



Model for thermal convection with uniform volumetric energy sources

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A theoretical model is derived to predict the heat fluxes at the upper and lower horizontal surfaces of an internally heated (IH) convection cell by extending the well-known Grossmann & Lohse (*Phys. Rev. Lett.*, vol. 86, 2001, pp. 3316–3319) theory. The approach of Creyssels (*J. Fluid Mech.*, vol. 900, 2020, p. A39) is generalized for a fluid heated internally and uniformly, confined between top and bottom plates of equal temperature. For each plate, a Nusselt number is defined and an analytical formula is given to predict its variations with the Rayleigh and Prandtl numbers. The turbulent flow produced in the upper half of the IH convection cell is very similar to that observed in standard Rayleigh–Bénard convection. On the contrary, the lower plate is swept by the large scale flow that circulates through the entire cell. The corresponding boundary layer is therefore modelled by a laminar boundary layer of the Blasius type. These predictions are consistent with the independent theoretical scalings proposed by Wang *et al.* (*Geophys. Res. Lett.*, vol. 48, issue 4, 2021, p. e2020GL091198) and they are confirmed by the numerical results obtained by Goluskin & van der Poel (*J. Fluid Mech.*, vol. 791, 2016, p. R6) and Wang *et al.* (*Geophys. Res. Lett.*, vol. 48, issue 4, 2021, p. e2020GL091198).

Key words: Bénard convection

1. Introduction

Natural convection induced by internal heat production is a phenomenon that occurs very often in geophysical, astrophysical and engineering systems (Goluskin 2015). By following the analysis recently given by Creyssels (2020), the purpose of this paper is to extend the many theoretical, numerical and experimental results of Rayleigh–Bénard (RB) convection to two internally heated (IH) convection systems shown in figures 1(a) and 1(b). In RB convection, convective flow is produced by thermal boundary conditions that cause heat to enter through the lower hot plate and exit through the upper cold plate. In this case,

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Figure 1. Two convective systems are theoretically studied here. Configuration (*a*) is closer to standard RB experiments since the two hot and cold plates play the same role and the same heat flux (Φ_0) crosses them. But the heat flux at the centre of the cell is lower or greater than Φ_0 , depending on the sign of q: $\bar{\Phi}_a(z=h/2) = \Phi_0 - qh/2$. Configuration (*b*) is more complex because the upper half-cell is very similar to the upper half-cell of RB experiments whereas the flow is stratified in the lower part of the bottom half-cell. The heat flux at the bottom plate is negative $(-q_{bot}h)$ and is not equal in absolute value to the heat flux at the top plate $(q_{top}h)$. Besides, $q_{top} > q_{bot}$. For configuration (*b*), the mean temperature profile $\bar{T}(z)$ was measured by Goluskin & van der Poel (2016). The upper and lower thermal boundary layers of thickness δ_{top} and δ_{bot} are also displayed for each configuration.

convection is controlled by the temperature difference between the plates (ΔT) and the height of the cell (*h*), or by the Rayleigh number defined as

$$Ra = \frac{g\beta\Delta Th^3}{\nu\kappa},\tag{1.1}$$

where g is the uniform gravitational acceleration, β , ν and κ are, respectively, the coefficient of thermal expansion, the kinematic viscosity and the thermal diffusivity of the fluid. On the contrary, for IH convection, the flow is produced by a volumetric source of internal heating (q in W m⁻³), itself produced by chemical or nuclear reactions, or by radiation. Instead of using *Ra*, the Rayleigh–Roberts number (Roberts 1967) is adopted as follows

$$Rr = \frac{g\beta qh^{5}}{\lambda\nu\kappa},\tag{1.2}$$

where λ is the thermal conductivity of the fluid. A first approach of IH convection is to take q constant and uniform throughout the volume. Using h^2/κ as the unit of time, h as the unit of length and qh^2/λ as the unit of temperature, the dimensionless Boussinesq equations governing the velocity (\tilde{u}) , pressure (\tilde{p}) and temperature (\tilde{T}) are the incompressibility condition $(\nabla \cdot u)$ and

$$\widetilde{\partial}_{t}\widetilde{\boldsymbol{u}} + \widetilde{\boldsymbol{u}} \cdot \widetilde{\boldsymbol{\nabla}}\widetilde{\boldsymbol{u}} = -\widetilde{\boldsymbol{\nabla}}\widetilde{p} + Pr\widetilde{\boldsymbol{\nabla}}^{2}\widetilde{\boldsymbol{u}} + Pr\,Rr\,\widetilde{T}\,\hat{\boldsymbol{e}}_{z},\tag{1.3}$$

$$\widetilde{\partial}_t \widetilde{T} + \widetilde{\boldsymbol{u}} \cdot \widetilde{\nabla} \widetilde{T} = \widetilde{\nabla}^2 \widetilde{T} + 1, \tag{1.4}$$

where dimensionless variables and operators are designated by tildes, \hat{e}_z is the vertical unit vector and *Pr* is the Prandtl number. With regard to thermal boundary conditions, we consider hereafter those used by most of previous studies i.e. fixed and equal temperature conditions (*T*₀) at the top and bottom plates. By adopting this condition, modelling an

IH convection experiment becomes a great challenge because in a single cell, there are both positive and negative vertical mean temperature gradients (see figure 1b). This leads to both turbulent convection and a stably stratified lower boundary layer. Therefore, the experiments on RB and IH convection have some similarities but also some differences. For both cases, the mean temperature is almost constant in the middle of the cell and we call it the temperature of the 'bulk flow' (\overline{T}_b) . At high Ra numbers and assuming negligible non-Boussinesq effects, \overline{T}_b is equal to the average of the top and bottom plate temperatures for a RB experiment, while \bar{T}_{h} is the maximum of the mean temperature in an IH convection cell (see figure 1b). Besides, the two thermal boundary layers located near the top and bottom plates have the same thickness for a RB experiment. On the contrary, heating in volume leads to an asymmetry between the two thermal boundary layers. The mean temperature profile was measured by Goluskin & van der Poel (2016) and is shown in figure 1(b). Therefore, the upper boundary layer is similar to that observed in a RB experiment, whereas in the lower boundary layer, the mean temperature gradient tends to stop convective flows produced in the upper region of the cell. Consequently, in an IH convection cell, the thickness of the lower boundary layer is greater than that of the upper boundary layer ($\delta_{bot} > \delta_{top}$). In addition, the difference in thickness must increase as the Rayleigh number increases. Likewise, at a fixed Ra number, the mean vertical heat flux is constant for a RB experiment and is given by the Nusselt number as $\Phi_{RB} = Nu_{RB}(\lambda \Delta T/h)$. In contrast, in an IH convection cell, the heat produced inside the fluid is evacuated through both lower and upper boundaries, leading to a mean vertical heat flux that changes sign from the bottom plate to the top plate (see figure 1b). As the mechanisms that drive the two thermal boundary layers are different, the heat fluxes through the upper and lower plates are not equal in absolute value. A coefficient α can be defined to quantify this down-up asymmetry. Indeed, the fraction of heat produced inside the fluid flowing outwards from the bottom plate can be written as

$$q_{bot} = \left(\frac{1}{2} - \alpha\right)q,\tag{1.5}$$

leading to a heat flux at the bottom plate equal to $-q_{bot}he_z$. In steady state, energy conservation yields to a heat flux that leaves through the top plate as $q_{top}he_z$, with

$$q_{top} = \left(\frac{1}{2} + \alpha\right)q. \tag{1.6}$$

Then, for each plate, we can define one Nusselt number as

$$Nu_{bot} = \frac{q_{bot}h^2}{\lambda\Delta T},\tag{1.7}$$

$$Nu_{top} = \frac{q_{top}h^2}{\lambda\Delta T}.$$
(1.8)

To be consistent with the definition of the Nusselt number adopted for RB convection experiments, the characteristic temperature difference ΔT used in (1.7) and (1.8) is $\Delta T = 2(\bar{T}_b - T_0)$. There is therefore a factor 2 by comparing the definitions of Nu_{bot} and Nu_{top} given in previous studies on IH convection (Goluskin 2015). Note also that other definitions of Nusselt numbers are given in the literature. Instead of using \bar{T}_b , the mean fluid temperature ($\langle \bar{T} \rangle$, where angle brackets denote an average over the entire volume) can be chosen as temperature reference in (1.7) and (1.8). Figure 6(*a*) shows that the difference of temperature $\bar{T}_b - \langle \bar{T} \rangle$ becomes negligible only for very high Rayleigh–Roberts numbers.

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Reference	Pr	Rr range	Nutop fit	Nubot fit
Laboratory experiments				
Kulacki & Goldstein (1972)	≈ 6	$5 \times 10^4 - 2 \times 10^7$	$0.18 Rr^{0.23}$	$0.66 Rr^{0.10}$
Jahn & Reineke (1974)	≈ 7	$10^{5} - 10^{9}$	$0.39 Rr^{0.20}$	$1.07 Rr^{0.10}$
Mayinger et al. (1975)	≈ 7	$10^{5} - 10^{11}$	$0.22Rr^{0.23}$	$0.68 Rr^{0.10}$
Ralph, McGreevy & Peckover (1977)	\approx 7	$4 \times 10^8 - 10^{12}$	$0.10 Rr^{0.24}$	$0.14 Rr^{0.17}$
Lee, Lee & Suh (2007)	≈ 0.7	$10^{10}-4 \times 10^{11}$	$0.09 Rr^{0.24}$	$0.71 Rr^{0.11}$
Simulations (3D DNS)				
Wörner, Schmidt & Grötzbach (1997)	7	$10^{5} - 10^{8}$	$0.19 Rr^{0.23}$	$0.65 Rr^{0.10}$
Goluskin & van der Poel (2016)	1	$5 \times 10^{7} - 2 \times 10^{10}$	$0.18 Rr^{0.23}$	$0.38 Rr^{0.14}$

Table 1. Previous results giving Nusselt numbers as a function of *Rr* for experiments and numericalsimulations of three-dimensional IH convection (Goluskin 2015).

Kulacki & Goldstein (1972), Jahn & Reineke (1974), Mayinger et al. (1975), Ralph et al. (1977), Lee et al. (2007), Wörner et al. (1997) and Goluskin & van der Poel (2016) have measured experimentally and numerically the Nusselt numbers Nu_{top} and Nu_{bot} as a function of the control parameter Rr. As is usual for RB experiments, they presented their results as power-law fits. The value of Nu_{top} has been found to increase with Rr at rates between $Rr^{0.20}$ and $Rr^{0.24}$ (see table 1). The value of Nu_{bot} increases more slowly with Rr, at rates between $Rr^{0.10}$ and $Rr^{0.17}$. Up to now, no theory has been able to predict these exponents whereas, for RB convection, the theory developed by Grossmann & Lohse (2000, 2001) (henceforth the GL theory, (2.18) and (2.19)) describes well the behaviour of $Nu_{RB}(Ra, Pr)$ and $Re_{RB}(Ra, Pr)$. Creyssels (2020) has recently developed a simple theoretical model to predict the Ra and Pr-dependent Nusselt number for a modified RB experiment in which heat is injected by volume but only in the lower thermal boundary layer. At the same time, the upper boundary layer is cooled with the same rate in order to have a constant energy in the convection cell, as in a standard RB cell. Note that Lepot, Aumaître & Gallet (2018) and Bouillaut et al. (2019) presented an experimental method to bypass the cooling boundary layer in order to perform this modified RB experiment. For configuration shown in figure 1(b), the problem is much more complex since the symmetry between the two plates is broken. Because of this, GL theory is first extended to the modified RB configuration (a) (see figure 1a) for which the fluid is confined between two horizontal plates of different temperature ($\Delta T > 0$ is fixed). To preserve the symmetry between the two plates, the lower part of the cell for configuration (a) is cooled with the same volumetric power as the upper part is heated

$$q_{v}(z) = \begin{cases} -q, & \text{for } 0 \le z \le h/2, \\ +q, & \text{for } h/2 \le z \le h. \end{cases}$$
(1.9)

In steady state, energy conservation yields the following variation of the heat flux averaged over a horizontal cross-section:

$$\bar{\Phi}_a(z) = \begin{cases} \Phi_0 - qz, & \text{for } 0 \le z \le h/2, \\ \Phi_0 - q(h-z), & \text{for } h/2 \le z \le h, \end{cases},$$
(1.10)

where Φ_0 is the heat flux that crosses each horizontal plate. For configuration (a), the Nusselt number can then be defined as for standard RB experiment using

$$Nu_a = \frac{\Phi_0 h}{\lambda \Delta T}.$$
(1.11)

The Nusselt (Nu_a) and Reynolds (Re_a) numbers depend on three non-dimensional parameters: Ra, Pr and Q, where Q represents the non-dimensional form of the volumetric heating or cooling sources

$$Q = \frac{qh^2}{\lambda\Delta T}.$$
(1.12)

Note that: (i) $Nu_a = Nu_{RB}(Ra, Pr)$ when Q = 0; (ii) Q can be either positive or negative; (iii) a priori, $Nu_a < Nu_{RB}$ and $Re_a < Re_{RB}$ when Q > 0 whereas $Nu_a > Nu_{RB}$ and $Re_a > Nu_{RB}$ and $Nu_{RB} > Nu_{RB}$ and $Nu_{RB} > Nu_{RB}$ and $Nu_{RB} > Nu_{RB}$ and $Nu_{RB} > Nu_{RB} > Nu_{RB} > Nu_{RB}$ and $Nu_{RB} > Nu_{RB} >$ Re_{RB} if Q < 0; (iv) $\bar{\Phi}_a(z)$ must be positive to avoid the weakening of the turbulent flow in the centre of the cell and the appearance of a stratified flow. The condition $\Phi_a(h/2) =$ $\Phi_0 - q(h/2) > 0$ implies that $Q < 2Nu_a$. Finally, for Q > 0, the upper boundary layer and the turbulent flow observed in the upper half of the cell of configuration (a) are similar to those observed in the upper half of a cell of configuration (b). However, Q is a control parameter in configuration (a) while Q must be measured or predicted in configuration (b) because $\Delta T = 2(\bar{T}_b - T_0)$ is a function of *Rr* and *Pr*. For configuration (b), using the Grossmann & Lohse (2000) ansatz based on the decomposition of the kinetic and thermal dissipation rates, Wang, Shishkina & Lohse (2021) have theoretically modelled the inverse of Q ($\tilde{\Delta} \approx 1/Q$) by power laws of Rr and Pr (see table 1 in Wang *et al.* (2021). In §2, the GL theory is adapted to configuration (a) and the two (2.20) and (2.21) predict the evolution of Nu_a and Re_a as a function of Ra, Pr and Q. Under the same conditions (same *Ra*, *Pr* and *Q*), we assume in § 3.1 that both Nu_{top} and Re_{h} measured in configuration (b) are given by $Nu_a(Ra, Pr, Q)$ and $Re_a(Ra, Pr, Q)$ (3.1) and (3.2). In § 3.2, a theoretical model based on Prandtl–Blasius–Pohlhausen theory is derived to predict Nu_{bot} as a function of the Reynolds number (3.10). Finally, these predictions for the Ra and Pr dependence of Nutop and Nubot are compared with three-dimensional experimental and numerical results in §4. A detailed comparison with the two-dimensional results of Goluskin & Spiegel (2012) and Wang *et al.* (2021) is given in Appendix A.

2. Grossmann & Lohse (2001) theory for RB experiment with volumetric energy sources (configuration a)

To predict the variations of Nusselt numbers with Ra and Pr for internal heating and cooling convection experiments, Creyssels (2020) proposed the assumption that, for high Ra numbers, the dynamical structure of the convective flow be considered to be the same as that in standard RB experiments. Notably, he assumed that the 5 dimensionless parameters (a, c_1-c_4) defined within the framework of the GL theory do not depend on the way the heat is injected and extracted in the experiment. Besides, the thickness of each thermal boundary layer is assumed to be only controlled by the difference of temperature between the bulk flow and the corresponding plate using the following equation ((2.1) in Creyssels 2020):

$$\frac{\delta_T}{h} = \frac{1}{2Nu_{RB}(Ra, Pr)}.$$
(2.1)

Then, the central idea of the GL theory is to split both mean kinetic energy and thermal dissipation rates into two contributions each, one from the bulk (Bu) and one from the boundary layers (BLs) as

$$\langle \epsilon_u \rangle = \langle \epsilon_u \rangle_{Bu} + \langle \epsilon_u \rangle_{BL}, \qquad (2.2)$$

$$\langle \epsilon_T \rangle = \langle \epsilon_T \rangle_{Bu} + \langle \epsilon_T \rangle_{BL},$$
 (2.3)

and each contribution is modelled as follows:

$$\langle \epsilon_u \rangle_{Bu} \sim \frac{U_a^2}{h/U_0} \left(1 - \frac{\delta_u}{h} \right) \approx \frac{\nu^3}{h^4} R e_0^3,$$
 (2.4)

$$\langle \epsilon_T \rangle_{Bu} \sim \frac{(\Delta T)^2}{h/U_a^{edge}} \left(1 - \frac{\delta_T}{h}\right) \approx \kappa \left(\frac{\Delta T}{h}\right)^2 Re_a Prf\left(\frac{\delta_u}{\delta_T}\right),$$
 (2.5)

$$\langle \epsilon_u \rangle_{BL} \sim \nu \left(\frac{U_a}{\delta_u}\right)^2 \frac{\delta_u}{h} = \frac{\nu^3}{h^4} R e_a^2 \frac{h}{\delta_u},$$
 (2.6)

$$\langle \epsilon_T \rangle_{BL} \sim \kappa \left(\frac{\Delta T}{\delta_T}\right)^2 \frac{\delta_T}{h} = \kappa \left(\frac{\Delta T}{h}\right)^2 \frac{h}{\delta_T}.$$
 (2.7)

In (2.4)–(2.6), U_a represents the characteristic velocity of the bulk flow. Grossmann & Lohse (2001) introduced the function $0 \le f \le 1$ because the relevant velocity at the edge between the thermal BL and the thermal bulk can be less than U_a , depending on the ratio δ_u/δ_T . They gave $f(x) = (1 + x^n)^{-1/n}$, with n = 4 as an example of function f. Besides, they assumed that the velocity BLs are Blasius like, with a thickness of

$$\frac{\delta_u}{h} = \frac{a}{\sqrt{Re_a}}.$$
(2.8)

For the mean kinetic energy and thermal dissipation rates, the balances of the turbulent kinetic energy and of the thermal variance give the following two exact relations (Creyssels 2020):

$$\langle \epsilon_u \rangle = \frac{g\beta}{h} \left[\int_0^h \frac{\bar{\Phi}_a(z)}{\rho c_p} dz - \frac{\lambda \Delta T}{\rho c_p} \right], \qquad (2.9)$$

$$\langle \epsilon_T \rangle = \frac{1}{h} \int_0^h \bar{T}(z) \frac{q_v(z)}{\rho c_p} dz + \frac{\Delta T \Phi_0}{\rho c_p h}.$$
(2.10)

Using (1.10), (1.1), (1.12) and (1.11), (2.9) becomes

$$\langle \epsilon_u \rangle = \frac{\nu^3}{h^4} \frac{\left(Nu_a - 1 - \frac{Q}{4}\right)Ra}{Pr^2}.$$
(2.11)

As the GL theory is based on Prandtl–Blasius–Pohlhausen laminar boundary layers (Grossmann & Lohse 2000), the mean temperature can be written as (Creyssels 2020)

$$2\frac{\bar{T}(z) - T_b}{2} = \begin{cases} 1 - \Theta_P\left(\frac{z}{\delta_T}\right), & \text{for } z \le h/2, \end{cases}$$
(2.12*a*)

$$\Delta T \qquad \left(\begin{array}{c} \Theta_P\left(\frac{h-z}{\delta_T}\right) - 1, \quad \text{for } h/2 \le z \le h, \end{array} \right.$$
(2.12b)

with Θ_P the Pohlhausen temperature profile, which is assumed to be independent of the Prandtl number. In particular, $\Theta_P(0) = 0$ and $\Theta_P(\eta) \to 1$ when $\eta \gg 1$. Using (1.9), (2.12),

(1.12) and (1.11), (2.10) then becomes

$$\langle \epsilon_T \rangle = \kappa \left(\frac{\Delta T}{h}\right)^2 \left[Nu_a - \lambda_d \frac{\delta_T}{h}Q\right],$$
(2.13)

where $\lambda_d = \int_0^\infty [1 - \Theta_P(\eta)] d\eta$ denotes the displacement thickness of the mean temperature profile ($\lambda_d \approx 0.57$ for a Blasius profile).

From decomposition of the two global dissipation rates (2.2) and (2.3), four regimes of convection (I, II, III and IV) can be defined depending on whether the bulk or the BL contributions dominate the global dissipations. For regimes *II* and *IV*, the kinetic energy dissipation rate is dominated by its bulk contribution whereas, for regimes *I* and *III*, $\langle \epsilon_u \rangle \sim \langle \epsilon_u \rangle_{BL}$. Combining (2.4) and (2.11), or (2.6) and (2.11), and using (2.8), we obtain

$$\frac{\left(Nu_a - 1 - \frac{Q}{4}\right)Ra}{Pr^2} \sim (Re_a)^{\theta_i},\tag{2.14}$$

whereas, for standard RB convection, we have

$$\frac{(Nu_{RB}-1)Ra}{Pr^2} \sim (Re_{RB})^{\theta_i},\tag{2.15}$$

with $\theta_{II} = \theta_{IV} = 3$ and $\theta_I = \theta_{III} = 5/2$.

Regimes *III* and *IV* are obtained for high *Ra* numbers for which thermal dissipation rate is dominated by its bulk contribution whereas, for regimes *I* and *II*, $\langle \epsilon_T \rangle \sim \langle \epsilon_T \rangle_{BL}$. For standard RB convection, Grossmann & Lohse (2000) predicted

$$(Nu_{RB})^{\phi_i} \sim Re_{RB} Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{RB}}}\right),$$
 (2.16)

with $\phi_{III} = \phi_{IV} = 1$ and $\phi_I = \phi_{II} = 2$. For internal heating and cooling convection, combining (2.5) and (2.13), or (2.7) and (2.13), and using (2.1), yields

$$Nu_{a} - \frac{\lambda_{d}}{2} \frac{Q}{Nu_{RB}} \sim \begin{cases} Re_{a} Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{a}}}\right) & \text{(regimes III and IV),} \\ Nu_{RB} \sim \sqrt{Re_{RB} Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{RB}}}\right)} & \text{(regimes I and II).} \end{cases}$$
(2.17)

As the 4 previous regimes can only be observed experimentally and numerically for extreme values of the Ra and Pr numbers, Grossmann & Lohse (2001) proposed describing RB convection at any Ra and Pr numbers as a mixture of these 4 regimes. Equations (2.15) and (2.16) are then generalized as

$$(Nu_{RB} - 1)RaPr^{-2} = \frac{c_1}{2a}Re_{RB}^{5/2} + c_2Re_{RB}^3,$$
(2.18)

$$Nu_{RB} = c_3 \sqrt{Re_{RB}Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{RB}}}\right) + c_4Re_{RB}Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{RB}}}\right)}.$$
(2.19)



Figure 2. Results of the extension of GL theory for turbulent convection with uniform volumetric sources. Values of Nu_a/Nu_{RB} (*a*) and Re_a/Re_{RB} (*b*) as a function of *Q* normalized by $Nu_{RB}(Ra, Pr)$ for Pr = 1 and for 3 different Ra numbers: $Ra = 10^6$ (solid line), $Ra = 10^{10}$ (dashed line) and $Ra = 10^{14}$ (dash-dotted line). The black dotted line represents regime IV_u (2.22) i.e. the limit of system of (2.20) and (2.21) when $Ra \to \infty$ and $Pr \ge 1$. Values of Nu_a/Nu_{RB} (*c*) and Re_a/Re_{RB} (*d*) as a function of Pr for $Ra = 10^9$ and from top to bottom: $Q/Nu_{RB} = -1, -0.5, 0.5$ and 1.

By applying this idea to internal heating convection, (2.14) and (2.17) are generalized using the two following equations:

$$\left(Nu_a - 1 - \frac{Q}{4}\right)\frac{Ra}{Pr^2} = \frac{c_1}{2a}Re_a^{5/2} + c_2Re_a^3,$$
(2.20)

$$Nu_{a} - \frac{\lambda_{d}}{2} \frac{Q}{Nu_{RB}} = c_{3} \sqrt{Re_{RB} Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{RB}}}\right) + c_{4}Re_{a}Prf\left(\frac{2aNu_{RB}}{\sqrt{Re_{a}}}\right)}.$$
 (2.21)

Equations (2.20) and (2.21) give the Nu_a and Re_a numbers as functions of the 3 parameters Ra, Pr and Q. Figures 2(*a*) and 2(*b*) show the ratios Nu_a/Nu_{RB} and Re_a/Re_{RB} as a function of Q compensated by $Nu_{RB}(Ra, Pr)$ for Pr = 1 and for 3 different Ra numbers: $Ra = 10^6$ (solid line), $Ra = 10^{10}$ (dashed line) and $Ra = 10^{14}$ (dash-dotted line). As expected, heating the lower part of the cell (Q < 0) increases both Nu_a and Re_a while heating the upper part of the cell (Q > 0) decreases both Nu_a and Re_a . At fixed ratio Q/Nu_{RB} and when Ra increases, both ratios Nu_a/Nu_{RB} and Re_a/Re_{RB} increase for Q < 0 and they decrease for Q > 0. This can easily be explained by the fact that convection is described by a mixture of regimes II and IV for Pr = 1 and the higher Ra, the lower the

portion corresponding to regime *II*. Besides, $Nu_a \approx Nu_{RB}$ for pure regime *II* whereas we have

$$\frac{Nu_a}{Nu_{RB}} = \left(\frac{Re_a}{Re_{RB}}\right)^{3/2} = \frac{1}{2} + \frac{1}{2}\sqrt{1 - \frac{Q}{Nu_{RB}}}$$
(2.22)

for pure regime *IV* and $\delta_u > \delta_T$ (or $Pr \ge 1$). To obtain (2.22), the system of (2.20) and (2.21) is solved with $c_1 = c_3 = 0$ (regime *IV*), assuming that $Nu_a \gg 1 > (\lambda_d/2)(Q/Nu_{RB})$ and for $\delta_u > \delta_T [f(\delta_u/\delta_T) \approx \delta_T/\delta_u]$. Figures 2(*a*) and 2(*b*) show that ratios Nu_a/Nu_{RB} and Re_a/Re_{RB} tend to the solution given by (2.22) when $Ra \to \infty$ and Pr = 1.

Finally, when fixing Ra and Q, and for $Pr \ge 1$, ratios Nu_a/Nu_{RB} and Re_a/Re_{RB} depend very little on Pr (see figure 2c,d), in agreement with the solution of pure regime IV_u (2.22). On the contrary, when Pr decreases from 1 to 0 and even if Ra is held constant, both $Re_{RB}Pr$ and Re_aPr decrease. As a result, the importance of regime II in the mixture of regimes represented in (2.21) increases and thus ratio Nu_a/Nu_{RB} approaches 1 (figure 2c,d).

These results are used in the following section to study the configuration shown in figure 1(*b*) for which both non-dimensional numbers Ra and Q are not control parameters. Indeed, they both depend on the temperature of the bulk flow (\bar{T}_b) and thus \bar{T}_b need to be either measured or predicted. However, for configuration (b), using definitions (1.1) and (1.12), we have: Rr = Ra Q, with Rr the Rayleigh number adopted for IH convection experiment (1.2).

3. Theory for IH convection experiment (configuration b)

3.1. Model for the top Nusselt number

The main assumption of this model is to assume that the flow in the upper half of the configuration shown in figure 1(*b*) is similar to the flow seen in the upper half of the cell of configuration (a). By analogy with RB convection for which $Nu_{RB} \sim Ra^{1/3}$ means that the two boundary layers behave independently of each other, we assume here that the upper boundary layer is not sensitive to the flow structure that controls the lower boundary layer. Therefore, the thickness of the top BL (δ_{top}), the heat flux at the top plate ($q_{top}h$) and the characteristic velocity (U_b) are only controlled by the following parameters: $\Delta T = 2(\bar{T}_b - T_0)$, q, h and Pr. In non-dimensional form, it is assumed that Nu_{top} and Re_b are given by the Nusselt and Reynolds numbers calculated for configuration (a) under the same conditions

$$Nu_{top}(Rr, Pr) = Nu_a(Ra, Pr, Q),$$
(3.1)

$$Re_b(Rr, Pr) = Re_a(Ra, Pr, Q).$$
(3.2)

Besides, modelling both Nu_{bot} and Nu_{top} is sufficient to give Ra and Q as a function of the two control parameters Rr and Pr. Indeed, using (1.2)–(1.8), (1.1) and (1.12) become

$$Q(Rr, Pr) = Nu_{bot} + Nu_{top}, \tag{3.3}$$

$$Ra(Rr, Pr) = \frac{Rr}{Nu_{bot} + Nu_{top}}.$$
(3.4)

3.2. Model for the bottom Nusselt number

Since the mean temperature gradient is positive throughout the lower boundary layer (see figure 1b), forced convection is the main mechanism that drives the heat flux at the

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lower plate. Assuming static large scale flow, also called 'wind flow', a laminar velocity boundary layer develops against the bottom plate in the same way as a Blasius boundary layer. Using the theory of Prandtl–Blasius–Pohlhausen (Schlichting 1979), the thickness of the thermal boundary layer is given by

$$\frac{\delta_{T,x}}{x} = \frac{1}{A \,\mathcal{F}(Pr) \,Re_x^{1/2}},\tag{3.5}$$

with x the horizontal coordinate, $A \approx 0.33$ and $\mathcal{F}(1) = 1$. Here, δ_T is defined as $(\delta_T)^{-1} = (d\Theta_P/dz)_{z=0}$, with $\Theta_P = (T - T_0)/(\bar{T}_b - T_0)$ the non-dimensional difference of temperature. The Prandtl-dependent function \mathcal{F} varies as $Pr^{1/3}$ when $Pr \ge 0.5$, whereas, when Pr decreases, the exponent of the power-law behaviour of \mathcal{F} increases up to 1/2. In the range $10^{-3} \le Pr \le 10^3$, \mathcal{F} can be approximated by $\mathcal{F}(Pr) = d_1Pr^{1/2}f(d_2Pr^{1/6})$ with a precision of ± 1 %, using $f(x) = (1 + x^n)^{-1/n}$, n = 4, $d_1 = 1.68$ and $d_2 = 1.63$.

At the bottom plate (z = 0), the local heat flux is the sum of two contributions

$$\Phi_x = \Phi_{x,P} + \Phi_{x,q}. \tag{3.6}$$

Here, $\Phi_{x,P}$ and $\Phi_{x,q}$ represent, respectively, the heat flux given by the theory of Prandtl–Blasius–Pohlhausen and the heat flux due to the presence of volumetric heat sources inside the thermal boundary layer

$$\Phi_{x,P} = A \mathcal{F}(Pr) R e_x^{1/2} \frac{\lambda(\bar{T}_b - T_0)}{x}, \qquad (3.7)$$

$$\Phi_{x,q} = B \, q \delta_{T,x}. \tag{3.8}$$

In (3.8), *B* is a numerical constant to be determined using experiments or numerical simulations. In the case of $\Phi_{x,P} = 0$ (pure conductive state), B = 0.5.

Assuming an aspect ratio of 1 for the cell and using (3.7), (3.8) and (3.5), the integration of the local heat flux Φ_x (3.6) between x = 0 and x = h gives

$$\bar{\Phi}_{bot} = \frac{1}{h} \int_0^h \Phi_x = \frac{2A \,\mathcal{F}(Pr)}{h} \lambda (\bar{T}_b - T_0) R e_b^{1/2} + \frac{2Bqh}{3A \,\mathcal{F}(Pr) R e_b^{1/2}}.$$
(3.9)

Using (3.9) and (1.12), the bottom Nusselt number can be calculated as

$$Nu_{bot} = \frac{\bar{\Phi}_{bot}h}{2\lambda(\bar{T}_b - T_0)} = A \,\mathcal{F}(Pr) \,Re_b^{1/2} + \frac{2B}{3A \,\mathcal{F}(Pr)} \frac{Q}{Re_b^{1/2}}.$$
(3.10)

Likewise, from (3.5), the mean bottom boundary layer thickness can be written as

$$\frac{\bar{\delta}_{bot}}{h} = \frac{1}{h^2} \int_0^h \delta_{T,x} \, \mathrm{d}x = \frac{2}{3A \,\mathcal{F}(Pr) \, Re_h^{1/2}}.$$
(3.11)

3.3. *Mean fluid temperature* $\langle \bar{T} \rangle$

Both models presented in §§ 3.1 and 3.2 are based on Blasius profiles for the z-evolution of $\overline{T}(z)$ in the two boundary layers. In addition, $\overline{T}(z) \approx \overline{T}_b$ in the bulk flow i.e. for

 $\bar{\delta}_{bot} \ll z \ll h - \delta_{top}$. Thus, for configuration (b), using (2.12b), $\bar{T}(z)$ can be expressed as

$$\frac{\bar{T}_b - \bar{T}(z)}{\bar{z}} = \begin{cases} 1 - \Theta_P\left(\frac{z}{\bar{\delta}_{bot}}\right), & \text{for } z \le h/2, \end{cases}$$
(3.12*a*)

$$\bar{T}_b - T_0 = \left[1 - \Theta_P \left(\frac{h - z}{\bar{\delta}_{top}} \right), \quad \text{for } h/2 \le z \le h.$$
 (3.12b)

The thicknesses of the bottom and top boundary layers ($\bar{\delta}_{bot}$ and $\bar{\delta}_{top}$) are given by (2.1) and (3.11), respectively. Using (3.12), the mean fluid temperature can be calculated yielding

$$\frac{\bar{T}_b - \langle \bar{T} \rangle}{\bar{T}_b - T_0} = \lambda_d \frac{\bar{\delta}_{bot} + \bar{\delta}_{top}}{h},\tag{3.13}$$

$$=\lambda_d \left(\frac{2}{3A \mathcal{F}(Pr) Re_b^{1/2}} + \frac{1}{2Nu_{RB}}\right),\tag{3.14}$$

with $Re_b(Rr, Pr)$ (3.2) and $Nu_{RB}(Ra, Pr)$ with Ra(Rr, Pr) (3.4).

4. Comparison between theories and numerical results

Theoretical predictions for Nu_{top} , Nu_{bot} and mean fluid temperature are tested below using experimental and numerical results on three-dimensional IH convection (see table 1 and Goluskin (2015) for a literature survey). A comparison between the theories and two-dimensional simulations of IH convection (Goluskin & Spiegel 2012; Goluskin & van der Poel 2016; Wang *et al.* 2021) is given in Appendix A.

4.1. Top Nusselt number

Figures 3(*a*) and 3(*b*) show the variations of Nu_{top} (filled squares) defined by (1.8) and given by Goluskin & van der Poel (2016) for IH convection simulations. The variations of the Nusselt number obtained for standard RB simulations are also shown (data from Pandey & Verma (2016) and Shishkina *et al.* (2017), open symbols). Clearly, at constant *Ra* and *Pr*, Nu_{top} is slightly lower than Nu_{RB} . Dashed and solid lines represent, respectively, GL theory for RB convection (2.18) and (2.19) and for RB convection with volumetric energy sources (2.20) and (2.21). For the latter case, *Q* and *Ra* are calculated using (3.3), (3.4) and the data of Goluskin & van der Poel (2016). A good agreement between the numerical simulations and the extension of GL theory is observed both by plotting Nu_{top} as a function of Ra for Pr = 1 (figure 3*a*) and by considering the variations of Nu_{top} with Pr for $Ra \approx 1.7 \times 10^6$ (figure 3*b*).

4.2. Bottom Nusselt number

The theory developed in § 3.2 to predict the variations of Nu_{bot} is based on forced convection in the lower half of the cell and the theory of Blasius. Therefore, to get Nu_{bot} with good precision, a very good estimate of the Reynolds number is necessary (see (3.10)). However, to date there are no experimental or numerical measurements for either $Re_b(Rr, Pr)$ or $Re_a(Ra, Pr, Q)$ in three dimensions. Even for RB convection, experimental and numerical measurements of the Reynolds number have larger uncertainties than those of the Nusselt number because they require measurement of velocity fluctuations



Figure 3. The compensated Nusselt number for standard RB convection (open symbols) and for IH convection (filled squares) as a function of Ra for Pr = 1 (a) and against Pr for $Ra \approx 1.7 \times 10^6$ (b). For IH convection, Ra values are calculated using (3.4). Dashed lines: Grossmann & Lohse (2001) theory with prefactors given by Stevens *et al.* (2013) (2.18) and (2.19). Solid blue lines: extension of GL theory for convection with volumetric energy sources (2.20) and (2.21).



Figure 4. The compensated Reynolds number for standard RB convection as a function of Ra for Pr = 1 (a) and against Pr for $Ra \approx 1.7 \times 10^6$ (b) (symbols as figure 3). Also shown: the predictions of the GL theory (dashed lines, (2.18) and (2.19)) and those of its extension to convection with volumetric heat source (solid lines, (2.20) and (2.21)). For the second case, Ra and Q are calculated using (3.4) and (3.3) and the data of Goluskin & van der Poel (2016). The horizontal lines represent power-law behaviours (upper line: (4.1), lower line: (4.2)).

throughout the cell. In addition, several definitions of *Re* are given in the literature (root mean square of the velocity, root mean square of the vertical velocity, maximum of the velocity fluctuations, maximum mean velocity along the heated plate). Likewise, the GL theory is less effective in predicting Re_{RB} than Nu_{RB} . Figures 4(*a*) and 4(*b*) show the variations of Re_{RB} as a function of Ra and Pr given by Pandey & Verma (2016) and Shishkina *et al.* (2017). For both cases i.e. Pr = 1, $10^5 \le Ra \le 5 \times 10^8$ (figure 4*a*) and $Ra = 1.7 \times 10^6$, $0.1 \le Pr \le 10$, Re_{RB} can be fitted by a power law of Ra and Pr as

$$Re_{RB} \approx 0.15 \, Ra^{1/2} Pr^{-\gamma},\tag{4.1}$$

with $\gamma \approx 0.86$. As no data on Re_a are available, the same power-law behaviour will be used for Re_a . In addition, the extension of the GL theory (see § 2) predicts that $Re_a/Re_{RB} \approx$ 0.8 in the range of Ra, Pr and Q investigated by Goluskin & van der Poel (2016)



Figure 5. Value of Nu_{bot} vs Ra(a) and Pr(b). Symbols: data from Goluskin & van der Poel (2016). Solid lines: (3.10) with A = 0.21 and B = 0.4. Dashed lines: first term of (3.10) ($\propto \sqrt{Re} \propto Ra^{1/4}$). Dotted lines: second term of (3.10) ($\propto Q/\sqrt{Re}$).



Figure 6. Difference between the temperature of the bulk flow (\bar{T}_b) and the mean fluid temperature $(\langle \bar{T} \rangle)$ as a function of Ra (a) and Pr (b). Symbols: data from Goluskin & van der Poel (2016). Solid lines: (3.14) with A = 0.21 and $\lambda_d = 0.57$. Dashed lines: first term of (3.14) i.e. $\lambda_d (\bar{\delta}_{bot}/h) \propto Re_b^{-1/2} \propto Ra^{-1/4}$. Dotted blue lines: second term of (3.14) i.e. $\lambda_d (\bar{\delta}_{top}/h) \propto 1/Nu_{RB} \propto Ra^{-\theta_{RB}}$ with $\theta_{RB} \approx 0.3$ (see § 4.4).

(see figure 4a,b). Thus, the following equation will be use to give Re_a or Re_b as a function of Ra and Pr:

$$Re_b(R, Pr) = Re_a(Ra, Pr, Q) = 0.12 Ra^{1/2} Pr^{-\gamma}.$$
(4.2)

Using (4.2), (3.10) gives Nu_{bot} as a function of Ra, Pr and Q. Note that Ra and Q are calculated from the data of Goluskin & van der Poel (2016) and using (3.3) and (3.4). Figure 5(*a*,*b*) shows that the (3.10) describes well both the evolution of Nu_{bot} with Ra and with Pr, using A = 0.21 (not far away from 0.33) and B = 0.4 (close to the value 0.5 which characterizes a purely conductive boundary layer).

4.3. Mean fluid temperature $\langle \bar{T} \rangle$

The model presented in § 3 also predicts the evolution with Ra and Pr of the difference between the mean fluid temperature $(\langle \bar{T} \rangle)$ and the temperature of the bulk flow (\bar{T}_b) . Figure 6(a,b) shows a good agreement between (3.14) and the numerical results of Goluskin (2015) using no new adjustable parameters. Indeed, as for figure 5(a,b),

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A is taken equal to 0.21 whereas the displacement thickness of the mean temperature profile is given by the Blasius theory ($\lambda_d = 0.57$). As shown by (3.13), the dimensionless temperature difference $(\bar{T}_b - \langle \bar{T} \rangle)/(\bar{T}_b - T_0)$ is equal to the sum of the thicknesses of the two boundary layers. For $Ra > 10^5$, as $\bar{\delta}_{bot} \gg \bar{\delta}_{top}$ and $Re_b \propto Ra^{1/2}$ (see (4.2)), (3.14) can be approximated by

$$\frac{\bar{T}_b - \langle \bar{T} \rangle}{\bar{T}_b - T_0} \approx \lambda_d \frac{\bar{\delta}_{bot}}{h} \propto R a^{-1/4}$$
(4.3)

where $\bar{\delta}_{bot}/h \propto Re_b^{-1/2}$ (see (3.11)) and $Re_b \propto Ra^{1/2}$ (4.2). As the Rayleigh number increases, the thicknesses of the upper and lower boundary layers decrease, which also results in a decrease of $\bar{T}_b - \langle \bar{T} \rangle$.

4.4. Approximation of Nutop and Nubot by power laws of Rr

For $10^5 \le Ra \le 10^9$ and $Pr \ge 1$, numerical simulations (Goluskin & van der Poel 2016; Pandey & Verma 2016; Shishkina *et al.* 2017) and GL theory show that both Nu_{RB} and Nu_{top} can be approximated by a power law of Ra with an exponent $\theta_{RB} \approx 0.3$ (see also figure 3*a*). Neglecting Nu_{bot} in front of Nu_{top} , (3.4) yields $Ra \sim Rr^{1/(1+\theta_{RB})}$. Consequently, we get

$$Nu_{top} \approx 0.18 \, Rr^{\theta_{RB}/(1+\theta_{RB})} \approx 0.18 \, Rr^{0.23}.$$
 (4.4)

This scaling for Nu_{top} is consistent with the independent theoretical findings of Wang *et al.* (2021). Indeed, for regimes $I_{\infty}^{<}$ and IV_u , Wang *et al.* (2021) have shown that $Q \approx \tilde{\Delta}^{-1} \sim Rr^{1/4}$ whereas for regimes I_u and I_l , $Q \approx \tilde{\Delta}^{-1} \sim Rr^{1/5}$.

Using (4.2) and assuming again $Q \approx Nu_{top}$, (3.10) becomes

$$Nu_{bot} \approx 0.11 \, Pr^{-\Gamma} \, Rr^{1/4(1+\theta_{RB})} + 0.42 \, Pr^{\Gamma} \, Rr^{(\theta_{RB}-1/4)/(1+\theta_{RB})}, \tag{4.5}$$

$$\approx 0.11 \, Pr^{-0.10} \, Rr^{0.19} + 0.42 \, Pr^{0.10} \, Rr^{0.04}. \tag{4.6}$$

Indeed, for $Pr \ge 1$, $\mathcal{F}(Pr) \approx Pr^{1/3}$ and $\Gamma = \gamma/2 - \frac{1}{3} \approx 0.10$. Equation (4.6) shows that Nu_{bot} can be expressed by the sum of two power laws of Rr with exponents 0.04 and 0.19 (see also figure 7). This explains the high variability of the measured exponents when the experimental or numerical results are fitted by a single power law (see table 1). Since the experiments of Ralph *et al.* (1977) were performed for higher Rr numbers $(4 \cdot 10^8 - 10^{12})$ than the other works, the corresponding exponent (0.17) is larger than the exponents calculated using the other results (~0.10). Of course, the observed differences are also due to experimental uncertainties, boundary conditions and the way the thermal power q_v is imposed in volume in the fluid.

5. Conclusions

The Grossmann & Lohse (2001) theory has been extended here to RB convection with volumetric energy sources in the fluid (configuration a). The two equations of the GL theory (2.18) and (2.19) have been modified to take account of the effect of the new non-dimensional parameter $Q = qh^2/(\lambda\Delta T)$. Equations (2.20) and (2.21) predict the evolution of Nu_a and Re_a as a function of Ra, Pr and Q. Besides, these two equations can also describe the top Nusselt number Nu_{top} and the Reynolds number of an IH convection experiment (configuration b). In this case, the non-dimensional parameter Rr = Ra Q is the control parameter of the IH convection system whereas Ra and Q are theoretically



Figure 7. Values of Nu_{top} (squares) and Nu_{bot} (circles) vs Rr for Pr = 1 (data from Goluskin & van der Poel 2016). Solid upper line: $Nu_{top} \propto Rr^{0.23}$ (4.4). Solid lower line: (3.10) with A = 0.21 and B = 0.4. Dashed line: $Nu_{bot} = 0.11 Rr^{0.19} + 0.42 Rr^{0.04}$ (4.6). Dotted line: $Nu_{bot} = 0.11 Rr^{0.19}$.

modelled using $Q = Nu_{top} + Nu_{bot}$, with $Nu_{top} = Nu_a(Ra, Pr, Q)$ and $Nu_{bot}(Re_a, Pr, Q)$. The bottom Nusselt number is the sum of two terms (3.10). The first term can be modelled using the Prandtl–Blasius–Pohlhausen theory and is proportional to square root of the Reynolds number. The second term comes from the presence of volumetric energy sources and is proportional to Q/\sqrt{Re} . In agreement with the independent theoretical findings of Wang *et al.* (2021), for high *Rr* numbers, Nu_{top} scales as $Rr^{\theta_{RB}/(1+\theta_{RB})}$, with θ_{RB} the exponent of the power-law behaviour of Nu_{RB} with *Ra*. These predictions are confirmed by experimental and numerical results and show that natural and forced convections are the two mechanisms that control the heat fluxes in an IH convection experiment.

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Appendix A. Comparison between theories and numerical results for twodimensional IH convection

For two-dimensional IH convection, the predictions of the theoretical work presented in § 3 can be tested thanks to the numerical investigations of Goluskin & Spiegel (2012) and Wang *et al.* (2021). Figure 8(*a*,*b*) shows the variations of Nu_{top} and Re as a function of the Rayleigh number ($Ra = Rr/(Nu_{top} + Nu_{bot})$), for Pr = 1. For comparison purposes, the measured Nusselt and Reynolds numbers for RB convection are also displayed. We observe very little difference between Nu_{top} and Nu_{RB} whereas the Reynolds number is approximately 65 % lower for IH convection than for RB convection. With good precision, Nu_{top} and Re_b can be approximated by power laws of Ra as

$$Nu_{top} \approx Nu_{RB} \approx 0.25 \, Ra^{0.25},\tag{A1}$$

$$Re_b \approx 0.65 \, Re_{RB} \approx 0.038 \, Pr^{0.86} \, Ra^{0.58}.$$
 (A2)

The results for two-dimensional IH convection are therefore slightly different from those obtained for three-dimensional IH convection. In three dimensions, numerical data of



Figure 8. Results for two-dimensional IH and RB convection. (*a*) Values of Nu_{top} and Nu_{bot} compensated by $Ra^{1/3}$ as a function of Ra. Data from Goluskin & Spiegel (2012) (diamonds) and Wang *et al.* (2021) (squares), Pr = 1. The results for RB convection ($Nu_{RB}Ra^{-1/3}$) are also shown: Johnston & Doering (2009) (asterisks), Zhu *et al.* (2018) (stars) and Zhang, Zhou & Sun (2017) (plusses: Pr = 0.7, crosses: Pr = 5.3). Blue dashed line: (A1). Red solid line: (3.10) with $A_{2D} = 0.28$ and $B_{2D} = 0.4$. (*b*) Compensated Reynolds number ($RePr^{0.86}/Ra^{1/2}$) vs Ra. Blue solid line: (A2).

Goluskin (2015) and the extension of the GL theory show that $0.83 \le Nu_{top}/Nu_{RB} \le 0.91$ and $Re_b \approx 0.8 Re_{RB}$ (using GL theory) whereas $Nu_{top} \approx Nu_{RB}$ and $Re_b \approx 0.65 Re_{RB}$ in two dimensions. However, beyond these small differences, the two- and three-dimensional numerical results validate the theoretical approach chosen here to separate the IH convection cell into two distinct parts. The upper half of the cell behaves like the upper half of a RB convection cell ($Nu_{top} \sim Nu_{RB}$) and controls the Reynolds number. The heat transfer at the bottom plate is controlled by the Reynolds number and, using the theory of Blasius, Nu_{bot} is given by (3.10). Indeed, figure 8(*a*) shows a good agreement between the results of Wang *et al.* (2021) (red squares) and (3.10) (red solid line) using $A_{2D} = 0.28$ and $B_{2D} = B_{3D} = 0.4$. In two dimensions, the parameter *A* is therefore slightly higher than in three dimensions ($A_{3D} \approx 0.22$) and closer to the value for a Blasius boundary layer ($A \approx 0.33$). A greater difference is observed between the data from Goluskin & Spiegel (2012) (red diamonds) and (3.10). This is probably due to the fact that the equation that gives the Reynolds number (A2) is less well verified when the Rayleigh number is low.

The up–down asymmetry that turbulent convection induces is quantified measuring the coefficient α . From definitions given by (1.5)–(1.7), α is directly linked to the ratio of Nu_{top} over Nu_{bot}

$$2\alpha = \frac{\frac{Nu_{top}}{Nu_{bot}} - 1}{\frac{Nu_{top}}{Nu_{bot}} + 1}.$$
(A3)

Both for two- and three-dimensional IH convection, figure 9 shows that the theory presented in § 3 describes well the evolution of the asymmetry coefficient with *Ra*. However, the evolution of α with *Ra* is different in two dimensions than in three dimensions. In particular, for two dimensions, α saturates and even decreases when *Ra* increases. This behaviour is well described by (3.10). For high *Ra* numbers, (3.10) can be approximated by $Nu_{bot} \sim Re^{1/2}$. Using (A2) (two dimensional) or (4.2) (three dimensions), we obtain $Nu_{bot} \sim Ra^{0.29}$ for two dimensions while $Nu_{bot} \sim Ra^{0.25}$



Figure 9. Asymmetry coefficient α vs Ra for two-dimensional (red open symbols) and three-dimensional IH convection (blue circles: data of Goluskin & van der Poel 2016). Red solid line: (A3) and using (A1), (A2) and (3.10). Blue dashed line: (A3) and using (4.4), (4.2) and (3.10). Lower and upper dotted lines: (A3) with $Nu_{top}/Nu_{bot} \sim Ra^{-0.04}$ and $Nu_{top}/Nu_{bot} \sim Ra^{0.05}$, respectively.

for three dimensions. As $Nu_{top} \sim Ra^{0.25}$ for two dimensions (A1) and $Nu_{top} \sim Ra^{0.3}$ for three dimensions (see § 4.4), Nu_{top}/Nu_{bot} varies as $Ra^{-0.04}$ for two dimensions and as $Ra^{0.05}$ for three dimensions IH convection. That is why, for $Ra \gtrsim 10^7$, α decreases with Ra in two dimensions while α continues to increase with Ra in three dimensions (see figure 9).

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