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Turbulent mixing and layer formation in double-diffusive convection: 3D numerical simulations and theory

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ABSTRACT

Double-diffusive convection, often referred to as semi-convection in astrophysics, occurs in thermally and compositionally stratified systems which are stable according to the Ledoux-criterion but unstable according to the Schwarzschild criterion. This process has been given relatively little attention so far, and its properties remain poorly constrained. In this paper, we present and analyze a set of three-dimensional simulations of this phenomenon in a Cartesian domain under the Boussinesq approximation. We find that in some cases the double-diffusive convection saturates into a state of homogeneous turbulence, but with turbulent fluxes several orders of magnitude smaller than those expected from direct overturning convection. In other cases the system rapidly and spontaneously develops closely-packed thermo-compositional layers, which later successively merge until a single layer is left. We compare the output of our simulations with an existing theory of layer formation in the oceanographic context, and find very good agreement between the model and our results. The thermal and compositional mixing rates increase significantly during layer formation, and increase even further with each merger. We find that the heat flux through the staircase is a simple function of the layer height. We conclude by proposing a new approach to studying transport by double-diffusive convection in astrophysics.

Subject headings: convection – hydrodynamics – planets and satellites:general – stars:interior
1. Introduction

1.1. Convection, double-diffusive convection (semi-convection) and fingering convection (thermohaline convection)

One of the longest-standing problems in stellar and planetary astrophysics is that of modeling the transport of heat and chemical species within turbulent regions. The best-studied and most ubiquitously relevant case is that of overturning convection through a chemically homogeneous gas layer. There, the well-known Schwarzschild criterion is used to determine the extent of the convective region, while the transport properties through the layer are commonly modeled using mixing-length theory (Biermann 1932). The success of well-calibrated mixing-length models in explaining many observable properties of stars is quite remarkable.

However, much less is known about convection in the presence of additional factors such as strong rotation, strong magnetic fields and strong compositional gradients (Spiegel 1972). In all cases, the linear stability of the system is well-understood (Chandrasekhar 1961), but characterizing its fully-nonlinear transport properties remains the subject of ongoing research. In this work, we focus on the case of convection in the presence of a strong stabilizing compositional gradient, but in the absence of rotation or magnetic field. This regime is often called “semi-convection” in astrophysics (Schwarzschild & Härm 1958), although we prefer to use the terminology “double-diffusive convection” commonly used in oceanography to clarify the true nature of the instability responsible for the turbulence.

It has long been known that the relevant criterion for instability to overturning convection in the presence of a compositional gradient is not the Schwarzschild criterion,

\[ \nabla - \nabla_{\text{ad}} = \left( \frac{\partial \ln T}{\partial \ln p} \right) - \left( \frac{\partial \ln T}{\partial \ln p} \right)_{\text{ad}} > 0, \]  

but the Ledoux criterion (Ledoux 1947):

\[ \nabla - \nabla_{\text{ad}} > \nabla_{\mu}, \]

\[ \Leftrightarrow \left( \frac{\partial \ln T}{\partial \ln p} \right) - \left( \frac{\partial \ln T}{\partial \ln p} \right)_{\text{ad}} > \left( \frac{\partial \ln \mu}{\partial \ln p} \right), \]

where \( T \) is the temperature, \( \mu \) the mean molecular weight, \( p \) the pressure, and where the subscript “ad” expresses a derivative at constant specific entropy. In fact, both of these criteria merely express the same property when written in terms of the density stratification:

\[ \left( \frac{\partial \rho}{\partial p} \right)_{\text{ad}} > \left( \frac{\partial \rho}{\partial p} \right). \]
A system is unstable to overturning convection if the density of a parcel of fluid, raised adiabatically and in pressure equilibrium from its original position, is lower than that of its new surroundings.

The question of what happens to regions which are stable according to the Ledoux criterion but unstable according to the Schwarzchild criterion was first raised by Schwarzchild & Härm (1958). It was later found that this regime is in fact also linearly unstable (Walin 1964; Kato 1966), but through a double-diffusive instability, i.e. an instability which cannot occur unless the thermal diffusivity of the fluid, $\kappa_T$, is larger than its compositional diffusivity $\kappa_\mu$. This condition is however automatically satisfied in stellar and planetary interiors where the diffusivity ratio $\tau = \kappa_\mu / \kappa_T$ (often called the inverse Lewis number), can be as low as $10^{-7}$.

As a result, a wide range of situations arise in which double-diffusive convection occurs and controls transport within the object. A commonly studied case is that of semi-convection at the edge of core-convective stars (Ledoux 1947; Tayler 1954; Schwarzchild & Härm 1958; Merryfield 1995). In moderately massive stars for example, a mean molecular weight gradient develops over time at the edge of the core, and eventually begins to affect convection. When it is strong enough to stabilize the fluid, the fully-convective region shrinks in size, leaving behind a “semi-convective” region in which transport is controlled by double-diffusive processes instead. Other related examples are reviewed by Merryfield (1995).

The possible role of double-diffusive convection in regulating thermal and compositional transport in the interior of giant planets was recognized later, and has been discussed in the context of convective planetary envelopes where the stabilizing component is helium (Stevenson & Salpeter 1977), higher-metallicity material at the edge of a rocky core (see Stevenson 1985 and in particular Figure 2 of his paper), or methane (Gierasch & Conrath 1987). Double-diffusive convection has also been proposed to explain the abnormally large radii of some transiting exoplanets (Chabrier & Baraffe 2007). Finally, it has recently been invoked as a new mechanism for driving pulsations in white dwarfs (Shibahashi 2007; Kurtz et al. 2008).

Before we move on to describe existing work on double-diffusive convection, we note that it should not be confused with that arising from the related double-diffusive “fingering” instability (Stern 1960). The latter also occurs in Ledoux-stable systems, but in the opposite situation when the more rapidly diffusing thermal field is stably stratified while the slowly-diffusing compositional field is un unstably stratified. Its turbulent manifestation is often referred to as “thermohaline convection” in astrophysics, by analogy with the oceanic case in which the compositional gradient is due to salt. We prefer a more general terminology and use the alternative name “fingering convection”. Figure 2 illustrates for clarity the various regimes of convective instability. See Traxler et al. (2010a) for a study of fingering
convection in the astrophysical regime.

1.2. Previous work on double-diffusive convection

Very little is known about mixing by double-diffusive convection, despite its obvious importance in stellar and planetary astrophysics. Linear stability reveals that the unstable modes take the form of overstable gravity waves (Walin 1964; Kato 1966; Baines & Gill 1969). What governs the saturation of the instability in the astrophysically-relevant parameter regime, however, remains essentially unknown. Note that by “astrophysically-relevant” we imply a low diffusivity ratio, $\tau \ll 1$, and a low Prandtl number, $\text{Pr} = \nu/\kappa \ll 1$, where $\nu$ is the viscosity. Both numbers are typically of the order of $10^{-5} – 10^{-7}$ in stellar and planetary interiors.

To add to the complexity of the problem, double-diffusive convection is known in some cases to lead to thermo-compositional layering, i.e. to the development of stacks of well-mixed fully-convective layers separated by strongly stratified interfaces. This double-diffusive layering is commonly observed for example in the arctic ocean (Neal et al. 1969; Toole et al. 2006) where cool fresh water lies on top of warmer, saltier water. It has been studied extensively in laboratory experiments (Turner 1968; Huppert & Linden 1979; Huppert & Turner 1980; Turner 1985). An important result of these studies is that turbulent mixing in the presence of layers is significantly enhanced compared with that of a system which has the same overall contrast in temperature and composition, but where the stratification is everywhere much smoother.

Whether layer formation occurs in double-diffusive convection at low Prandtl number actually remains to be determined – it is usually assumed (Spruit 1992; Chabrier & Baraffe 2007), by analogy with the high-Prandtl number oceanographic case. It is important to realize, however, that such analogies can be misleading. This was recently demonstrated by Traxler et al. (2010a) in the case of fingering convection. Similar thermohaline staircases are ubiquitously observed in fingering-unstable regions of the ocean (Schmitt et al. 2005), and have been shown to form spontaneously through a secondary linear instability of homogeneous fingering convection (Radko 2003; Traxler et al. 2010b; Stellmach et al. 2010) (see §5 too). However, Traxler et al. (2010a) demonstrated that this secondary instability does not happen in the astrophysical context and concluded that thermo-compositional layers are not expected in that case. In other words, given that the analogy with the heat-salt system doesn’t hold in the fingering regime, one should be extremely cautious about using it a priori in the double-diffusive regime.
To summarize, it is known that “homogeneous” and “layered” double-diffusive convection have rather different mixing properties. A good mixing parametrization therefore needs to incorporate both cases, and must include a criterion to decide whether the system considered lies in one or the other. Existing parametrizations, by contrast, have so far either ignored the possible effect of layering (Schwarzschild & Härn 1958; Langer et al. 1983), or relied on it (Spruit 1992).

A few numerical simulations of double-diffusive convection have been performed to date to address the problem. The first of this kind (to our knowledge), were presented by Merryfield (1995). He ran a series of two-dimensional (2D) anelastic simulations, horizontally-periodic, and bounded in the vertical direction by two plates. Double-diffusive convection was forced through the imposed boundary conditions, which maintained a given overall temperature and compositional contrast across the domain. Merryfield (1995) focused on understanding how the mechanism responsible for the saturation of the initial double-diffusive instability depends on the governing parameters, and in particular the strength of the thermal driving, the Prandtl number and the diffusivity ratio. He also compared the outcome of his simulations with existing parametrizations of double-diffusive convection both in the absence of layers, and in the presence of initially forced layers. One of the main difficulties encountered was the development of numerical instabilities in cases with strong thermal driving, which prevented him from drawing definite conclusions on the long-term statistical properties of the turbulence. In addition, in many of the runs the flows were eventually influenced by the presence of domain boundaries.

In subsequent years, two additional attempts at modeling double-diffusive convection were made. Biello (2001), as part of his PhD thesis, ran a series of 2D fully-compressible simulations which complement those of Merryfield (1995). There, the system was also confined between two plates, but the boundary conditions were “fixed flux” conditions. Biello (2001) was interested in studying more specifically the layer formation process, and his experimental setup was similar to that of heat-salt laboratory experiments (e.g. Huppert & Linden 1979). For this purpose, the simulations were initialized with stable uniform gradients in temperature and composition, but destabilized at \( t = 0 \) by increasing the heat flux at the bottom boundary. He found that the first bottom layer easily forms, but did not observe any subsequent layer formation. He analyzed the dynamics of the interface, and concluded that interfacial transport was dominated by wave-breaking rather than by diffusive processes as is often assumed. However, boundary effects also began to influence the results of his simulations after some time. Bascoul (2007), also as part of his PhD thesis, studied a similar 2D time-dependent system, in which the initial background state had a homogeneous composition and neutrally stable temperature gradient, and where the system was destabilized by an imposed heat and mean-molecular weight flux through the bottom boundary. He
also observed the growth of a convective layer near the bottom boundary and studied its development, for high Prandtl number (heat-salt regime) and for low Prandtl number (“astrophysical” regime). He was unable, however, to run his simulations long enough to achieve statistical equilibrium.

In this paper, we present a new series of three-dimensional (3D) numerical experiments to study mixing by double-diffusive convection. We approach the problem from a different but complementary angle, and try to address some of the inherent shortcomings of the experimental setup used in previous studies: we focus specifically on measuring the quasi-steady statistical properties of double-diffusive turbulence, and use a numerical setup which minimizes the effects of domain boundaries.

We discuss the model setup and briefly summarize its linear stability properties in §2. The numerical algorithm and the selection of the experimental parameters are described in §3. In each case, as described in §4, we extract values for the transport coefficients while the system is in a state of homogeneous turbulence. However, we find that for more unstable systems a secondary instability leads to the formation of thermo-compositional layers. The layers continue evolving and successively merge, and each merger is accompanied with a significant increase in the transport coefficients. These results are discussed in §5, and compared with a recent theory of layer formation in the oceanographic context (Radko 2003). We find good agreement between this theory and our numerical results, which enables us to deduce a general criterion for the spontaneous formation of layers in double-diffusive convection in the astrophysical context. We discuss our results in §6.

2. Model description and linear analysis

In this section we present the governing equations and briefly summarize known results on the linear stability of the problem. Our formalism is overall very similar to the one used by Traxler et al. (2010a) for fingering convection.

2.1. Governing equations and boundary conditions

As we demonstrate later (see §2.2), the typical lengthscale of the unstable motions (in the absence of layers) is of the order of a few kilometers at most in parameter regimes typical of stellar and planetary interiors. It is therefore justified to neglect the effect of curvature entirely and work in a local Cartesian reference frame \((x, y, z)\). Here, gravity defines the vertical direction: \( g = -g e_z \). We ignore the possible presence of magnetic fields
for simplicity, and neglect the effect of rotation. The latter is justified whenever the mean rotation rate is much smaller than the buoyancy (Brunt-Väisälä) frequency, which is often the case.

In all that follows, we simplify the problem by using the Boussinesq approximation [Spiegel & Veronis 1960]. This approximation is a regular asymptotic limit of the primitive governing equations (for mass, momentum and energy conservation) when the domain height $L_z$ is much smaller than the density scaleheight $D_\rho$ (i.e. $L_z/D_\rho \to 0$), and when the typical flow velocity $u$ is much smaller than the sound speed $c_s$ (i.e. $u/c_s \to 0$). It is therefore particularly relevant here since the typical lengthscale of the perturbations is much smaller than a pressure or density scaleheight (unless the region considered is very close to the photosphere), and the typical velocities are always substantially subsonic (unless the system is very close to the onset of overturning convection).

Since we consider a small fluid region within a much larger system, we can assume that the background composition and temperature gradients $\mu_{0z}$, $T_{0z}$ and $T^{\text{ad}}_{0z}$ are constant. Here, the index 0 denotes a background field, and the index $z$ denotes a derivative with respect to the vertical coordinate $z$. We restrict our study to the case of double-diffusive convection by choosing the background stratification such that $\nabla \mu > \nabla - \nabla^{\text{ad}} > 0$, or equivalently $\mu_{0z} < T_{0z} - T^{\text{ad}}_{0z} < 0$.

In the Boussinesq approximation the mass conservation equation is replaced by the continuity equation
\begin{equation}
\nabla \cdot \mathbf{u} = 0 ,
\end{equation}
where $\mathbf{u} = (u, v, w)$ is the velocity field, and the (dimensional) thermal energy equation is approximated (e.g. Ulrich 1972) by
\begin{equation}
\frac{\partial T}{\partial t} + \mathbf{u} \cdot \nabla T + (T_{0z} - T^{\text{ad}}_{0z})w = \kappa_T \nabla^2 T ,
\end{equation}
where $T$ now represents the dimensional temperature perturbation (all background quantities being denoted by the subscript 0 instead), and we have assumed for simplicity that the thermal diffusivity is constant. This equation is derived from an energy conservation principle, noting that in the Boussinesq approximation temperature and entropy perturbations are proportional. The term $T_{0z} - T^{\text{ad}}_{0z}$ therefore models the advection of the background entropy gradient.

In the Boussinesq approximation, the density, temperature and mean molecular weight ($\rho$, $T$ and $\mu$ respectively) are related via
\begin{equation}
\frac{\rho}{\rho_0} = -\alpha T + \beta \mu ,
\end{equation}
where $\alpha$ and $\beta$ are constants.
where \( \rho_0 \) is the mean density of the region considered, and \( \alpha \) and \( \beta \) are the coefficients of thermal expansion and compositional contraction respectively (e.g. \( \alpha = 1/T_0 \) and \( \beta = 1/\mu_0 \) for a perfect gas, where \( T_0 \) and \( \mu_0 \) are the mean temperature and mean-molecular weight in the region considered).

We construct our numerical model in such a way as to minimize the effects of the computational domain boundaries. For this purpose we use a triply-periodic box of size \( (L_x, L_y, L_z) \), in which convection is permanently forced by the aforementioned background stratification. This approach has recently been used with success in modeling and studying the formation of thermostaline staircases in the oceanographic context by [Radko (2003)] and [Stellmach et al. (2010)], and is discussed in more detail in these papers. In this framework, the temperature and mean molecular weight fields can be written as the sum of a background variation plus perturbations which are triply-periodic functions of \((x, y, z)\) such that

\[
T(x, y, z, t) = T(x + L_x, y, z, t) = T(x, y + L_y, z, t) = T(x, y, z + L_z, t) ,
\]

and similarly for \( \mu \). The pressure perturbation \( p \) and velocity field \( \mathbf{u} \) are also assumed to be triply-periodic functions in the same way.

We non-dimensionalize the equations using the anticipated lengthscale of the fastest growing modes of linear instability, which is a thermal diffusion scale (e.g. [Baines & Gill 1969]):

\[
d = \left( \frac{\kappa_T \nu}{\alpha g |T_{0z} - T_{0z}^{ad}|} \right)^{1/4} = \left( \frac{\kappa_T \nu}{N^2} \right)^{1/4} ,
\]

where \( N \) is the buoyancy frequency. Very roughly, in typical stellar interiors, \( \nu = O(10) \text{cm}^2/\text{s} \), \( \kappa_T = O(10^7) \text{cm}^2/\text{s} \) and \( N^2 = O(10^{-6}) \text{s}^{-2} \) so \( d = O(10^{3.5}) \text{cm} \), or in other words, \textit{a few hundreds of meters} only.

Note that with this definition, the thermal Rayleigh number defined on the finger scale is exactly one, while the global Rayleigh number,

\[
Ra_T = \frac{\alpha g |T_{0z} - T_{0z}^{ad}| L_z^4}{\kappa_T \nu} = \left( \frac{L_z}{d} \right)^4
\]

is a function of the dimensionless height of the domain only. The unit timescale is taken to be the diffusion timescale across \( d \), namely \( [t] = d^2/\kappa_T \), and the velocity scale is \( [v] = \kappa_T/d \). The unit temperature is \( [T] = d |T_{0z} - T_{0z}^{ad}| \), and the unit mean-molecular weight is \( [\mu] = (\alpha/\beta) |T_{0z} - T_{0z}^{ad}| d \). The resulting non-dimensional governing equations are

\[
\frac{1}{Pr} \left( \frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} \right) = -\nabla \tilde{p} + (\tilde{T} - \tilde{\mu}) \mathbf{e}_z + \nabla^2 \mathbf{u} ,
\]
\[
\frac{\partial \tilde{T}}{\partial t} + \tilde{u} \cdot \nabla \tilde{T} - \tilde{w} = \nabla^2 \tilde{T},
\]
\[
\frac{\partial \tilde{\mu}}{\partial t} + \tilde{u} \cdot \nabla \tilde{\mu} - R_0^{-1} \tilde{w} = \tau \nabla^2 \tilde{\mu},
\]
\[
\nabla \cdot \tilde{\mathbf{u}} = 0 ,
\]
where quantities with tildes now represent the dimensionless, triply-periodic perturbations. The non-dimensionalization introduces three parameters, namely the aforementioned Prandtl number \( \text{Pr} = \nu/\kappa_T \) and diffusivity ratio \( \tau = \kappa_\mu/\kappa_T \), as well as the so-called density ratio \( R_0 = \frac{\alpha |T_0z - T_0^\text{ad}|}{\beta |\mu_0z|} = \frac{\nabla - \nabla_\text{ad}}{\nabla_\mu} \).

Finally, note that for reasons described in §2.3, it is common and preferable to work with the inverse density ratio \( R_0^{-1} = 1/R_0 \) as a governing parameter.

2.2. Linear stability

The linear stability of this problem is well-understood, thanks to the works of Walin (1964), Veronis (1965) and Baines & Gill (1969) in the oceanographic context, and Kato (1966) in the astrophysical context. The salient points are summarized here for completeness and clarity, as some of them will be used later.

To analyze the stability of the governing equations we first linearize them around \( \tilde{T} = \tilde{\mu} = 0 \) and \( \tilde{\mathbf{u}} = 0 \), and assume normal forms for all perturbations as \( q = \hat{q} e^{ilx + imy + ikz + \lambda t} \) where hatted quantities are the mode amplitudes, \( l, m, \) and \( k \) are horizontal and vertical wavenumbers respectively, and \( \lambda \) is the growth rate. This procedure yields a cubic equation for \( \lambda \) in terms of the wavevector \((l,m,k)\) and of the non-dimensional parameters:

\[
\left( \frac{\lambda}{\text{Pr}} + K^2 \right) (\lambda + K^2)(\lambda + \tau K^2) \left( \frac{K^2}{l^2 + m^2} \right) - (\lambda + \tau K^2) + R_0^{-1}(\lambda + K^2) = 0 ,
\]
where \( K^2 = l^2 + m^2 + k^2 \).

It can be shown that the most unstable mode always occurs when the vertical wavenumber, \( k \), is equal to zero. This corresponds to an "elevator mode", similar to the ones found in the related problem of fingering convection and of homogeneous Rayleigh-Bénard convection. Furthermore, without loss of generality, we may select \( l \) or \( m \) to be zero by reorienting the Cartesian frame so that the fastest growing modes can be described using one horizontal wavenumber only. The cubic equation (12) then becomes:

\[
\left( \frac{\lambda}{\text{Pr}} + l^2 \right) (\lambda + l^2)(\lambda + \tau l^2) - (\lambda + \tau l^2) + R_0^{-1}(\lambda + l^2) = 0 .
\]
There are typically one real and two complex roots of this equation. The real root is always negative, achieving a maximum of zero when \( l = 0 \). However, the complex root yields a positive growth rate when the inverse density ratio lies in the interval \( R_0^{-1} \in (1, \frac{Pr + 1}{Pr + \tau}) \). The oscillation frequency is close to the buoyancy frequency. The unstable modes can therefore be viewed as overstable gravity waves, as mentioned earlier.

By maximizing the real part of \( \lambda \) over all horizontal wavenumbers, we can find the most unstable mode in the system. Figure 1 shows its wavenumber and growth rate for \( Pr = \tau = 1/3 \). By and large, the most rapidly growing mode has wavelength of the order of 20\( d \) regardless of \( R_0^{-1} \) for the parameters selected, which implies a lengthscale of a few kilometers as described in §2.1.

Fig. 1.— Growth rate (solid line) and horizontal wavenumber \( l \) (dotted line) of the most rapidly growing gravity wave (i.e. the double-diffusive mode of instability), as a function of the inverse density ratio, for \( Pr = \tau = 1/3 \). All quantities plotted are in the units used in this paper (see §2.1).

2.3. \( R_0 \) vs. \( R_0^{-1} \)

As mentioned previously, it is standard in oceanographic studies of double-diffusive convection to use the parameter \( R_0^{-1} = \beta \mu_0 z/\alpha T_0 z \), while \( R_0 = \alpha T_0 z/\beta \mu_0 z \) is used in studies
of fingering convection. The reason for this change in convention is to emphasize the symmetry between the two regimes, which is apparent in Figure 2. Indeed, in that case, one can write that overturning convection occurs for

\[
R_0 < 1 \text{ fingering regime,} \\
R_0^{-1} < 1 \text{ diffusive regime,}
\]

(14)

the double-diffusive regime occurs for

\[
1 \leq R_0 \leq \frac{1}{\tau} \text{ fingering regime,} \\
1 \leq R_0^{-1} \leq \frac{\Pr + 1}{\Pr + \tau} \text{ diffusive regime,}
\]

(15)

and finally that the system is stable for

\[
R_0 > \frac{1}{\tau} \text{ fingering regime,} \\
R_0^{-1} > \frac{\Pr + 1}{\Pr + \tau} \text{ diffusive regime.}
\]

(16)

The symmetry is even more striking at low Prandtl number, where the critical value for complete stability is much larger than one in both cases. For the diffusive regime, this implies that the system is fully convective up to the Ledoux-stability criterion, and then double-diffusively convective between Ledoux-stability and Schwarzchild-stability. A summary of the various regimes of instability, for the compositionally homogeneous case, for the fingering case and for this diffusive case is presented in Figure 2.

3. Numerical Experiments

3.1. Description of the experiments

We solve the governing equations and boundary conditions presented in §2 using a high performance spectral code developed by S. Stellmach (Traxler et al. 2010b; Stellmach et al. 2010; Traxler et al. 2010a), specifically designed for the study of fingering convection in the oceanographic context. The code can be used “as is” to model double-diffusive convection simply by reversing the sign of both background gradients. Since it was developed for

\[1 \text{ For added confusion, studies of double-diffusive convection often call } \beta \mu_0 z / \alpha T_0 z \text{ the density ratio and denote it as } R_0 \text{ although we will not use this naming convention here.}\]
studying oceanic convection, the code does not consider entropy separately from temperature. As a result, it intrinsically assumes that $T_{0z}^{\text{ad}} = 0$, so that $R_0 = \alpha T_{0z}/\beta \mu_{0z}$. By virtue of being non-dimensional, however, the results can nevertheless straightforwardly be applied to the astrophysical context simply by interpreting $R_0$ as being defined by $11$.

Our goals are threefold: (a) to characterize transport by homogeneous double-diffusive convection (i.e. in the absence of layers), (b) to determine if, under which conditions, and through which process thermo-compositional layers may form and (c) to characterize transport by layered double-diffusive convection if appropriate.

For this purpose, we ran a sequence of exploratory numerical experiments setting $\Pr = 1/3$ and $\tau = 1/3$. We selected $\Pr$ and $\tau$ below unity to be in the “low Prandtl number, low diffusivity ratio” regime, but not too small so that we could remain in a numerically-tractable region of parameter space. While it is possible to run simulations for lower values of these parameters in 3D (see Traxler et al. 2010a, for example), these are computationally much more demanding. In this first analysis, we wanted to be able to run a uniform set of simulations across the whole instability range, and integrate some of them for a significant length of time to observe the layer formation and merger process if and when it occurs. We discuss the applicability of our results to lower Prandtl number and lower diffusivity ratio environments in §6.

In order to get statistically meaningful measurements of the turbulent fluxes in this system, to address points (a) and (c) raised above, we must ensure that the computational box contains at least a few wavelengths of the most unstable mode in the horizontal directions (Traxler et al. 2010b). As found in Figure 1, these are of the order of about $20d$ so we select a domain size with $L_x = L_y = 100d$ in all simulations. We use an aspect ratio of 1 and choose $L_z = 100$, which corresponds to $\text{Ra}_T = 10^8$. For comparison, we also ran a series of simulations with narrower domains ($L_x = L_y = 50d$) at $\text{Ra}_T = 10^8$ for all $R_0^{-1}$, as well as one taller-domain simulations using $L_z = 178d$ (equivalently, $\text{Ra}_T = 10^9$) for $R_0^{-1} = 1.2$. See Table 1 for a comparison of these runs. In dimensional terms, the domain height is of the order of a few tens of kilometers using the estimates presented in §2.1 for typical stellar interiors, which is much smaller than a pressure or density scaleheight and therefore fully justifies the use of the Boussinesq approximation.

In terms of spatial resolution, we use a sufficient number of Fourier modes in all simulations to resolve the typical size of the composition and velocity perturbations (which are roughly of the same size when $\tau = \Pr$). Similar runs with different spatial resolutions are presented in Table 1, and show consistent results.
Table 1: Summary of runs performed and measured turbulent fluxes in the homogeneous phase.

<table>
<thead>
<tr>
<th>$R_0^{-1}$</th>
<th>$N_x, N_y$</th>
<th>$N_z$</th>
<th>$t_{\text{start}}$</th>
<th>$t_{\text{end}}$</th>
<th>Nu$_T$</th>
<th>Nu$_\mu$</th>
<th>1/$\gamma_1$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1.1</td>
<td>192</td>
<td>192</td>
<td>260</td>
<td>370</td>
<td>5.7±3.9</td>
<td>11.2±4.2</td>
<td>.63±.10</td>
</tr>
<tr>
<td>1.1$^{(a)}$</td>
<td>192</td>
<td>192</td>
<td>270</td>
<td>460</td>
<td>5.0±1.2</td>
<td>9.6±3.5</td>
<td>.61±.10</td>
</tr>
<tr>
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<td>192</td>
<td>300</td>
<td>630</td>
<td>3.9±7</td>
<td>6.7±2.3</td>
<td>.58±.11</td>
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<tr>
<td>1.15$^{(a)}$</td>
<td>96</td>
<td>192</td>
<td>290</td>
<td>510</td>
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<td>96</td>
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Note. — For each value of the inverse density ratio $R_0^{-1}$, the second and third columns show the horizontal and vertical number of grid-points used. The times $t_{\text{start}}$ and $t_{\text{end}}$ define the interval over which the turbulent fluxes, measured via Nu$_T$ and Nu$_\mu$, were averaged. The quantity $\gamma_1$ is the primary way of measuring the total flux ratio $\gamma^{\text{tot}}$: see text for detail. Errors quoted indicate the rms fluctuations of Nu$_T(t)$, Nu$_\mu(t)$ and 1/$\gamma_1(t)$ about the respective means. Unless otherwise specified (see footnotes), simulations were run at domain sizes of 100$d \times 100d \times 100d$. Note that the simulation for $R_0^{-1} = 2.1$ was integrated for significant time, but by the end of the run the perturbations had not grown to high-enough amplitudes to induce a statistically significant heat or compositional flux.

$^{a}50d \times 50d \times 100d$

$^{b}44d \times 44d \times 178d$

$^{c}89d \times 89d \times 178d$
3.2. Qualitative description of the results

We ran a first set of cubic-domain simulations \((L_x = L_y = L_z = 100d)\) for \(R_0^{-1}\) varying across the whole instability range, which for our selected parameters corresponds to \(1 < R_0^{-1} < 2.167\). We found that the initial behavior of the system is qualitatively similar for all density ratios: the perturbations first grow exponentially, and then saturate into a homogeneous turbulent state.

The initial exponential growth is well-approximated by linear theory. This is illustrated in Figure 3 for example, which shows the early temporal evolution of the rms velocity for the \(R_0^{-1} = 1.2\) run, and compares it with the growth of the fastest-growing mode according to linear theory. The fit is very good – the small discrepancy at early times can be attributed to the fact that more than one mode are excited, but the most rapidly growing mode then quickly takes over. We confirmed that the initial instability is independent of the Rayleigh number (i.e. the domain height) by comparing these results with those of a taller-domain simulation \((L_z = 178)\) at the same \(R_0^{-1}\).

The perturbations saturate once nonlinear effects become important. We find that the level of saturation of the turbulence is also independent of the domain height (alternatively, of the Rayleigh number), as shown in Figure 3. However, it depends sensitively on the value of the density ratio. Figure 4 shows the temporal evolution of the thermal Nusselt number \(\text{Nu}_T\) as a function of the inverse density ratio. A Nusselt number is the ratio of the total flux (diffusive + turbulent) to the diffusive flux, so we define the thermal Nusselt number as

\[
\text{Nu}_T = \frac{-\kappa T T_0 z + wT}{-\kappa T T_0 z} = 1 + \langle \tilde{w} \tilde{T} \rangle, \tag{17}
\]

where the angular brackets denote a spatial average over the entire domain. We also define the equivalent compositional Nusselt number \(\text{Nu}_\mu\) as

\[
\text{Nu}_\mu = \frac{-\kappa \mu \mu_0 z + w\mu}{-\kappa \mu \mu_0 z} = 1 + \frac{R_0}{\tau} \langle \tilde{w} \tilde{\mu} \rangle. \tag{18}
\]

In each case, in the second expression the turbulent fluxes are expressed in non-dimensional form recalling that \(T_0 z < 0\) and \(\mu_0 z < 0\) while \(T\) and \(\mu\) are non-dimensionalized using \(|T_0 z|\).

In all simulations presented in Figure 4 the thermal Nusselt number increases exponentially until saturation, and remains approximately constant during the early saturated phase. After saturation, however, simulations which were run using a lower \(R_0^{-1}\) behave in a fundamentally different way from those at higher \(R_0^{-1}\). In latter case, (for \(R_0^{-1} \geq 1.35\)), the Nusselt number at saturation remains statistically steady for the entire duration of the run. By contrast, for \(R_0^{-1} < 1.35\), the Nusselt number later continues to increase.
When visualizing the results, we find that this second increase in the turbulent transport properties of the system corresponds to the formation of well-mixed fully convective layers separated by thin stably stratified interfaces (see Figure 5 for example). The Nusselt number continues to increase as the layers merge, until a single layer is left. We have therefore established that layers can indeed form in low-Prandtl number double-diffusive convection, and that, as in the high-Prandtl number regime, a layered system transports heat more efficiently than a homogeneous system with the same overall temperature and compositional gradient. We now study both the homogeneous phase and the layered phase in more detail in §4 and §5 respectively.

4. Homogeneous Double-Diffusive Convection

We focus here on the homogeneous phase, prior to the formation of the first set of layers, and measure the transport properties of the turbulence via the respective Nusselt numbers defined in (17) and (18). Note that the time period between the initial saturation of the double-diffusive instability and the onset of layer formation, when it occurs, varies with $R_0^{-1}$ (see Figure 4). Table 1 indicates, for each simulation, the time interval over which the system is in this homogeneously turbulent phase and during which we average the instantaneous Nusselt numbers.

The mean Nusselt numbers thus extracted are presented in Table 1 and illustrated in Figure 6a. The errors quoted denote the rms of the fluctuations around the respective means. For most values of $R_0^{-1}$, we ran a series of simulations with different resolution, or different box size, or both. The measured Nusselt numbers are always consistent within the errorbars.

As expected, turbulent mixing is negligible close to marginal stability, i.e. when $R_0^{-1} \rightarrow (Pr + 1)/(Pr + \tau)$. It increases as $R_0^{-1}$ decreases through the instability range, and grows rapidly close to the onset of overturning convection (i.e. as $R_0^{-1} \rightarrow 1$). However, we find that it remains fairly weak, with $\text{Nu}_T$ of the order of a few and $\text{Nu}_\mu$ of the order of ten, even for our lowest $R_0^{-1}$ run ($R_0^{-1} = 1.1$). By contrast, homogeneous Rayleigh-Bénard convection in the absence of compositional gradient, in the same parameter regime ($Pr = 1/3$, $Ra_T = 10^8$ and aspect ratio one), would have a thermal Nusselt number of the order of several thousands (Garaud et al. 2010). This shows that while turbulent mixing is not negligible in this homogeneous double-diffusive regime, it is nevertheless much smaller than that induced by standard convection. Presumably, there exists a very narrow range of inverse density ratio close to unity, but above it, across which turbulent mixing rapidly but continuously increases towards the fully convective value. This will be the subject of a subsequent study.
Since this quantity is crucial to the theory of layer formation in the fingering regime (Radko 2003; Traxler et al. 2010a), we also compute the so-called total flux ratio \( \gamma_{\text{tot}} \), defined as the ratio of the total buoyancy flux due to heat transport to that due to compositional transport:

\[
\gamma_{\text{tot}} = \frac{F_{T}^{\text{tot}}}{F_{\mu}^{\text{tot}}} = \frac{R_{0} \text{Nu}_{T}}{\tau \text{Nu}_{\mu}}.
\]  

Equation (19)

There are two different ways of computing \( \gamma_{\text{tot}} \) from our numerical results: the first and preferable method involves calculating \( \gamma_{\text{tot}}(t) \) at every timestep and then taking its mean during the homogeneous turbulent phase. The second involves using the measured mean values of \( \text{Nu}_{T} \) and \( \text{Nu}_{\mu} \) directly into (19). These two versions of \( \gamma_{\text{tot}} \) are denoted as \( \gamma_{1} \) and \( \gamma_{2} \) respectively, in Figure 6b. They are consistent within the errorbars.

Note that we are actually plotting the function \( \gamma_{\text{tot}}^{-1}(R_{0}^{-1}) \) in Figure 6b. The reason for showing it rather than \( \gamma_{\text{tot}}(R_{0}^{-1}) \) will be revealed in §5.2 and is related to the layer formation mechanism. As we shall see, the fact that \( \gamma_{\text{tot}}^{-1} \) is a decreasing function of \( R_{0}^{-1} \) in the interval \([1, 1.35]\), i.e. the same interval in which layers are observed to emerge, is not a coincidence.

5. Layered convection

As discussed in §3.2, in all simulations with \( R_{0}^{-1} < 1.35 \) we find that the system does not remain for long in a state of homogeneous double-diffusive convection, but spontaneously develops thermo-compositional layers instead. We study this process in more detail in this section, focusing on the \( R_{0}^{-1} = 1.2 \) run. We choose this value of \( R_{0}^{-1} \) rather than one closer to unity, based on the results of Figure 4. Indeed, for \( R_{0}^{-1} = 1.2 \) layer formation and mergers are “slow enough” to be studied, but proceed much more rapidly for lower \( R_{0}^{-1} \).

5.1. General considerations

The results presented in Figure 4 are for a cubic box with height \( L_{z} = 100 \). In that run, we observe that two layers initially form, then merge into one a little while later. In order

---

2Note that \( F_{T}^{\text{tot}} \) and \( F_{\mu}^{\text{tot}} \) are dimensionless here. The interpretation of \( \gamma_{\text{tot}} \) as a buoyancy flux is more apparent when we go back to the dimensional quantities \( F_{T}^{\text{tot, dim}} \) and \( F_{\mu}^{\text{tot, dim}} \), since \( F_{T}^{\text{tot}} = F_{T}^{\text{tot, dim}} / \kappa |T_{0z}| \) and \( F_{\mu}^{\text{tot}} = F_{\mu}^{\text{tot, dim}} / (\alpha/\beta) \kappa |T_{0z}| \).

3To be precise, we calculate \( \gamma_{\text{tot}}^{-1}(t) \) at every timestep, take the average of this function (shown in Figure 6b), and then take its inverse to get the mean \( \gamma_{\text{tot}} \). The reason for doing this in two steps is that there are very occasional events where \( \text{Nu}_{\mu}(t) < 0 \), at which point \( \gamma_{\text{tot}}(t) \to \infty \) but the inverse remains well-defined.
to minimize the influence of the finite domain height on the layer formation and merger process, we use from here on the taller-domain simulation, for which $L_z = 178$ ($Ra_T = 10^9$). The appearance and successive merger of layers observed in that run is shown in Figure 5. We find that twice as many layers initially form in this nearly-twice-as-tall domain, and appear roughly at the same time as they did in the cubic box run. This shows that layer formation depends only on local processes, and knows about the thermal scale $d$ rather than the domain scale. The initial layer height, in each case, is of the order of $45d – 50d$.

Figure 7 shows the evolution of the thermal and compositional Nusselt numbers in the tall-domain simulation. We see quite clearly the stepwise increase in transport which accompanies the layer formation and successive mergers: there are five fairly well-defined plateaus, corresponding to the homogeneous phase, (up to about $t = 1000$, see Figure 5a), the four-layer phase (up to about $t = 1200$, see Figure 5b), the three-layer phase (up to $t = 1450$, see Figure 5c), the two-layer phase (up to $t = 1650$, see Figure 5d), and the final one-layer phase, see Figure 5e. It is also interesting to note how the heat and compositional fluxes follow each other closely throughout the entire simulation.

A useful way of studying the formation and structure of the layers was presented by Stellmach et al. (2010), and consists in looking at Fourier modes of the density perturbation:

$$\tilde{\rho}(x, y, z, t) = \sum_{l, m, k} \tilde{\rho}_{l, m, k}(t)e^{ilx + imy + ikz}, \quad (20)$$

where $k$ is an integer multiple of $2\pi/L_z$ and similarly for $l$ and $m$. These Fourier modes are straightforwardly extracted from our numerical solutions since our code is spectral.

By definition, the $\tilde{\rho}_{0,0,k}$ modes are the vertical Fourier modes of the horizontally-averaged density profile. Since thermo-compositional staircases are also density staircases (i.e. with a nearly uniform density within the layers separated by sharp pycnoclines), a staircase with $n$ layers has a dominant vertical wavenumber $k_n = 2\pi n/L_z$. This is seen most clearly in Figure 8 which shows the norm of $\tilde{\rho}_{0,0,k}$ for the four gravest non-zero modes, and illustrates how $k_4, k_3, k_2$ and $k_1$ successively take over as the dominant mode as the layers form and merge.

A rather striking feature of Figure 8 however, is that the $k_4$ and $k_3$ modes actually begin to grow as early as $t = 500$, long before the layers appear in visual inspection of the temperature and composition field (as in Figure 5 for example). An even more striking result is that this growth is well-approximated by an exponential. This strongly suggests that layer formation arises through a secondary linear instability of homogeneous double-diffusive convection (see next section) rather than through stochastic overturning events of the growing gravity waves, as is commonly assumed (e.g. Stevenson 1979; Spruit 1992).
We now study in more detail, successively, the layer formation process, and the evolution of the fluxes through the staircase as the mergers proceed.

5.2. The $\gamma$−instability

5.2.1. The $\gamma$−instability of fingering convection

Recently, significant progress has been made in understanding the spontaneous formation of layers in fingering convection, thanks to the groundbreaking work of Radko (2003) in the oceanographic context. Radko discovered that homogeneous fingering convection in that regime is linearly unstable to a secondary large-scale instability, which takes the form of exponentially-growing horizontally-invariant perturbations in the density (equivalently, temperature and composition) profile. The perturbations grow through positive feedback between the perturbed stratification and the modulated fingering fluxes. Upon reaching a critical amplitude, the modulated density profile becomes unstable to direct overturning convection, and the system rapidly transitions into a fully-formed, regularly-spaced staircase. A sufficient condition for this instability to occur is that the ratio of the total buoyancy flux due to heat to the total buoyancy flux due to salt, the quantity referred to as $\gamma_{\text{tot}}$ in §4, should be a strictly decreasing function of the density ratio. For this reason, this new instability was called the $\gamma$−instability and the associated perturbations, the $\gamma$−modes.

Radko’s theory was validated first through two-dimensional simulations (Radko 2003), and more crucially through 3D simulations (Stellmach et al. 2010). Stellmach et al. (2010) were the first to find spontaneous layer formation in 3D simulations of fingering convection in the “oceanic” parameter regime. They analyzed the growth and structure of the emergent layers using the vertical Fourier modes of the density profile, as we have done in the previous section. They found, as we do in Figure 8, that the Fourier mode which corresponds to the number of layers of the emerging staircase began to grow long before the layers form, and that its growth rate could be predicted very accurately by Radko’s $\gamma$−instability theory (see their Figure 6).

Traxler et al. (2010a) extended Radko’s theory for layer formation in fingering convection to a parameter regime more relevant of the astrophysical context. Their results suggest that at low Prandtl number and low diffusivity ratio $\gamma_{\text{tot}}$ is always an increasing function of the density ratio, so that $\gamma$−modes are stable. They concluded that spontaneous layer formation is unlikely in astrophysical fingering convection.
5.2.2. The $\gamma$–instability of double-diffusive convection

Radko’s theory is quite general, and can be applied with only minor modifications to the case of double-diffusive convection. Let us consider a system in a homogeneously turbulent state, with a background density ratio $R_0$. We know through the results of §4 that this system drives a non-zero total vertical heat flux $F_T^{\text{tot}}$ and a non-zero total vertical compositional flux $F_\mu^{\text{tot}}$ with

$$F_T^{\text{tot}} = \text{Nu}_T(R_0) \text{ and } F_\mu^{\text{tot}} = \frac{F_T^{\text{tot}}}{\gamma^{\text{tot}}(R_0)} .$$

(21)

As long as the background is homogeneous, however, $F_T^{\text{tot}}$ and $F_\mu^{\text{tot}}$ are constant in the domain, and do not affect the temperature or chemical composition of the system.

Let us now assume that this homogeneously turbulent system is modulated by a large-scale, horizontally-invariant perturbation, so that the horizontally-averaged temperature and mean-molecular weight profiles, $\bar{T}$ and $\bar{\mu}$ can be written as

$$\bar{T}(z,t) = \hat{T} e^{ikz + \Lambda t} ,$$

(22)

and similarly for $\bar{\mu}$. These large-scale perturbations change the local density ratio, which we now write as $R_\rho(z,t) = R_0 + R'(z,t)$. Since the turbulent fluxes are functions of $R_\rho$, as shown in Figure 6a, then the perturbations also induce a spatial modulation of the turbulent fluxes. In adequate circumstances, the convergence/divergence of the modulated fluxes reinforce the original temperature and compositional perturbations, and close the feedback loop. We show in Appendix A that the growth rate of the perturbations, $\Lambda$, is the solution of the following quadratic:

$$\Lambda^2 + \Lambda k^2 \left[ A_2 \left( 1 - R_0 \frac{1}{\gamma_0} \right) + \text{Nu}_0 (1 - A_1 R_0) \right] - A_1 k^4 R_0 \text{Nu}_0 = 0 ,$$

(23)

where

$$A_1 = R_0 \left. \frac{d(1/\gamma^{\text{tot}})}{dR_\rho} \right|_{R_0} , \quad A_2 = R_0 \left. \frac{d\text{Nu}_T}{dR_\rho} \right|_{R_0} ,$$

$$\text{Nu}_0 = \text{Nu}_T(R_0) , \quad \gamma_0 = \gamma^{\text{tot}}(R_0) .$$

(24)

(25)

This quadratic is exactly the same as that of Traxler et al. (2010a) and by proxy that of Radko (2003) provided his $\gamma$ is interpreted as the total flux ratio $\gamma^{\text{tot}}$, see Appendix A for detail. As discussed by Radko (2003) there is a positive real root when $A_1 > 0$, i.e. when $\gamma^{\text{tot}}$ decreases with increasing density ratio, or alternatively, when $\gamma^{-1}$ decreases with $R_0^{-1}$. Exactly as in the case of fingering convection, a necessary condition for layer formation in double-diffusive convection is that $\gamma^{\text{tot}}$ should be a decreasing function of $R_0$. 
Based on the results of Figure 6b, we can now straightforwardly explain, thanks to this theory, the dichotomy between the cases with $R^{-1}_0 > 1.35$, for which we do not expect (and do not observe) layer formation, and the cases with $R^{-1}_0 < 1.35$ for which we do. The case $R^{-1}_0 = 1.35$ is unclear given our measurement errors on $\gamma_{tot}$. Furthermore, since the mode growth rate increases with $A_1$, we can also explain, at least qualitatively, why the staircase forms much more rapidly at lower $R^{-1}_0$ (i.e. because the $\gamma_{tot}$ curve is steeper in that regime). In what follows, we now compare theory and simulations more quantitatively.

5.3. Comparison of the $\gamma$–instability theory with numerical results

The tall-domain simulation described in §5.1 shows the emergence of a four-layer staircase. Based on the discussion of the $\gamma$–instability of §5.2, we need to compare the growth rate of the $k_4$ mode shown in Figure 8 to the solution of (23) with $k = k_4$. In order to do this, we first estimate the coefficients $A_1$ and $A_2$. The derivatives of $Nu_T$ and $\gamma_{tot}$ with respect to $R_\rho$, at $R_\rho = R_0 = 1.2$, are calculated using the fluxes obtained in simulations at neighboring values of the inverse density ratio. We find that

$$A_1 = 0.453, A_2 = 12.9.$$ (26)

We also use the value of $Nu_0$ (and associated errorbar) given in Table 1 for $R^{-1}_0 = 1.2$, and the average of the two $\gamma_{tot}$ values for $\gamma_0$.

We find that the dominant $\gamma$–mode has a dimensionless growth rate

$$\Lambda(k_4) = 9.42 \times 10^{-3}.$$ (27)

Similarly, we find that

$$\Lambda(k_3) = \Lambda(k_4) \frac{k_3^2}{k_4^2} = 5.30 \times 10^{-3}.$$ (28)

Figure 8 compares these theoretical predictions with the numerical results for the $k_4$ and $k_3$ modes. We find that they over-predict the growth rate of the dominant $k_4$ mode, but provide a fairly accurate estimate of the growth rate of the $k_3$ mode.

The fact that the $k_4$ mode grows somewhat slower than predicted is actually expected, and leads us to discuss an important caveat of the $\gamma$–instability theory. The derivation of (23) is fundamentally based on an assumption of scale separation between the small-scale turbulence, which drives the heat and compositional fluxes, and the large-scale temperature and compositional perturbations $\bar{T}$ and $\bar{\mu}$. However, solutions of (23) satisfy the similarity law $\Lambda \propto k^2$, implying that for any given $\gamma$–mode there exists, in theory, another more rapidly
growing one with smaller wavelength. This “ultra-violet catastrophe” problem was discussed by [Radko (2003)], and is clearly an artefact of the model. In practice, the $\gamma$–instability theory should only be applied for $k$ significantly smaller than the wavenumber of the fastest growing mode of the primary instability, in order to satisfy the required separation of scales.

It is therefore interesting to see that the mode which emerges as the dominant $\gamma$–mode in Figure 8 has a wavelength of about $45d$, which is only about twice as large as the wavelength of the most rapidly growing primary gravity wave (about $20d$, see §2.2). As such, it appears to be the smallest-scale mode for which the $\gamma$–instability theory remains applicable, since it still grows exponentially, albeit with a growth rate somewhat slower than predicted. This behavior is very similar to the one found by [Stellmach et al. (2010)] in the case of fingering convection. It is reassuring to see that, by comparison, the larger-scale $k_3$ mode grows as well, and this time with a growth rate which is much closer to the predicted one.

### 5.4. Staircase formation

The dominant $\gamma$–mode grows in amplitude until it causes regularly spaced local inversions in the density profile (see Figure 9). The critical amplitude $\hat{\rho}_{\text{crit}}$, for which the total density gradient $\rho_0 + d\rho/dz = 0$ somewhere in the domain, depends on the mode wavenumber and on the background density ratio (e.g. [Stellmach et al. 2010]):

$$\hat{\rho}_{\text{crit}} = 1 - R_0^{-1} k. \quad (29)$$

As the mode grows beyond this amplitude (shown as a horizontal line in Figure 8), progressively larger regions of the domain are unstable to direct overturning convection, and a fully-formed staircase rapidly appears.

Once layers have formed, the $\gamma$–instability theory no longer applies and the mode stops growing. The subsequent evolution of the staircase through mergers is caused by subtle differences in the fluxes through the interfaces and through the convective regions. While a full study of the merger dynamics is beyond the scope of this paper, we present in the next section an analysis of the heat and compositional fluxes though the staircase as a function of the layer height.
5.5. Layer mergers and flux increase

The time of each merger roughly corresponds, in Figure 8, to the point where the amplitude of the $k_n$ mode overtakes that of the $k_{n+1}$ mode. Shortly afterward, the new dominant mode stops growing, at which point the staircase has reached a new temporary $n-$layered equilibrium. We can then measure the transport properties of the system while it is in this $n-$layered state. The results are presented in Table 2 and illustrated in Figure 7.

Assuming that the sum of the thicknesses of all the interfaces is small compared with the height of the domain, we deduce the layer height $H_L(n) = L_z/n = 178d/n$. We can then construct a Rayleigh number based on the layer height rather than the domain height,

$$Ra_L = \left( \frac{H_L}{d} \right)^4.$$  \hspace{1cm} (30)

Figure 10 shows the variation of $Nu_T$ with $Ra_L$. Rather remarkably, we find that within the errorbars the heat flux through the system is more or less consistent with the standard scaling law for convection between two bounded plates separated by a distance $H_L$ (see Garaud et al. 2010, and references therein),

$$Nu_T \simeq 0.06Ra_L^{1/3}.$$  \hspace{1cm} (31)

Note that no fitting was involved here. This scaling would be consistent with an interpretation of the interfaces as impermeable, diffusive boundary layers.

The compositional Nusselt number is about twice the thermal Nusselt number, in the layered phase and in the homogeneous phase. A possible explanation for this result is the following. The turbulent flux ratio $\gamma = <\tilde{w}\tilde{T}> / <\tilde{w}\tilde{\mu}>$ is typically of order unity for turbulence induced by double-diffusive instabilities (Radko 2003). As a result, using (17) and (18) we have:

$$Nu_\mu \simeq 1 + \frac{R_0}{\tau} <\tilde{w}\tilde{T}> = 1 + \frac{R_0}{\tau} (Nu_T - 1) = 1 + 2.5(Nu_T - 1)$$  \hspace{1cm} (32)

\hspace{1cm} $^4$A plausible reason for this is the following, as argued by Radko (2003). As $R_0^{-1} \to 1$, the compositional field acts more and more like a passive tracer. As a result, the turbulent diffusivities for heat and composition tend to one-another, and since $R_0 = 1$, so do the induced turbulent buoyancy fluxes. As $R_0^{-1} \to (Pr+\tau)/(Pr+1)$, $\gamma$ tends to one for a different reason. Since the instability is driven by the conversion of potential energy into kinetic energy, near the marginal stability limit this available potential energy must vanish. Hence, the turbulent buoyancy flux due to heat must exactly equal that due to composition for the potential energy available to vanish. Finally, since $\gamma$ has to tend to one in both limits, it cannot deviate significantly away from one inbetween.
for our selected parameters. For large enough Nusselt numbers, this implies \( \text{Nu}_\mu \simeq 2.5 \text{Nu}_T \). Figure 10 shows (32) in comparison with the data. Again, the fit is satisfactory within the errorbars, although not quite as compelling as that of \( \text{Nu}_T \). Note that a similar argument was invoked by Traxler et al. (2010a) to explain the relationship between their measured scaling laws for \( \text{Nu}_T \) and \( \text{Nu}_\mu \) in fingering convection.

It is also interesting to compare our numerical results with the work of Spruit (1992), who proposed the following parametrization for heat transport by layered convection:

\[
\text{Nu}_T \simeq 0.5(\text{PrRa}_L)^{1/4}. \tag{33}
\]

His estimate of the turbulent compositional diffusivity (see his equation (44)) yields the following compositional Nusselt number:

\[
\text{Nu}_\mu \propto \tau^{-1/2} \nabla \nabla \mu \text{Nu}_T = R_0 \tau^{-1/2} \text{Nu}_T \tag{34}
\]

in this Boussinesq case, where the proportionality constant is of order unity. Both estimates are shown in Figure 10 as well, and are also more-or-less consistent with our flux measurements within the errorbars. This time, \( \text{Nu}_\mu \) seems to be better accounted for than \( \text{Nu}_T \).

Further simulations will be needed to determine which (if any) of these scaling laws best explains the data. Wider domains will be needed to improve our statistics and reduce the variability of the mean fluxes. In addition, taller domains will be necessary to allow for larger layer heights. But beyond the details of the power laws or the prefactors, our simulations clearly show that the overall transport through a staircase depends only on the layer height, and that the heat and compositional fluxes are roughly proportional to each other for given \( R_0^{-1}, \text{Pr} \) and \( \tau \).

6. Discussion and conclusion

6.1. Summary of the results

In this work, we have studied a set of numerical simulations of double-diffusive convection, in a triply-periodic domain, for Prandtl number \( \text{Pr} = \nu/\kappa_T = 1/3 \) and diffusivity ratio \( \tau = \kappa_\mu/\kappa_T = 1/3 \). We have explored the entire instability range, varying the inverse density ratio \( R_0^{-1} \) between 1 (the onset of direct overturning convection) and \( (\text{Pr} + 1)/(\text{Pr} + \tau) \) (the marginal stability limit). Our simulations were performed in a “small” domain spanning, in the horizontal direction, about five wavelengths of the fastest-growing double-diffusive
mode (i.e. \( L_x = L_y = 100d \) where \( d \) is a thermal diffusion lengthscale), and in the vertical direction, \( L_z = 100d \) or \( L_z = 178d \) depending on the runs.

In all cases we initialized a double-diffusively unstable system with infinitesimal perturbations, and found that these first grow exponentially according to linear theory, then saturate into a state of homogeneous double-diffusive convection. In that state, the turbulent contribution to thermal and compositional transport is significant but much smaller than that expected from standard convection, ranging from 5-10 times the diffusive rate near the onset of direct convective instability, and rapidly dropping towards zero as \( R_0^{-1} \) increases towards marginal stability (see §4).

For small \( R_0^{-1} \), however, the system does not remain in this homogeneously convecting state. Instead, thermo-compositional layers rapidly appear, and transport through the system strongly increases. We showed that the layer formation process is governed by Radko’s \( \gamma \)-instability theory ([Radko 2003; Stellmach et al. 2010; Traxler et al. 2010a]), both qualitatively and quantitatively. In particular, it explains why our simulations with \( R_0^{-1} < 1.35 \) transition into layers while those with \( R_0^{-1} > 1.35 \) do not. The key factor is the variation of the total buoyancy flux ratio \( \gamma_{\text{tot}} \) with density ratio (see §5.2): layers can only form when \( \gamma_{\text{tot}} \) decreases with \( R_0 \).

In the layered phase, we found that the flux through the staircase depends sensitively on the mean layer height \( H_L \). Given the large variability of the measured fluxes during the layered phase, our results are roughly consistent both with Spruit’s theory ([Spruit 1992]) and with heat transport between two solid plates (as given by equation (31)). Further simulations will be needed to help distinguish between these two possibilities – or perhaps suggest an alternative one. Finally, note that in our small-domain simulations, the mergers always proceed until a single layer is left. In that sense, the dynamical evolution of the system is always eventually influenced by the domain size.

6.2. Discussion of the applicability of our results to real systems

Our initial goals were threefold: (a) to characterize transport by homogeneous double-diffusive convection (i.e. in the absence of layers), (b) to determine if, under which conditions, and through which process thermo-compositional layers may form and (c) to characterize transport by layered double-diffusive convection when appropriate.

To answer part (a) in detail, a much larger number of simulations will be needed, using progressively smaller Prandtl numbers and diffusivity ratios. These are the subject of an ongoing investigation. We hope to find similar scaling laws for \( \text{Nu}_T \) and \( \text{Nu}_\mu \) as functions of
Pr and τ as the one found by Traxler et al. (2010b) for fingering convection.

By contrast with the case of fingering convection, however, we now know that thermocompositional staircases can form spontaneously from double-diffusive convection. Radko’s criterion for layer formation, namely that $\gamma_{\text{tot}}$ should decrease with $R_0$, is the answer to part (b) of our goals, but does require knowledge of the function $\gamma_{\text{tot}}(R_\rho)$ to be applied in practice. The latter must be determined separately for each parameter set ($\Pr, \tau$).

Finally, our findings have provided some insight into part (c). Since transport through a staircase depends on the layer height only (for given fluid parameters and overall stratification), the problem shifts to estimating actual layer heights in astrophysical objects. The layers we observe in our simulations have a strong tendency to merge, which suggests two possible outcomes: the mergers continue indefinitely, until the scale of the equilibrium layers is commensurate with the system size; or the mergers stop for other reasons (see below), with an equilibrium layer height significantly smaller than the system size. It is of course crucial to know which of these two scenarios is correct, as they imply vastly different transport rates through the staircase.

Unfortunately, our simulations were not able to provide a definitive answer to this question. In the two cases studied, the final layer height was equal to the domain height, but this should not be interpreted as a result in favor of the first scenario since this could simply mean that our domain was too small to “contain” the intrinsic equilibrium layer height of the second scenario.

Radko (2005) proposed a theory supporting the idea that mergers stop before layers reach the system size, and deduced a means of estimating the equilibrium layer height. Starting from an initial staircase with uniform “jumps” in temperature and chemical composition across the interfaces, he studied how the staircase evolves if it is perturbed slightly, by making some of the jumps larger and some of the jumps weaker. He concluded that the staircase is unstable to mergers if the total buoyancy flux ratio through the interfaces is a decreasing function of the density ratio across the interfaces – a criterion very similar to the $\gamma$–instability criterion.

We have tried to test Radko’s merger theory against our simulations, but this has unfortunately proven to be difficult. The statistical fluctuations in the measured turbulent fluxes (see Figure 7 for example) are too large to detect a significant variation of the interfacial flux ratio as the mergers proceed. We would need a much larger domain to improve the signal to “noise” ratio to a point where our results could be compared with his theory. We would also need a much taller domain (at least a few times taller than the equilibrium layer height) to see if the merger process indeed stops as expected. Finally, we would need to integrate
the simulation long enough to establish convincingly that the mergers have indeed stopped. Unfortunately, running equivalent simulations in a much larger domain, and for long enough to observe the layer formation and merger process, is impossible within current numerical limitations.

6.3. Future prospects

The preliminary findings presented in this paper still enable us to lay out a clear path towards obtaining better parametrizations of mixing by double-diffusive convection in the near future, using currently available computational resources:

- Firstly, we must gain a better understanding of the instability saturation mechanism at low Prandtl number and low diffusivity ratio, in order to determine the flux laws $\text{Nu}_T(\text{Pr}, \tau, R_0)$ and $\text{Nu}_\mu(\text{Pr}, \tau, R_0)$ for homogeneous double-diffusive convection. These flux laws are needed to determine when layers are expected to form, and can be used “as is” to parametrize mixing otherwise. They can be measured using “small-domain” simulations similar to the ones we have presented here, at least for values of $\text{Pr}$ and $\tau$ as low as about 0.01 or so. Semi-analytical weakly-nonlinear models will then be helpful to guide extrapolations to the much lower parameter values appropriate of the astrophysical regime.

- Secondly, we must gain a general understanding of mixing in the layered case, at low Prandtl number and low diffusivity ratio. In order to do this, we need to determine how interfacial transport depends on the fluid parameters ($\text{Pr}, \tau$) and on the interfacial density ratio (i.e. a density ratio based on the difference in temperature and composition across the layers). We must also understand how transport scales within the convective layers, as a function of the same parameters but also as a function of the layer height. This can be done today using simulations in which a single layer is pre-seeded, to bypass the rather lengthy layer formation and merger phases. Using this information, we will be able to test the basic flux laws which are central to Radko’s merger theory more quantitatively \cite{Radko2005}. If this theory holds, then one can straightforwardly deduce the equilibrium layer height for a given parameter set, and ultimately quantify the staircase transport properties.

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A. Derivation of the $\gamma$–instability in the diffusive case.

Following Radko (2003) and Traxler et al. (2010a), we begin with the general non-dimensional governing equations (10), and average them over several wavelengths of the fastest growing mode of the primary instability. We get

$$\frac{1}{Pr} \left( \frac{\partial u}{\partial t} + u \cdot \nabla u \right) = -\nabla p + (T - \mu)e_z + \nabla^2 u - \frac{1}{Pr} \nabla \cdot R, \quad (A1)$$

$$\frac{\partial T}{\partial t} - w + u \cdot \nabla T = -\nabla \cdot F_T^{tot}, \quad (A2)$$

$$\frac{\partial \mu}{\partial t} - \frac{1}{R_0} w + u \cdot \nabla \mu = -\nabla \cdot F_\mu^{tot}, \quad (A3)$$

where $R$ is the Reynolds stress, $F_T^{tot}$ and $F_\mu^{tot}$ are the total heat and compositional fluxes respectively, and $T$, $\mu$, $p$, and $u$ now denote large-scale fields only.

The $\gamma$–instability drives horizontally-invariant perturbation with zero mean flow (Radko 2003). We can therefore neglect the momentum equation, set $u = 0$, and ignore all horizontal derivatives. The mean temperature and composition equations simplify to:

$$\frac{\partial T}{\partial t} = -\frac{\partial F_T^{tot}}{\partial z}, \quad (A4)$$

$$\frac{\partial \mu}{\partial t} = -\frac{\partial F_\mu^{tot}}{\partial z}. \quad (A4)$$

Finally, we assume that $Nu_T$ and $\gamma_{tot}$ depend only on the local value of the density ratio $R_\rho$. Note that $R_\rho$ is no longer constant, but varies with $z$ as a result of the large-scale background temperature and compositional perturbations, as

$$R_\rho = \frac{\alpha(T_{0z} + T_{\dim}^{z})}{\beta(\mu_{0z} + \mu_{\dim}^{z})} = \frac{R_0(1 - T_z)}{1 - R_0 \mu_z}, \quad (A5)$$

where, for clarity, we first expressed $R_\rho$ as the ratio of dimensional quantities and then as the ratio of non-dimensional quantities.

We now linearize equations (A2) and (A3) around a state of homogeneous turbulent convection in which $T = 0 + T'$, $\mu = 0 + \mu'$, and $R_\rho = R_0 + R'$ where linearization of (A5) yields:

$$R' = R_0(1 - T_z + R_0 \mu_z). \quad (A6)$$
Using the fact that $F_T^{\mathrm{tot}} = \mathrm{Nu}_T(R_\rho)(1 - T_z)$ and $F_\mu^{\mathrm{tot}} = F_T^{\mathrm{tot}}/\gamma_{\mathrm{tot}}$ the linearized temperature equation becomes

$$
\frac{\partial T'}{\partial t} = A_2 \left( \frac{\partial^2 T'}{\partial z^2} - R_0 \frac{\partial^2 \mu'}{\partial z^2} \right) + \mathrm{Nu}_0 \frac{\partial^2 T'}{\partial z^2}, \quad (A7)
$$

while the linearized composition equation is

$$
\frac{\partial \mu'}{\partial t} = \frac{1}{\gamma_0} \frac{\partial T'}{\partial t} + A_1 \left( \frac{\partial^2 T'}{\partial z^2} - R_0 \mathrm{Nu}_0 \frac{\partial^2 \mu'}{\partial z^2} \right), \quad (A8)
$$

where we have used the following notation for simplicity:

$$
A_1 = R_0 \frac{d(1/\gamma_{\mathrm{tot}})}{dR_\rho} \bigg|_{R_0}, \quad A_2 = R_0 \frac{d\mathrm{Nu}_T}{dR_\rho} \bigg|_{R_0}, \quad \mathrm{Nu}_0 = \mathrm{Nu}_T(R_0), \quad \gamma_0 = \gamma_{\mathrm{tot}}(R_0).
$$

Assuming normal modes of the form $T', \mu' \sim e^{ikz + \Lambda t}$, we get the quadratic

$$
\Lambda^2 + \Lambda k^2 \left[ A_2 \left( 1 - \frac{R_0}{\gamma_{\mathrm{tot}}} \right) + \mathrm{Nu}_0(1 - A_1 R_0) \right] - A_1 k^4 R_0 \mathrm{Nu}_0 = 0. \quad (A9)
$$

This quadratic is exactly the same as the one obtained in the fingering case. In hindsight, this result is trivial, and could be obtained immediately had we allowed ourselves to non-dimensionalize $T$ and $\mu$ using negative dimensions (in which case the governing equations and all definitions are exactly the same as in the fingering case).

**REFERENCES**


Biermann, L. 1932, ZAp, 5, 117


Kato, S. 1966, PASJ, 18, 374
Radko, T. 2003, J. Fluid Mech., 497, 365
—. 2005, J. Fluid Mech., 523, 79
Stern, M. 1960, Tellus, 12, 172
—. 1985, Icarus, 62, 4
Toole, J., et al. 2006, EOS Transactions, 87, 434


Traxler, A., Stellmach, S., Garaud, P., Radko, T., & Brummell, N. 2010b, ArXiv e-prints

Turner, J. 1985, Annu. Rev. Fluid Mech., 17, 11

Turner, J. S. 1968, Journal of Fluid Mechanics, 33, 183


Walin, G. 1964, Tellus, 16, 389
Fig. 2.— Summary of the various regimes of convective instability. The top line corresponds to the chemically homogeneous case, with $\nabla \mu = 0$. The system is unstable to direct overturning convection if $\nabla > \nabla_{\text{ad}}$ or equivalently $T_{0z} < T_{0z}^{\text{ad}} < 0$. The Schwarzchild criterion (dotted line) appropriately marks the stability boundary (solid line). In the presence of an unstable mean molecular weight gradient (middle line), the region unstable to convective overturning extends into the subadiabatic regime, and is stabilized only when $\nabla - \nabla_{\text{ad}} = \nabla_{\mu}$ (Ledoux criterion, vertical dashed line). The Schwarzchild criterion in this regime is not relevant. Beyond the Ledoux limit, the system can still be unstable, this time to fingering convection. In the case where the system has a stable mean molecular weight gradient (bottom line), the region of parameter space unstable to overturning convection shrinks according to the Ledoux criterion. The system can still be unstable to double-diffusive convection in a subset of the interval between Ledoux-stability and Schwarzchild stability.
Fig. 3.— Temporal evolution of the non-dimensional rms velocity in simulations with \( R_0^{-1} = 1.2 \), for two different computational domain heights: \( L_z = 100 \) corresponding to \( Ra_T = 10^8 \) (dashed line), and \( L_z = 178 \) corresponding to \( Ra_T = 10^9 \) (solid line). The straight dotted line shows an estimate of the early exponential growth based on the growth rate of the most rapidly growing mode only (see §2.2). The two simulations saturate at the same level, confirming that the dynamics of the system in the saturated phase are independent of the domain size (for large enough domains).

Table 2: Variation of the Nusselt numbers as a function of the number of layers, for the \( R_0^{-1} = 1.2, L_z = 178d \) run.

<table>
<thead>
<tr>
<th>( n )</th>
<th>( H_L )</th>
<th>( t_{\text{start}} )</th>
<th>( t_{\text{end}} )</th>
<th>( Nu_T )</th>
<th>( Nu_\mu )</th>
</tr>
</thead>
<tbody>
<tr>
<td>4</td>
<td>44.5</td>
<td>1020</td>
<td>1240</td>
<td>10.5 ± 1.7</td>
<td>20.6 ± 4.1</td>
</tr>
<tr>
<td>3</td>
<td>59.3</td>
<td>1240</td>
<td>1480</td>
<td>16.4 ± 2.9</td>
<td>33.5 ± 8.4</td>
</tr>
<tr>
<td>2</td>
<td>88.9</td>
<td>1480</td>
<td>1660</td>
<td>25.8 ± 5.0</td>
<td>55.0 ± 12.6</td>
</tr>
<tr>
<td>1</td>
<td>177.8</td>
<td>1660</td>
<td>2190</td>
<td>46.0 ± 14.5</td>
<td>99.0 ± 33.3</td>
</tr>
</tbody>
</table>

Note. — The first column shows the number of layers, and the second column an estimate of the layer depth in units of \( d, H_L = L_z / n \). The times \( t_{\text{start}} \) and \( t_{\text{end}} \) mark the interval of the time over which the turbulent fluxes, measured via \( Nu_T \) and \( Nu_\mu \), were averaged. Errorbars represent the rms fluctuations about the mean.
Fig. 4.— Evolution of the thermal Nusselt number for seven of the simulations presented in Table 1. In all cases, Pr = τ = 1/3, Ra_T = 10^8 (L_z = 100), and the aspect ratio is one. Nu_T − 1 is shown to emphasize the exponential growth phase. The results are also staggered for clarity, so each curve actually shows f(Nu_T − 1) where the multiplicative factor f is 1, 10, 100, 1000, 10^4, 10^5, and 10^6 respectively for R_0^{-1} = 1.85, 1.6, 1.5, 1.35, 1.2, 1.15 and 1.1. A straight horizontal line of the same color in each case marks the point at which Nu_T − 1 = 1 for reference (i.e. when turbulent and diffusive fluxes are equal to one another). Note how runs with R_0^{-1} ≥ 1.35 remain in a quasi-steady saturated state, while runs with R_0^{-1} < 1.35 show a subsequent increase in transport. The R_0^{-1} = 1.35 run was actually integrated until t = 2500, but was found to remain at the same saturated level.
Fig. 5.— Volume-rendered visualization of the mean molecular weight perturbation, for $R_0^{-1} = 1.2$, using the tall-domain simulation ($L_z = 178d$). Shown are five snapshots taken at different times, (a) in the homogeneous phase at $t = 400$, (b) in the four-layer phase at $t = 1100$, (c) three-layer phase at $t = 1350$, (d) two-layer phase at $t = 1550$ and (e) single-layer phase at $t = 1850$. The color scale is adjusted in each panel to emphasize the perturbations, so that $\tilde{\mu} \in [-0.1, 0.1] \mu_0 L_z$ in (a), $\tilde{\mu} \in [-0.25, 0.25] \mu_0 L_z$ in (b) and (c), $\tilde{\mu} \in [-0.4, 0.4] \mu_0 L_z$ in (d) and $\tilde{\mu} \in [-0.5, 0.5] \mu_0 L_z$ in (e).
Fig. 6.— Left: Mean Nusselt numbers in the homogeneously turbulent phase (see Table 2), as a function of the inverse density ratio. The Nu$_T$ measurements are shown as (+) symbols, and the Nu$_\mu$ measurements as (×) symbols. The turbulent contribution to transport, represented by Nu − 1 in both cases, goes from a few to zero over the instability range. The $R_0^{-1} = 2.1$ point was left out since we measured Nu$_T$ = Nu$_\mu$ = 1 in that run. Right: Inverse of the total buoyancy flux ratio $\gamma_{\text{tot}}^{-1}(R_0^{-1})$ as measured in our simulations, using two different methods (see main text for detail). Note how $\gamma_{\text{tot}}^{-1}$ shows a pronounced minimum around $R_0^{-1} = 1.4$. In all cases, the errorbars represent rms fluctuations of the respective functions around the mean.
Fig. 7.— Evolution of $\text{Nu}_T$ (solid line) and $\text{Nu}_\mu$ (dashed line) for the $R_0^{-1} = 1.2$ tall domain run ($L_z = 178$). This plot shows the step-wise increase in transport through the various phases. In the layered phase in particular, the heat transport through the staircase depends on the layer height. The horizontal lines indicate the mean compositional Nusselt number measured in the four-, three-, two- and one-layer state in the tall-domain run (see §5.5 and Table 2 for detail). The solid part of each line indicates the interval of time over which the averages were measured. Equivalent lines for $\text{Nu}_T$ are left out to avoid cluttering the plot. Note how closely the two curves follow each other throughout the entire run.
Fig. 8.— Evolution of the norm of the Fourier coefficient of each of the four gravest Fourier modes of the vertical density perturbation profile, $|\hat{\rho}_{0,0,k}|^2 = \hat{\rho}_{0,0,k}\hat{\rho}_{0,0,k}^*$. The figure illustrates the initial exponential growth of the $k_4$ and $k_3$ modes, and compares them with the prediction from the $\gamma$–instability theory (see §5.2), shown as the same-color straight lines. The mode grows until it reaches the critical amplitude for overturning (horizontal black line), see main text for detail. Shortly afterward, four equally spaced layers appear (around $t = 1000$). The layers later merge successively, which can be seen here easily as the $k_3$, $k_2$ and $k_1$ mode respectively take over as the dominant mode.
Fig. 9.— Evolution of the density profile of a $k_{4}$ $\gamma$–mode. For simplicity, the density is normalized to the density difference across the domain height, and the vertical coordinate is in units of the domain height. The dotted line shows the background density profile with constant gradient $\rho_{0z}$, the dashed line the background + perturbation for intermediate perturbation amplitude, while the solid line shows the background + perturbation at the critical amplitude for the onset of overturning convection given by (29). Note the existence of four specific points where the total density gradient $\rho_{0z} + d\rho/dz$ is zero, which will become the middle of the emergent layers.
Fig. 10.— Variation of the two Nusselt numbers in the layered phase, for the tall-domain $R_0^{-1} = 1.2$ run, with the Rayleigh number defined with layer height, $Ra_L$. The symbols correspond to the measurements presented in Table 2, with $Nu_T$ shown as (+) symbols and $Nu_\mu$ shown as ($\times$) symbols. The errorbars show the standard deviation of the measured fluxes about the mean. The various lines correspond to different possible scaling laws discussed in the main text: the blue lines are from Spruit (1992), for $Nu_T$ (solid line) and $Nu_\mu$ (dashed line), and the green lines are from (31) for $Nu_T$ (solid line) and (32) for $Nu_\mu$ (dashed line).