,4}$ are treated explicitly with the forward Euler scheme¹, to give

$$(\rho c_p(\frac{T^{i+1} - T^i}{\Delta t} + \boldsymbol{u}^{i+1} \cdot \nabla T^{i+1}), s) + (\lambda \nabla T^{i+1}, \nabla s) = (\chi^i, s)$$

+
$$\sum_{K \in \mathcal{T}_h} [(\tau_3^i \mathcal{R}_T^{i+1}, \boldsymbol{u}^{i+1} \cdot \nabla s)_K + (\tau_4^i \mathcal{R}_T^{i+1}, \zeta^i \nabla T^i \cdot \nabla s)_K],$$
(15)

¹⁹⁵ with residual

$$-\mathcal{R}_T^{i+1} = \rho c_p \left(\frac{T^{i+1} - T^i}{\Delta t} + \boldsymbol{u}^{i+1} \cdot \nabla T^{i+1}\right) - \chi^i \,. \tag{16}$$

We solve equations (13)-(15) with an in-house VMS solver whose flexibility, accuracy and reliability 196 is assessed in a series of previous papers to which the reader is referred for further information, see 197 in particular [55, 60] for the detailed mathematical formulation of the IVM in the context of finite 198 element VMS methods. The ability of the IVM to handle the abrupt conductivity change across 199 the fluid/solid interface is documented in [53, 61, 62]. Excellent agreement with reference solutions 200 available from the literature and in-house data obtained enforcing proper thermal conditions at the 201 boundary of body-fitted meshes is reported for several time-dependent conjugate heat transfer test 202 cases (e.g., mixed convection in a plane channel flow, combined convection in square enclosures 203 and conduction/radiation heat transfer, all in two dimensions). Ref. [61] also reports favorable 204 agreement between the IVM and in-house experimental data pertaining to a three-dimensional test 205 case representative of an industrial cooling system, which provides strong evidence of relevance for 206 the intended application. 207

²⁰⁸ 3. Deep reinforcement learning and proximal policy optimization

209 3.1. Neural networks

A neural network (NN) is a collection of artificial neurons, i.e., connected computational units that can be trained to arbitrarily well approximate the mapping function between input and output spaces. Each connection provides the output of a neuron as an input to another neuron. Each neuron performs a weighted sum of its inputs, to assign significance to the inputs with regard to the task the algorithm is trying to learn. It then adds a bias to better represent the part of the output that is actually independent of the input. Finally, it feeds an activation function that determines whether and to what extent the computed value should affect the outcome. As sketched in figure 1,

¹That is, with respect to T, but the velocity is modeled after its latest computed approximation u^{i+1} wherever needed.

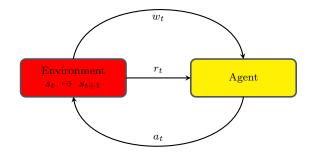


Figure 2: RL agent and its interactions with its environment.

a fully connected network is generally organized into layers, with the neurons of one layer being
connected solely to those of the immediately preceding and following layers. The layer that receives
the external data is the input layer, the layer that produces the outcome is the output layer, and
in between them are zero or more hidden layers.

The design of an efficient neural network requires a proper optimization of the weights and biases, together with a relevant nonlinear activation function. The abundant literature available on this topic points to a relevant network architecture (e.g., type of network, depth, width of each layer), finely tuned hyper parameters (i.e., parameters whose value cannot be estimated from data, e.g., optimizer, learning rate, batch size) and a sufficiently large amount of data to learn from as being the key ingredients for success; see, e.g., Ref. [63] and the references therein.

227 3.2. Deep reinforcement learning

Deep reinforcement learning (DRL) is an advanced branch of machine learning in which deep neural networks train in solving sequential decision-making problems. It is a natural extension of reinforcement learning (RL), in which an agent (the neural network) is taught how to behave in an environment by taking actions and by receiving feedback from it under the form of a reward (to measure how good or bad the action was) and information (to gauge how the action has affected the environment). This can be formulated as a Markov Decision Process, for which a typical execution goes as follows (see also figure 2):

- assume the environment is in state $s_t \in S$ at iteration t, where S is a set of states,
- the agent uses w_t , an observation of the current environment state (and possibly a partial subset of s_t) to take action $a_t \in \mathcal{A}$, where \mathcal{A} is a set of actions,
- the environment reacts to the action and transitions from s_t to state $s_{t+1} \in S$,
- the agent is fed with a reward $r_t \in \mathcal{R}$, where \mathcal{R} is a set of rewards, and a new observation w_{t+1} ,

This repeats until some termination state is reached, the succession of states and actions defining a trajectory $\tau = (s_0, a_0, s_1, a_1, ...)$. In any given state, the objective of the agent is to determine the action maximizing its cumulative reward over an episode, i.e., over one instance of the scenario in which the agent takes actions. Most often, the quantity of interest is the discounted cumulative reward along a trajectory defined as

$$R(\tau) = \sum_{t=0}^{T} \gamma^t r_t \,, \tag{17}$$

where T is the horizon of the trajectory, and $\gamma \in [0, 1]$ is a discount factor that weighs the relative importance of present and future rewards (the agent being short-sighted in the limit where $\gamma \to 0$, since it then cares solely about the first reward, and far-sighted in the limit where $\gamma \to 1$, since it then cares equally about all rewards).

There exist two main types of RL algorithms, namely model-based methods, in which the 250 agent tries to build a model of how the environment works to make predictions about what the 251 next state and reward will be before taking any action, and model-free methods, in which the agent 252 conversely interacts with the environment without trying to understand it, and are prominent in 253 the DRL community. Another important distinction to be made within model-free algorithms 254 is that between value-based methods, in which the agent learns to predict the future reward of 255 taking an action when provided a given state, then selects the maximum action based on these 256 estimates, and policy-based methods, in which it optimizes the expected reward of a decision policy 257 mapping states to actions. Many of the most successful algorithms in DRL (including proximal 258 policy optimization, whose assessment for flow control and optimization purposes is the primary 259 motivation for this research) proceed from policy gradient methods, in which gradient ascent is 260 used to optimize a parameterized policy with respect to the expected return, as further explained 261 in the next section. The reader interested in a more thorough introduction to the zoology of RL 262 methods (together with their respective pros and cons) is referred to Ref. [64]. 263

²⁶⁴ 3.3. From policy methods to Proximal policy optimization

This section intended for the non-specialist reader briefly reviews the basic principles and assumptions of policy gradient methods, together with the various steps taken for improvement.

Policy methods. A policy method maximizes the expected discounted cumulative reward of a
 decision policy mapping states to actions. It resorts not to a value function, but to a probability
 distribution over actions given states, that fully defines the behavior of the agent. Since policies
 are most often stochastic, the following notations are introduced:

• $\pi(s, a)$ is the probability of taking action a in state s under policy π ,

• $Q^{\pi}(s, a)$ is the expected value of the return of the policy after taking action a in state s (also termed state-action value function or Q-function)

$$Q^{\pi}(s,a) = \mathbb{E}_{\pi} \left[R(\tau) | s, a \right], \tag{18}$$

where we use \mathbb{E}_{π} for the expected value \mathbb{E} under policy π .

• $V^{\pi}(s)$ is the expected value of the return of the policy in state *s* (also termed value function or V-function)

$$V^{\pi}(s) = \mathbb{E}_{\pi} \left[R(\tau) | s \right].$$
⁽¹⁹⁾

²⁷⁸ The V and Q functions are therefore such that

$$V^{\pi}(s) = \sum_{a} \pi(s, a) Q^{\pi}(s, a) , \qquad (20)$$

so $V^{\pi}(s)$ can also be understood as the probability-weighted average of discounted cumulated rewards over all possible actions in state s.

- Policy gradient methods. A policy gradient method aims at optimizing a parametrized policy π_{θ} , where θ denotes the free parameters whose value can be learnt from data (as opposed to the hyper parameters). In practice, one defines an objective function based on the expected discounted cumulative reward

$$J(\theta) = \mathbb{E}_{\pi_{\theta}} \left[R(\tau) \right], \tag{21}$$

and seeks the parameterization θ^* maximizing $J(\theta)$, hence such that

$$\theta^* = \arg\max_{\alpha} \mathbb{E}_{\pi_{\theta}} \left[R(\tau) \right], \tag{22}$$

which can be done on paper by plugging an estimator of the policy gradient $\nabla_{\theta} J(\theta)$ into a gradient ascent algorithm. This is no small task as one is looking for the gradient with respect to the policy parameters, in a context where the effects of policy changes on the state distribution are unknown (since modifying the policy will most likely modify the set of visited states, which will in turn affect performance in some indefinite manner). One commonly used estimator, derived in [64] using the log-probability trick, reads

$$\nabla_{\theta} J(\theta) = \mathbb{E}_{\pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \left(\pi_{\theta}(s_t, a_t) \right) R(\tau) \right] \sim \mathbb{E}_{\pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \left(\pi_{\theta}(s_t, a_t) \right) \widehat{A}^{\pi}(s_t, a_t) \right], \quad (23)$$

where \hat{A}^{π} is some biased estimator (here its normalization to zero mean and unit variance) of the advantage function

$$A^{\pi}(s,a) = Q^{\pi}(s,a) - V^{\pi}(s), \qquad (24)$$

that measures the improvement (if $A^{\pi} > 0$, otherwise the lack thereof) associated with taking action *a* in state *s* compared to taking the average over all possible actions. This is because the value function does not depend on θ , so taking it off changes neither the expected value, nor the gradient, but it does reduce the variance, and speeds up the training. Furthermore, when the policy π_{θ} is represented by a neural network (in which case θ simply denotes the network weights and biases to be optimized), the focus is rather on the policy loss defined as

$$L(\theta) = \mathbb{E}_{\pi_{\theta}} \left[\sum_{t=0}^{T} \log \left(\pi_{\theta}(a_t | s_t) \right) \widehat{A}(s_t, a_t) \right], \qquad (25)$$

whose gradient is equal to the (approximated) policy gradient (23) (since the gradient operator acts only on the log-policy term, not on the advantage) and is computed with respect to each weight and bias by the chain rule, one layer at the time, using the back-propagation algorithm [10].

- Trust regions. The performance of policy gradient methods is hurt by the high sensitivity to the 304 learning rate, i.e., the size of the step to be taken in the gradient direction. Indeed, small learning 305 rates are detrimental to learning, but large learning rates can lead to a performance collapse if the 306 agent falls off the cliff and restarts from a poorly performing state with a locally bad policy. This 307 is all the more harmful as the learning rate cannot be tuned locally, meaning that an above average 308 learning rate will speed up learning in some regions of the parameter space where the policy loss 309 is relatively flat, but will possibly trigger an exploding policy update in other regions exhibiting 310 sharper variations. One way to ensure continuous improvement is by imposing a trust region con-311 straint to limit the difference between the current and updated policies, which can be done by 312 determining first a maximum step size relevant for exploration, then by locating the optimal point 313 within this trust region. We will not dwell on the intricate details of the many algorithms developed 314 315 to solve such trust region optimization problems, e.g., natural policy gradient (NPG [65]), or trust region policy optimization (TRPO [66]). Suffice it to say that they use the minorize-maximization 316 algorithm to maximize iteratively a surrogate policy loss (i.e. a lower bound approximating locally 317 the actual loss at the current policy), but are difficult to implement and can be computationally 318 expensive, as they rely on an estimate of the second-order gradient of the policy log probability. 319 320

³²¹ - *Proximal policy optimization*. Proximal policy optimization (PPO) is another approach with ³²² simple and effective heuristics, that uses a probability ratio between the two policies to maximize ³²³ improvement without the risk of performance collapse [22]. The focus here is on the PPO-clip ³²⁴ algorithm², that optimizes the surrogate loss

$$L(\theta) = \mathbb{E}_{\pi_{\theta}} \left[\min \left(\frac{\pi_{\theta}(a|s)}{\pi_{\theta_{old}}(a|s)}, g(\epsilon, \widehat{A}^{\pi}(s, a)) \right) \widehat{A}^{\pi}(s, a) \right],$$
(26)

 $^{^{2}}$ There is also a PPO-Penalty variant which uses a penalization on the average Kullback–Leibler divergence between the current and new policies, but PPO-clip performs better in practice.

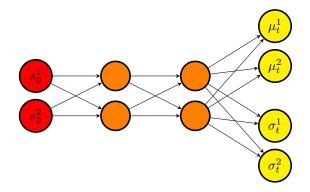


Figure 3: Agent network example used to map states to policy. The input state s_0 , here of size 2, is mapped to a mean μ and a standard deviation σ vectors, each of size 2. All activation functions are ReLu, except for that of the last layer, which are linear for the μ output, and softplus for the σ output. Orthogonal weights initialization is used throughout the network.

325 where

$$g(\epsilon, A) = \begin{cases} 1+\epsilon & A \ge 0, \\ 1-\epsilon & A < 0, \end{cases}$$
(27)

and $\epsilon \in [0.1, 0.3]$ is the clipping range, a small hyper parameter defining how far away the new policy is allowed to go from the old. The general picture is that a positive (resp. negative) advantage increases (resp. decreases) the probability of taking action a in state s, but always by a proportion smaller than ϵ , otherwise the min kicks in (26) and its argument hits a ceiling of $1 + \epsilon$ (resp. a floor of $1 - \epsilon$). This prevents stepping too far away from the current policy, and ensures that the new policy will behave similarly.

There exist more sophisticated PPO algorithms (e.g., Trust region PPO [67], that determines first a maximum step size relevant for exploration, then adaptively adjusts the clipping range to find the optimal within this trust region), but standard PPO has simple and effective heuristics. Namely, it is computationally inexpensive, easy to implement (as only the first-order gradient of the policy log probability is needed to calculate the clipped surrogate), and remains regarded as one of the most successful RL algorithms, achieving state-of-the-art performance across a wide range of challenging tasks.

339 3.4. Single-step PPO

We now come to single-step PPO, a "degenerate" version of PPO introduced in [30] and intended 340 for situations where the optimal policy to be learnt by the neural network is state-independent, as 341 is notably the case in optimization and open-loop control problems (closed-loop control problems 342 conversely require state-dependent policies for which standard PPO is best suited). The main 343 difference between standard and single-step PPO can be summed up as follows: where standard 344 PPO seeks the optimal set of actions a^* yielding the largest possible reward, single-step PPO seeks 345 the optimal mapping f_{θ^*} such that $a^* = f_{\theta^*}(s_0)$, where θ denotes the network free parameters 346 and s_0 is some input state (usually a vector of zeros) consistently fed to the agent for the optimal 347 policy to eventually embody the transformation from s_0 to a^* . The agent initially implements a 348 random state-action mapping f_{θ_0} from s_0 to an initial policy determined by the free parameters 349 initialization θ_0 , after which it gets only one attempt per learning episode at finding the optimal 350 (i.e., it interacts with the environment only once per episode). This is illustrated in figure 4 showing 351 the agent draw a population of actions $a_t = f_{\theta_t}(s_0)$ from the current policy, and being returned 352 incentives from the associated rewards to update the free parameters for the next population of 353 actions $a_{t+1} = f_{\theta_{t+1}}(s_0)$ to yield larger rewards. 354

In practice, the agent outputs a policy parameterized by the mean and variance of the probability density function of a *d*-dimensional multivariate normal distribution, with *d* the dimension of the action required by the environment. Actions drawn in $[-1,1]^d$ are then mapped into relevant physical ranges, a step deferred to the environment as being problem-specific. The resolution

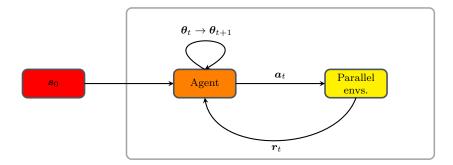


Figure 4: Action loop for single-step PPO. At each episode, the input state s_0 is provided to the agent, which in turn provides n actions to n parallel environments. The latter return n rewards, that evaluate the quality of each action taken. Once all the rewards are collected, an update of the agent parameters is made using the PPO loss (26).

essentially follows the process described in section 3.3, only a normalized averaged reward substitutes for the advantage function. This is because classical PPO is actor-critic, i.e., it improves the learning performance by updating two different networks, a first one called actor that controls the actions taken by the agent, and a second one called critic, that learns to estimate the advantage from the value function as

$$A(s_t, a_t) = r_t + \gamma V(s_{t+1}) - V(s_t).$$
(28)

In single-step PPO, the trajectory consists of a single state-action pair, so the discount factor can be set to $\gamma = 1$ with no loss of generality. In return, the advantage reduces to the whitened reward since the two rightmost terms cancel each other out in (28). This means that the approach can do without the value-function evaluations of the critic network, i.e., it is not actually actor-critic.

368 3.5. Numerical implementation

The present workflow relies on the online PPO implementation of Stable Baselines, a toolset 369 of reinforcement learning algorithms dedicated to the research community and industry [68], for 370 which a custom OpenAI environment has been designed using the Gym library [69]. Hyperbolic 371 tangent is used as default activation function. The instant reward r_t used to train the neural 372 network is simply the quantity subjected to optimization (modulo a plus or minus sign to tackle 373 both maximization and minimization problems). A moving average reward is also computed on the 374 fly as the sliding average over the 100 latest values of r_t (or the whole sample if it has insufficient 375 size). All other relevant hyper parameters are documented in the next sections, with the exception 376 of the discount factor (set to $\gamma = 1$). 377

In practice, actions are distributed to multiple environments running in parallel, each of which 378 executes a self-contained MPI-parallel CFD simulation and feeds data to the DRL algorithm (hence, 379 two levels of parallelism related to the environment and the computing architecture). The algorithm 380 waits for the simulations running in all parallel environments to be completed, then shuffles and 381 splits the rewards data set collected from all environments into several buffers (or mini-batches) 382 used sequentially to compute the loss and perform a network update. The process repeats for 383 several epochs, i.e., several full passes of the training algorithm over the entire data set (so the 384 policy network ends up being trained on samples generated by older policies, which is customary in 385 standard PPO operation). This simple parallelization technique is key to use DRL in the context 386 of CFD applications, as a sufficient number of actions drawn from the current policy must be 387 evaluated to accurately estimate the policy gradient. This comes at the expense of computing 388 the same amount of reward evaluations, and yields a substantial computational cost for high-389 dimensional fluid dynamics problems (typically from a few tens to several thousand hours for the 390 steady-state optimization problems considered herein). In the same vein, it should be noted that 391 the common practice in DRL studies to gain insight into the performances of the selected algorithm 392 by averaging results over multiple independent training runs with different random seeds is not 393 tractable, as it would trigger a prohibitively large computational burden. The same random seeds 394 have thus been deliberately used over the whole course of study to ensure a minimal level of 395 performance comparison between cases. 396

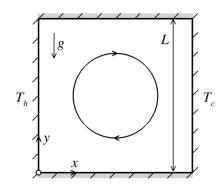


Figure 5: Schematic of the two-dimensional Rayleigh-Bénard set-up.

³⁹⁷ 4. Control of natural convection in 2-D closed cavity

398 4.1. Case description

We address first the control of natural convection in the two-dimensional differentially heated 300 square cavity schematically illustrated in figure 5(a). This is a widely studied benchmark system for 400 thermally-driven flows, relevant in nature and technical applications (e.g., ocean and atmospheric 401 convection, materials processing, metallurgy), that is thus suitable to validate and compare nu-402 merical solution algorithms while enriching the knowledge base for future projects in this field. 403 404 A Cartesian coordinate system is used with origin at the lower-left edge, horizontal x-axis, and vertical y-axis. The cavity has side L, its top and bottom horizontal walls are perfectly insulated 405 from the outside, and the vertical sidewalls are isothermal. Namely, the right sidewall is kept at 406 a constant, homogeneous "cold" temperature T_c , and the left sidewall is entirely controllable via a 407 constant in time, varying in space "hot" distribution $T_h(y)$ such that 408

$$\langle T_h \rangle > T_c \,, \tag{29}$$

where the brackets denote the average over space (here over the vertical position along the sidewall). In the following, we neglect radiative heat transfer ($\chi = 0$) and consider a Boussinesq system driven by buoyancy, hence

$$\boldsymbol{\psi} = \rho_0 \beta (T - T_c) g \boldsymbol{e}_y \,, \tag{30}$$

where g is the gravitational acceleration parallel to the sidewalls, β is the thermal expansion coefficient, and we use the cold sidewall temperature as Boussinesq reference temperature. By doing so, the pressure featured in the momentum equation (2) and related weak forms must be understood as the pressure correction representing the deviation from hydrostatic equilibrium. The governing equations are solved with no-slip conditions u = 0 on $\partial\Omega$ and temperature boundary conditions

$$\partial_y T(x,0,t) = \partial_y T(x,L,t) = 0, \qquad T(0,y,t) = \langle T_h \rangle + \ddot{T}_h(y), \qquad T(L,y,t) = T_c,$$
(31)

where \tilde{T}_h is a zero-mean (in the sense of the average over space) distribution of hot temperature fluctuations subjected to optimization, whose magnitude is bounded by some constant ΔT_{max} according to

$$|\tilde{T}_h(y)| \le \Delta T_{max} \,, \tag{32}$$

to avoid extreme and nonphysical temperature gradients. All results are made non-dimensional
using the cavity side, the heat conductivity time, and the well-defined, constant in time difference
between the averaged sidewall temperatures. The retained fluid properties yield values of the
Rayleigh and Prandtl numbers

$$\operatorname{Ra} = \frac{g\beta(\langle T_h \rangle - T_c)L^3}{\nu\alpha} = 10^4, \qquad \operatorname{Pr} = \frac{\nu}{\alpha} = 0.71, \qquad (33)$$

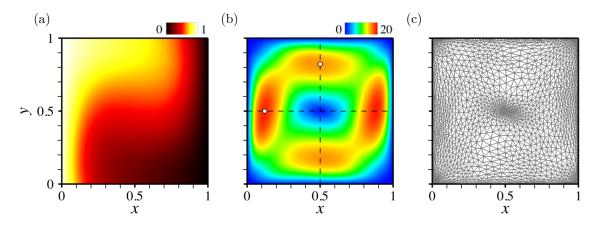


Figure 6: Iso-contours of the uncontrolled steady state (a) temperature and (b) velocity magnitude. (c) Adapted mesh. The circle symbols in (b) mark the positions of the maximum horizontal and vertical velocity along the centerlines reported in table 1.

		Present	Ref. [70]	Ref. [71]	Ref. [72]	Ref. [73]	Ref. [74]
	Nu	2.267	2.238	2.245	2.201	2.245	2.245
	$\max u(0.5, y)$	16.048	16.178	16.179	_	16.262	16.178
$Ra = 10^4$	y_{max}	0.823	0.823	0.824	0.832	0.818	0.827
	$\max v(x, 0.5)$	19.067	19.617	19.619	_	19.717	19.633
	x_{max}	0.120	0.119	0.121	0.113	0.119	0.123

Table 1: Comparison of the present numerical results in the absence of control with reference benchmark solutions from the literature.

425 where $\alpha = \lambda/(\rho c_p)$ is the thermal diffusivity.

In order to assess the accuracy of the numerical framework, the uncontrolled solution has 426 been computed by performing 60 iterations with time step $\Delta t = 0.5$ to march the initial solution 427 (consisting of zero velocity and uniform temperature, except at the hot sidewall) to steady state. 428 At each time step, an initially isotropic mesh is adapted under the constraint of a fixed number 429 of elements $n_{el} = 4000$ using a multiple-component criterion featuring velocity and temperature, 430 but no level-set. This is because the case is heat transfer but not conjugate heat transfer, as the 431 solid is solely at the boundary $\partial \Omega$ of the computational domain, where either the temperature is 432 known, or the heat flux is zero. It is thus implemented without the IVM and without a level set 433 (although accurate IVM numerical solutions have been obtained in [53] using thick sidewalls with 434 high thermal conductivity). The solution shown in figure 6(a,b) features a centered roll confined 435 by the cavity walls, consistently with the fact that Ra exceeds the critical value $Ra_c \sim 920$ for the 436 onset of convection (as extrapolated from the near-critical benchmark data in [70]) by one order 437 of magnitude, and heat transfer is thus driven by both conduction and convection. This shows in 438 the Nusselt number, i.e., the non-dimensional temperature gradient averaged over the hot sidewall 439

$$\mathrm{Nu} = -\langle \partial_x T \rangle \,, \tag{34}$$

whose present value Nu = 2.27 (as computed from 68 points uniformly distributed along the 440 sidewall) exceeds that Nu = 1 of the purely conductive solution, and exhibits excellent agreement 441 with benchmark results from the literature. This is evidenced in table 1 where we also report the 442 magnitude and position of the maximum horizontal velocity u (resp. the vertical velocity v) along 443 the vertical centerline (resp. the horizontal centerline). The corresponding adapted mesh shown in 444 figure 6(c) stresses that all boundary layers are sharply captured via extremely stretched elements, 445 and that the adaptation strategy yields refined meshes near high temperature gradients and close 446 to the side walls. Note however, the mesh refinement is not only along the boundary layers but also 447 close to the recirculation regions near the cavity center, while the elements in-between are coarse 448 and essentially isotropic. 449

450 4.2. Control

The question now being raised is whether DRL can be used to find a distribution of temperature 451 fluctuations T_h capable of alleviating convective heat transfer. To do so, we follow [39] and train a 452 DRL agent in selecting piece-wise constant temperature distributions over n_s identical segments, 453 each of which allows only two pre-determined states referred to as hot or cold. This is intended 454 to reduce the complexity and the computational resources, as large/continuous action spaces are 455 known to be challenging for the convergence of RL methods [28, 75]. Simply put, the network 456 action output consists of n_s values $\hat{T}_{hk \in \{1...n_s\}} = \pm \Delta T_{max}$, mapped into the actual fluctuations 457 according to 458

$$\tilde{T}_{hk} = \frac{\hat{T}_{hk} - \langle \hat{T}_{hk} \rangle}{\max_l \{1, \frac{|\hat{T}_{hl} - \langle \hat{T}_{hl} \rangle|}{\Delta T_{max}} \}},$$
(35)

to fulfill the zero-mean and upper bound constraints.³ Ultimately, the agent receives the reward $r_t = -Nu$ to minimize the space averaged heat flux at the hot sidewall.

All results reported herein are for $\Delta T_{max} = 0.75$ (so the hot temperature varies in the range [0.25; 1.75]) and $n_s = 10$ segments, as [39] report that $n_s = 20$ was computationally too demanding for their case, and that $n_s = 5$ yielded poor control efficiency. The agent is a fully-connected network with two hidden layers, each holding 2 neurons. The resolution process uses 8 environments and 2 steps mini-batches to update the network for 32 epochs, with learning rate 5×10^{-3} , and PPO loss clipping range $\epsilon = 0.2$.

467 4.3. Results

For this case, 120 episodes have been run, each of which follows the exact same procedure as 468 above and performs 60 iterations with time step $\Delta t = 0.5$ to march the zero-initial condition to 469 steady state. This represents 960 simulations, each of which is performed on 4 cores and lasts 470 20s, hence 5h of total CPU cost. We present in figure 7 representative iso-contours of the steady-471 state temperature and velocity magnitude computed over the course of the optimization. The 472 latter exhibit strong temperature gradients at the hot sidewall, together with a robust steady roll-473 shaped pattern accompanied by a small corner eddy at the upper-left edge of the cavity, whose size 474 and position depends on the specifics of the temperature distribution. The corresponding meshes 475 are displayed in figure 7(c) to stress the ability of the adaptation procedure to handle well the 476 anisotropy of the solution caused by the intrinsic flow dynamics and the discontinuous boundary 477 conditions. 478

We show in figure 8 the evolution of the controlled averaged Nusselt number, whose moving 479 average decreases monotonically and reaches a plateau after about 90 episodes, although we notice 480 that sub-optimal distributions keep being explored occasionally. The optimal computed by averag-481 ing over the 10 latest episodes (hence the 800 latest instant values) is $\langle Nu \rangle^* \sim 0.57$, with variations 482 ± 0.01 computed from the root-mean-square of the moving average over the same interval (which 483 is a simple yet robust criterion to assess qualitatively convergence a posteriori). Interestingly, 484 the optimized Nusselt number is almost twice as small as the purely conductive value (Nu = 1), 485 meaning that the approach successfully alleviates the heat transfer enhancement generated by the 486 onset of convection, although it does not alleviate convection itself, as evidenced by the consistent 487 roll-shaped pattern in figure 9. Similar results are reported in [39], for a different set-up in which 488 the horizontal cavity walls are isothermal and control is applied at the bottom at the cavity (hence 489 a different physics because of buoyancy), albeit with lower numerical and control efficiency since 490 the authors report an optimal Nusselt number Nu ~ 1 using up to 512 DRL environments with 491 learning rate of 2.5×10^{-4} . The reason for such discrepancies probably lies in different ways of 492 achieving and assessing control, as we use single-step PPO to optimize the steady-state Nusselt 493

³Another possible approach would have been to penalize the reward passed to the DRL for those temperature distributions deemed non-admissible (either because the average temperature is non-zero or the temperature magnitude is beyond the threshold). However, this would have made returning admissible solutions part of the tasks the network is trained on (not to mention that non-zero average temperatures amount to a change in the Rayleigh number), which would likely have slowed down learning substantially.

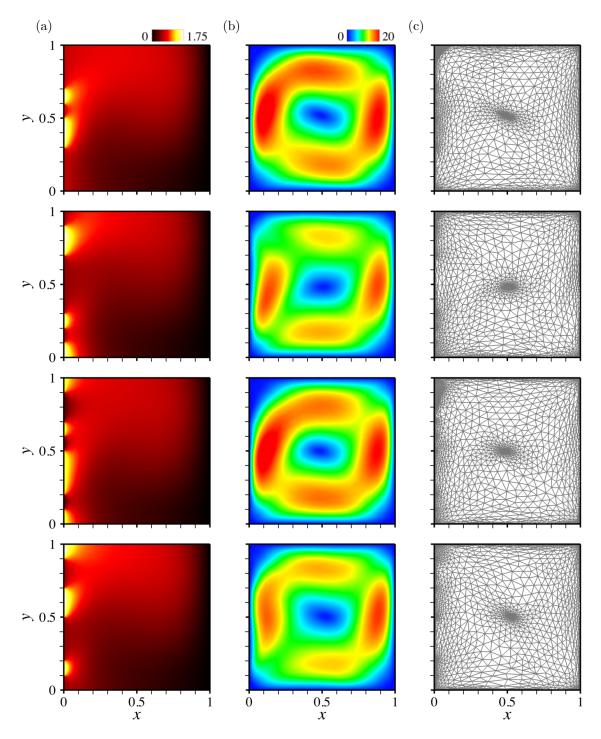


Figure 7: (a,b) Steady-state (a) temperature and (b) velocity magnitude against zero-mean temperature distributed at the left sidewall. (c) Adapted meshes.

⁴⁹⁴ number via a time-independent control, which requires choosing a sidewall temperature, marching ⁴⁹⁵ the controlled solution to steady state, then computing the reward. The problem considered in [39] ⁴⁹⁶ is more intricate, as classical PPO is used to optimize the reward accumulated over time via a ⁴⁹⁷ time-dependent control temperature updated with a certain period scaling with the convection ⁴⁹⁸ time in the cavity (the so-determined optimal control being ultimately time-independent for the ⁴⁹⁹ considered value of Ra, but truly time-dependent for Rayleigh numbers above ~ 10⁵).

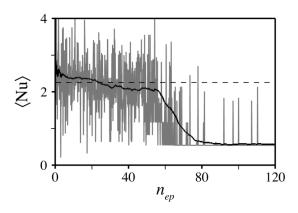


Figure 8: Evolution per learning episode of the instant (in grey) and moving average (in black) Nusselt number. The horizontal dashed line marks the uncontrolled value.

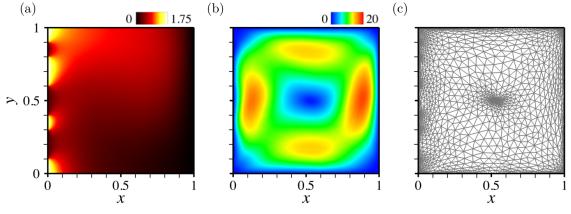


Figure 9: (a,b) Steady-state (a) temperature and (b) velocity magnitude for the optimal zero-mean temperature distribution. (c) Adapted mesh.

500 5. Control of forced convection in 2-D open cavity

501 5.1. Case description

This second test case addresses the control of actual conjugate heat transfer in a model setup 502 for the cooling of a hot solid by impingement of a fluid; see figure 10(a). A Cartesian coordinate 503 system is used with origin at the center of mass of the solid, horizontal x-axis, and vertical y-axis. 504 The solid has rectangular shape with height h and aspect ratio 2:1, and is initially at the hot 505 temperature T_h . It is fixed at the center of a rectangular cavity with height H and aspect ratio 506 4:1, whose walls are isothermal and kept at temperature T_w . The top cavity side is flush with n_j 507 identical holes of width e_i whose distribution is subjected to optimization, each of which models 508 the exit plane of an injector blowing cold air at velocity V_i and temperature T_c , and is identified 509 by the horizontal position of its center $x_{k \in \{1...n_i\}}$. Hot air is released through the cavity sidewalls, 510 blown with two identical exhaust areas of height e_o , and identified by the vertical position of their 511 center $(e_0 - H)/2$. 512

For this case, both buoyancy and radiative heat transfer are neglected ($\psi = 0$ and $\chi = 0$), meaning that temperature evolves as a passive scalar, similar to the mass fraction of a reactant in a chemical reaction. All relevant parameters are provided in Table 2, including the material properties used to model the composite fluid, that yield fluid values of the Reynolds and Prandtl numbers

$$\operatorname{Re} = \frac{\rho V_i e}{\mu} = 200, \qquad \operatorname{Pr} = 2. \tag{36}$$

Note the very high value of the solid to fluid viscosity ratio, that ensures that the velocity is zero in the solid domain and that the no-slip interface condition is satisfied. By doing so, the convective

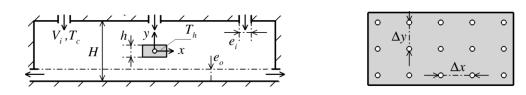


Figure 10: (a) Schematic of the 2-D forced convection set-up. (b) Sensors positions in the solid domain.

 Н	h	e_i	e_0	V_i	T_w	T_c	T_h	μ	ρ	λ	c_p	
1	0.2	0.2	0.2	1	10	10	150	0.001	1	0.5	1000	Fluid
 -	0	0.2	0.2	-	10	10	100				300	

Table 2: Numerical parameters used in the 2-D forced convection problem. All values in SI units, with the exception of temperatures given in Celsius.

terms drop out in the energy equation, that reduces to the pure conduction equation for the solid. The governing equations are solved with no-slip isothermal conditions $\boldsymbol{u} = \boldsymbol{0}$ and $T = T_w$ on $\partial\Omega$, except at the injection exit planes ($\boldsymbol{u} = -V_i \boldsymbol{e}_y$, $T = T_c$), and at the exhaust areas, where a zeropressure outflow condition is imposed ($p = \partial_x u = \partial_x T = 0$). No thermal condition is imposed at the interface, where heat exchange is implicitly driven by the difference in the individual material properties.

526 5.2. Control

(a)

The quantity being optimized is the distribution of the injectors center positions $x_{k \in \{1...n_j\}}$. Several control strategies are assessed in the following, whose ability to manage increasing design complexity translates into less constrained operation when it comes to optimizing a practically meaningful device. In practice, each injector is forced to sit in an interval $[x_k^-; x_k^+]$ whose edge values are determined beforehand of recomputed on the fly (depending on the control strategy), and bounded according to

$$|x_k^{\pm}| \le x_m \,, \tag{37}$$

where we set $x_m = 2H - 0.75e_i$ to avoid numerical issues at the upper cavity edges. The network action output therefore consists of n_j values $\hat{x} \in [-1; 1]$, mapped into the actual positions according to

$$x_k = \frac{x_k^+(\hat{x}_k+1) - x_k^-(\hat{x}_k-1)}{2} \,. \tag{38}$$

In order to compute the reward passed to the DRL, we distribute uniformly 15 probes in the solid domain, into $n_x = 5$ columns and $n_y = 3$ rows with resolutions $\Delta x = 0.09$ and $\Delta y = 0.075$, respectively; see figure 10(b). Selected tests have been carried out to check that the outcome of the learning process does not change using $n_y = 5$ rows of $n_x = 5$ probes (not shown here). The magnitude of the tangential heat flux is estimated by averaging the norm of the temperature gradient over all columns and rows, i.e., *i*-th column (resp. the *j*-th row) as

$$\langle ||\nabla_{\parallel}T||\rangle_{i} = \frac{2}{n_{y}-1} |\sum_{j\neq 0} \operatorname{sgn}(j)||\nabla T||_{ij}|, \qquad \langle ||\nabla_{\parallel}T||\rangle_{j} = -\frac{2}{n_{x}-1} |\sum_{i\neq 0} \operatorname{sgn}(i)||\nabla T||_{ij}|, \quad (39)$$

where subscripts i, j and ij denote quantities evaluated at $x = i\Delta x$, $y = j\Delta y$ and $(x, y) = (i\Delta x, j\Delta y)$, respectively, and symmetrical numbering is used for the center probe to sit at the intersection of the zero-th column and row. The numerical reward $r_t = -\langle ||\nabla_{\parallel}T||\rangle$ fed to the DRL agent deduces ultimately by averaging over all rows and columns, to give

$$\langle ||\nabla_{\parallel}T||\rangle = \frac{1}{n_x + n_y} \sum_{i,j} \langle ||\nabla_{\parallel}T||\rangle_i + \langle ||\nabla_{\parallel}T||\rangle_j , \qquad (40)$$

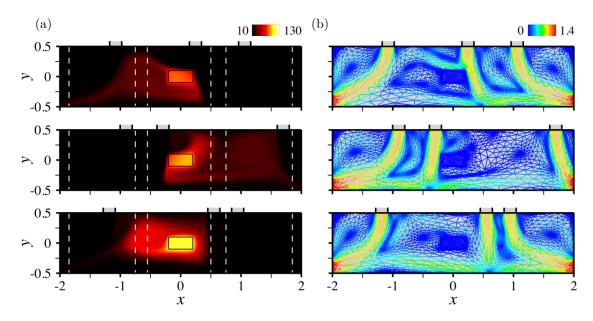


Figure 11: (a) Steady-state temperature against arrangements of 3 injectors, with admissible values under the fixed domain decomposition strategy S_1 delimited by the dashed lines. (b) Adapted meshes colored by the magnitude of velocity.

which especially yields $r_t = 0$ for a perfectly homogeneous cooling.

All results reported in the following are for $n_j = 3$ injectors. The agent is a fully-connected network with two hidden layers, each holding 2 neurons. The resolution process uses 8 environments and 2 steps mini-batches to update the network for 32 epochs, with learning rate set to 5×10^{-3} , and PPO loss clipping range to $\epsilon = 0.3$.

551 5.3. Results

552 5.3.1. Fixed domain decomposition strategy

We consider first the so-called fixed domain decomposition strategy S_1 in which the top cavity wall is split into n_j equal subdomains, and each injector is forced to sit in a different subdomain (a somehow heavily constrained optimization problem if n_j is not to small, relevant for cases where the design is rigid and the practitioner has limited freedom to act). The edge values for the position x_k of the k-th injector read

$$x_{k}^{-} = -x_{m} + (k-1)\frac{2x_{m} + e_{i}}{n_{j}}, \qquad x_{k}^{+} = x_{k}^{-} + \frac{2x_{m} - (n_{j} - 1)e_{i}}{n_{j}}.$$
(41)

It can be checked that $x_k^- = x_{k-1}^+ + e_i$, so, it is possible to end up with two side-by-side injectors, 558 which is numerically equivalent to having $n_j - 1$ injectors, $n_j - 2$ of width e_i plus one of width $2e_i$. 559 For this case, 60 episodes have been run, each of which performs 1500 iterations with time step 560 $\Delta t = 0.1$ to march the same initial condition (consisting of zero velocity and uniform temperature, 561 except in the solid domain) to steady state, using the level set, velocity and temperature as multiple-562 component criterion to adapt the mesh (initially pre-adapted using the sole level set) every 5 563 time steps under the constraint of a fixed number of elements $n_{el} = 15000$. This represents 480 564 simulations, each of which is performed on 8 cores and lasts 10mn, hence 80h of total CPU cost. 565

It is out of the scope of this work to analyze in details the many flow patterns that develop when 566 the blown fluid travels through the cavity. Suffice it to say that the outcome depends dramatically 567 on the injectors arrangement, and features complex rebound phenomena (either fluid/solid, when 568 a jet impinges on the cavity walls or on the workpiece itself, or fluid/fluid, when a deflected jet 569 meets the crossflow of another jet), leading to the formation of multiple recirculation varying in 570 number, position and size. Several such cases are illustrated in figure 11 via iso-contours of the 571 steady-state temperature distributions, together with the corresponding adapted meshes colored by 572 the magnitude of velocity to illustrate the ability of the numerical framework to capture accurately 573 all boundary layers and shear regions via extremely stretched elements. 574

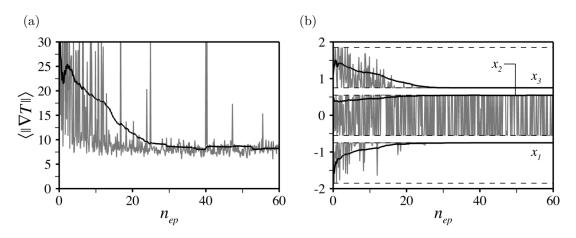


Figure 12: (a) Evolution per learning episode of the instant (in grey) and moving average (in black) rewards under the fixed domain decomposition strategy S_1 . (b) Same as (a) for the injectors center positions, with admissible values delimited by the dashed lines.

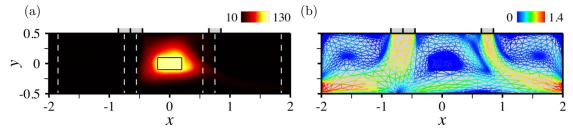


Figure 13: Same as figure 11 for the optimal arrangement of 3 injectors under the fixed domain decomposition strategy S_1 .

One point worth mentioning is that the individual position signals are best suited to draw robust 575 quantitative conclusion, as there is noise in the reward signal shown in figure 12(a). We believe 576 the issue to be twofold: on the one hand, the reward is approximated from point-wise temperature 577 data (similar to experimental measurements) that are more sensitive to small numerical errors 578 (e.g., the interpolation error at the probes position) than an integral quantity. On the other hand, 579 the mesh adaptation procedure is not a deterministic process, as the outcome depends on the 580 processors and number of processors used, and any initial difference propagates over the course 581 of the simulation because the meshes keep being adapted dynamically. In return, two exact same 582 control parameters can thus yield different rewards on behalf of different interpolation errors at 583 the probes position. This likely slows down learning and convergence, but we show in figure 12(b)584 that the moving average distribution does converge to an optimal arrangement after roughly 25 585 episodes. The latter consists of an injector at the right-end of the left subdomain $(x_1^{\star} = -0.75)$ 586 and two side-by-side injectors sitting astride the center and right subdomains ($x_2^* = 0.55$ and 587 $x_3^{\star} = 0.75$), that enclose the workpiece in a double-cell recirculation; see figure 13. These values 588 have been computed by averaging the instant positions of each injector over the 10 latest episodes, 589 with variations ± 0.002 computed from the root-mean-square of the moving average over the same 590 interval, a procedure that will be used consistently to assess convergence for all cases reported in the 591 following. The efficiency of the control itself is estimated by computing the magnitude of tangential 592 heat flux averaged over the same interval, found to be $\langle ||\nabla_{\parallel}T|| \rangle^{\star} \sim 8.3$. Note, the position x_2^{\star} is 593 actually obtained by averaging the absolute value of the instant position x_2 (although the true, 594 signed value is depicted in the figure), as the center injector keeps oscillating between two end 595 positions ± 0.55 on behalf of reflectional symmetry with respect to the vertical centerline. 596

597 5.3.2. Follow-up strategy

⁵⁹⁸ A less constrained problem is considered here using the so-called follow-up strategy S_2 , in which ⁵⁹⁹ all injectors are distributed sequentially the ones with respect to the others. The corresponding

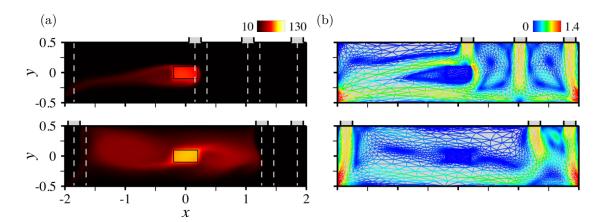


Figure 14: (a) Steady-state temperature against arrangements of 3 injectors, with admissible values under the follow-up strategy S_2 delimited by the dashed lines. (b) Adapted meshes colored by the magnitude of velocity.

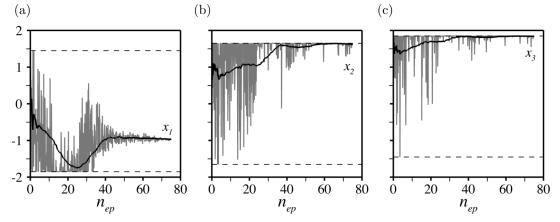


Figure 15: Evolution per learning episode of the instant (in grey) and moving average (in black) injectors center positions under the follow-up strategy S_2 , with admissible values delimited by the dashed lines.

600 edge values

$$x_1^- = -x_m, \qquad x_1^+ = x_m - (n_j - 1)e_i, \qquad (42)$$

$$x_k^- = x_{k-1}^+ + e_i, \qquad x_k^+ = x_m - (n_j - k)e_i, \qquad (43)$$

readily express that the k-th injector is forced to sit between the k-1-th one and the upper-right cavity edge while leaving enough space to distribute the remaining $n_j - k$ injectors, which increases the size of the control parameter space while again leaving the possibility for side-by-side injectors (since $x_k^- = x_{k-1}^+ + e_i$ by construction). 75 episodes have been run for this case following the exact same procedure as above, i.e., marching the zero-initial condition in time up to t = 150 with $\Delta t = 0.1$, hence 600 simulations, each of which is performed on 8 cores and lasts 10mn, hence 100h of total CPU cost.

The computed flow patterns closely resemble those obtained under the previous fixed domain 608 decomposition strategy, although figure 14 exhibits increased dissymmetry when two or more in-609 jectors move simultaneously to the same side of the cavity. We show in figure 15 that the moving 610 average distribution converges after roughly 60 episodes, with the optimal arrangement consisting 611 of one injector roughly midway between the left cavity sidewall and the workpiece ($x_1^{\star} = -0.96$), 612 and two side-by-side injectors at the right end of the cavity $(x_2^* = 1.65 \text{ and } x_3^* = 1.85)$. The 613 variations over the same interval are by ± 0.006 ; see also figure 16 for the corresponding flow pat-614 tern. Convergence here is much slower than under S_1 , as the search for an optimal is complicated 615 by the fact that all injector positions are interdependent the ones on the others and it is up to the 616 network to figure out exactly how. Another contingent matter is that the agent initially spans a 617 fraction of the control parameter space because the large values of x_1 considered limit the space 618 available to distribute the other two injectors. This is all the more so as such configurations turn 619

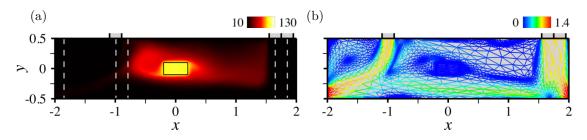


Figure 16: Same as figure 14 for the optimal arrangement of 3 injectors under the follow-up strategy S_2 .

to be far from optimality, for instance the magnitude of tangential heat flux is $\langle ||\nabla_{\parallel}T|| \rangle \sim 41.3$ for $x_1 = 1.45, x_2 = 1.65$ and $x_3 = 1.85$, but $\langle ||\nabla_{\parallel}T|| \rangle^* \sim 6.3$ at optimality. The latter value is smaller than the optimal achieved under S_1 , consistently with the fact that all positions spanned under S_1 are admissible under S_2 , hence the S_1 optimal is expected to be a S_2 sub-optimal.

624 5.3.3. Free strategy

We examine now a third strategy S_3 referred to as the free strategy, in which all injectors are independent and free to move along the top cavity wall (a mildly constrained optimization problem, relevant for cases where the design is flexible and the practitioner has great freedom to act). The edge values for the position x_k of the k-th injector read

$$x_k^- = -x_m, \qquad x_k^+ = x_m, \qquad (44)$$

⁶²⁹ so two injectors can end up side-by side and even overlapping one another if $|x_l - x_m| < e_i$. If ⁶³⁰ so, we implement a single injector of width $e_i + |x_l - x_m|$ and maintain the blowing velocity (not ⁶³¹ the flow rate) for the purpose of automating the set-up design process, meaning that having n_j ⁶³² injectors, two of which overlap exactly (i.e., $|x_l - x_m| = 0$) is rigorously equivalent to having $n_j - 1$ ⁶³³ injectors. 60 episodes have been run for this case following the exact same procedure as above.

All flow patterns are reminiscent of those obtained under the previous fix decomposition S_1 634 and follow-up S_2 strategies, even when two injectors overlap; see figure 17. Other than that, we 635 show in figures 18 that the moving average distribution converges to an optimal consisting of two 636 injectors almost perfectly overlapping one another at the left end of the cavity $(x_1^{\star} = -1.85 \text{ and}$ 637 $x_2^* = -1.82$), and a third injector at the right end of the cavity ($x_3^* = 1.85$). The variations over 638 the same interval are by ± 0.007 , and the associated flow pattern shown in figure 19 is symmetrical 639 and features two large recirculation regions on either side of the workpiece. Convergence occurs 640 after roughly 40 episodes, i.e., faster than under S_2 (consistently with the fact that there is no 641 need to learn anymore about how the network outputs depend the ones on the others) but slower 642 than under S_1 (consistently with the fact that the size of the control parameter space has increased 643 substantially). It is worth noticing that the system is invariant by permutations of the network 644 outputs, meaning that there exist $2^{n_j} - 2$ distributions (hence 6 for $n_j = 3$) associated with the 645 same reward. Nonetheless, a single optimal is selected, which is essentially fortuitous since the 646 agent does not learn about symmetries under the optimization process (otherwise S_1 would have 647 similarly selected a single optimal). The magnitude of tangential heat flux is $\langle || \nabla_{\parallel} T || \rangle^{\star} \sim 11.2$ at 648 optimality, i.e., larger than that achieved under S_2 . This can seem surprising at first, because all 649 positions spanned under S_2 are admissible under S_3 , and the S_2 optimal is thus expected to be a 650 S_3 sub-optimal. However, the argument does not hold here because the overlap in the S_3 optimal 651 reduces the flow rate to that of a two-injectors set-up, so the comparison should be with the S_2 652 optimal with $n_j = 2$. 653

654 5.3.4. Inverse strategy

Finally, we propose here to make the most of the numerical framework flexibility to solve a different optimization problem consisting in selecting first an injector distribution, then in finding the position x_0 of the solid center of mass minimizing the magnitude of tangential heat flux (which is relevant for cases where the practitioner simply cannot act on the design). The so-called inverse strategy S_4 considered herein features two injectors at each end of the cavity ($x_1 = -1.85$ and $x_2 = 1.85$), identical to the optimal arrangement of 3 injectors under the free strategy S_3 . The

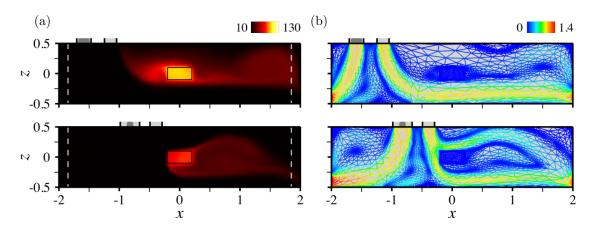


Figure 17: (a) Steady-state temperature against arrangements of 3 injectors, with admissible values under the free strategy S_3 delimited by the dashed lines and overlaps marked by the dark grey shade. (b) Adapted meshes colored by the magnitude of velocity.

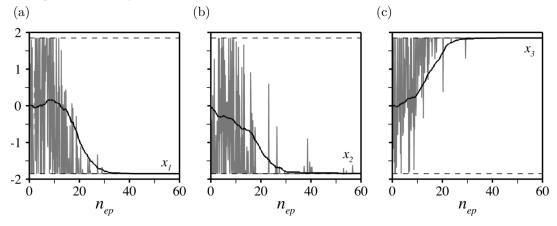


Figure 18: Evolution per learning episode of the instant (in grey) and moving average (in black) injectors center positions under the free strategy S_3 , with admissible values delimited by the dashed lines.

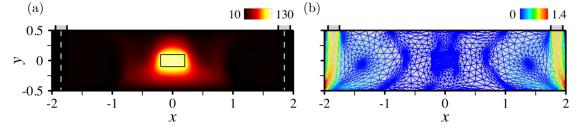


Figure 19: Same as figure 17 for the optimal arrangement of 3 injectors under the free strategy S_3 .

center of mass can take any value in $[-x_{0m}; x_{0m}]$ where we set $x_{0m} = 2(H-h)$ to avoid numerical issues at the sidewalls. The same coordinate system as above is used, but with reference frame attached to the cavity, not the moving solid (hence all results obtained under the previous strategies pertain to $x_0 = 0$ in the new system).

⁶⁶⁵ A total of 60 episodes have been run for this case using the exact same DRL agent, the only ⁶⁶⁶ difference being in the network action output, now made up of a single value $\hat{x}_0 \in [-1; 1]$, mapped ⁶⁶⁷ into the actual position using

$$x_0 = x_{0m} \hat{x}_0 \,. \tag{45}$$

⁶⁶⁸ A large variety of flow patterns is obtained by doing so, that closely resemble those computed under ⁶⁶⁹ the previous strategies, only the outcome is now also altered by the width of the gap between the ⁶⁷⁰ cavity sidewalls and the workpiece, as illustrated in figure 20. We show in figure 21 that the position ⁶⁷¹ of the solid center of mass converges to an optimal $x_0^* = 0.42$ (the variations over the same interval

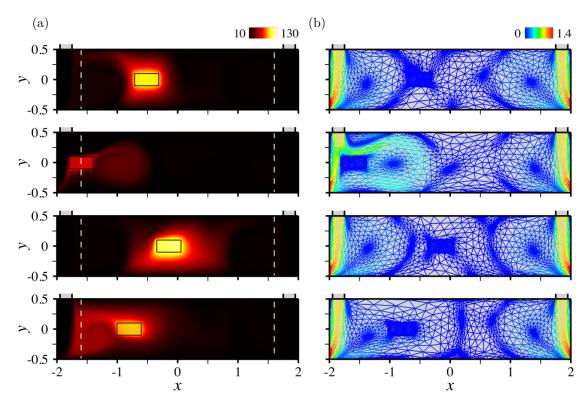


Figure 20: (a) Steady-state temperature against solid center of mass position, with admissible domains under the inverse strategy S_4 marked by the dashed lines. (b) Adapted meshes colored by the norm of velocity.

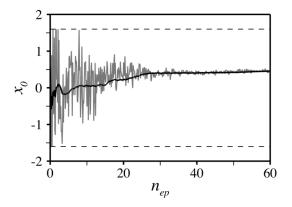


Figure 21: Evolution per learning episode of the instant (in grey) and moving average (in black) center of mass positions under the inverse strategy S_4 , with admissible values delimited by the dashed lines.

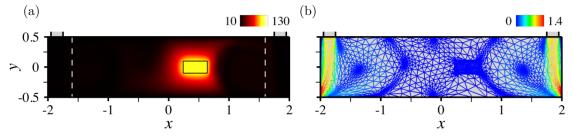


Figure 22: Same as figure 20 for the optimal center of mass position under the inverse strategy S_4 .

⁶⁷² being by ± 0.005), the associated magnitude of tangential heat flux $\langle ||\nabla_{\parallel}T|| \rangle^* \sim 4.1$, being smaller ⁶⁷³ than that achieved under S_3 using a centered workpiece. The fact that the optimal position is ⁶⁷⁴ offset from the vertical centerline is a little surprising at first, because intuition suggests that the

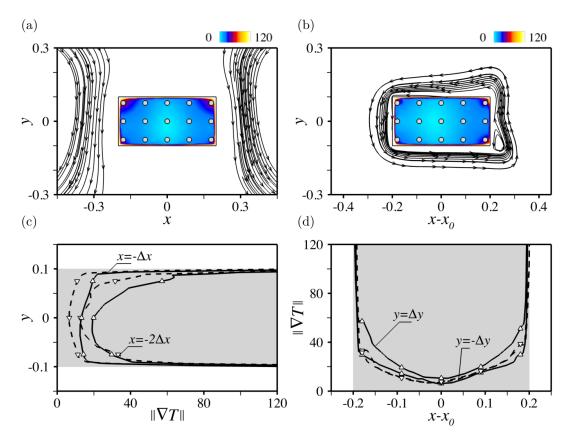


Figure 23: (a,b) Norm of the temperature gradient in the solid domain with superimposed streamlines of the underlying velocity field, as computed for (a) $x_0 = 0$, and (b) $x_0^* = 0.45$, i.e., the optimal position selected under the inverse strategy S_4 . (c) Cuts along the two leftmost columns of probes. The solid and dashed lines refer to $x_0 = 0$ and $x_0^* = 0.42$, respectively, and the symbols mark the probe values. (d) Same as (c) for cuts along the lower and upper rows of columns.

simplest way to achieve homogeneous heat transfer is by having symmetrically distributed injectors. 675 Nonetheless, examining carefully the norm of the temperature gradient in the solid domain shows 676 that $x_0 = 0$ achieves close to perfect horizontal symmetry but vertical asymmetry, owing to the 677 formation of two large-scale, small velocity end vortices entraining heat laterally downwards; see 678 figure 23(a). Conversely, for $x \sim x_0^*$, the workpiece it almost at the core of the closest recirculation 679 region, hence the surrounding fluid particles have small velocities and wrap almost perfectly around 680 its surface, as illustrated in figure 23(b). This restores excellent vertical symmetry, as evidenced 681 by relevant cuts along the two leftmost columns of probes in figure 23(c), and along the lower and 682 upper rows in figure 23(d), which explains the improved the reward. 683

684 5.4. Discussion

Figure 24 reproduces the optimal temperature distributions computed under the various strate-685 gies considered above. For benchmarking purposes, we also provide in table 3 relevant convergence 686 data computed over the 10 latest episodes. To recap, the most homogeneous cooling is achieved 687 under the follow-up strategy S_2 , but the DRL agent seems more easily trained under the fixed 688 decomposition domain strategy S_1 and the free strategy S_3 . Another interesting point is the ex-689 tent to which the workpiece is actually cooled, for which S_2 seems more relevant, on behalf of the 690 dissymmetry in the left and right flow rates that creates order one velocities at the bottom of the 691 cavity. This stresses S_2 as a possible compromise to achieve efficient and homogeneous cooling, 692 although a true optimal with this regard can be computed rigorously by applying the same ap-693 proach to compound functionals weighing, e.g., the magnitude of the tangential heat flux and the 694 solid center temperature (which we defer to future work). 695

These results provide a basis for future self-assessment of the method and identifies potential for improvement regarding the convergence efficiency. The approach can certainly benefit from

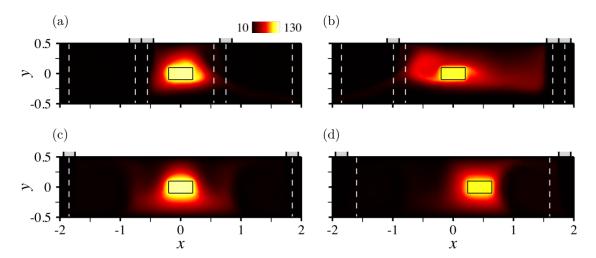


Figure 24: (a-c) Optimal arrangements of 3 injectors under the (a) fixed decomposition domain strategy S_1 , (b) follow-up strategy S_2 and (c) free strategy S_3 . (b) Optimal position of the workpiece under the inverse strategy S_4 .

	n_j	n_{ep}	x_0	x_1	x_2	x_3	$\langle \nabla_{\parallel}T \rangle$
S_1	3	60	0	-0.75	± 0.55	0.75	8.3
S_2	3	75	0	-0.96	1.65	1.85	6.3
S_3	3	60	0	-1.85	-1.82	1.85	11.2
S_4	2	60	0.42	-1.85	1.85	—	4.1

Table 3: Numerical data for the optimal arrangements computed under strategies S_{1-4} . All values computed by averaging the instant signal over the 10 latest learning episodes.

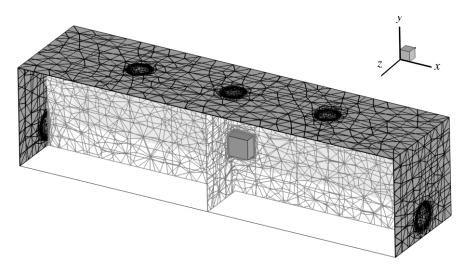


Figure 25: Schematic of the 3-D forced convection set-up.

a fine tuning of the reward computation, as having sufficient spatial resolution on the relevant state of the system is an obvious requirement to allow a successful control. Adjusting the trade-off between exploration and exploitation is also worth consideration to better handle the existence of multiple global optima (whether they stem from symmetries of from the topology of the reward itself) which could be done using non-normal probability density functions.

 Η	h	d_i	d_0	δ_0	V_i	T_w	T_c	T_h	μ	ρ	λ	c_p	
1	0.2	0.2	0.24	0.16	1	10	10	150	0.01	1	0.5	1000	Fluid
-	0.2	0	0.21	0.10	-	10	10	100	1000	100	15	300	Solid

Table 4: Numerical parameters used in the 3-D forced convection problem. All values in SI units, with the exception of temperatures given in Celsius.

703 6. Extension to 3-D forced convection

704 6.1. Case description

The model cooling set up considered in section 5 is extended here to 3-D to assess the extent to 705 which the approach carries over to three-dimensional conjugate heat transfer. The main differences 706 between 2-D and 3-D are as follows: a Cartesian coordinate system is used with origin at the center 707 of mass of the solid, horizontal x-axis, vertical y-axis, and the z-axis completes the direct triad; see 708 figure 25. The solid is a rectangular prism with aspect ratio 2:1:1, and is fixed at the center of a 709 rectangular cavity with height H and aspect ratio 4:1:1. We consider n_i circular-shaped injectors 710 with diameter d_i , whose exit planes are forced to be symmetrical with respect to z = 0, hence each 711 injector is identified by the horizontal position of its center $x_{k \in \{1...n_i\}}$. We also use circular-shaped 712 exhaust areas with diameter d_o , offset by a distance δ_o from the bottom of the cavity, and whose 713 exit planes are also symmetrical with respect to z = 0, hence each exhaust area is identified by the 714 vertical position of its center $(d_0 + \delta_o - H)/2$. The governing equations are solved with the exact 715 same boundary conditions as in section 5. All parameters are provided in Table 4, including the 716 material properties used to model the composite fluid, that yield fluid values of the Reynolds and 717 Prandtl numbers 718

Re =
$$\frac{\rho V_i d_i}{\mu} = 20$$
, Pr = 20. (46)

719 6.2. Control strategy

We keep here the same control objective and compute the reward fed to the DRL from 45 probes arranged symmetrically into $n_z = 3$ transverse layers with resolution $\Delta z = 0.075$, each of which distributes uniformly 15 probes into $n_x = 5$ columns and $n_y = 3$ rows with resolutions $\Delta x = 0.09$ and $\Delta y = 0.075$. In practice, the 3-D reward is simply the average over z of the 2-D reward defined in section 5, hence $r_t = -\langle ||\nabla_{\parallel}T|| \rangle$ with

$$\langle ||\nabla_{\parallel}T||\rangle = \frac{1}{(n_x + n_y)n_z} \sum_{i,j,k} \langle ||\nabla_{\parallel}T||\rangle_{ik} + \langle ||\nabla_{\parallel}T||\rangle_{jk} , \qquad (47)$$

725 with

$$\langle ||\nabla_{\parallel}T||\rangle_{ik} = \frac{2}{n_y - 1} |\sum_{j \neq 0} \operatorname{sgn}(j)||\nabla T||_{ijk}|, \qquad \langle ||\nabla_{\parallel}T||\rangle_{jk} = -\frac{2}{n_x - 1} |\sum_{i \neq 0} \operatorname{sgn}(i)||\nabla T||_{ijk}|, \quad (48)$$

and the subscripts ik, jk and ijk denote quantities evaluated at $(x, z) = (i\Delta x, k\Delta z)$, $(y, z) = (i\Delta y, k\Delta z)$ and $(x, y, z) = (i\Delta x, j\Delta y, k\Delta z)$, respectively.

All results reported in the following are for $n_j = 3$ injectors. The edge values needed to map the 728 network action output into the actual injectors positions deduce straightforwardly from (41)-(44) 729 substituting the diameter d_i of the 3-D injectors for the length e_i of the 2-D injectors. The same 730 DRL agent is used, that consists of two hidden layers, each holding 2 neurons, and the resolution 731 process uses 8 environments and 2 steps mini-batches to update the network for 32 epochs. Each 732 environment performs 1250 iterations with time step $\Delta t = 0.1$ to march the same initial condition 733 (consisting of zero velocity and uniform temperature, except in the solid domain) to steady state, 734 using the level set, velocity and temperature as multiple-component criterion to adapt the mesh 735 (initially pre-adapted using the sole level set) every 10 time steps under the constraint of a fixed 736

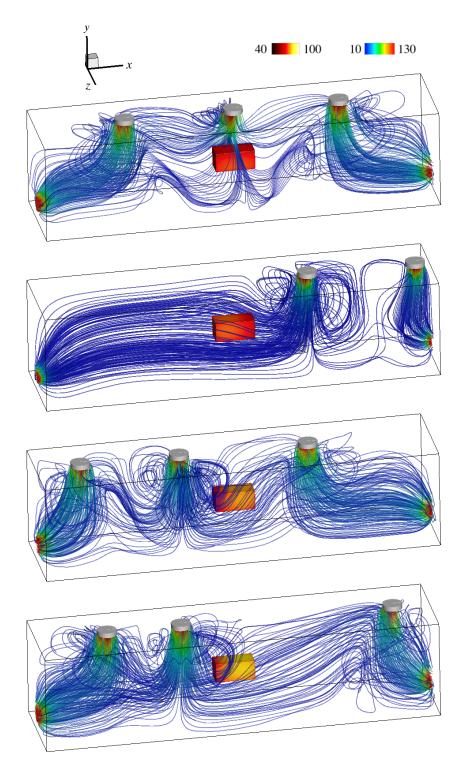


Figure 26: Representative steady-state temperature distributions at the solid/fluid interface together with 3-D streamlines colored by the magnitude of velocity.

⁷³⁷ number of elements $n_{el} = 120000$. This is likely insufficient to claim true numerical accuracy, but ⁷³⁸ given the numerical cost (320 3-D simulations per strategy, each of which is performed on 8 cores ⁷³⁹ and lasts 2h30, hence 800h of total CPU cost), we believe this is a reasonable compromise to assess ⁷⁴⁰ feasibility while producing qualitative results to build on.

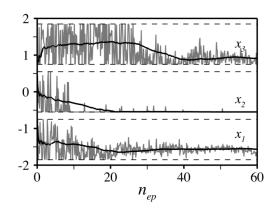


Figure 27: Evolution per learning episode of the instant (in grey) and moving average (in black) injectors center positions under the three-dimensional fixed domain decomposition strategy S_1 , with admissible values delimited by the dashed lines.

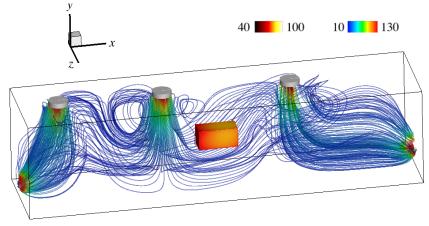


Figure 28: Optimal 3 injector arrangement under the three-dimensional fixed decomposition domain strategy S_1 .

741 6.3. Results

Only the fixed domain decomposition S_1 strategy (in which the top cavity wall is split into n_j equal subdomains and each injector is forced to sit in a different subdomain) and the free S_3 strategy (in which the injectors are entirely independent and free to move along the top cavity wall) are considered here to save computational resources, as learning has been seen to be slower in 2-D under the follow-up S_2 strategy.

A total of 60 episodes have been run under the fixed domain decomposition strategy S_1 . Several 747 representative flow patterns computed over the course of optimization are shown in figure 26 via 748 iso-contours of the steady-state temperature at the fluid-solid interface and 3-D streamlines colored 749 by the magnitude of velocity, to put special emphasis on transverse inhomogeneities and display 750 the increased degree of complexity due to the formation of large-scale horseshoe vortices wrapped 751 around the nozzle jets. We show in figure 27 that the distribution slowly converges to an optimal 752 arrangement consisting of one injector at the left end of the left subdomain $(x_1^* = -1.63)$, another 753 one at the left end of the center subdomain $(x_2^{\star} = -0.55)$, and a third one at the left end of 754 the right subdomain $(x_3^{\star} = 0.87)$, as has been determined by averaging the instant positions 755 of each injector over the latest 10 learning episodes, with variations by roughly ± 0.04 computed 756 from the root-mean-square of the moving average over the same interval. This is larger by one 757 order of magnitude than the variations reported in 2-D, as the agent keeps exploring slightly 758 sub-optimal positions of the lateral injectors, which likely simply reflects the challenging nature 759 of performing three-dimensional optimal control. The 3-D S_1 optimal somehow resemble its 2-D 760 counterpart, namely the center injector is at the exact same position, while the lateral injectors 761 (especially the leftmost one) have been pushed towards the cavity sidewalls. The associated flow 762 pattern is reported in figure 28. The associated optimal reward computed over the same interval 763

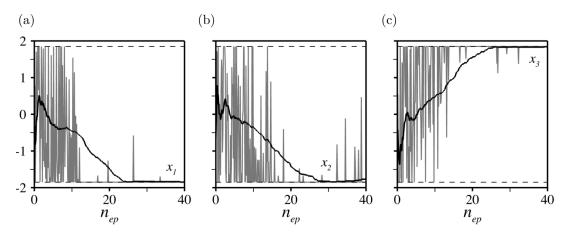


Figure 29: Evolution per learning episode of the instant (in grey) and moving average (in black) injectors center positions under the three-dimensional free strategy S_3 , with admissible values delimited by the dashed lines.

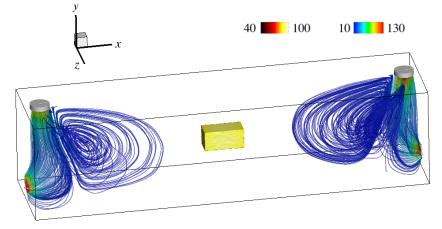


Figure 30: Optimal 3 injector arrangement under the three-dimensional free strategy S_3 .

	n_j	n_{ep}	x_0	x_1	x_2	x_3	$\langle \nabla_{\parallel}T \rangle$
$\overline{S_1}$	3	60	0	-1.63	-0.55	0.87	19.5
S_3	3	40	0	-1.83	-1.82	1.83	4.7

Table 5: Numerical data for the optimal arrangements computed in three-dimensions under strategies S_1 and S_3 . All values computed by averaging the instant signal over the 10 latest learning episodes.

⁷⁶⁴ is $\langle ||\nabla_{\parallel}T|| \rangle^{\star} \sim 19.5$, i.e. twice as large than in 2-D, although it is difficult to compare further ⁷⁶⁵ because of the difference in the Reynolds and Prandtl number.

Another 40 episodes have been run under the free strategy S_3 , for which the results are al-766 most identical to their 2-D counterparts, as the distribution converges in figure 29 to an optimal 767 arrangement consisting of two overlapping injectors at the left end of the cavity ($x_1^{\star} = -1.83$ and 768 $x_2^{\star} = -1.82$), and a third injector at the right end ($x_3^{\star} = 1.83$), with variations by with ± 0.01 769 for the lateral injectors, but ± 0.03 for the center injector, for which the agent keeps occasionally 770 exploring sub-optimal positions. The corresponding flow pattern shown in figure 30 is thus again 771 symmetrical with two large, 3-D recirculation regions on either side of the workpiece. The asso-772 ciated optimal reward computed over the same interval is $\langle ||\nabla_{\parallel}T|| \rangle^{\star} \sim 4.7$ substantially smaller 773 than that achieved under the 3-D S_1 strategy, which again demonstrates the feasibility to improve 774 performances by allowing overlaps. All relevant numerical data are reported in table 5 for the sake 775 of completeness. 776

777 7. Conclusion

Optimization of conjugate natural and forced heat transfer systems is achieved here training a fully connected network with a novel single-step PPO deep reinforcement algorithm, in which it gets only one attempt per learning episode at finding the optimal. The numerical reward fed to the network is computed with a finite elements CFD environment solving stabilized weak forms of the coupled Navier–Stokes and heat equations with a combination of variational multi-scale modeling, immerse volume method, and multi-component anisotropic mesh adaptation.

Convergence is assessed by alleviating the natural convection induced enhancement of heat 784 transfer in a two-dimensional, differentially heated square cavity controlled by piece-wise constant 785 fluctuations of the sidewall temperature. The approach is also relevant to forced convection prob-786 lems, as single-step PPO shows capable of improving the homogeneity of temperature across the 787 surface of two and three-dimensional hot workpieces under impingement cooling. Several control 788 strategies are considered, in which the position of multiple cold air injectors is optimized relative 789 to a fixed workpiece position, each of which mimics a different levels of design constraint. The 790 flexibility of the numerical framework also allows solving the inverse problem, i.e., optimizing the 791 workpiece position relative to a fixed injector distribution, which is relevant in situations where the 792 design cannot be changed. The approach is beneficial in two important respects: first, it is effi-793 cient, even though the parameter spaces are large and it may be costly to identify optimal control 794 parameters from simple parametric searches. Second, and more significantly, it is capable of deter-795 mining additional optimal configurations, as the results of the inverse problem under symmetrical 796 actuation indicate that the workpiece is best positioned offset from the symmetry axis, which had 797 not been anticipated. Such results clearly stress that single-step PPO (and DRL in general) can 798 be effective to explore and discover new solutions from unforeseen parameter combinations. 799

Fluid dynamicists have just begun to gauge the ability of DRL to design optimal control strate-800 gies. The efforts for developing single-step PPO are ongoing and remain at an early stage, so we 801 do not expect the approach to compete right away with more established methods, for instance 802 Evolution strategies (ES), a popular class of algorithms imitating principles of organic evolution 803 processes as rules for black-box optimum seeking. ES rely on a stochastic description of the vari-804 ables to optimize, i.e., they consider probability density functions, not deterministic variables. 805 Simply put, at each generation (or iteration) new candidate solutions are sampled isotropically 806 by variation of the current parental individuals according to a multivariate normal distribution. 807 Recombination and mutation transformations are applied (that amount respectively to changing 808 the mean and adding a random, zero-mean perturbation), after which the individuals with the 809 highest cost function are selected to become the parents in the next generation. Improved variants 810 include the covariance matrix adaptation evolution strategy (CMA-ES), that speeds up conver-811 gence by updating its full covariance matrix (which amounts to learning a second-order model of 812 the objective function). In present form, single-step PPO can be thought as an evolutionary-like 813 algorithm with simpler heuristics (i.e., without an evolutionary update strategy, as the optimal 814 model parameters are learnt via gradient ascent), so it is our guess that the performance should 815 be comparable to that of standard ES algorithms with isotropic covariance matrix. Besides con-816 solidating the acquired knowledge, future research should thus aim at improving efficiency (by 817 818 fine-tuning the hyper parameters, or using pre-trained deep learning models) and convergence (by coupling with a surrogate model trained on-the-fly, using non-normal probability density functions, 819 or modifying the balance between exploration and exploitation, as PPO prevents large updates of 820 the policy to avoid the issue of performance collapse). For complex configurations representative of 821 industrial applications, the implementation of properly designed numerical rewards (under partial 822 state information) and noise reduction techniques is another issue that deserves consideration, as 823 pointed out in [38]. 824

Scope is another key ingredient to keep pushing forward the state of the art. The next step is to tackle more complex test cases exhibiting flow unsteadiness and turbulence, which the CFD environment is perfectly suited to do via a combination of Reynolds-averaged Navier–Stokes modeling [76, 77] and second-order, semi-implicit time discretization [78]. We believe that this will highlight even more clearly the relevance of the methodology, as [42] speculates that DRL should be able to handle chaotic systems without suffering from the shortcomings and limitations of the adjoint method, and it is shown in [39] to outperform a canonical linear proportional-derivative controller in controlling turbulent natural convection. The long-term objective would be to enrich the description of the test cases using multi-physics modeling (e.g., radiative heat transfer, phase transformation) in order to pave the way toward flexible, ready-to-use control of industrially relevant applications, such as thermal comfort for building design or manufacturing processes.

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