

Tidal Heating via Internal Gravity Waves in White Dwarf Binaries

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Accepted XXX. Received YYY; in original form ZZZ

ABSTRACT

In white dwarf binaries tidal torques are expected to raise a train of internal gravity waves in the white dwarf interior. These outward-propagating waves can then grow to be nonlinear and break via hydrodynamical instabilities, transferring energy and angular momentum from the binary orbit to the white dwarf. We perform 2D numerical simulations of nonlinear wave breaking of outwards-propagating internal gravity waves in an incompressible, isothermal atmosphere with an exponentially decaying density stratification. We find that tidal synchronization, after an initial transient phase, proceeds inwards from the surface in thin layers. We argue the thickness of these layers is limited by the Kelvin-Helmholtz Instability. We provide simple analytical formulae for the location and thickness of tidal synchronization over time that are in good agreement with our simulations.

Key words: Wave Breaking, Dynamical Tides, Internal Gravity Waves

1 INTRODUCTION

Compact white dwarf (WD) binary systems, with orbital periods in the range of minutes to hours, are important for a range of astrophysical problems. They are the most important sources of gravitational waves (GWs) for the Laser Interferometric Space Antenna (LISA) (Nelemans (2009)). They are also thought to produce interesting optical transients such as underluminous supernovae (Perets et al. (2010)), Ca-rich fast transients (García-Berro et al. (2017)), and tidal novae (Fuller & Lai (2012b)). Most importantly, they have been proposed as the likely progenitors of type Ia supernovae (e.g. Iben Jr & Tutukov (1984); Webbink (1984) or more recently Gilfanov & Bogdán (2010); Maoz et al. (2010)). While presently only a few tens of compact WD binaries are known (Korol et al. (2017)), *Gaia* (currently gathering data) is expected to expand the catalog to a few hundreds (Korol et al. (2017)) (results based on *Gaia*’s second data release have already begun to appear (Shen et al. (2018); Kilic et al. (2018))), and the Large Synoptic Survey Telescope (LSST, first light scheduled for 2020) will likely detect a few thousand more (Korol et al. (2017)). These observations will significantly advance the understanding of WD binaries and their evolution. My proposed theoretical and computational research is well-timed to take advantage of these new advances.

In spite of the broad importance of WD binaries, the evolution of these systems prior to their final mergers is not well understood. Much of this uncertainty comes from our imprecise understanding of tidal interactions, which play an important role during a compact WD binary’s inspiral (Fuller & Lai (2012a)). Previous studies have shown that these interactions manifest as tidal excitation of internal gravity waves (IGW), waves in the WD fluid restored by the buoyancy force due to density stratification (Fuller & Lai (2011)). As these waves propagate outwards towards the WD surface, they grow in amplitude until they break, as do ocean waves on a shore, and

transfer both energy and angular momentum from the binary orbit to the outer envelope of the WD (Fuller & Lai (2011, 2012a)).

Previous works have found that the dissipation of IGW can generate significantly more energy than thermal radiation from the isolated WD surface and is thus a major contributor to the WD energy budget (Fuller & Lai (2012a, 2013)). However, these works parameterized the wave breaking process in an ad hoc manner. The details of dissipation, namely the location and spatial extent of the wave breaking, affect the observable outcome: dissipation near the surface of the WD can be efficiently radiated away and simply brightens the WD, while dissipation deep in the WD envelope causes an energy buildup that results in energetic flares (Fuller & Lai (2012b)). Works in other fields based on numerical simulations show that strongly nonlinear wave breaking behaves differently than predictions based in linear and weakly nonlinear theory (Winters & D’Asaro (1994); Barker & Ogilvie (2010)). Such fully nonlinear numerical simulations have not been performed for WDs.

In section 2, we will describe the system of equations we will use to analyze IGW breaking. In section 3, we discuss relevant analytical results. In section 4 we present the results of numerical simulations. Finally, in section 5 we discuss the results of the preceding section.

2 PROBLEM DESCRIPTION

We consider a two-dimensional incompressible, isothermal fluid, representative of degenerate matter in WDs. We neglect temperature variations and assume a barotropic equation of state as a first approximation. As we are interested in dynamics far from the center of the WD, we approximate the gravitational field as uniform. We model the background density stratification as

$\rho_0(x, z) = \rho_0(z) = \rho_0(z=0)e^{-z/H}$ for some reference density $\rho_0(z=0) = \rho_0(z)|_{z=0}$.

The Euler equations for an incompressible, barotropic fluid in a uniform gravitational field are

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (1a)$$

$$\frac{D\rho}{Dt} = 0, \quad (1b)$$

$$\frac{D\vec{u}}{Dt} + \frac{\vec{\nabla}P}{\rho} + g\hat{z} = 0. \quad (1c)$$

$\frac{D}{Dt} = \frac{\partial}{\partial t} + (*)\vec{u} \cdot \vec{\nabla}$ is the Lagrangian or material derivative. \vec{u}, ρ, P denote the velocity field, density and pressure respectively. We denote $-g\hat{z}$ constant gravitational acceleration. Note that at hydrostatic equilibrium $\frac{\partial}{\partial t} = 0$ we have $\vec{\nabla}P_0 = -\rho_0 g\hat{z}$ and so $P_0 = \rho_0 gH$. A shear flow $u_x(x, z, t) = u_x(z)$ is permitted in hydrostatic equilibrium, but we will initially consider no shear flow in equilibrium.

In practice, it is convenient to introduce coordinate $\Upsilon = \ln \frac{\rho}{\rho_0}$. This both identically enforces $\rho > 0$ and avoids numerical issues if ρ is small. We also define reduced pressure $P = \frac{P}{\rho}$. Then, we may rewrite the second two equations of Equation 1 as

$$\frac{D\Upsilon}{Dt} + u_z \frac{\partial \ln \rho_0}{\partial z} = 0, \quad (2a)$$

$$\frac{D\vec{u}}{Dt} + \vec{\nabla}P + P\vec{\nabla}\Upsilon + g\hat{z} = 0. \quad (2b)$$

Note that in the new coordinates, hydrostatic equilibrium corresponds to $\Upsilon_0 = 1, P_0 = gH$.

3 INTERNAL GRAVITY WAVE THEORY

3.1 Analytical Properties: Linear

In the small perturbation limit, where flow velocities are small compared to the characteristic space and time scales $\frac{\partial}{\partial t} \gg \vec{u} \cdot \vec{\nabla}$, we may linearize Equation 2. We ignore the advective components of the material derivative and consider small deviations $\Upsilon_1 = \Upsilon - \Upsilon_0, P_1 = P - P_0$ about hydrostatic equilibrium:

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (3a)$$

$$\frac{\partial \Upsilon_1}{\partial t} - \frac{u_z}{H} = 0, \quad (3b)$$

$$\frac{\partial \vec{u}}{\partial t} + \vec{\nabla}P_1 + (*)1 + \Upsilon_1 g\hat{z} = 0. \quad (3c)$$

This can be solved to obtain solutions up to undetermined amplitude A the well-known result (Drazin (1977); Dosser & Sutherland (2011b)):

$$u_z(*)x, z, t = Ae^{z/2H} e^{i(k_{0x}x + k_{0z}z - \omega_0 t)}, \quad (4a)$$

$$\omega_0^2 = \frac{N^2 k_{0x}^2}{k_{0x}^2 + k_{0z}^2 + \frac{1}{4H^2}}, \quad (4b)$$

where

$$N^2 \equiv g^2(*) \frac{d\rho}{dP} - \frac{\partial \rho}{\partial P} \Big|_{ad} = \frac{g}{H}, \quad (5)$$

the Brunt-Väisälä frequency is constant

In the weak stratification limit $k_{0z}H \gg 1$, the solution exhibits the following characteristics:

- The amplitude of the wave grows like $e^{z/2H}$. Thus, the linear approximation is violated for sufficiently large z .

- The phase and group velocities can be computed respectively:

$$\vec{c}_{ph} = (*) \frac{\hat{x}}{k_{0x}} + \frac{\hat{z}}{k_{0z}} \omega, \quad (6)$$

$$\begin{aligned} \vec{c}_g &= \frac{N(*)k_{0x}k_{0z}\hat{x} - (*)k_{0z}^2 + \frac{1}{4H^2}\hat{z}}{(*)k_{0x}^2 + k_{0z}^2 + \frac{1}{4H^2}} \\ &\approx \frac{\omega k_{0z}}{|*|\vec{k}_0^{3/2}} (*)k_{0x}\hat{x} - k_{0z}\hat{z}. \end{aligned} \quad (7)$$

We recover the usual result $\vec{c}_{ph} \cdot \vec{c}_g = 0$ (e.g. Drazin (1977); Dosser & Sutherland (2011a)). Note that an IGW transporting energy and momentum upwards $c_{g,z} > 0$ has $k_{0z} < 0$.

- The time-averaged total x -momentum flux in the \hat{z} direction can be computed. Since the linear solution is separable as $f(x, z, t) = f(z)e^{i(k_{0x}x - \omega_0 t)}$, x averaging and time averaging are equivalent, so we may write

$$\begin{aligned} S_{px,z} &\equiv \langle (*) \rho u_x u_{z,x} \rangle \equiv \frac{1}{L_x} \int_0^{L_x} \rho u_x u_z dx \\ &\approx -\frac{A^2}{2} \rho_0(z=0) \frac{k_{0z}}{k_{0x}}, \end{aligned} \quad (8)$$

where $\langle (*) \dots \rangle_x$ denotes x -averaging.

3.2 Wave Generation

To model a continuous IGW wave train excited deep in the WD interior propagating towards the surface, we use a volumetric forcing term to excite IGW near the bottom of the simulation domain¹. Our forcing excites both IGWs propagating upwards, imitating a wave tidally excited deeper in the WD, and downwards, which are damped away by the damping layers described in subsection 4.2.

As not to interfere with the incompressibility constraint, we force the system on the density equation. This constitutes replacing Equation 2a with

$$\frac{D\Upsilon}{Dt} + u_z \frac{\partial \ln \rho_0}{\partial z} = F e^{-\frac{(z-z_0)^2}{2\sigma^2}} \cos(*)k_{0x}x - \omega_0 t. \quad (9)$$

Using a narrow Gaussian profile excites a broad k_z wavenumber spectrum, and only the k_{0z} satisfying dispersion relation Equation 4b for the given $k_{0x}, \omega_0(k_{0x}, k_{0z})$ will propagate.

In the linearized system, the effect of this forcing can be solved analytically up to good accuracy: we first approximate the driving term using $e^{-\frac{(z-z_0)^2}{2\sigma^2}} \approx \sqrt{2\pi\sigma^2} \delta(z - z_0)$, the $\sigma \rightarrow 0$ limit². This system is solved exactly by matching the two homogeneous solutions above and below z_0 . We may then approximately relax the solution to nonzero σ : an extra factor of $e^{-\frac{(k_{0z}\sigma)^2}{2}}$ arises compared to the δ -function solution (evaluating the Fourier Transform of $e^{-\frac{(z-z_0)^2}{2\sigma^2}}$

¹ Interfacial forcing at the bottom boundary incurs strong stability limits, as we use a Chebyshev polynomial basis along the z axis in our spectral method which has very small grid spacing at the boundaries.

² In practice, σ must be large enough to be numerically resolved by the spectral code.

at $k_z = k_{0z}$), and we obtain

$$u_{1z}(x, z, t) = \frac{Fgk_{0x}^2}{\omega_0^2} \frac{1}{2ik_{0z}} \frac{e^{-\frac{(k_{0z}\sigma)^2}{2}}}{\sqrt{2\pi\sigma^2}} \times \begin{cases} e^{\frac{z-z_0}{2H}} e^{i(*)k_{0x}x + k_{0z}(z-z_0) - \omega_0 t + \frac{1}{2k_{0z}H}} & z > z_0 \\ e^{\frac{z-z_0}{2H}} e^{i(*)k_{0x}x - k_{0z}(z-z_0) - \omega_0 t + \frac{1}{2k_{0z}H}} & z < z_0 \end{cases}. \quad (10)$$

The $z > z_0$ solution models an upwards-propagating IGW wavetrain inbound on the simulation domain from below. The solution for $u_{1x}(x, z, t)$ is correspondingly $u_{1x} \approx \frac{k_{0z}u_{1z}}{k_{0x}} e^{\frac{i}{2k_{0z}H}}$.

3.3 Wave Breaking

In developing the results above, we have neglected advective terms $\frac{\partial}{\partial t} \gg \vec{u} \cdot \vec{\nabla}$. However, since $|*|\vec{u}| \propto e^{z/2H}$, any nonzero amplitude IGW will eventually violate the linearity criterion. This regime is relevant in practice since the IGW may grow to nonlinear amplitudes before it reaches the surface of the WD (Fuller & Lai (2011, 2012a)). We can coarsely estimate the height at which this happens by when the Lagrangian displacement $\vec{\xi}$ of a fluid parcel satisfies

$$\vec{\xi} \cdot \vec{k} \gtrsim 1. \quad (11)$$

The general understanding of nonlinear IGW breaking is that the waves' interactions with the mean flow of the fluid transfer horizontal momentum from the waves into the fluid; such a process has been conjectured to be responsible for tidal synchronization in stellar binaries (Zahn (1975); Goldreich & Nicholson (1989)) as well as the quasi-biennial oscillation (Lindzen & Holton (1968)). The details of this process have been laid out in a few key papers: IGWs are globally unstable to resonant three-wave interactions (Drazin (1977)), causing energy transfer out of the IGW to daughter modes. These steeper daughter modes facilitate wave breaking, depositing horizontal momentum in the fluid's mean flow (Klostermeyer (1991)). Once the mean flow reaches the horizontal phase velocity of the parent IGW, a critical layer forms at which the frequency of the IGW is Doppler-shifted to zero. The interaction of the parent IGW with the mean flow was first studied in the inviscid, linear regime in Booker & Bretherton (1967), which found nearly complete absorption. This result was reproduced with nonzero viscosity (Hazel (1967)), but weakly nonlinear theory (Brown & Stewartson (1982)) and fully nonlinear simulations (Winters & D'Asaro (1994)) suggest that nonlinear effects can induce reflection.

An estimate of where these mean flow effects become important can be made using the wave-induced mean flow (Andrews & McIntyre (1976); Dosser & Sutherland (2011b)):

$$\langle * \rangle u_x = \bar{U}_0 \equiv \frac{1}{L_x} \int u_x dx = \frac{\langle * \rangle u_x u_{zx}}{c_{g,z}}, \quad (12)$$

where $\langle * \rangle \dots_x$ denotes horizontal averaging. Setting $\bar{U}_0 = c_{ph,x}$ the critical layer condition, we recover $\frac{u_x k_x}{\omega} = \xi_x k_x \sim 1$, roughly equivalent to the nonlinearity criterion (Equation 11).

4 INTERNAL GRAVITY WAVE NUMERICAL SIMULATION

To study IGW breaking as a weak perturbation grows to nonlinear amplitudes, we perform direct numerical simulation using Dedalus (Burns et al. (2016)) a spectral numerical method. Our simulation domain is smaller than our full physical domain, so we

adopt reflection-suppressing boundary conditions. We use periodic boundary conditions in the x direction and damping layers (described in subsection 4.2) in the z direction to damp perturbations exiting the top and bottom of the domain. We use a Fourier basis in the x direction and either a Chebyshev or Fourier basis in the z direction; our results were robust to either set of basis functions. We varied the number of x, z modes (denoted N_x, N_z respectively), and used 3/2 dealiasing (Boyd (2001)).

The time integration uses a split implicit-explicit first-order scheme where certain terms are treated implicitly using forward Euler, and the remaining terms are treated explicitly using backward Euler. A third-order, four-stage DIRK-ERK scheme (Ascher et al. (1997)) is used with adaptive timesteps taken to be the advective Courant-Friedrichs-Lewy (CFL) time given by $0.7 \min(\Delta x/u_x, \Delta z/u_z)$. The minimum is taken over every grid point in the domain and $\Delta x, \Delta z$ are the grid spacings in the x and z directions respectively.

4.1 Numerical Parameters

We nondimensionalize by taking $H = N = \rho_0(z = 0) = 1$ in Equation 1.

Denote L_x, L_z to be the physical dimensions of the simulation domain. The physics of our simulation is fixed by six parameters: L_x, L_z , and k_{0x}, ω_0, F, ν . We describe and choose these parameters below:

- L_z : Since the wave amplitude grows $\sim e^{z/2H}$, we choose $L_z = 10H$ to give $\sim e^4$ amplitude growth between the damping zones.
- L_x : We want similar grid spacing $\frac{L_x}{N_x} \sim \frac{L_z}{N_z}$. For computational savings, we fix $N_z = 4N_x$. Expecting wave breaking to be mostly in the z direction, we use $L_x = 4H$.
- k_{0x} : Astrophysical IGWs in stars are generally excited by the $l = 2$ component of the tidal potential, for which $k_\perp \sim \frac{1}{R}$ where R is the radius of the star. To best emulate this, we use $k_{0x} = \frac{2\pi}{L_x}$ the smallest permitted wavenumber permitted by periodic boundary conditions.
- ω_0 : We choose ω_0 by evaluating dispersion relation $\omega_0(*)k_{0x}, k_{0z}$ for some desired k_{0z} .
- Astrophysical IGWs also generally satisfy $\omega \ll N$, or equivalently $\frac{k_z}{k_\perp} \sim k_r R \gg 1$. Since $H \lesssim R$, we aim to study $k_{0z}H \gg 1$. However, to satisfy this well is computationally expensive: $L_z \gg H$ is required to give waves ample room to grow within the simulation domain, but we want k_{0z} to be sufficiently separated from the grid spacing $\sim L_z/N_z$ that nonlinear effects can be seen in later simulations (see subsection 4.5). Thus, N_z sets how well we can satisfy $k_{0z}H \gg 1$; we choose $k_{0z}H = 1$.
- F : We first choose F forcing strength such that $\vec{\xi} \cdot \vec{k} \ll 1$ is satisfied everywhere throughout the domain. This becomes a constraint on F by using Equation 10 and $|*|\vec{u}| \propto e^{z/2H}$.
- ν : Nonlinear effects transfer wave energy from \vec{k}_0 to smaller scales. Since well-resolved simulations using spectral methods have no inherent numerical viscosity, energy will accumulate at grid scales in the absence of artificial dissipation. We introduce Navier-Stokes viscosity ν while ensuring our equations of motion continue to conserve horizontal momentum (see section A for details). When $\vec{\xi} \cdot \vec{k} \ll 1$, we may set $\nu = 0$.

4.2 Damping Layers

To imitate an infinite fluid using a finite simulation domain, we use periodic boundary conditions in the x direction and damping layers

at the top and bottom of the z direction. This damps waves that reach the edge of the simulation domain without inducing nonphysical reflection. Specifically, these are implemented by introducing linear damping terms to Equation 2:

$$\begin{aligned} \frac{D\Upsilon}{Dt} + u_z \frac{\partial \ln \rho_0}{\partial z} &= -\Gamma(z)\Upsilon, \\ \frac{D\vec{u}}{Dt} + \vec{\nabla}P + P\vec{\nabla}\Upsilon + g\hat{z} &= -\Gamma(z)\vec{u}, \\ \frac{1}{2\tau} [*] 2 + \tanh \frac{z - z_T}{\Delta z} + \tanh \frac{z_B - z}{\Delta z} &= \Gamma(z), \end{aligned} \quad (13)$$

where $z_B = 0.05L_z$, $z_T = 0.95L_z$ are the boundaries of the damping zone. This strongly damps perturbations below z_B and above z_T with damping time τ , negligibly affects dynamics between z_B , z_T and has transition width governed by Δz ; we use $\Delta z = 0.025L_z$. This prescription, used in similar studies (Lecoanet et al. 2016), has the advantage of being smooth, important for spectral methods. The full implementation of the fluid equations in Dedalus is described in section A.

4.3 Linear Regime

We first test our numerical simulation with a “linear” simulation, where F is chosen such that $\vec{\xi} \cdot \vec{k} \ll 1$ everywhere. We emphasize that our “linear” simulations solve the full nonlinear equations, just in the weak perturbation limit.

4.4 Mean Flow Critical Layer Absorption

A Navier-Stokes numerical viscosity ν is used to damp high wavenumbers and regularize the nonlinear cascade at near grid resolution: $\nu \sim 0.1 \frac{\omega_0}{|k_{0z}|} \frac{L_z}{2\pi N_z}$ was found to be suitable for $N_z = 2048$.

In past studies of IGWs in WDs (cite Fuller & Lai), $\xi_z \equiv \frac{u_z}{\omega_0}$ the Lagrangian fluid displacement was often used towards wave breaking criterion $k_{0z}\xi_z \gtrsim 1$. We argue that the wave’s self-interaction via its generated mean flow \bar{U}_0 induces total absorption when the mean flow exceeds critical value

$$\bar{U}_c = \frac{\omega_0}{k_{0x}}. \quad (14)$$

This is consistent with the picture put forth in e.g. Goldreich and Nicholson (cite).

A purely horizontal shear flow $\bar{U}_0(z)\hat{x}$ can be seen in Equation 1 to have the effect of modifying time derivatives ∂_t to their frequency in the comoving frame of the fluid $\partial_t - \bar{U}_0(z)\partial_x$. For a critical value $\omega_0 - \bar{U}_c k_{0x} = 0$, the frequency of the linear wave in the fluid’s frame of reference vanishes and critical behavior is observed. In a linear theory or a theory where small scales are viscosity rather than advection dominated, the incident wave has amplitude reflection and transmission coefficients

$$\mathcal{R} = e^{-2\pi\sqrt{\text{Ri}-\frac{1}{4}}}, \quad \mathcal{T} = e^{-\pi\sqrt{\text{Ri}-\frac{1}{4}}}, \quad (15)$$

where we have defined Richardson number $\text{Ri} \equiv \frac{N^2}{(*)\frac{\partial \bar{U}_0}{\partial z}} \bigg|_{z_c}$ at the

critical layer z_c : $\bar{U}_0(z_c) = \frac{\omega_0}{k_{0x}}$. For most shear flows, $\text{Ri} \gg 1$ and so $\mathcal{R}, \mathcal{T} \ll 1$ and the incident wave is absorbed.

When the fluid absorbs the incident wave, it absorbs the incident horizontal momentum flux as well, which is converted into additional horizontal momentum of the shear flow. Since the shear flow cannot exceed \bar{U}_c the horizontal phase velocity of the incident

wave, the critical layer must thus propagate downwards (towards the wave source) to accommodate the incident momentum. In other words, the total horizontal momentum of the shear flow obeys conservation equation

$$\frac{\partial}{\partial t} \int_0^{L_z} \rho(z) \bar{U}_0(z, t) dz - S_{px} = 0. \quad (16)$$

Treating $\bar{U}_0(z > z_c) = \bar{U}_c$, $\bar{U}_0(z < z_c) = 0$ gives us exactly

$$\rho \bar{U}_c \frac{\partial z_c}{\partial t} = -S_{px}. \quad (17)$$

For constant S_{px} in space and $\rho \approx \rho_0$, this has analytical solution

$$z(t) = -H \ln t - H \ln \frac{H \rho_0(z=0) c_{ph,x}}{S_{px}}. \quad (18)$$

4.5 Numerical Simulation

We use the same k_{0x}, ω_0 as subsection 4.3. Our other parameters are:

- We choose $\nu = 0.35 \frac{\omega_0}{k_{0z} k_{z,\max}}$, where $k_{z,\max} = \frac{2\pi N_z}{L_z}$. Note that $\nu = \frac{\omega_0}{k_{0z} k_{z,\max}}$ corresponds to the advective term $\vec{u} \cdot \vec{\nabla}$ being of the same order as the time derivative ∂_t for flow velocities $\vec{u} \sim \frac{\omega_0}{k_{0z}}$ at the grid spacing.

Two representative snapshots from our simulation are provided in TODO after the critical layer has had time to form. We may note that the critical layer, where S_{px} is absorbed and $\bar{U}_0 = \bar{U}_c$, travels downwards as predicted.

4.6 Propagating Critical Layer

For the simulation in TODO, we may define $z_c = \arg\max_z \frac{\partial \bar{U}_0}{\partial z}$. Computing $\frac{\partial z_c}{\partial t}$ using the analytical flux Equation 8 allows comparison to Equation 17, which we exhibit in TODO. We see overall good agreement, though some small deviation is expected since S_{px} is not perfectly conserved owing to numerical viscosity (and YUBONOTE S_{px} is misestimated?).

However, recalling Equation 15, we only expect complete critical layer absorption when $\text{Ri} \gg \frac{1}{4}$. We may plot Ri over the same time interval in TODO. We may observe that the Richardson number is initially decreasing but eventually levels out and begins to increase again. This corresponds to an increasingly sharp critical layer transition then subsequently a decreasingly sharp critical layer transition.

We argue that the Richardson number is bounded from below by viscosity. A repeated simulation with a larger viscosity is shown in TODO, where the Richardson number does not go nearly as low.

5 DISCUSSION

Discussion foo

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APPENDIX A: EQUATION IMPLEMENTATIONS

We denote $x \in [0, L_x], z \in [0, L_z]$ the simulation domain and N_x, N_z the number of spectral modes in the respective dimensions. We perform direct numerical simulation of Equation 9 with the open-source pseudo-spectral code Dedalus (CITE).

Numerically, the nonlinear $\frac{\vec{\nabla}P}{\rho}$ term is problematic: we desire a system where the fluid fields are not divided by one another. We introduce $\varpi = \frac{P}{\rho}$ instead, then mandate ρ_0, ϖ_0 background fields satisfy hydrostatic equilibrium $\vec{\nabla}\varpi_0 + \varpi_0\vec{\nabla}\rho_0 + g\hat{z} = 0$. Taking isothermal stratification, we find $\varpi_0 = gH$. We further change variables to $\Upsilon = \ln \rho - \ln \rho_0$ and $\varpi_1 = \varpi - \varpi_0$ deviations from the background state to obtain a system of equations at most quadratic in fluid fields:

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (\text{A1a})$$

$$\frac{\partial \Upsilon}{\partial t} + (*)\vec{u} \cdot \vec{\nabla}\Upsilon - \frac{u_z}{H} = 0, \quad (\text{A1b})$$

$$\frac{\partial u_x}{\partial t} + (*)\vec{u} \cdot \vec{\nabla}u_x + \frac{\partial \varpi_1}{\partial x} + gH\frac{\partial \Upsilon}{\partial x} + \varpi_1\frac{\partial \Upsilon}{\partial x} = 0, \quad (\text{A1c})$$

$$\frac{\partial u_z}{\partial t} + (*)\vec{u} \cdot \vec{\nabla}u_z + \frac{\partial \varpi_1}{\partial z} + gH\frac{\partial \Upsilon}{\partial z} + \varpi_1\frac{\partial \Upsilon}{\partial z} - \frac{\varpi_1}{H} = 0. \quad (\text{A1d})$$

It bears noting that these equations are exactly equivalent to the original Euler equations and hence conserve horizontal momentum.

A1 Artificial Dissipation

The nonlinear terms in the above equations will transfer energy from lower wavenumbers to higher wavenumbers. Since spectral codes have no numerical dissipation, artificial dissipation must be added. To ensure the dissipative system conserves horizontal momentum, we begin by adding dissipative terms to the flux-conservative form of the Euler fluid equations Equation 1 (we use stress tensor $\tau_{ij} = P\delta_{ij}$):

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (\text{A2a})$$

$$\partial_t \rho + \vec{\nabla} \cdot (\rho \vec{u} - \nu \vec{\nabla}(\rho - \rho_0)) = 0, \quad (\text{A2b})$$

$$\partial_t (\rho \vec{u}) + \vec{\nabla} \cdot (\rho \vec{u} \vec{u} + \text{diag}(\rho \varpi) - \nu \rho \vec{\nabla} \vec{u}) + \rho g \hat{z} = 0. \quad (\text{A2c})$$

The same ν is used for both the diffusive and viscous term, though this is not required. Since the dissipation is not physical and is purely used for numerical stability, we choose it such that hydrostatic equilibrium is not modified. Some algebraic manipulation to re-cast it in the form of Equation A1 gives

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (\text{A3a})$$

$$\partial_t \Upsilon + (*)\vec{u} \cdot \vec{\nabla}\Upsilon - \frac{u_z}{H} - \nu(*)\nabla^2 \Upsilon + (*)\vec{\nabla}\Upsilon \cdot (*)\vec{\nabla}\Upsilon - \frac{2}{H}\partial_z \Upsilon + \frac{1 - e^{-\Upsilon}}{H^2} = 0, \quad (\text{A3b})$$

$$\begin{aligned} \partial_t \vec{u} + (*)\vec{u} \cdot \vec{\nabla}\vec{u} + \vec{\nabla}\varpi + \varpi\vec{\nabla}\Upsilon - \nu\nabla^2 \vec{u} + \vec{u}\nu(*)\nabla^2 \Upsilon + (*)\vec{\nabla}\Upsilon \cdot (*)\vec{\nabla}\Upsilon - \frac{2}{H}\partial_z \Upsilon + \frac{1 - e^{-\Upsilon}}{H^2} \\ - 2\nu(*)\vec{\nabla}\Upsilon \cdot \vec{\nabla}\vec{u} - \frac{1}{H}\partial_z \vec{u} - \frac{\varpi_1}{H} = 0. \end{aligned} \quad (\text{A3c})$$

Hydrostatic equilibrium is still $\vec{\nabla}\varpi_0 + g\hat{z} = 0$ where $\rho = \rho_0, \vec{u} = 0$. Including the damping layers and forcing terms as described in subsection 4.1, we finally obtain the full system of equations as simulated in Dedalus:

$$\vec{\nabla} \cdot \vec{u} = 0, \quad (\text{A4a})$$

$$\begin{aligned} \partial_t \Upsilon - \frac{u_z}{H} = \nu(*)\nabla^2 \Upsilon + (*)\vec{\nabla}\Upsilon \cdot (*)\vec{\nabla}\Upsilon - \frac{2}{H}\partial_z \Upsilon + \frac{1 - e^{-\Upsilon}}{H^2}, \\ - (*)\vec{u} \cdot \vec{\nabla}\Upsilon - \Gamma(z)\Upsilon + \frac{F}{\rho_0(z)}e^{-\frac{(z-z_0)^2}{2\sigma^2}}\cos(*)k_x x - \omega t, \end{aligned} \quad (\text{A4b})$$

$$\begin{aligned} \frac{\partial u_x}{\partial t} + \frac{\partial T}{\partial x} + gH\frac{\partial \Upsilon}{\partial x} = \nu\nabla^2 u_x - u_x\nu(*)\nabla^2 \Upsilon + (*)\vec{\nabla}\Upsilon \cdot (*)\vec{\nabla}\Upsilon - \frac{2}{H}\partial_z \Upsilon + \frac{1 - e^{-\Upsilon}}{H^2} \\ + 2\nu(*)\vec{\nabla}\Upsilon \cdot \vec{\nabla}u_x - \frac{1}{H}\partial_z u_x - \Gamma(z)u_x - (*)\vec{u} \cdot \vec{\nabla}u_x - T_1\frac{\partial \Upsilon}{\partial x}, \end{aligned} \quad (\text{A4c})$$

$$\begin{aligned} \frac{\partial u_z}{\partial t} + \frac{\partial T}{\partial z} + gH\frac{\partial \Upsilon}{\partial z} - \frac{T_1}{H} = \nu\nabla^2 u_z - u_z\nu(*)\nabla^2 \Upsilon + (*)\vec{\nabla}\Upsilon \cdot (*)\vec{\nabla}\Upsilon - \frac{2}{H}\partial_z \Upsilon + \frac{1 - e^{-\Upsilon}}{H^2} \\ + 2\nu(*)\vec{\nabla}\Upsilon \cdot \vec{\nabla}u_z - \frac{1}{H}\partial_z u_z - \Gamma(z)u_z - (*)\vec{u} \cdot \vec{\nabla}u_z - T_1\frac{\partial \Upsilon}{\partial z}. \end{aligned} \quad (\text{A4d})$$