



Experimental study of the convection in a thin cylindrical gas layer with imposed bottom and top fluxes and imposed side temperature

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We investigate convection in a thin cylindrical gas layer with an imposed flux at the bottom and a fixed temperature along the side, using a combination of direct numerical simulations and laboratory experiments. The experimental approach allows us to extend by two orders of magnitude the explored range in terms of flux Rayleigh number. We identify a scaling law governing the root-mean-square horizontal velocity and explain it through a dimensional analysis based on heat transport in the turbulent regime. Using particle image velocimetry, we experimentally confirm, for the most turbulent regimes, the presence of a drifting persistent pattern consisting of radial branches, as identified by Rein *et al.* (2023, *J. Fluid Mech.* **977**, A26). We characterise the angular drift frequency and azimuthal wavenumber of this pattern as functions of the Rayleigh number. The system exhibits a wide distribution of heat flux across various time scales, with the longest fluctuations attributed to the branch pattern and the shortest to turbulent fluctuations. Consequently, the branch pattern must be considered to better forecast important wall heat flux fluctuations, a result of great relevance in the context of nuclear safety, the initial motivation for our study.

Key words: convection

1. Introduction

The present study follows up the numerical study of Rein *et al.* (2023a) and is motivated by nuclear safety issues. During a severe accident (SA) in a nuclear power plant, the radioactive fuel and metallic components of the reactor melt, forming a fluid known as corium. The corium relocates from the reactor core to the lower plenum of the reactor vessel, where non-miscible oxidic and metallic phases separate. The oxide phase contains

the majority of radioactive elements. It thus heats from below the less dense and thinner liquid metal phase floating on its surface, which then concentrates the heat towards the vessel wall. This phenomenon is referred to as the ‘focusing effect’ in nuclear safety literature. Understanding heat transfer through the top metal layer is crucial for predicting vessel failure or ensuring its integrity when in-vessel retention is employed as a SA management strategy (Theofanous *et al.* 1997; Carénini *et al.* 2018).

A previous study of the metal layer dynamics (Rein *et al.* 2023a) highlighted the interplay between the Rayleigh–Bénard convection triggered by the bottom heating and the vertical convection resulting from the lateral cooling. Both configurations have been the subject of numerous studies; see, e.g. Ahlers *et al.* (2009b); Chapman & Proctor (1980); Otero *et al.* (2002); Verzicco & Sreenivasan (2008); Johnston & Doering (2009); Fantuzzi (2018); Batchelor (1954); Churchill & Chu (1975); Bejan & Tien (1978); George & Capp (1979); Wells & Worster (2008); Ng *et al.* (2015); Shishkina (2016), respectively. In integral SA codes, such as ASTEC (accident source term evaluation code) developed by the Institut de Radioprotection et de Sûreté Nucléaire (IRNS, Chatelard *et al.* 2014), the entire process of a reactor core meltdown accident is simulated, from initiating events to the release of radioactive materials. Different modules address different aspects of the accident, the corium behaviour in the lower plenum of the vessel being one of them (Carénini *et al.* 2014). In such codes, the focusing effect evaluation is based on a simplified approach proposed by Theofanous *et al.* (1997) (denoted the zero-dimensional (0-D) model), which combines correlations from both Rayleigh–Bénard and natural convection to characterise the fluid by a single, mean surface temperature. It is assumed that the fluid in the bulk is thoroughly mixed and that the vertical heat transfer is symmetrical, meaning that the temperature difference between the bottom and the bulk is equal to the temperature difference between the bulk and the top.

The 0-D model was experimentally tested using a rectangular geometry and water as a simulant for the metal layer. Initially, assumptions regarding a well-mixed fluid and symmetry were verified through the MELAD experiment (Theofanous *et al.* 1997). Further examination was conducted in the BALI-metal experiment, focusing on higher aspect ratios (Bonnet & Seiler 1999). Findings indicated that the initial model might overestimate the focusing effect for shallow layer thickness. Recent HELM (Li *et al.* 2015) and HELM-LR (Li *et al.* 2021) experiments, employing a cylindrical geometry but maintaining water as the simulant, aimed at refining these conclusions for various aspect ratios. Additionally, numerical simulations of the metal layer were carried out (Shams *et al.* 2020) and revealed a significant impact of the fluid properties (water versus metal layer) on the overall phenomenology, particularly due to differences in the Prandtl number, substantially lower for the metal layer compared with water (0.1 versus 7). The lateral heat flux and concentration factor were found to be up to 50 % higher with steel compared with water under similar boundary conditions. The validity of results obtained with water as a simulant for the metallic layer, as well as the derived model for SA integral codes, are therefore questionable. Note that although the Prandtl number in liquid metals is generally lower than 0.1, for example, close to 0.026 in liquid gallium (Grannan *et al.* 2022), the present metal layer is composed of a mixture of zirconium, metallic uranium and steel. In this configuration the expected Prandtl value is in the range [0.07–0.2] (Le Guennic *et al.* 2020). Hence, a representative intermediate value of 0.1 is considered.

Rein *et al.* (2023b) proposed an improvement of the 0-D model, incorporating a description of the radial temperature profile at the top surface based on direct numerical simulations (DNS) and scaling laws (called the one-dimensional (1-D) model in the following). However, the applicability of this 1-D model was only tested numerically up to a flux Rayleigh number $Ra_\phi = 10^8$ (see the exact definition (3.5) below), while the system

can reach up to $Ra_\phi = 10^{14}$ during SAs. Even for thin metal layers (typically below 20 cm in depth), for which the focusing effect is strongest, the Rayleigh number can reach up to $Ra_\phi = 10^{10}$.

In addition, all existing models focus on the mean value of the heat flux only, and overlook possible large amplitude fluctuations on the side boundary. Rein *et al.* (2023a) highlighted the presence of a three-dimensional (3-D) drifting thermal structure taking the form of radial hot branches, potentially playing a crucial role in those heat transfer fluctuations, but up to now unaccounted for.

This paper presents an experimental study of the convection in a thin cylindrical gas layer with a Prandtl number $Pr = 0.7$, alongside 3-D DNS of the experimental set-up using both the experimental working fluid and a fluid with a Prandtl number $Pr = 0.1$ representative of the nuclear application (Carénini *et al.* 2018). The objectives are to: (i) experimentally confirm and expand the scope of scaling laws identified through our previous numerical simulations; (ii) experimentally identify the pattern of drifting branches and analyse its properties depending on the input parameters; and (iii) investigate the heat flux fluctuations at the wall and quantify the influence of the drifting branch pattern on heat transfer. The paper is divided into four sections. The next section delves into the experimental set-up, detailing the instrumentation and the experimental protocol. In § 3 we outline the governing equations and the numerical simulation tool employed. In § 4 we present both experimental and numerical results, beginning with an analysis of mean values within the system and describing scaling laws through dimensional analysis arguments. Additionally, we examine the angular drift velocity and the azimuthal wavenumber of the drifting branch pattern in relation to the Rayleigh number. We also investigate the distribution of heat flux fluctuations and explore the role of the drifting pattern in heat flux transport along the side. The concluding section discusses the implications of our findings for nuclear safety and outlines potential avenues for future research.

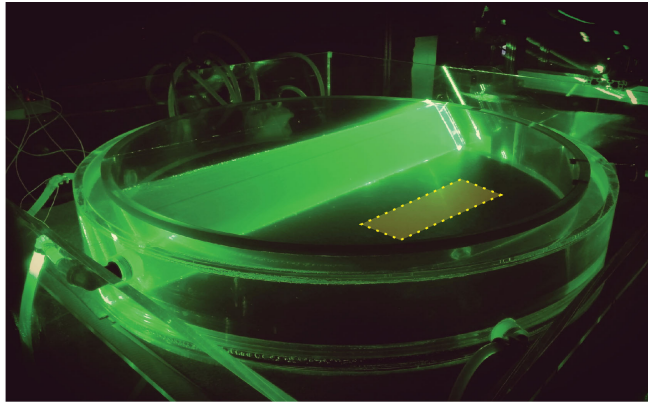
2. Experimental method

2.1. Experimental set-up: FOCUS

The experimental set-up, named FOCUS and depicted in figure 1, consists of an upright acrylic cylindrical enclosure with a vertical main axis and an aspect ratio $\Gamma = R/H = 4$, with R the tank radius and H its height. Inside this enclosure, a gas (air or air/SF₆ mixture) is heated from below and cooled along the side.

The choice of a gas as a simulant was motivated by the Prandtl number characteristic of the metal layer in the nuclear application, which is less than 1 ($Pr = \nu/\kappa \simeq 0.1$, with ν the kinematic viscosity and κ the thermal diffusivity). When $Pr < 1$, the viscous boundary layer is nested within the thermal boundary layer. This structure plays a major role in heat transfers (Rein *et al.* 2023a). Using gases with $Pr \simeq 0.7 < 1$ ensures maintaining a similar configuration, thus preserving similar heat transfer mechanisms, without the cost and many difficulties of using liquid metals (including, for instance, their opacity or the necessity to work at higher than ambient temperature). With SF₆ being denser than air, it allows more turbulent regimes to be reached by significantly increasing the Rayleigh number: indeed, the thermal conductivity and the dynamic viscosity of air and SF₆ are mostly similar, so the Rayleigh number (see the exact definition (3.5)) scales as the gas density squared. Note that such gases have already been successfully used to study high-Rayleigh-number thermal convection in the past; see, e.g. Ashkenazi & Steinberg (1999); Ahlers *et al.* (2009a); and more recently Cierpka *et al.* (2019).

(a)



(b)

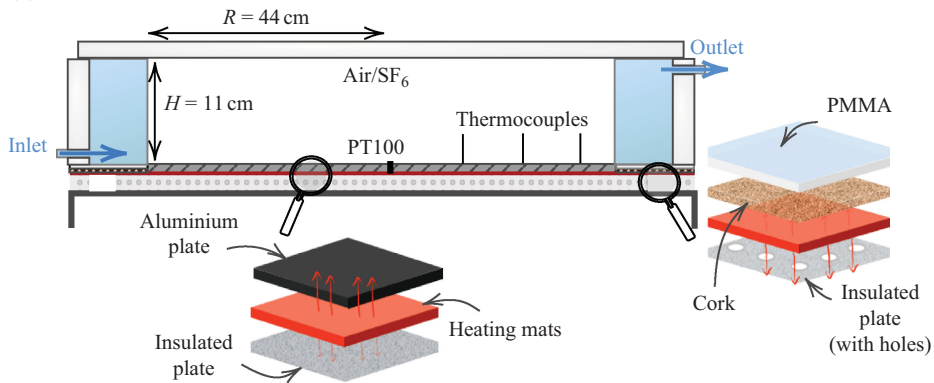


Figure 1. (a) Photo of the experimental set-up with the laser sheet inclined for better visualisation (it is horizontal during acquisition). The dotted yellow rectangular region shows the PIV camera field of view. (b) Sketch of a vertical cross-section of the experimental set-up with magnified views of the bottom plate configuration below the working tank and below the side boundary, respectively.

In our set-up (see [figure 1](#)), the gas layer is heated from below through eight silicone heating elements ($450 \text{ mm} \times 250 \text{ mm}$ each) with a combined maximum power of 7.2 kW , resting on a cylindrical aluminium plate with a radius of $R = 44 \text{ cm}$, for better heat flux homogenisation. These heating elements are positioned on a 5 cm thick calcium silicate insulating plate to limit heat losses and to concentrate heat into the aluminium plate. Cooling on the side is provided by a cylindrical polymethyl methacrylate (PMMA) torus with a rectangular section of 11 cm in height and 40 mm in width, through which water flows at an average rate of 15 L min^{-1} . The water temperature is regulated by a thermostatic bath with a power of 1.2 kW . Three water inlets and outlets located at the bottom and top of the torus, arranged in a staggered pattern and spaced at 60° intervals, promote water mixing through natural and forced convection. The thickness of the inner wall of the torus is only 4 mm (minimum thickness achievable through manufacturing processes), so as to ensure a temperature as constant as possible at the wall. Note that on the edges of the bottom plate, a thin 3 mm cork plate was added to prevent direct contact between the heating elements and the bottom of the cooling torus, while holes were drilled in the insulating plate to dissipate heat. Note also that insulated walls can introduce side wall

effects, especially when gas is used (Roche *et al.* 2001). In our set-up however, we expect that the temperature-regulated water inside the tank, controlled by the powerful thermal bath, largely compensates for any residual heat flux.

A transparent polycarbonate lid with a thickness of 35 mm is used to maintain the gas in the tank and to constrain the outgoing heat flux, while allowing flow visualisation from above. This outgoing heat flux through the lid is a parameter of our study and accounts for approximately 50 % of the input heat flux across all experiments (see Appendix A.1). The choice of using a polycarbonate cover was guided by the temperature range that can be reached with the experimental set-up, in order to prevent any changes in lid opacity and ensure optimal flow visualisation.

Finally, to minimise the SF₆ mixing with the external environment, the entire set-up is enclosed in a large square PMMA tank measuring (1220 mm × 1220 mm × 470 mm).

The operational temperature range of the device has been limited to less than 1000 °C, reaching a maximal bulk temperature of 64 °C. Indeed, since we use gas as a simulant for the incompressible metal layer, we want to avoid excessive temperature differences that would result in significant and undesirable compressibility effects. For an ideal gas, the density variation can be estimated as the ratio of the largest temperature change between the initial and statistically stationary states to the mean temperature in the statistically stationary state. Throughout all the experiments, we measured an average relative density variation of 5 %, peaking at 11.5 % for the highest Rayleigh number obtained.

2.2. Instrumentation

2.2.1. Heat flux sensor

In order to measure heat flux fluctuations through the side wall, we use a thermal flux sensor ('gSkin' manufactured by 'greenTEG') measuring 18 mm × 18 mm × 0.5 mm and positioned on the inner wall of the torus. It has a response time of the order of 0.7 s and its sensitivity depends on temperature via the relation

$$s = S_0 + (\theta_0 - \theta_S)S_c, \quad (2.1)$$

with $S_0 = 39.92 \mu\text{V W}^{-1} \text{m}^2$, $\theta_S = 22.5 \text{ }^\circ\text{C}$, $S_c = 0.0499 \mu\text{V W}^{-1} \text{m}^2 \text{ }^\circ\text{C}^{-1}$. Here θ_0 corresponds to the temperature value of the substrate on which the flux sensor is placed, hence, the cooling temperature here. During the measurement campaign, the sensor was mainly placed at mid-height of the torus, using thermal paste to avoid any parasitic thermal resistance and to ensure good adhesion to the wall over the entire surface despite the local curvature.

2.2.2. PT100 and thermocouples

We also deployed a set of temperature measurements inside the set-up, as sketched in figure 1(b). Six thermocouples are positioned at mid-height of the tank, distributed along two radii offset by 90°, and located at $r = 10 \text{ cm}$, 20 cm and 30 cm from the tank centre. This specific arrangement aims at characterising the dynamics of the thermal branch pattern, particularly by measuring its drift frequency. Additionally, we have inserted and levelled a PT100 sensor in the centre of the aluminium bottom plate. Two other PT100 sensors are used to measure the inlet and outlet temperatures of the water in the torus to estimate the extracted heat flux. To complete the set-up, two additional thermocouples are positioned above and below the lid at its centre. All signals are acquired using National Instruments modules (NI9212, NI9211 and NI9217 modules for the acquisition of thermocouples, flux sensor and PT100 sensors, all connected to an NI cDAQ-9174

rack). These measurements contribute to a comprehensive thermal characterisation of the system. In particular, they allow for quantifying the thermal losses and the distribution of heat flux between the lateral and upper surfaces, as described in Appendix A.3.

2.2.3. Particle Image Velocimetry

We use time-resolved particle image velocimetry (PIV) to determine the horizontal velocities in an horizontal laser sheet positioned 6 cm above the aluminium plate. This sheet is produced by a 10 W Asur light systems 488–532 nm continuous wave laser with a Powell lens. Finding particles acting as passive scalars and persisting during a long time in our closed gas system proved particularly challenging. Initially, the use of fine water droplets seemed promising, but in an environment subjected to high temperatures the lifespan of these droplets was too short due to rapid evaporation. This issue was resolved by adding a small fraction of UCON oil to the water droplets. However, a fine balance had to be struck, as a substantial addition of oil leads to an increase in the typical size of the droplets and, thus, to rapid sedimentation, altering their behaviour as passive scalars. An oil concentration of 3.5 % by volume was identified as an optimal compromise (Dorel, Le Gal & Le Bars 2023), allowing the droplets to persist for a period of 30–40 min while faithfully tracking flow motions. These particles are produced and introduced into the system via a 1 cm diameter conduit on the side of the cover (within its thickness) using an atomizer. Their average size is of the order of 10 μm . An estimate of the viscous relaxation time of the droplet motion shows that it is much smaller than the Kolmogorov time, indicating that the droplets behave as passive tracers. Furthermore, we verified that the volume of liquid introduced into the system is too small to have a significant thermal impact on the gas properties.

The droplets motion is tracked using a camera positioned above the set-up (Blackfly U3-51S5M FLIR with a Fujinon 12.5 mm lens). We narrowed the camera's field of view to achieve a sampling frequency ranging from 100 fps to 180 fps, resulting in images of 1500 pixels \times 700 pixels. The observation area forms a rectangle of 34 cm \times 16 cm, offset by 18 cm from the centre (see figure 1a). The velocity fields are derived from these images using the DPIVSoft program developed by Meunier & Leweke (2003). We consider 24 pixels \times 24 pixels boxes with 50 % overlap on 1500 pixels \times 700 pixels images, resulting in a velocity field of 125 \times 58 vectors. Throughout the results, the percentage of spurious vectors in the velocity field never exceeded an average of 5 %.

2.3. Experimental protocol

We now detail the experimental protocol implemented during the measurement campaign. We begin by initiating the acquisition system for temperature and flux. In the case of SF₆ usage, plastic sheets are placed above the set-up and fixed to the outer basin to contain the gas. The outer basin is first filled completely with SF₆ from the bottom, until gas is detected on the surface of the sheets using a gas detector. Then, the experimental set-up is filled with SF₆ while aiming to maximise density. There, we temporarily use a densimeter (WIKIA Tech, Avenisense NorthDome) placed inside the system, with its sensor positioned 2 cm above the aluminium plate, near the edge opposite to the flux sensors.

Water circulation and water cooling via the thermostated bath are initiated. The bath temperature set point is chosen to maintain a difference of approximately 5 °C with the ambient temperature, aiming to keep the same external heat losses between each experiment (see details in Appendix A.3). Finally, the bottom heating system is activated by selecting the electrical power to be supplied to the heating elements.

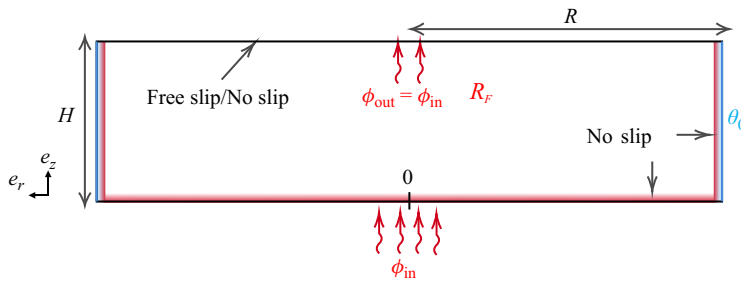


Figure 2. Sketch of the modelled fluid layer in a vertical plane through the cylinder. The top boundary condition is no slip for the gas experiment simulations, representing the lid used in the set-up, while a free-slip condition is applied to model the metal layer in the nuclear application.

We now enter the transient regime. When using SF_6 , it is expected that the density inside the set-up decreases as it is not gas tight, and SF_6 diffuses into the ambient air. If the density drops below the desired threshold, more SF_6 is injected into the system until the desired density is reached. The gas in the system is then a mixture of air and SF_6 . The mass fraction of SF_6 , denoted y , is determined via a density measurement further detailed in Appendix A.2.

During this transient regime, the characteristic time and thermal losses are evaluated (see Appendix A.3). The steady-state regime is achieved when the heating time exceeds 3 characteristic times, i.e. in about approximately 9 h. At this stage, it must be ensured that the temporal variation of temperature is less than $1\text{ }^\circ\text{C h}^{-1}$.

Once the steady state is reached, flux data are collected for 3–6 diffusive times (approximately 1 h). During this period, the necessary elements for PIV measurements are prepared, including the water/UCON oil mixture to generate PIV particles. After the flux data acquisition is completed, the seeding phase begins. A low-power horizontal laser sheet (0.2 W) is set up, an air/ SF_6 injection flow rate through the atomizer is fixed, and the set-up is seeded for a predefined duration to control the quantity of injected PIV particles, ensuring experimental reproducibility. The injection of these particles disturbs the flow and the thermal fluxes involved. It is necessary to wait for the particle jet to homogenise and the thermal signals to return to steady state before starting the flow video acquisition. The laser power, gain, sampling frequency and exposure time are then adjusted to identify and track the movements of the PIV particles. A series of video acquisitions are then conducted, each for a duration of approximately one minute (a duration for which we verify a constant sampling frequency).

3. Mathematical and numerical formulation

3.1. Governing equations

Following Rein *et al.* (2023a), we model the experimental set-up considering an incompressible fluid in the Boussinesq approximation, confined in a cylinder of thickness H and radius R , with gravity directed downward (see figure 2). Heating from below is applied with a uniform heat flux ϕ_{in} , while cooling from above occurs with a uniform outgoing flux ϕ_{out} . Note that this upper boundary condition is an approximation, since the experimental outgoing flux is dynamically constrained by the convective heat transfers within the system and may change radially. Nonetheless, the thermal budget described in Appendix A.3 confirms that the surface average measurement and the local, centre plate

measurement of the outgoing flux have close values, hence validating this assumption with an outgoing flux ϕ_{out} ranging between 40 % and 60 % of the incoming one. The study hence focuses on cases where ϕ_{in} and ϕ_{out} are unequal, resulting in residual heat flux escaping through the side boundary, whose temperature is fixed at θ_0 . No-slip conditions are applied to the bottom and side boundaries, while two configurations are considered for the top boundary in the DNS: a no-slip condition to represent the experimental lid and a free-slip condition to model the liquid metal layer in the nuclear application.

Lengths are rescaled using the height of the cylinder H , while time is rescaled using the vertical diffusive dash-dot time scale H^2/κ . In our set-up, the two main control parameters are the bottom heat flux and the side temperature, which is the only imposed temperature. Hence, the non-dimensional temperature T is defined relative to the imposed side temperature and is rescaled using the imposed bottom flux ϕ_{in} ,

$$T = \frac{k}{\phi_{in}H}(\theta - \theta_0), \tag{3.1}$$

where k is the thermal conductivity. The non-dimensional conservation equations of momentum, mass and energy are then

$$\frac{1}{Pr} \left(\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} \right) = -\nabla P + Ra_\phi T \mathbf{e}_z + \nabla^2 \mathbf{u}, \tag{3.2}$$

$$\nabla \cdot \mathbf{u} = 0, \tag{3.3}$$

$$\frac{\partial T}{\partial t} + \mathbf{u} \cdot \nabla T = \nabla^2 T. \tag{3.4}$$

Here \mathbf{u} and P are the non-dimensional velocity and pressure (including the hydrostatic contribution) fields. The problem is characterised by four non-dimensional parameters: the aspect ratio Γ , the flux ratio R_F , the Rayleigh number Ra_ϕ based on the heat flux imposed at the bottom ϕ_{in} , the Prandtl number Pr set to 0.1 to mimic the properties of the metal layer or 0.7 to replicate those of the gas used in the experiment. They are defined by

$$\Gamma = \frac{R}{H}, \quad R_F = \frac{\phi_{out}}{\phi_{in}}, \quad Ra_\phi = \frac{\beta g \phi_{in} H^4}{k\nu\kappa}, \quad Pr = \frac{\nu}{\kappa} = \begin{cases} 0.1 \text{ (application),} \\ 0.7 \text{ (experiment),} \end{cases} \tag{3.5}$$

where β is the thermal expansion coefficient, assumed to be constant. The non-dimensional boundary conditions can be written as

$$\begin{aligned} \mathbf{u}(z=0) &= \mathbf{0} \quad \text{and} \quad \left. \frac{\partial T}{\partial z} \right|_{z=0} = -1, \\ \mathbf{u}(r=\Gamma) &= \mathbf{0} \quad \text{and} \quad T(r=\Gamma) = 0, \\ \left\{ \begin{array}{l} \text{Application: } \left. \frac{\partial u}{\partial z} \right|_{z=1} = \left. \frac{\partial v}{\partial z} \right|_{z=1} = w(z=1) = 0 \\ \text{Experiment: } \mathbf{u}(z=1) = \mathbf{0} \end{array} \right. & \quad \text{and} \quad \left. \frac{\partial T}{\partial z} \right|_{z=1} = -R_F, \end{aligned} \tag{3.6}$$

where $\mathbf{u} = (u, v, w)$ are the velocity components in cylindrical coordinates along the unit vectors $\mathbf{e}_r, \mathbf{e}_\phi, \mathbf{e}_z$.

3.2. Numerical approach

The governing equations (3.2)–(3.4) with corresponding boundary conditions (3.6) are computationally solved using the spectral element code [Nek5000](#) (Fischer 1997; Deville, Fischer & Mund 2002). The cylindrical geometry is discretised with up to $\mathcal{E} = 9984$

hexahedral elements, refined near all boundaries to accurately resolve viscous and thermal boundary layers, with typically 15 grid points across the viscous boundary layers, reaching a minimum of 10 grid points for the highest Rayleigh number.

Numerical simulations are initiated with a quiescent fluid and a uniform temperature field $T = 0$ throughout the domain. Small temperature perturbations of amplitude 10^{-3} are introduced, leading to thermal convection growth during a transient phase lasting approximately 5 vertical diffusive times, with longer durations observed for higher aspect ratios (Γ). Once the system reaches a statistically stationary state, various spatio-temporal averages are computed. Temporal and volume average operators $\langle \cdot \rangle$ are defined over the entire fluid domain volume V and time τ as

$$\langle \cdot \rangle = \frac{1}{\tau V} \int_{t_0}^{t_0+\tau} \int_V \cdot dV dt, \quad (3.7)$$

where τ typically ranges between 2 and 0.1 diffusive times for the lowest and highest Rayleigh numbers, respectively. Additionally, adding specific variables as a subscript means that an average along those specific directions is made. We always consider temporal averages during the statistically steady state so that the time variable is never explicitly written. For instance, $\langle \cdot \rangle_\phi$ indicates an average in time and along the azimuthal direction only.

Although most results are derived from DNS, some extreme cases necessitated filtered simulations, as described in Fischer & Mullen (2001). A viscous dissipation criterion was employed to distinguish between DNS and filtered simulations. For further details, readers are directed to Rein *et al.* (2023a).

4. Experimental and numerical results

Detailed information for all experiments and simulations is available in tables 1 and 2 of Appendix A.1, respectively. The experimental data set comprises 12 independent experiments, including three with the SF₆ mixture where the average density ranged from 2.9 to 4.1 kg m⁻³. With the heating power varied from 75 to 400 W, we have explored the range $Ra_\phi \in [3.2 \times 10^7 : 7.6 \times 10^9]$. This experimental data set is completed by eight 3-D DNS of the experimental set-up (with a no-slip condition on the upper surface and $Pr = 0.7$), exploring the range $Ra_\phi \in [10^5 : 10^8]$. We also performed 70 DNS with a free-slip condition on the upper surface and $Pr = 0.1$, corresponding to the nuclear application, whose results have already been reported in Rein *et al.* (2023a).

4.1. Scalings

In figure 3(a) the non-dimensional mean temperature is plotted against the Rayleigh number. To do so, we averaged in time each thermocouple signal and then averaged over the six thermocouples within the system. Using DNS, we compute the mean temperature following the same set of discrete locations as in the experiments, considering both $R_F = 0.5$ and $R_F = 0.6$ (which encompasses most of the experimental values for the flux ratio; see table 1), and spanning a range of Ra_ϕ from 10^5 to 10^8 .

Overall, a satisfactory agreement between DNS and experimental measurements is observed. One can notice here the slight influence of the flux ratio. Indeed, the two lowest experimental measurement points in terms of Ra_ϕ (close to $Ra_\phi \sim 10^7$) correspond to flux ratios $R_F \simeq 0.5$ (see Appendix A.3) and align well with the numerical trend observed for $R_F = 0.5$. Similarly, the air experimental measurements near $Ra_\phi \sim 10^8$, associated with flux ratios close to 0.6, show excellent agreement with the numerical trend at $R_F = 0.6$.

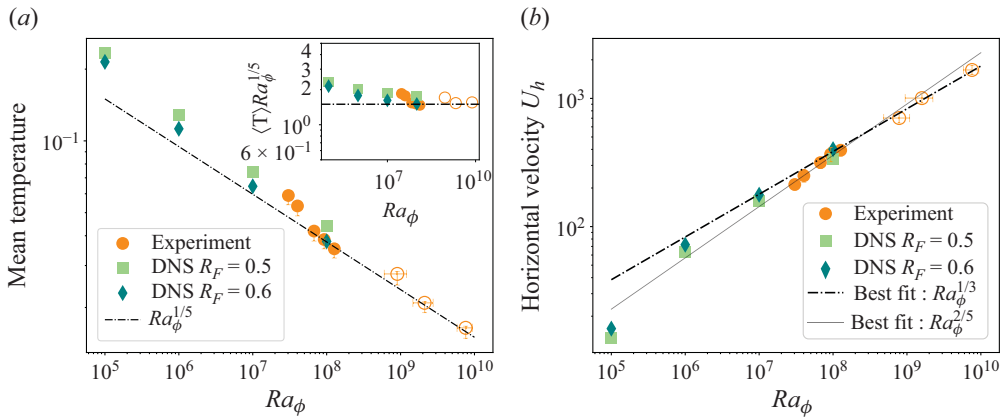


Figure 3. (a) Non-dimensional temperature averaged in time and over the six thermocouples in the bulk of the experiment, as a function of Ra_ϕ . The dash-dot line shows the $Ra_\phi^{-1/5}$ asymptotic scaling from Rein *et al.* (2023a) (the inset shows the compensated plot). (b) Non-dimensional horizontal r.m.s. velocity in the (x, y) plane at height $z = 6$ cm as a function of Ra_ϕ . The dash-dot line and continuous line show a $1/3$ and $2/5$ best fit power law based on experimental data. For both plots, square (\square), diamond (\diamond) and circle (\circ) symbols represent respectively data from DNS simulations at $R_F = 0.5$ and $R_F = 0.6$, and the experiments. Empty circles indicate when SF_6 gas is used. For all DNS simulations, $\Gamma = 4$ and $Pr = 0.7$.

The experimental measurements conducted with SF_6 have a ratio close to 0.6, and extend the numerical trend by nearly two orders of magnitude in terms of Ra_ϕ .

Also shown is the slope of the asymptotic scaling in $Ra_\phi^{-1/5}$ predicted by Rein *et al.* (2023a) and which corresponds to the classical scaling of vertical convection predicted by Batchelor (1954); the inset graph shows the results compensated by this scaling with a good agreement above $Ra_\phi = 10^8$. The scaling, as detailed in Rein *et al.* (2023a), comes from the balance at a statistically stationary state between the top/bottom flux mismatch, and the lateral outgoing diffusive flux through the thermal boundary layer at the vertical boundary. As the Rayleigh number increases, the system becomes more turbulent leading to more efficient bulk mixing limited by the diffusive boundary layer, and hence, a better agreement with the scaling law.

In figure 3(b) we plot the norm of the non-dimensional horizontal velocity value U_h as a function of the Rayleigh number for numerical simulations at $R_F = 0.5$ and 0.6, and the experimental data. This velocity is computed as follows:

$$U_h = \left[\langle u^2 \rangle_{r\varphi} + \langle v^2 \rangle_{r\varphi} \right]^{1/2}. \quad (4.1)$$

The averaging procedure along the azimuthal and radial directions involves the same spatial domain for the experiments and DNS (i.e. the PIV field of view; see § 2.2.3). The time averaging is processed during a characteristic time of 1 min for the experiments (0.1 diffusive time for air, 0.02 for SF_6) and over 0.1 diffusive time for DNS. Each point represents a measurement of the characteristic horizontal fluid velocity localised in the PIV measurement zone. First, it is worth noting the excellent agreement between the experimental measurements and DNS over the range of Rayleigh numbers common to both approaches. In particular, the experimental points between $Ra_\phi = 10^7$ and $Ra_\phi = 10^8$, conducted in air, compare favourably with the numerical trend obtained at $Ra_\phi = 10^8$. Additionally, the experimental measurements involving the use of SF_6 extend the trend

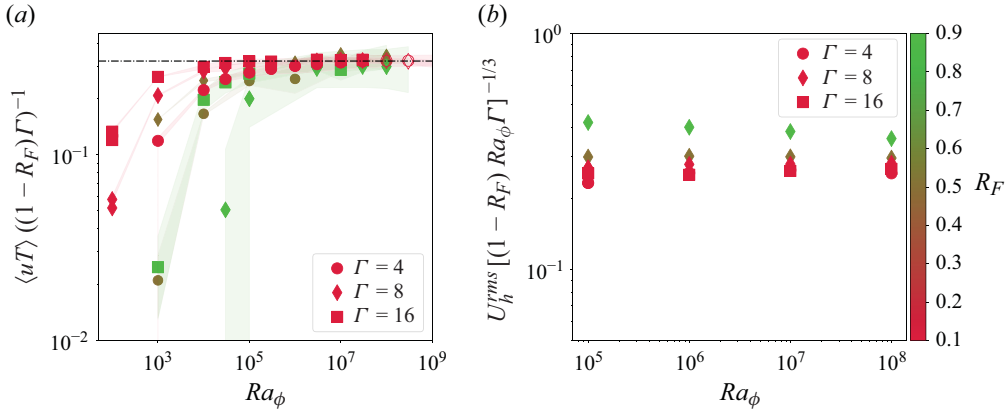


Figure 4. (a) Compensated horizontal advective flux following (4.4) as a function of Ra_ϕ . (b) Compensated radial r.m.s. velocity following (4.5) as a function of Ra_ϕ . For both plots, different symbols correspond to different aspect ratios Γ and colour variation from red to green involves R_F variation from 0.1 to 0.9. Data are from DNS with $Pr = 0.1$ and a free-slip upper boundary, for which our data set is more complete in terms of Γ and R_F systematics, but the same behaviour is expected with the experimental configuration. Empty/full symbols indicate respectively filtered/DNS simulations. The shadings represent three times the standard deviation of the horizontal advective flux time series at the statistically stationary state.

over almost two orders of magnitude in Ra_ϕ . A power law relationship compatible with $Ra_\phi^{1/3}$ is observed, with a best fit estimate of $Ra_\phi^{0.36}$ using the least squares method based on the experimental points. Note that in the numerical simulations, the root-mean-square (r.m.s.) velocity exhibits minimal dependence on the flux ratio. This observation aligns with our previous findings using $Pr = 0.1$ and free-slip boundary conditions applied to the upper surface (not shown). In the following section we use dimensional analysis to gain insight into the physics underlying this scaling law.

4.2. Dimensional analysis

Let us focus on the heat transport by considering the following non-dimensional energy equation derived from (3.4) using cylindrical coordinates:

$$\frac{\partial T}{\partial t} + \frac{1}{r} \frac{\partial}{\partial r} \left[r \left(uT - \frac{\partial T}{\partial r} \right) \right] + \frac{1}{r} \frac{\partial}{\partial \varphi} \left[vT - \frac{1}{r} \frac{\partial T}{\partial \varphi} \right] + \frac{\partial}{\partial z} \left[wT - \frac{\partial T}{\partial z} \right] = 0. \quad (4.2)$$

The horizontal transport of the heat flux at a radius r is expressed by averaging the energy equation (4.2) in time, vertical and azimuthal directions, which leads to

$$\langle uT \rangle_{\varphi z} - \left\langle \frac{\partial T}{\partial r} \right\rangle_{\varphi z} = \frac{(1 - R_F)}{2} r. \quad (4.3)$$

This equation illustrates that the difference in heat flux between the top and bottom (i.e. the right-hand side of (4.3)) is transported horizontally through a combination of diffusion ($\langle \partial_r T \rangle_{\varphi z}$) and advection ($\langle uT \rangle_{\varphi z}$) processes. If we now assume the system to be fully turbulent, the radial diffusive gradient matters only within the thermal boundary layer. Within the bulk, we expect the balance $\langle uT \rangle_{\varphi z} \sim ((1 - R_F)/2)r$. By averaging along the radial direction, the horizontal advective flux scales as follows:

$$\langle uT \rangle = \frac{1}{\Gamma} \int_0^\Gamma \langle uT \rangle_{\varphi z} dr \sim (1 - R_F)\Gamma. \quad (4.4)$$

This scaling is confirmed in figure 4(a), where we use DNS data with Γ ranging from 4 to 16, R_F from 0.1 to 0.9 and $Pr = 0.1$. The higher the Rayleigh number, the better the agreement with the scaling given by (4.4). In particular, above $Ra_\phi \sim 10^5$, $\langle uT \rangle$ becomes independent of the Rayleigh number and all the data collapse on the same straight line.

We now seek to explain the scaling of the horizontal velocity. Following the determination of the free fall velocity scaling in classical Rayleigh–Bénard convection, we assume that the dimensional horizontal velocity scale is independent of viscosity and diffusivity, but depends on its driving buoyancy, i.e. the horizontal turbulent buoyancy flux $\beta g F_h \sim \beta g \phi / \rho c_p$, related to $\langle uT \rangle$ in non-dimensional form.

Scaling analysis gives $\hat{U}_h^{rms} \sim (\beta g F_h H)^{1/3}$, where $\hat{\cdot}$ denotes dimensional variable. In non-dimensional form, this becomes $U_h^{rms} \sim (Ra_\phi Pr \langle uT \rangle)^{1/3}$. Hence, with (4.4),

$$U_h^{rms} \sim (Ra_\phi Pr(1 - R_F)\Gamma)^{1/3}. \tag{4.5}$$

We recover the scaling law in $Ra_\phi^{1/3}$ identified in figure 3(b). We further validate this scaling law in figure 4(b), which shows the compensated r.m.s. horizontal velocity over a large range of (Γ, R_F, Ra_ϕ) at $Pr = 0.1$. From our DNS data set, we consider the inner part of the cylinder ($r \leq \Gamma/2$) in order to focus on the turbulent bulk, and we compute

$$U_h^{rms} = \left\langle (u - \langle u \rangle_{\varphi z r_f})^2 + (v - \langle v \rangle_{\varphi z r_f})^2 \right\rangle_{\varphi z r_f}^{1/2}, \tag{4.6}$$

where the r_f subscript indicates an average along the radial direction over $r \in [0, \Gamma/2]$. The agreement with (4.5) is very good with all the data converging at large Ra_ϕ . This is particularly clear for the lower flux ratios. As the value of R_F increases, the relevance of a characteristic horizontal velocity derived from horizontal heat transport becomes less meaningful. In the asymptotic case $R_F = 1$, this typical velocity scale effectively disappears since all the injected heat flux is evacuated at the top.

Note that in the context of vertical convection (flow between two differentially heated vertical walls), similar dimensional arguments have been put forward in the seminal work of George & Capp (1979). This scaling has been notably revisited and corroborated by subsequent studies such as Versteegh & Nieuwstadt (1999); Hölling & Herwig (2005); Ng, Chung & Ooi (2013). Note also that even if we observe the same 1/3 exponent in Ra_ϕ as for horizontal convection with imposed flux conditions (Mullarney, Griffiths & Hughes 2004), the velocity corresponds in this case to the horizontal velocity within the thermal boundary layers (Rossby 1965; Griffiths, Hughes & Gayen 2013), whereas in our case, it corresponds to the bulk velocity driven by the imposed buoyancy flux. Finally, we pointed out in our previous paper that our system shares similarity with the Grossmann & Lohse (2000) regime $I-l$ (Rein *et al.* 2023a). Hence, one might expect a typical r.m.s. velocity scaling like $Ra_\phi^{2/5}$ (i.e. the imposed flux version of the classical $Ra^{1/2}$ scaling in the imposed temperature convection). This does not appear to be consistent with our data, as show in figure 3(b). Presumably, the experimental range of the flux ratio observed ($R_F \in [0.4, 0.6]$) does not reach the asymptotically high value necessary to manifest the fully developed $I-l$ regime described by Grossmann & Lohse (2000), which would require larger Ra_ϕ and/or R_F .

4.3. Persistent drifting pattern in the turbulent regime

4.3.1. Characteristics

In Rein *et al.* (2023a) a drifting pattern was observed at high Rayleigh numbers for low flux ratios (see their figure 2(c) for example), but no detail was given about its

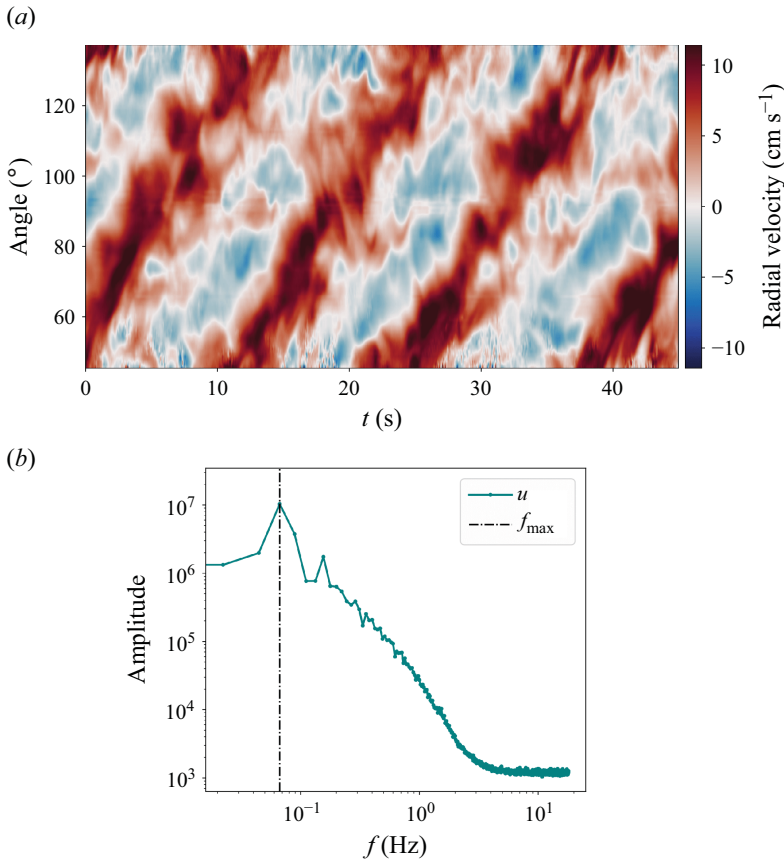


Figure 5. (a) Hovmöller (space–time) diagram of the radial velocity averaged along the radial direction. The radial velocity is measured at $z = 6$ cm (i.e. $z/H = 0.54$). (b) Average of the radial velocity time spectra from each PIV measurement location. The vertical dash–dot line shows the frequency ($f_{max} = 6.7 \times 10^{-2}$ Hz) of the peak amplitude. For both, data correspond to experiment 9 from table 1 where $Ra_\phi = 1.26 \times 10^8$, using air.

properties. During the experimental campaign discussed here, the system also showed the spontaneous emergence of such a branch pattern. A video of experiment 3 from table 1 conducted in air at $Ra_\phi \sim 10^7$ can be downloaded as supplementary material available at <https://doi.org/10.1017/jfm.2025.2>. In this video we have removed the average image spanning the entire duration of the footage and applied a moving average over five frames across all images. This adjustment allows for a clearer visualisation of the movement of the PIV particles, represented by white streaks. The video is in real time, except for the clear appearance of a branch where a slowdown was applied. It shows a branch of the pattern, visible as extended white streaks moving from right to left.

The drifting branches are also visible through the velocity fields obtained from PIV processing, as, for instance, on the Hovmöller diagram shown in figure 5. Thick bands associated with the highest radial velocities indicate the passage of the branches in the camera’s field of view. The global angular drift frequency of the whole pattern ω_p can be deduced from such diagrams by tracking a branch of the pattern. The observed drift can occur in both clockwise and anticlockwise directions in independent realisations of the same configuration. This indicates that there is no preferred rotation direction as

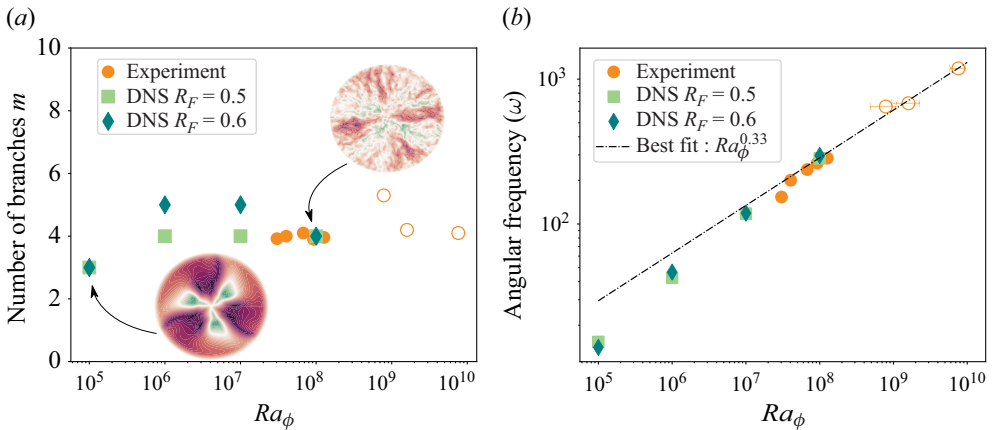


Figure 6. (a) Number of branches as a function of Ra_ϕ . The insets show two snapshots of the radial velocity maps from DNS at $z = 0.66$ for $Ra_\phi = 10^5$ and 10^8 , $R_F = 0.6$, $Pr = 0.7$ and $\Gamma = 4$. (b) Angular drift frequency as a function of Ra_ϕ . The dashed–dotted line represents a best fit based on the experimental data for $Ra_\phi > 5 \times 10^7$. In both graphs, square symbols (\square), diamond symbols (\diamond) and circle symbols (\circ) respectively represent data for DNS at $R_F = 0.5$ and $R_F = 0.6$, and experimental data. Empty circles indicate the use of SF₆. All DNS were performed with $Pr = 0.7$ and $\Gamma = 4$.

expected from the problem axisymmetry. A local frequency ω can also be computed from the temporal spectrum of the radial velocity at each spatial location. The spectra are then averaged over the PIV measurement area for experiments and over the whole domain for DNS. An example based on the PIV measurement from experiment 9 is shown figure 5(b).

The two determined frequencies are related by $\omega = m \omega_p$, where m is the number of branches or azimuthal wavenumber. Figure 6 shows m and ω as functions of the Rayleigh number for all our experimental and numerical results. We note a good overall qualitative agreement, even if DNS exhibit a stronger dependence of m on the flux ratio. Regarding ω , a power law emerges, indicating a dependence $\omega \sim Ra_\phi^{1/3}$. This relationship was determined through a best fit using the least squares method, applied to the experimental data while excluding values below $Ra_\phi = 5 \times 10^7$ to focus on the asymptotic behaviour at high Rayleigh numbers. This 1/3 exponent is reminiscent of the one identified for the horizontal velocity in figure 3. However, we argue that both are not directly related. Indeed, from the DNS at $R_F = 0.5$ and 0.6 , we can compute the mean angular velocity averaged in φ and z at $r = 3\Gamma/2$ (our conclusion remains qualitatively unchanged for other radii). For the low Ra_ϕ values of DNS, this advection angular velocity is significantly smaller than the measured angular drift velocity. Hence, we argue that the drifting pattern is not caused by the advection of some structure by a mean azimuthal velocity, but rather seems related to a wave-like phenomenon where the fluid, on average, does not undergo significant azimuthal movement. We also note from our systematic DNS study that this pattern appears only above a given threshold in Ra_ϕ . Initially, prograde and retrograde similar structures superimpose. By increasing Ra_ϕ , one of the two directions is randomly selected and then persists up to the largest Ra_ϕ investigated. A detailed study of the onset of this oscillatory instability is necessary. This will be the subject of future work, including comparisons with other large-scale structures observed in turbulent convection, such as the

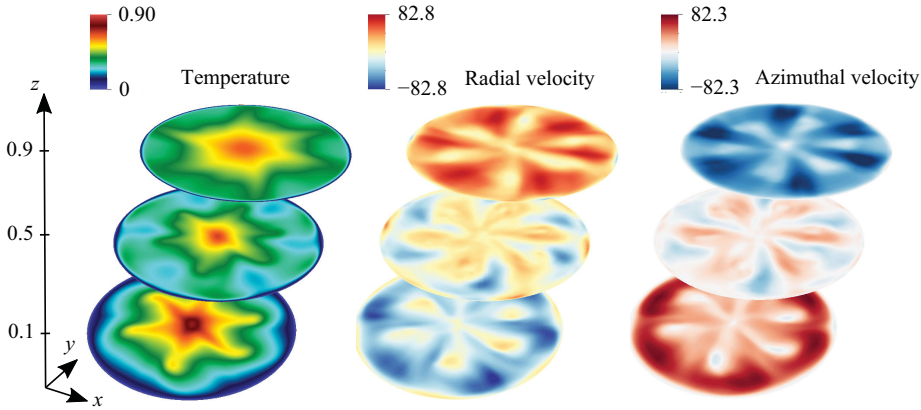


Figure 7. Phase-averaged temperature, radial velocity and azimuthal velocity fields, processed during 1 diffusive time. Results are presented for three horizontal slices at $z = 0.1, 0.5$ and 0.9 . Parameters are $\Gamma = 4$, $R_F = 0.1$, $Ra_\phi = 10^5$ and $Pr = 0.1$. In the present DNS, the pattern drifts anticlockwise.

jump rope vortex observed in liquid metal by Akashi *et al.* (2022); Cheng *et al.* (2022); Teimurazov *et al.* (2023).

4.3.2. Analysis in the reference frame of the drifting pattern

To gain deeper insights into the temperature and velocity field structure associated with the drifting pattern, we analyse our DNS in the frame of reference rotating with the pattern. To do so, we first determine the angular drift frequency (ω) and azimuthal wavenumber (m) from the time and spatial spectra of the azimuthal velocity. Then, we spectrally interpolate, at each time step, all variables on a grid $(X(t), Y(t), Z(t))$ rotating with the phase speed ω/m of the pattern. We finally introduce a phase-averaging operator denoted by $\langle \cdot \rangle_p$, which is effectively a time average over a time τ_p within the reference frame of the drifting pattern, defined as

$$\langle T \rangle_p = \frac{1}{\tau_p} \int_{t_0}^{t_0 + \tau_p} T(t, X, Y, Z) dt . \quad (4.7)$$

The typical value of τ_p ranges between 0.2 and 1 diffusive times for the largest and the lowest Rayleigh numbers, respectively. To ensure the convergence of the data obtained from this averaging, we verify that the probability density functions (PDF) of the radial temperature gradient and velocity fields remain consistent even when considering only half of the data.

Figure 7 shows an example of the obtained temperature, radial velocity and azimuthal velocity fields at three different depths. Thanks to the phase average, we now clearly identify the six branches pattern on all fields, while it would otherwise disappear when using a standard time average in the laboratory frame. Furthermore, in the upper part of the domain, we observe positive radial velocities moving from the centre towards the edge, while in the lower part, the radial velocities are negative, moving from the edge towards the centre. This aligns with the cooling effect applied at the edge, causing the fluid to cool, increasing its density, and subsequently falling down along the wall.

Concerning the azimuthal velocity, a clear shear is evident between the upper ($z > 0.5$) and lower ($z < 0.5$) regions. This shear was unnoticed without the phase average, because it was masked by temporal fluctuations of similar magnitude to the average flow (not

shown). Note that we observe a shear in the opposite direction compared with [figure 7](#) when the pattern drifts clockwise. In the pattern frame of reference, phase shifts between the fields are observed. The temperature and radial velocity seem to be in phase opposition, while the azimuthal velocity appears to be in phase quadrature with respect to both the temperature and radial velocity (not shown).

We see on the temperature field a well-defined six branches hot pattern at $z = 0.1$, which becomes more diffuse as z increases. Yet, for any given depth and radius, large azimuthal fluctuations are observed, including close to the outer wall. This thermal pattern thus plays a crucial role in the transport of the heat flux to the side, which will be the focus of the next section.

4.4. Wall flux fluctuations

In line with Rayleigh–Bénard studies (e.g. Shang *et al.* (2003, 2004); Gasteuil *et al.* (2007); Lakkaraju *et al.* (2012); Labarre, Fauve & Chibbaro (2023)), we want to quantify heat flux fluctuations on the lateral wall and to understand their connection with both turbulent fluctuations and the large-scale flow structures. To do so, we define the outgoing flux at the wall $r = \Gamma$ by

$$\Phi(t, \varphi, z) = -\frac{\partial T}{\partial r} \Big|_{r=\Gamma}. \quad (4.8)$$

Once the system reaches a statistically stationary state, the global flux balance (by integrating (3.4) over the volume) implies that the flux mismatch between the top and the bottom $(1 - R_F)\pi\Gamma^2$ has to be balanced on time average by the outgoing flux at the side denoted by $\Phi_{side}2\pi\Gamma$. Hence,

$$\Phi_{side} = \frac{(1 - R_F)\Gamma}{2}. \quad (4.9)$$

For the following subsections, the wall heat flux Φ is normalised by this mean side flux Φ_{side} .

4.4.1. The DNS results

In [figure 8](#) we analyse the flux fluctuations in DNS, considering all azimuthal and vertical positions along the side wall over a period of 0.2 thermal diffusive time. The PDFs of the heat flux are depicted for Ra_ϕ ranging from near the instability threshold at $Ra_\phi = 2 \times 10^4$ to the fully turbulent regime at $Ra_\phi = 10^8$. Note that the PDFs remain mostly unchanged when varying the aspect ratio from 4 to 16 at $Ra_\phi = 10^6$, $Pr = 0.1$ and $R_F = 0.1$ (not shown).

With increasing Ra_ϕ , more intense fluctuations across a wider range of values are observed. An exponential tail seems to emerge notably at $Ra_\phi = 10^8$. It is worth noting that Lakkaraju *et al.* (2012) highlighted a similar exponential tail pattern in the heat flux along the sidewall of a cylindrical Rayleigh–Bénard system, with the slope increasing with Ra_ϕ . Additionally, Shang *et al.* (2003, 2004), in a Cartesian Rayleigh–Bénard configuration, identified a similar exponential behaviour in the distribution of both vertical and horizontal bulk heat fluxes. This exponential tail is not observed close to the onset of instability, but requires reaching sufficiently turbulent conditions.

The large-scale drifting pattern is also responsible for intense fluctuations in heat flux near the wall. To illustrate this point, we performed a DNS at $Ra_\phi = 10^7$, $\Gamma = 4$, with $Pr = 0.1$ and the flux ratio at $R_F = 0.1$. [Figure 9\(a\)](#) shows a snapshot of the temperature field, seen from below. The largest temperatures are localised along four drifting branches,

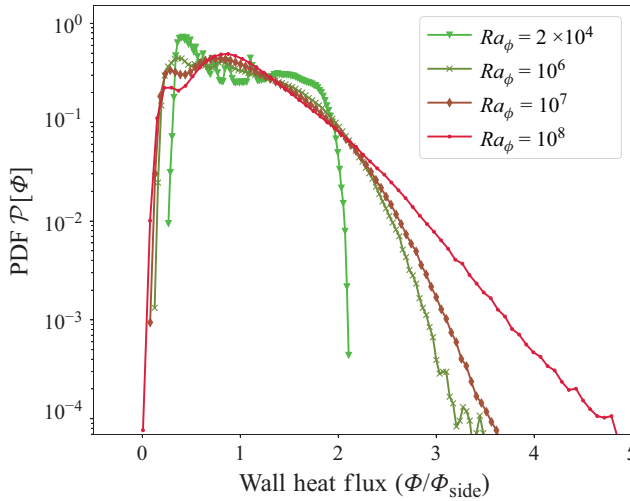


Figure 8. Wall heat flux PDF from DNS with different values of Ra_ϕ . The fluxes have been scaled by the time-averaged flux Φ_{side} . Input parameters are $Pr = 0.1$, $R_F = 0.1$ and $\Gamma = 4$.

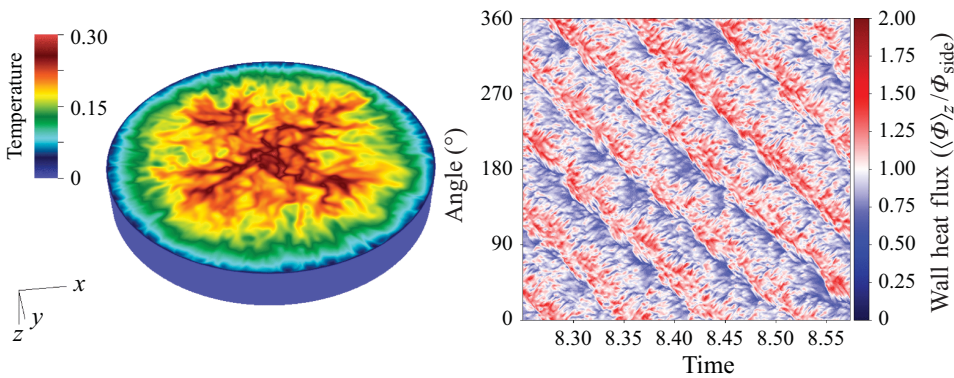


Figure 9. (a) Three-dimensional view of the temperature field at the bottom plate (note the upside-down axes system with z pointing downward) with $\Gamma = 4$, $Ra_\phi = 10^7$, $R_F = 0.1$ and $Pr = 0.1$. (b) Hovmöller (space-time) diagram from this simulation of the wall heat flux averaged in z and scaled by the mean wall flux Φ_{side} .

as more clearly identified in the space–time diagram in [figure 9\(b\)](#). Within these branches, intense fluctuations involve heat fluxes up to two times the average flux. Once scaled for our air experiment, the characteristic size and duration of one of these fluctuations are approximately 4 cm and 1.5 s, respectively.

We now want to disentangle the contributions to the side heat flux fluctuations from the large-scale drifting pattern and those from the small-scale turbulent convective fluctuations. [Figure 10\(a\)](#) shows the wall heat flux PDF after applying phase averaging: it thus focuses on the effect of the large-scale drifting pattern only. Note that similar analyses have been conducted in turbulent Rayleigh–Bénard systems in cylindrical geometry by [Lakkaraju *et al.* \(2012\)](#) to assess how the local heat flux varies depending on the measurement location relative to the orientation plane of the large-scale circulation.

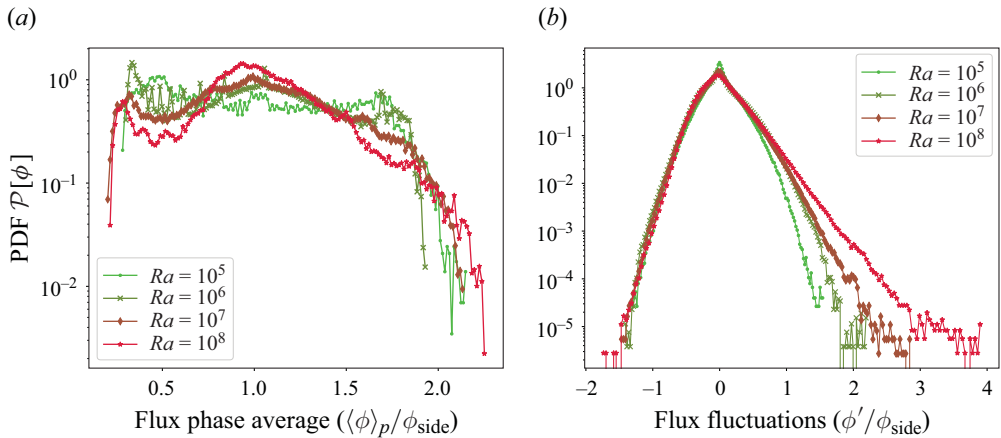


Figure 10. (a) The PDF of the phase-averaged wall heat flux as a function of Ra_ϕ . The other input parameters are $Pr = 0.1$, $R_F = 0.1$ and $\Gamma = 4$. All depths and azimuths along the side wall have been considered. (b) Same but for the wall heat flux fluctuations related to the phase average. For both plots, the fluxes have been scaled by the time-averaged flux Φ_{side} .

We observe that the flux distribution carried by the pattern extends over a range from 0.5 to 2 times the mean flux Φ_{side} . This range barely depends on the Rayleigh number. Indeed, the drifting pattern is essentially responsible for the transport of the mean flux (Φ_{side}) out of the system, which is independent of Ra_ϕ . Increasing the Ra_ϕ increases the rapid turbulent fluctuations, which will be considered next. However, the shape of the PDFs changes with Ra_ϕ . Indeed, the observed pattern evolves as Ra_ϕ increases, notably by varying the number of branches and increasing the skewness of the flux profile associated with a branch (not shown).

Figure 10(b) shows the PDF of the wall heat flux fluctuations related to the phase-averaged operator: we thus focus here on the turbulent fluctuations taking place around the phase average field. As the Rayleigh number increases, the range of heat flux fluctuations widens. We also see the emergence of the exponential tail at $Ra_\phi = 10^8$, previously identified in figure 8. This means that the largest and rarest heat flux events are associated with turbulent fluctuations, and not with the pattern itself. Their probability will keep increasing with Ra_ϕ . It is also worth noting that the left portion of the fluctuations, with the amplitude ranging from -2 to 0 , appears to be unaffected by changes in the Rayleigh number, showing a consistent overlap. Since $\phi' = \phi - \langle\phi\rangle_p$ and the heat flux is consistently positive for $R_F = 0.1$, negative ϕ' values suggest low heat fluxes ϕ and high heat fluxes caused by the pattern's drift $\langle\phi\rangle_p$. These occurrences indicate scenarios where the observed heat flux primarily arises from the pattern's drift. This finding aligns with the observation that the pattern's drift consistently induces similar fluctuations regardless of the Rayleigh number (refer to figure 10a).

In summary, the heat flux fluctuations within the system exhibit a broad distribution that can be attributed to two main contributions: the pattern contribution, which transports the mean flux Φ_{side} with a fixed amplitude independent of Ra_ϕ , and the fluctuations, which exhibit wider distributions and exponential tails as Ra_ϕ increases. In light of these

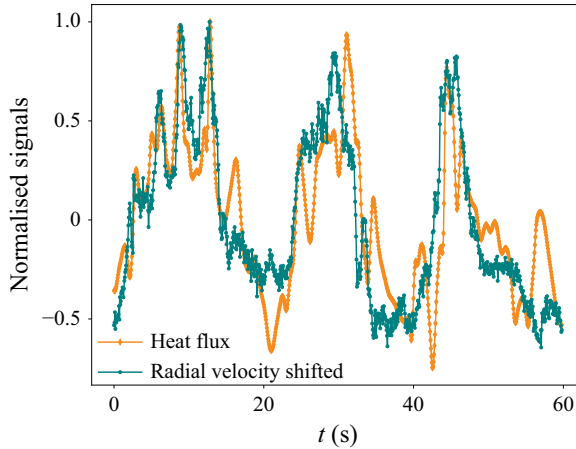


Figure 11. Temporal evolution of the wall heat flux measured by our probe spanning a width of 18 mm and a depth $z \in [4, 6 \text{ cm}, 6, 4 \text{ cm}]$ (orange), and radial velocity measured by PIV at $r = 34 \text{ cm}$, $z = 6 \text{ cm}$, in the direction aligned with the flux sensor (green). The time is shifted by 2.6 s for the velocity. For each signal, the mean temporal component has been removed to keep only the fluctuating part, which is then normalised by the maximum value of the signal. Data from experiment 3 in [table 1](#).

numerical observations, how does experimental analysis contribute to our understanding of the heat flux fluctuations?

4.4.2. Experimental results

In [figure 11](#) we plot the time evolution of the wall heat flux measured experimentally as well as the radial velocity at a location situated 10 cm from the flux sensor in the radial direction towards the centre. The purpose of this measurement is to explore the relationship between radial velocity and heat flux measured at the wall. Upon examination, we observe that when we shift the radial velocity by a characteristic time $\tau_{adv} = 2.6 \text{ s}$, the two rescaled signals reasonably match. Moreover, the characteristic time τ_{adv} aligns with an advective time scale, namely d/U , where d represents the distance between the flux sensor and the location of the velocity measurement and U corresponds to the time-average radial velocity at that specific location. This alignment suggests a correlation between the radial velocity and the wall heat flux. This is compatible with the numerical results shown in [figure 9\(b\)](#), and indicates that the drifting branches associated with higher radial velocity values act as pathways for transporting heat fluctuations towards the wall.

In [figure 12\(a\)](#) we compare the PDF of the wall heat flux derived from experimental measurements and from numerical simulations. The heat flux is measured experimentally at mid-height and averaged over the entire surface of the sensor. It is positioned halfway up the sidewall, extending end to end from $z = 4.6 \text{ cm}$ to $z = 6.4 \text{ cm}$, with a width of 1.8 cm. We consider experiment $9n^{o}9$ from [table 1](#), where the non-dimensional parameters have been estimated by $Ra_\phi = 1.26 \times 10^8$ and $R_F = 0.54$. Numerically, we fix $Ra_\phi = 10^8$, $\Gamma = 4$, $Pr = 0.7$ and consider the flux ratio $R_F = 0.6$. Given the numerical difficulty to get statistics over long periods, we assume system ergodicity. As a result, the numerical flux calculations consider all azimuthal events at each output snapshot, limited in height to match the depth of the flux sensor ($z \in [4.6 \text{ cm to } 6.4 \text{ cm}]$).

When comparing the PDFs, we initially note that the numerical PDF predicts much larger fluctuations than those measured experimentally. However, it should be noted that

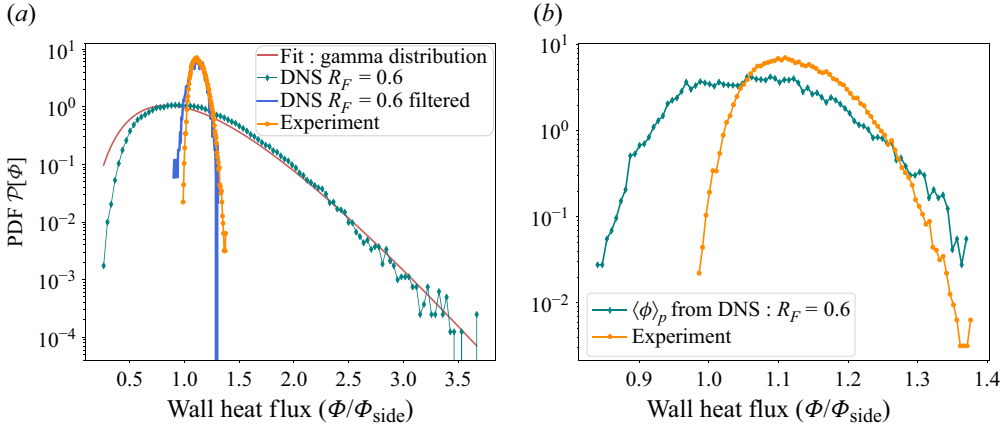


Figure 12. (a) Wall heat flux PDF for experiment 9n09 in table 1 (orange \circ) and for a DNS with $Ra_\phi = 10^8$, $\Gamma = 4$, $Pr = 0.7$, $R_F = 0.6$ (green \diamond). The continuous blue thick line represents the low-pass filtered DNS results. The thin red line represents a fit by a gamma distribution. The fluxes are all scaled by the time-averaged flux Φ_{side} . (b) Same for experiment 9n09 and for DNS heat flux fluctuations due to the drifting pattern only (see § 4.4.1).

the sensor has a response time (T_r) of 0.7 s, meaning that any event occurring over a period shorter than this will not be directly measured. To account for this experimental aspect, we apply a first-order low-pass temporal filter, with a cutoff frequency (f_c) corresponding to the flux sensor’s response time ($f_c = 1/T_r$). The application of this filter is done in Fourier space by multiplying the Fourier transform of the experimental signal by the transfer function of a first-order low-pass filter. We then note a good qualitative concordance between DNS and experimental results. *A priori*, a two-dimensional spatial filter should be considered, since the heat sensor averages over its entire surface. However, the good agreement suggests this is unnecessary, as short events correlate with small-scale phenomena. Note that we did the same numerical computation but for $R_F = 0.5$, and the same qualitative agreement was found. As a result, the most extreme events then appear to have a relatively brief duration, at least shorter than the response time associated with the sensor.

The shape of the numerical PDF of figure 12(a) is compatible with a gamma distribution, denoted $\mathcal{P}_\gamma[\Phi]$ and defined by

$$\mathcal{P}_\gamma[\Phi] = \frac{1}{\gamma(N)\Psi^N} \Phi^{N-1} e^{-\Phi/\Psi}, \tag{4.10}$$

where $\gamma(N)$ represents the gamma function with N and Ψ the two fitting parameters. Here N is an integer without dimensionality, while Ψ has a dimension of a heat flux. Since in our case the mean flux is imposed, Ψ is constrained to Φ_{side}/N . The distribution now has only one free parameter, N , which is set to 6 in figure 12(a), but depends on the input parameters Ra_ϕ , R_F , Pr and z due to vertical inhomogeneity. In figure 8 for example, the PDF accounting for all events along the vertical direction z at $R_F = 0.1$ and $Ra_\phi = 10^8$ is best fitted by a gamma distribution with $N = 4$.

In figure 12(b) we analyse the PDF of wall heat flux, comparing experiment 9 with the same DNS conducted at $Pr = 0.7$, $\Gamma = 4$, $R_F = 0.6$ and $Ra_\phi = 10^8$, but considering

only the heat flux distribution due to the drifting pattern. Both PDFs are similar. This makes sense since the drifting pattern is persistent and drifts slowly: it is hence easily caught by the sensor despite its limited response time. The heat sensor however misses the rapid, turbulent fluctuations described in § 4.4.1. The experimental PDF also appears to have fewer values with $\Phi/\Phi_{side} < 1$ compared with the DNS results. While a perfect match between the two PDFs is not expected, the remaining discrepancy likely stems from the fact that the flux sensor averages over its surface, a factor not accounted for in the numerical data.

5. Conclusion and future work

In conclusion, the experiments presented in the present paper allowed us to observe a $Ra_\phi^{-1/5}$ scaling law for the mean temperature, consistent with the scaling identified in Rein *et al.* (2023a,b), and extended over nearly two orders of magnitude in Rayleigh number, reaching $Ra_\phi \simeq 10^{10}$ values commensurate with those of the nuclear safety application that initially motivated our work. Additionally, a $Ra_\phi^{1/3}$ scaling law for the r.m.s. horizontal velocity has been identified. This behaviour is explained through dimensional analysis, suggesting that the scaling originates from horizontal heat transport in the turbulent regime. In addition to scaling laws for globally averaged quantities including the r.m.s. velocity and the temperature, experimental measurements have confirmed, even for the largest values of Ra_ϕ , the presence of a mesoscale, persistent branch pattern associated with large-scale, slow wall heat flux fluctuations. These superimpose to small-scale, rapid, high amplitude fluctuations associated with turbulent fluctuations.

Regarding the nuclear safety application, we are interested in understanding heat transfer through the studied metal layer towards the lateral metal wall, in order to predict its melting leading possibly to vessel failure. We argue that the turbulent fluctuations may be secondary: indeed, even if they involve events associated with the most extreme fluxes, they manifest in very localised spatial areas and for short durations, and hence, correspond energetically to weak events. On the other hand, the thermal branches, associated with slower and larger events, play a significant role in heat exchanges, as they cover substantial portions of the wall and persist for longer durations.

To further explore the influence of these fluctuations on the structural integrity of the wall, it would be valuable to examine a global melting scenario, incorporating both convective flow and melting phenomena. This sort of challenging problem has been investigated across different domains, spanning geophysics (Alboussière, Deguen & Melzani 2010; Labrosse *et al.* 2018) and industrial applications (Kalhori & Ramadhyani 1985; Fragnito *et al.* 2022). Published studies encompass Rayleigh–Bénard convection (Labrosse *et al.* 2018; Favier, Purseed & Duchemin 2019) as well as natural convection (Jany & Bejan 1988; Viskanta 1988; Lacroix & Arsenault 1993), particularly relevant to our case. The characteristic feature of our problem is the periodic heat forcing due to the pattern rotation, which may lead to various melting regimes (Ho & Chu 1993).

From a fundamental point of view, the next step is to tackle the instability at the origin of the drifting pattern. Physical insights into the underlying mechanism could indeed be useful to better forecast the heat flux fluctuations. This will be the focus of our future work.

Supplementary movie. Supplementary movie is available at <https://doi.org/10.1017/jfm.2025.2>.

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had no influence on the design, data collection, analysis, interpretation of results or the decision to publish. The content of this work remains the sole responsibility of the authors.

Appendix A

A.1. Summary of the experiments/simulations parameters

Tables 1 and 2 provide all relevant parameters for the performed experiments and DNS, respectively.

A.2. Estimation of the non-dimensional numbers within a gas mixture

For reaching the most extreme values of Ra_ϕ experimentally, we use SF_6 . However, SF_6 is not pure in the system; it is mixed with air. The evaluation of the Rayleigh and Prandtl numbers therefore requires a precise knowledge of the physical properties of the air/ SF_6 mixture under consideration. For a perfect gas mixture, the mass density (ρ_m) as well as the specific heat capacity (c_{p_m}) of the mixture are linearly expressed as a function of the mass fraction of SF_6 (denoted y), as follows:

$$\begin{cases} \rho_m = (1 - y)\rho_{air} + y\rho_{SF_6}, \\ c_{p_m} = (1 - y)c_{p_{air}} + yc_{p_{SF_6}}. \end{cases} \quad (A1)$$

Since the densities of SF_6 and air are tabulated as functions of temperature and pressure conditions, the experimental measurement of the mean density ρ_m enables the determination of the SF_6 mass fraction. This, in turn, allows for the estimation of the heat capacity of the gas mixture. Regarding dynamic viscosity and thermal conductivity, the kinetic theory of gases allows for the determination of these properties (Chapman & Cowling 1991), but it is difficult to apply because it involves numerous degrees of freedom. Simpler models allow for a relatively accurate estimation of these quantities. The model proposed by Wilke (1950) provides an estimation of the viscosity of a gaseous mixture (tested up to seven species) accurate to 1.9%. The viscosity of the air/ SF_6 mixture, denoted as μ_m , is expressed as

$$\mu_m = \frac{\mu_{air}}{1 + A \frac{x_{SF_6}}{x_{air}}} + \frac{\mu_{SF_6}}{1 + \frac{1}{A} \frac{x_{air}}{x_{SF_6}}}, \quad (A2)$$

where x_{air} and x_{SF_6} represent the mole fraction of air and of SF_6 . Here A is the Sutherland constant (Sutherland 1895), $A = ((M_{SF_6}/M_{air})(\mu_{air}/\mu_{SF_6}))^{1/2}$, where M_{SF_6} and M_{air} are the molar masses of SF_6 and air. Concerning thermal conductivity, a similar model has been proposed by Lindsay & Bromley (1950). The difference between the Wilke model (viscosity) and the Lindsay and Bromley model (conductivity) lies in a variation of the Sutherland constant.

Figures 13(a) and 13(b) show how the Rayleigh number and the Prandtl number evolve as a function of the system's density. This evolution is studied using two different models: one with linear models for conductivity and viscosity (similar to (A1)), and the other where the Wilke model (viscosity) and Lindsay/Bromley model (conductivity) are used to estimate these properties.

Significant deviation between these two estimates is observed for densities around 3 kg m^{-3} , with a deviation exceeding 50% for the Rayleigh number and approximately 20% for the Prandtl number. This large Rayleigh number deviation is mainly due to its quadratic dependency on conductivity. Henceforth, the Wilke and Lindsay/Bromley models are used for estimating the viscosity/conductivity of the air/ SF_6 mixture.

No	P (W)	Gas	ρ_{eff} (kg m ⁻³)	$\langle\theta\rangle$ (°C)	R_F	Ra_ϕ	Pr	Γ
1	75	Air	1.2	28.7	0.43	3.2×10^7	0.71	4
2	100	Air	1.2	30.5	0.44	4.0×10^7	0.71	4
3	100	Air	1.2	31.4	0.48	4.6×10^7	0.71	4
4	150	Air	1.2	35.6	0.57	6.8×10^7	0.71	4
5	150	Air	1.2	35.3	0.58	7.0×10^7	0.71	4
6	150	Air	1.2	36.6	0.54	6.75×10^7	0.71	4
7	200	Air	1.2	40.4	0.57	9.26×10^7	0.71	4
8	200	Air	1.2	42.3	0.54	9.88×10^7	0.71	4
9	250	Air	1.2	48.2	0.54	1.26×10^8	0.71	4
10	100	Air + SF ₆	2.9	32.1	0.54	8.02×10^8	0.87	4
11	170	Air + SF ₆	3.4	38.6	0.57	1.61×10^9	0.87	4
12	400	Air + SF ₆	4.1	61.5	0.56	7.6×10^9	0.85	4

Table 1. Experimental parameters for the 12 performed cases. Here P is the injected power in watts (W).

Ra_ϕ	Γ	Pr	R_F	DNS/filtered	\mathcal{E}	N	η_K/L	BC
10^2	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[2688; 9216; 33792]	8	[5.8; 4.4; 3.1]	FS
10^3	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[2688; 9216; 33792]	8	[5.8; 4.4; 3.1]	FS
10^3	[4; 8]	0.1	0.5	DNS	[2688; 9216]	8	[6.2; 3.7]	FS
10^4	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[2688; 9216; 33792]	10	[3.3; 2.7; 2.3]	FS
10^4	[4; 8]	0.1	0.5	DNS	[2688; 9216]	10	[3.3; 2.5]	FS
3×10^4	[4; 8; 16]	0.1	0.1	DNS	[2688; 9216; 33792]	10	[2.5; 2; 1.8]	FS
10^5	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[2688; 9216; 33792]	10	[1.88; 1.56; 1.52]	FS
10^5	4	0.7	[0.5; 0.6]	DNS	9216	10	4	NS
10^5	[4; 8]	0.1	0.5	DNS	[2688; 9216]	10	[1.87; 1.52]	FS
3×10^5	[4; 8; 16]	0.1	0.1	DNS	[2688; 9216; 33792]	10	[1.43; 1.35; 1.36]	FS
10^6	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[9984; 33608; 33792]	10	[4; 3.3; 1.22]	FS
10^6	4	0.7	[0.5; 0.6]	DNS	9216	10	2.2	NS
10^6	[4; 8]	0.1	0.5	DNS	[9216; 33608]	10	[3.7; 3.22]	FS
3×10^6	[4; 8; 16]	0.1	0.1	DNS	[9984; 33608; 33792]	10	[3; 2.5; 1.12]	FS
10^7	[4; 8; 16]	0.1	[0.1; 0.9]	DNS	[9984; 33608; 36608]	10	[2.2; 1.8; 1]	FS
10^7	4	0.7	[0.5; 0.6]	DNS	9984	10	1.3	NS
10^7	[4; 8]	0.1	0.5	DNS	[9984; 33608]	10	[2.3; 1.74]	FS
3×10^7	4	0.1	0.1	DNS	9984	10	1.8	FS
3×10^7	[8; 16]	0.1	[0.1; 0.9]	Filtered	[33792; 36608]	10	[1.44; 0.78]	FS
10^8	4	0.1	[0.1; 0.9]	DNS	9984	10	1.27	FS
10^8	[4; 8]	0.1	0.5	[DNS; filtered]	[9984; 33792]	10	[1.4; 1.3]	FS
10^8	4	0.7	[0.5; 0.6]	DNS	33608	10	2.5	NS
10^8	[8; 16]	0.1	[0.1; 0.9]	Filtered	[33792; 36608]	10	[1.33; 0.62]	FS
3×10^8	[4; 8]	0.1	[0.1]	Filtered	[9984; 33792]	10	[1.1; 1]	FS

Table 2. Simulations summary (DNS or filtered) according to the physical and numerical parameters. Here \mathcal{E} , N , BC and η_K/L represent the number of hexahedral elements, the polynomial spectral order, the top boundary conditions applied (FS denotes free slip and NS denotes no slip) and the ratio between the Kolmogorov dissipative scale (η_K) and the numerical grid size (L), respectively. For further details, readers are directed to Rein *et al.* (2023a).

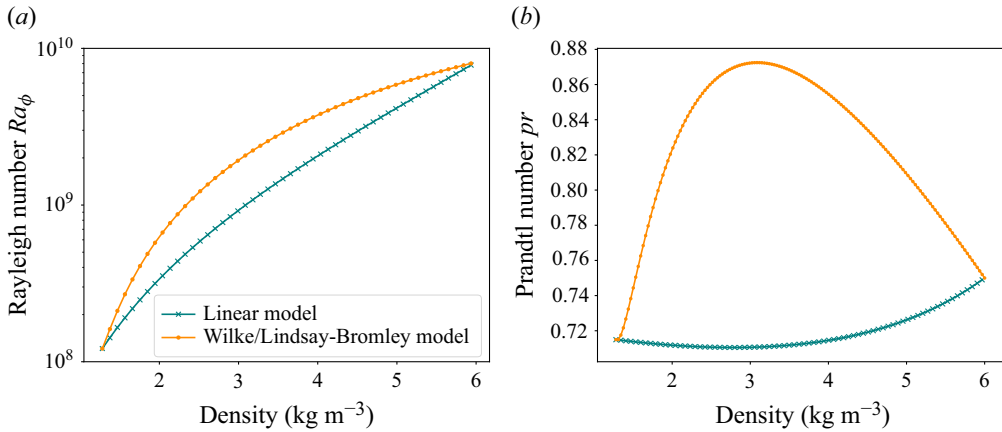


Figure 13. Evolution of (a) Ra_ϕ and (b) Pr as a function of the density of the air/SF₆ mixture, considering either a linear mixture model for the viscosity and thermal conductivity or the Wilke (1950)/Lindsay & Bromley (1950) model. In all cases, $\phi = 200 \text{ W m}^{-2}$ and $H = 11 \text{ cm}$.

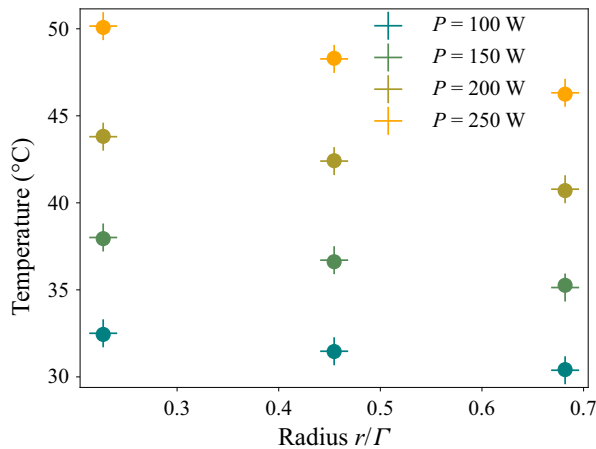


Figure 14. Average temperature as a function of radius for angular positions $\phi = 0^\circ$ (\bullet) and $\phi = 90^\circ$ ($+$), at different powers ranging from 100 W to 250 W.

A.3. Experimental setup: energy balance

In this section we evaluate the thermal balance of the experimental set-up in order to quantify both the thermal losses and the time-averaged ratio between the imposed heat flux and the one escaping through the lid (R_F). This last estimate is necessary for comparison with numerical simulations.

A.3.1. Input heat flux and losses

In figure 14(a) the temperature field from the camera, averaged over these 3 min, shows perfect uniformity, indicating a uniform heat flux. First, to ensure system azimuthal uniformity, we use thermal measurements from the thermocouples. In figure 14(b) the time-averaged temperature of each thermocouple is plotted against its location radius, for two specific angular positions $\phi = 0^\circ$ (\bullet) and $\phi = 90^\circ$ ($+$). These measurements were

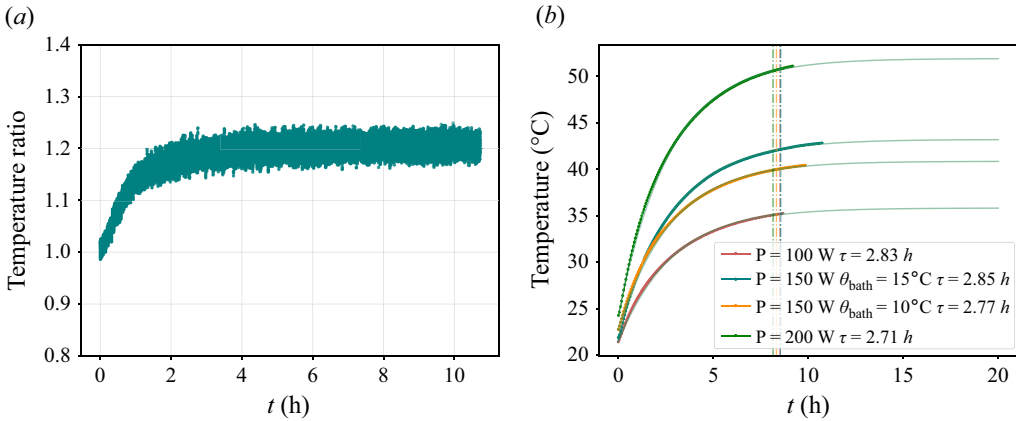


Figure 15. (a) Time evolution of the ratio between the temperature of the aluminium plate and the average temperature within the system, averaged over the six thermocouples (θ/θ_{gas}). The set power is $P = 150$ W and the bath set point temperature is $\theta_{bath} = 15$ °C. (b) Temporal evolution of the temperature at the centre of the aluminium plate measured via the PT 100 positioned just tangent to the plate, for various powers and bath temperatures (θ_{bath}). Vertical lines indicate when time reaches 3τ , and solid green lines represent the best fit of the data using (A4) as a model.

taken for different powers, ranging from 100 W to 250 W. The average temperatures are independent of their angular position, confirming the azimuthal homogeneity of the system.

Now, let us evaluate the heat losses in our set-up. When applying electrical power P to the eight heating elements, most of the flux is supplied to the aluminium plate, but a fraction escapes into the insulating plate. Let us introduce the heat efficiency coefficient, denoted χ and defined by $\chi = \phi_{in}/\phi_{set}$, where ϕ_{in} and ϕ_{set} represent the heat flux supplied to the aluminium plate and to the heat mats (calculated as $P/8S_{heat}$, with S_{heat} as the heat mat surface area). In the following, we outline two methods to estimate χ .

The first method, named the transient method, involves deriving the equation governing the time evolution of the aluminium plate temperature from an energy balance on the plate. The change in internal energy of the plate is due to the heat flux imposed by the heating elements, as well as the convective flux of the gas at the plate surface. This is written as

$$\rho_{Al}c_{pAl}e_{Al}\frac{d\theta}{dt} = \phi_{in} - h(\theta - \theta_{gas}), \tag{A3}$$

where ρ_{Al} , c_{pAl} , e_{Al} , θ respectively represent the density, specific heat capacity, thickness and temperature (assumed uniform) of the aluminium plate. Additionally, h and θ_{gas} represent the convective heat transfer coefficient between the gas and the plate, and the ambient gas temperature. It should be noted that θ_{gas} is not constant and increases during the heating period. However, after a short transient phase, figure 15(a) demonstrates that the ratio between the temperature of the aluminium plate measured at the centre and the average temperature within the system (averaged over the six thermocouples) remains globally constant over time, despite some fluctuations. Consequently, we consider θ_{gas} to be proportional to θ by introducing $(1 - \alpha)$, the proportionality factor, defined as $\theta_{gas} = (1 - \alpha)\theta$. Hence, (A3) leads to an exponential profile defined by the equation

$$\Delta\theta(t) = \frac{\phi_{in}}{\alpha h} (1 - e^{-t/\tau}), \tag{A4}$$

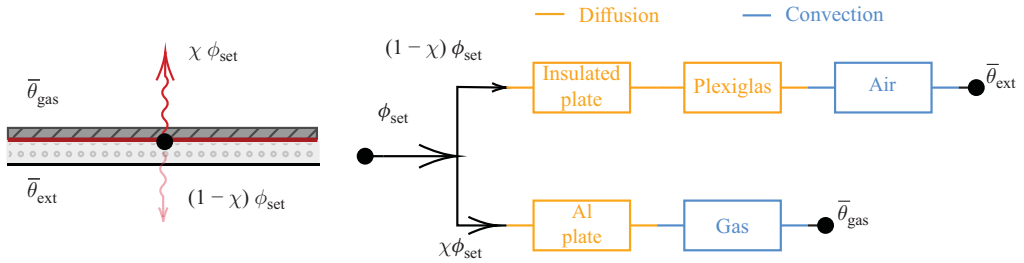


Figure 16. Left: sketch of the aluminium plate and the heat fluxes distribution due to the heating mats. Right: thermal circuit of the aluminium plate area, the orange and blue rectangles representing diffusive and convective thermal resistances, respectively.

where $\Delta\theta(t) = \theta(t) - \theta_{gas}(t = 0)$ and $\tau = \rho_{Al}c_{pAl}e_{Al}/\alpha h$ represents the time constant of the system. Figure 15(b) illustrates the temporal evolution of the temperature at the centre of the aluminium plate measured via the PT100 probe, for different powers and bath temperatures. The green line on these profiles corresponds to a fit performed using the least squares method based on the solution (A4). A very good agreement between this simple exponential model and the experimental measurements is observed. The fitting parameters $\Delta\theta_{max}$ (asymptotic value of $\Delta\theta(t)$ as $t \rightarrow \infty$) and τ then allow estimation of the effective heat flux supplied to the system as $\phi_{in} = \rho_{Al}c_{pAl}e_{Al} \Delta\theta_{max}/\tau$ and consequently the heat efficiency coefficient χ .

The second method, referred to as the steady method, involves conducting an energy balance under steady-state conditions. Therefore, a thermal circuit representation of the different fluxes and temperatures can be drawn (as illustrated in figure 16). The heat flux from the heating mats follows two paths. The main fraction conducts through the aluminium plate (thickness e_{Al} , conductivity k_{Al}), inducing convective flow in the system, characterised by the convective heat transfer coefficient h . In a steady state, h can be determined using the thermal circuit (figure 16) as $h = \chi\phi_{set}/(\bar{\theta} - \bar{\theta}_{gas})$, where $\bar{\theta}$ and $\bar{\theta}_{gas}$ represent the steady-state temperatures of the aluminium plate and the gas within the system, respectively. This path introduces a thermal resistance denoted as $R_{Al} = e_{Al}/k_{Al} + 1/h$. The remaining heat travels through the insulating plate (thickness e_{IP} , conductivity k_{IP}) and the plexiglass support (thickness e_{plex} , conductivity k_{plex}) before contributing to small convective motion in the ambient air, characterised by the convective heat transfer coefficient h_{air} (determined via the empirical free convection correlation of a hot horizontal plate heating from above; see, e.g. Lloyd & Moran (1974)). This path is characterised by a thermal resistance noted as $R_{ext} = e_{IP}/k_{IP} + e_{plex}/k_{plex} + 1/h_{air}$. Writing the global heat flux conservation, the heat efficiency coefficient can be expressed as

$$\chi = \frac{1}{1 + \frac{R_{Al}}{R_{ext}}} - \frac{\bar{\theta}_{gas} - \bar{\theta}_{ext}}{\phi_{set} (R_{Al} + R_{ext})}, \quad (A5)$$

with $\bar{\theta}_{ext}$ representing the room temperature. The first term on the right-hand side corresponds to heat distribution due to thermal resistance ratio, while the second term represents a heat loss due to the temperature difference between the system and the ambient environment. It reflects the ratio between the flux due to the temperature difference across all layers (aluminium plate, insulated plate, etc.) and the flux set by the heating mats. Notably, this term is relatively small compared with the first one, as $\phi_{set} \gg (\bar{\theta}_{gas} - \bar{\theta}_{ext})/(R_{Al} + R_{ext})$ and R_{ext} is roughly 10 times larger than R_{Al} , making the first term the dominant

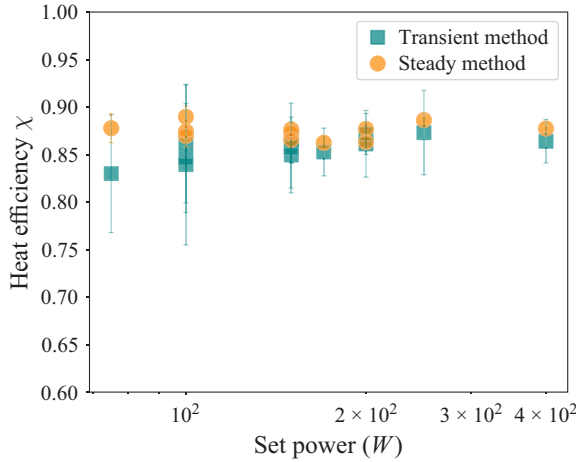


Figure 17. Heat efficiency coefficient $\chi = \phi_{in}/\phi_{set}$, as a function of the set power P using the transient method (\square) and the steady method (\circ).

factor. We measured $\bar{\theta}$ using a PT100 sensor on the aluminium plate and $\bar{\theta}_{gas}$ using the six thermocouples, during steady state (after 3τ) and averaged these readings over time. These measurements provided the heat efficiency χ . In figure 17, χ is plotted against the set power P using both transient and steady methods. Both methods show good agreement, with an average thermal loss of 15%. Error bars represent uncertainty propagation, calculated according to (A4) for the transient method and (A5) for the steady method, with contributions from the PT100 sensor uncertainty and the variances of the fit coefficients related to the aluminium plate temperature profiles. Larger discrepancies at lower power levels are due to difficulties in fitting the exponential profile, where small temperature variations have a greater impact at lower power, leading to increased uncertainty.

A.3.2. Flux ratio determination

In our numerical model, a uniform and constant flux escaping from the upper surface is imposed. In the experimental set-up, a rigid cover allows for a conductive heat flux, which is therefore non-uniform and fluctuating. This configuration resembles the real situation, where the resulting radiative flux is a dynamic consequence of the system's heat transfers. However, an average flux escaping from the upper surface can be estimated to evaluate a flux ratio R_F and compare the results with numerical simulations. In this subsection we outline two methods to estimate it. The room temperature is one of the parameters that can influence the top heat flux, similarly to the lid thickness and the top temperature in the system below the lid. The room temperature is not controlled, but measured for each experiment and does not significantly change during the period of the data acquisitions (approximately 1 h).

The first method, named the lid method, involves performing a thermal balance on the lid. To minimise disturbances to the flow, thin thermocouples (much thinner than PT-100) are carefully positioned above and just below the lid at the centre to take the necessary measurements. Through these measurements, the temperature difference between the bottom and top of the lid can be averaged and related to the flux escaping through conduction in the cover ($R_F\phi_{in}$) using the relation

$$R_F\phi_{in} = k_{Lid} \frac{\Delta\theta_{Lid}}{e_{Lid}}, \quad (\text{A6})$$

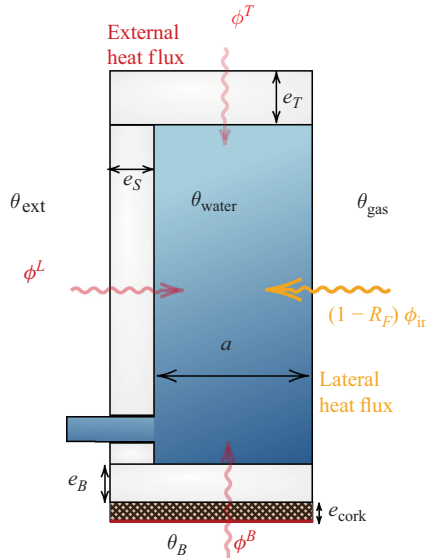


Figure 18. Sketch representing the energy balance of the water within the torus.

where k_{Lid} , e_{Lid} , $\Delta\theta_{Lid}$ represent the thermal conductivity, the thickness of the lid and the temperature difference between the bottom and top of the lid, respectively. Note that this flux ratio estimation is based on a local temperature measurement, so it does not account for system inhomogeneity.

The second method, called the bath method, relies on the thermal analysis of the cooling water on the lateral surface, providing a comprehensive measurement of the thermal fluxes involved in the system. The thermal balance of the water within the torus can be formulated as

$$\underbrace{\rho_e c_{pe} Q_v \Delta\theta_e}_{\text{Change in water internal energy}} = \underbrace{(1 - R_F)\phi_{in} S}_{\text{Lateral flux}} + \underbrace{\phi_{ext} S_{lat}}_{\text{External flux}}, \quad (\text{A7})$$

leading to the average flux ratio

$$R_F = 1 - \left[\frac{\rho_e c_{pe} Q_v \Delta\theta_e}{\phi_{in} S} - \frac{2\phi_{ext}}{\Gamma\phi_{in}} \right]. \quad (\text{A8})$$

Here, ρ_e , c_{pe} , Q_v , S_{lat} , S and $\Delta\theta_e$ represent the density, specific heat capacity, volumetric flow rate of water, the lateral ($S_{lat} = 2\pi RH$) and the heating surface area ($S = \pi R^2$), and the temperature difference between the water inlet and outlet in the bath, respectively.

Two PT 100 sensors enable the measurement of $\Delta\theta_e$. Additionally, a volumetric water meter measures the average flow rate of the system over the entire experiment duration (approximately 1 h). Equation (A7) indicates that the water undergoes a change in internal energy as it passes through the torus, due to the lateral heat flux of the system and the perturbative fluxes induced by the environment. In (A7) the lateral heat flux is expressed as the difference between the flux imposed on the bottom surface ($\phi_{in} S$) and that escaping from the cover ($R_F\phi_{in} S$). The term ϕ_{ext} denotes the perturbative heat flux from the external environment, heating the bath water. As sketched in figure 18, it decomposes into three contributions ϕ^L , ϕ^T and ϕ^B , representing the external heat fluxes from the lateral

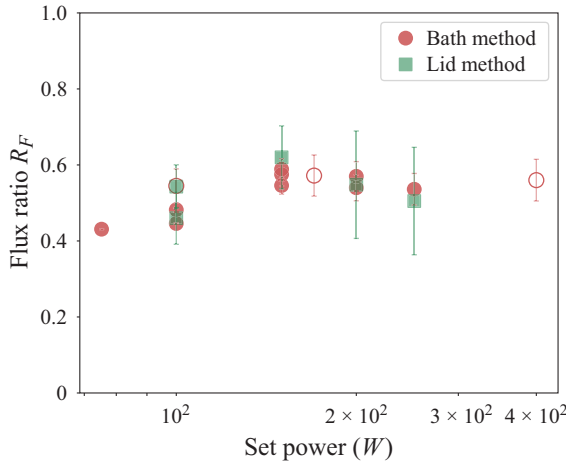


Figure 19. Estimated flux ratio R_F as a function of the set power, considering Lid (\square) and Bath (\circ) methods. The empty symbols indicate SF_6 gas is used.

surface, the upper surface of the torus and the lower surface of the torus due to the heating elements:

$$\phi_{ext} = \phi^L + (\phi^T + \phi^B) \frac{a}{H} \left(1 + \frac{a}{2R} \right). \quad (\text{A9})$$

Here, a represents the gap between the inner and outer walls of the torus. The perturbative fluxes can be expressed in terms of the torus characteristics and system temperatures as

$$\phi^L = \left(\frac{e_L}{k_{PMMA}} + \frac{1}{h_{ext}} \right)^{-1} (\theta_{water} - \theta_{ext}), \quad (\text{A10})$$

$$\phi^T = \left(\frac{e_T}{k_{PMMA}} + \frac{1}{h_{ext}} \right)^{-1} (\theta_{water} - \theta_{ext}), \quad (\text{A11})$$

$$\phi^B = \left(\frac{e_{cork}}{k_{cork}} + \frac{e_B}{k_{PMMA}} \right)^{-1} (\theta_{water} - \theta_B), \quad (\text{A12})$$

where k_{PMMA} , e_B , e_L , e_T , θ_{water} , θ_{ext} , θ_B represent respectively the thermal conductivity of PMMA, the thickness of the bottom, side and top of the torus, the temperature of the water, external environment and of the contact between the heating element and the cork. The convective heat transfer coefficient h_{ext} characterises the vertical convection generated by the external side wall and external air. While strictly speaking, separate h values apply to the vertical surface and the flat top, their estimated values are of the same order of magnitude (see, e.g. Fujii & Imura (1972)). Therefore, we use the same h value for both, based on the classical vertical convection correlation (Churchill & Chu 1975).

During the measurement campaign, consistent efforts were made to maintain a temperature difference of approximately 5 °C between the bath water and the external environment. This approach was intended to keep the same external perturbative flux across all experiments. An estimation of the perturbative term ϕ_{ext} was performed and represents an average of approximately 12 % of the imposed flux across all experiments.

In figure 19 the estimation of the flux ratio using either the global method (bath method) or the local method (lid method) is plotted against the power supplied to the heating elements. Error bars represent uncertainty propagation, calculated according to (A6) for

the lid method and (A7) for the bath method, with contributions from the PT100 sensor, thermocouple uncertainties, as well as ϕ_{in} uncertainty related to the variances of the fit coefficients from (A4). It is observed that the average flux ratio depends on the set power, ranging from approximately 0.4 to around 0.6. The global measurement using the bath provides a good experimental repeatability, with a measurement uncertainty of 0.02 for the measurement at 150 W. Furthermore, the local measurements using the lid are consistent and in good agreement with the global ones.

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