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A blended soundproof-to-compressible numerical model for small

to meso-scale atmospheric dynamics

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ABSTRACT

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A blended model for atmospheric flow simulations is introduced that enables seamless transition from fully-compressible to pseudo-incompressible dynamics. The model equations are written in non-perturbational form and integrated using a well-balanced second-order finite volume discretization. The semi-implicit scheme combines an explicit predictor for advection with elliptic corrections for the pressure field. Compressibility is implemented in the elliptic equations through a diagonal term. The compressible/pseudo-incompressible transition is realized by suitably weighting the term and provides a mechanism for removing unwanted acoustic imbalances in compressible runs.

As the gradient of the pressure is used instead of the Exner pressure in the momentum equation, the influence of perturbation pressure on buoyancy must be included to ensure thermodynamic consistency. With this effect included the thermodynamically consistent model is equivalent to Durran's original pseudo-incompressible model, which uses the Exner pressure.

Numerical experiments demonstrate quadratic convergence and competitive solution quality for several benchmarks. With the inclusion of an additional buoyancy term required for thermodynamic consistency, the " $p-\rho$ -formulation" of the pseudo-incompressible model closely reproduces the compressible results.

The proposed unified approach offers a framework for models that are largely free of
the biases which can arise when different discretizations are used. With data assimilation
applications in mind, the seamless compressible/pseudo-incompressible transition mechanism
is also shown to enable the flattening of acoustic imbalances in initial data for which balanced
pressure distributions are unknown.

₂₇ 1. Introduction

Physical processes in the atmosphere feature a wide range of spatio-temporal scales de-28 scribed by the fully-compressible non-hydrostatic flow equations. Accordingly, non-hydrostatic fully-compressible modelling approaches hold sway in atmospheric research codes and in op-30 erational dynamical cores, e.g., ICON (Zängl et al. 2014), NUMA (Kelly and Giraldo 2012), 31 DUNE (Brdar et al. 2013), the models in use at NCAR (Wong et al. 2014), ECMWF (Hortal 32 2002; Smolarkiewicz et al. 2014), the UK Met Office (Davies et al. 2005; Wood et al. 2013), 33 and others. 34 Despite very successful ongoing developments, the proper treatment of multiple charac-35 teristic time scales in atmospheric simulations remains a matter of scientific research. Two of 36 the biggest obstacles of multiple-scales simulations are (i) the discretization of fast processes 37 in the governing equations and (ii) balanced data assimilation. 38 Numerical stiffness is the source of the first remaining obstacle. Except for inertia-39

gravity waves of long wavelength, which are not considered here, quantities of meteorological interest propagate at low speed compared with sound waves. Sound modes are said to be nearly balanced and their effects are considered negligible for atmospheric dynamics. The difference between the sound and flow speeds stiffens the numerics of fully-compressible solvers rendering straightforward explicit schemes impractical due to severe stability-related time step constraints.

Filtering the data with respect to fast modes while minimally distorting the ensuing dynamics is the second remaining obstacle. Computational simulations never exactly track the evolution of the considered system. Hence, data assimilation is needed for exploiting observational data at regular time intervals to set up initial data for the next simulation period. However, importing observed field data from local weather stations directly to adjacent grid points would disregard the aforementioned balances of the fast modes. For example, in the presence of a low pressure system in the summer with high levels of convection, the local vertical velocities would project onto non-hydrostatic and compressible modes yielding

strongly unbalanced data on the numerical grid. Efficiently controlling such modes remains
 a challenge in data assimilation.

Numerical approaches aimed at overcoming the stiffness are split-explicit, semi-implicit, and fully implicit numerical time integrators for the fully-compressible flow equations. The first class of schemes subcycles a simplified discretization of the fast wave processes at short time steps and employs suitable synchronization procedures for coupling the results to large time steps of the slower modes (Skamarock and Klemp 1994, 2008; Jebens et al. 2009). Another option would be to adopt a fully implicit approach which even overcomes the time step restrictions associated with explicit discretizations of advection. Due to their computational expense these schemes have, to our knowedge, thus far not found widespread application in meteorology. A notable exception is the work by Reisner et al. (2005).

The focus of the present work lies instead on semi-implicit discretizations which invoke implicit integrators for the terms in the equations representing the fast wave modes while treating the slow modes explicitly. Many approaches to semi-implicit discretization for atmospheric flows have been reported, e.g., by Bonaventura (2000); Gatti-Bono and Colella (2006); Restelli and Giraldo (2009); Jebens et al. (2011); Durran and Blossey (2012); Giraldo et al. (2013); Wood et al. (2013); Smolarkiewicz et al. (2014); Weller and Shahrokhi (2014). For all-speed flow discretizations in computational fluid dynamics the reader is referred to Casulli and Greenspan (1984); Bijl and Wesseling (1998); Munz et al. (2003); Kwatra et al. (2009).

An alternative to these numerical approaches to overcoming the stiffness is to adopt a "soundproof" model. These reduced dynamical models include a diagnostic constraint on the velocity divergence and therefore do not support sound waves. The divergence constraint needs to be maintained numerically, which entails the solution of an elliptic pressure equation. Soundproof models suitable for atmospheric motions covering vertical distances comparable to the pressure scale height are the anelastic (Lipps and Hemler 1982; Bannon 1996) and pseudo-incompressible models (Durran 1989; Klein and Pauluis 2011).

Soundproof models have successfully been used to simulate small to meso-scale flows, and 81 their validity as slow-flow limit models has recently been established on theoretical grounds 82 (Klein et al. 2010; Achatz et al. 2010). However, their applicability to large-scale motions is 83 still under debate (Davies et al. 2003; Dukowicz 2013) despite recent successful large-scale simulations for atmospheric (Smolarkiewicz and Dörnbrack 2008; Smolarkiewicz et al. 2014) and astrophysical (Nonaka et al. 2010; Smolarkiewicz and Charbonneau 2013) applications. In line with these observations, one of our goals is to develop a numerical scheme for the 87 fully-compressible equations that defaults to the pseudo-incompressible limit for slow flows on small to meso scales. Such asymptotically adaptive schemes have a substantial history of studies (Klein 2000; Klein et al. 2001; Gatti-Bono and Colella 2006; Cullen 2007; Haack et al. 2012) in which the low Mach or low Froude number limits are discretely recovered through careful identification and separate discretization of the advection, acoustic, and/or buoyancy 92 terms in the fully-compressible equations. In the present work we suggest a particularly 93 straightforward approach of this type that is directly motivated by the theoretical framework 94 set out in Klein (2009, 2010). 95

More specifically, this paper documents the construction of a semi-implicit second-order 96 accurate numerical method for the simulation of weakly compressible atmospheric flows that shares the principal components of the discretization with the soundproof solver by Klein 98 (2009). The time integration for the fully compressible equations derives from that of the pseudo-incompressible model and the required adjustments amount to no more than adding 100 a diagonal term to the matrix of the elliptic pressure problem and synchronizing the cell-101 centered and node-based pressures. This is similar in spirit to parallel developments by 102 Smolarkiewicz et al. (2014) but technically different. In particular, these authors do not 103 address the possibility of a seamless blending of models and they work with perturbation 104 variables and with the Exner pressure in the momentum equation. 105

Besides constructing the compressible flow solver, we design the discretization such that it can be used directly to solve a continuous family of weakly compressible models that interpolate seamlessly between the fully-compressible and pseudo-incompressible ones. This
is realized by exploiting the close structural similarity of these two limiting models when
written in conservative, non-perturbational form for the densities of mass, momentum, and
potential temperature.

In the context of increasing computing resources and ever smaller scales accessible in 112 high-resolution weather and climate simulations, it is of arguable interest to operate differ-113 ent analytical formulations within a single numerical framework. Such a unified numerical 114 scheme becomes all the more desirable in the light of a recent study (Smolarkiewicz and 115 Dörnbrack 2008) that compared the errors made by using different numerical methods for 116 the same model equations with those made by considering different equation systems dis-117 cretized with nearly identical numerics. These authors found, somewhat surprisingly, that 118 the former errors exceeded the latter, and this underlines the importance of comparing flow 119 models within one and the same numerical framework. In an interesting investigation of this 120 type, Smith and Bannon (2008) compared anelastic and compressible models in a case of 121 localized instantaneous diabatic warming. 122

A second motivation for implementing the seamless model family lies in its potential use for balanced data assimilation. By adjusting the model interpolation parameter accordingly from zero to unity, such a "blended" scheme can be tuned to perform a few time steps in pseudo-incompressible mode and to then transition to its fully-compressible mode after a few further steps. As we will show, this effectively reduces initial acoustic imbalances. Considering the factors affecting predictability of the simulated precipitation field in cloud-resolving models, Hohenegger and Schär (2007) showed that uncontrolled small-scale acoustic perturbations may contribute to rapid error growth at the mesoscale.

The scheme we propose has more potentially attractive features. One of these features is the formulation in a non-perturbational form that does not rely on subtraction of a background state for accuracy. This is achieved for the present collocated finite volume method by a well-balanced discretization of the pressure gradient and gravity terms following Botta

et al. (2004); Klein (2009). Moreover, the scheme uses the gradient of the thermodynamic 135 instead of the Exner pressure, thereby allowing for a conservative discretization of the mo-136 mentum flux induced by the pressure force. In addition, as pointed out by Klein and Pauluis 137 (2011), Durran's original formulation of the pseudo-incompressible model using Exner pres-138 sure cannot be easily extended to general equations of state. One step towards overcoming 139 this obstacle is to adopt a formulation with pressure instead of Exner pressure in the momen-140 tum equation as done in this paper. Yet, this formulation is thermodynamically consistent only if first-order density perturbations are included in the gravity term in addition to Dur-142 ran's "pseudo-density". For an ideal gas with constant specific heat capacities, Durran's model and the present thermodynamically consistent formulation are equivalent as a short 144 calculation using the transformations $\pi_0 = (p_0/p_{\rm ref})^{R/c_p}$ and $\pi' = p'/(c_p P_0)$ shows. A second 145 step that is also necessary in extending to general equations of state, but which is not pursued 146 here, is a reformulation of the velocity divergence constraint. This step is needed because in 147 this case the pressure equation can no longer be easily cast into a simple conservation law 148 (Almgren et al. 2006a,b; Klein and Pauluis 2011). 149

Furthermore, the transition from the pseudo-incompressible via the blending to the compressible model is achieved by minimal code adjustments. These involve reassigning certain
weights in the grid stencil of the elliptic correction equations and applying a weighted superposition of pressure updates. These updates are calculated from the elliptic equations and
from the conservative balance of potential temperature.

The paper is structured as follows. Compressible, pseudo-incompressible, and blended models are presented in section 2. Section 3 summarizes the numerics. The results of numerical simulations in a number of two-dimensional test cases is documented in section 4. Grid convergence with the expected second-order rate is verified in a benchmark involving advection of a smooth axysimmetric vortex. For the standard test cases of a rising hot air thermal, density current and inertia-gravity waves, we compare the predictions obtained with the compressible and pseudo-incompressible models and demonstrate the importance

of the thermodynamic consistency correction within the pseudo-incompressible framework.

Usage of the blended model for filtering acoustic imbalances is demonstrated for both short sound-resolving time steps and for time steps corresponding to an advective CFL number of order unity. Section 5 provides a concluding discussion and an outline of open issues and future work.

₆₇ 2. Theoretical Framework

168 Fully-compressible equations

The dry, inviscid fully-compressible equations, henceforth referred to as "FC", describe conservation of mass, momentum, and energy under the influence of gravity. If we neglect rotational effects and use the transport equation for potential temperature to describe the energy balance, they read in conservative form and in the dry adiabatic case,

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) = 0, \tag{1a}$$

$$\frac{\partial \rho \mathbf{v}}{\partial t} + \nabla \cdot (\rho \mathbf{v} \circ \mathbf{v} + p \mathbf{I}) = -\rho g \mathbf{k}, \tag{1b}$$

$$\frac{\partial P}{\partial t} + \nabla \cdot (P\mathbf{v}) = 0. \tag{1c}$$

Here, ρ denotes the fluid density, \mathbf{v} the velocity vector, \circ the tensor product, g the acceleration of gravity, \mathbf{k} the vertical unit vector, and \mathbf{I} the identity tensor. As in Klein (2009), we have introduced the equation of state

$$P = \rho \theta = \frac{p_{\text{ref}}}{R} \left(\frac{p}{p_{\text{ref}}}\right)^{\frac{1}{\gamma}},\tag{2}$$

where potential temperature is defined as

$$\theta = T \left(\frac{p}{p_{\text{ref}}}\right)^{\frac{1-\gamma}{\gamma}}$$
 and $T = \frac{p}{\rho R}$ (3)

is the temperature. R is the gas constant for dry air, γ is the isentropic exponent, respectively.

Hereafter, we take $\gamma = 1.4$ and R = 287 N m kg⁻¹ K⁻¹ throughout. For smooth flows,

(1c) can equivalently replace total energy conservation in a finite volume discretization, 179 which is common in numerical meteorology, but which would not be adequate for flows 180 with shocks (LeVeque 2002). Together, (1a) and (1c) describe mass conservation and the 181 advection of potential temperature, while (1c) is equivalent to the pressure evolution equation 182 $p_t + \mathbf{v} \cdot \nabla p + \gamma p \nabla \cdot \mathbf{v} = 0$. Thus, a discretization of (1c) directly controls the pressure evolution, 183 and this is central to the blended compressible—soundproof formulation to be presented below. 184 The system is closed by appropriate initial and boundary conditions which we will specify 185 in conjunction with specific test cases below. 186

For later reference, using (2), we compute

$$\frac{\partial P}{\partial p} = \frac{1}{R\gamma} \left(\frac{p}{p_{\text{ref}}} \right)^{\frac{1}{\gamma} - 1} = \frac{1}{R\gamma} \left(\frac{PR}{p_{\text{ref}}} \right)^{1 - \gamma}. \tag{4}$$

The pseudo-incompressible approximation

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The pseudo-incompressible model (Durran 1989) is commonly derived from a compressible model that formulates the pressure gradient term in the momentum equation using the
Exner pressure,

$$\pi = \left(\frac{p}{p_{\text{ref}}}\right)^{\frac{\gamma - 1}{\gamma}} \tag{5}$$

so that, in view of (3), one finds

$$\frac{1}{\rho} \nabla p \equiv c_p \theta \nabla \pi \,. \tag{6}$$

To retain flexibility of the developed code, in particular with respect to generalizations of the equation of state, we adopt the $p-\rho$ formulation here (Klein and Pauluis 2011). When written in the latter form, extra care must be taken in formulating the momentum equation to ensure that it retains the influences of the pressure perturbation up to first order.

As in Durran (1989) we start our derivations by assuming that the pressure does not vary much from its hydrostatic background value and can be written as

$$p = p_0(z) + p'(\mathbf{x}, t), \tag{7}$$

where $p'/p_0 \ll 1$ and

$$\frac{\partial p_0}{\partial z} = -\rho_0 g. \tag{8}$$

Using (7) in the equation of state (2) gives, with a Taylor expansion,

$$\rho = \frac{1}{\theta} \frac{p_{ref}}{R} \left(\frac{p_0 + p'}{p_{ref}} \right)^{1/\gamma} \approx \frac{1}{\theta} \frac{p_{ref}}{R} \left(\frac{p_0}{p_{ref}} \right)^{1/\gamma} \left(1 + \frac{p'}{\gamma p_0} \right) = \rho^* \left(1 + \frac{p'}{\gamma p_0} \right)$$
(9)

where ρ^* is called the "pseudo-density" and is defined as the density calculated at the background pressure but using the full potential temperature, i.e.

$$\rho^* = \frac{1}{\theta} \frac{p_{ref}}{R} \left(\frac{p_0}{p_{ref}}\right)^{1/\gamma} = \rho(p_0, \theta). \tag{10}$$

To filter sound waves we suppress the effect of pressure pertubations on density to obtain

$$(\rho^*)_t + \nabla \cdot (\rho^* \mathbf{v}) = 0. \tag{11}$$

However, in the momentum equation we want to keep the effect of the pressure perturbations up to first order. Using an expansion as in (10) we re-write (1b) in non-conservative form

$$\mathbf{v}_t + \mathbf{v} \cdot \nabla \mathbf{v} + \frac{1}{\rho^*} \left(1 - \frac{p'}{\gamma p_0} \right) \nabla \left(p_0 + p' \right) = -g \mathbf{k} \,. \tag{12}$$

206 Keeping terms in (12) up to first order in the pressure perturbation and re-arranging we get

$$\mathbf{v}_t + \mathbf{v} \cdot \nabla \mathbf{v} + \frac{1}{\rho^*} \nabla \left(p_0 + p' \right) = -\left(1 + \frac{1}{\rho^*} \frac{\rho_0}{\gamma p_0} p' \right) g \mathbf{k}. \tag{13}$$

We re-write (13) in conservative form by multiplying by ρ^* and using (11),

$$(\rho^* \mathbf{v})_t + \nabla \cdot (\rho^* \mathbf{v} \circ \mathbf{v}) + \nabla p = -\left(\rho^* + \frac{\rho_0}{\gamma p_0} p'\right) g \mathbf{k}.$$
 (14)

Lastly, we redefine P as

$$P \approx \rho^* \theta = \frac{p_{ref}}{R} \left(\frac{p_0}{p_{ref}}\right)^{1/\gamma} = P_0 \tag{15}$$

209 and (1c) becomes

$$(P_0)_t + \nabla \cdot (P_0 \mathbf{v}) = \nabla \cdot (P_0 \mathbf{v}) = 0. \tag{16}$$

In (16) we have used that P is now a function of p_0 only which allows us to drop the time derivative term and the evolution equation becomes a divergence constraint. This constraint enforces the pseudo-incompressible form of the density equation in (11) thereby filtering the effect of pressure perturbations on the density and thus filtering sound waves.

The complete pseudo-incompressible governing equations are given by

$$(\rho^*)_t + \nabla \cdot (\rho^* \mathbf{v}) = 0 \tag{17a}$$

$$(\rho^* \mathbf{v})_t + \nabla \cdot (\rho^* \mathbf{v} \circ \mathbf{v}) + \nabla p = -\left(\rho^* + \frac{\rho_0}{\gamma p_0} p'\right) g\mathbf{k}$$
 (17b)

$$\nabla \cdot (P_0 \mathbf{v}) = 0 \tag{17c}$$

Klein (2009) showed agreement between (17a)-(17c) and the original formulation of Dur-215 ran (1989) to leading and first order in a perturbation expansion for small pressure variations. Moreover, if Exner pressure variables are introduced so that $\pi_0 = (p_0/p_{\rm ref})^{R/c_p}$ and $\pi' = p'/(c_p P_0)$, a straightforward calculation shows that the original formulation of Durran 218 (1989) and the present $\mathrm{PI}^{\mathrm{tc}}_{\rho,p}$ formulation are actually equivalent at the level of the partial 219 differential equations. An advantage of our formulation is that it is more easily extended to 220 incorporate more complex equations of state and that it is "thermodynamically consistent". 221 This notion refers to the existence of well-defined thermodynamic potentials describing the 222 proper increase/decrease of an entropy variable in the diabatic case (Klein and Pauluis 2011). 223 Note, however, that completing the extension to general equations of state also requires a 224 reformulation of the divergence constraint (Almgren et al. 2006a,b; Klein and Pauluis 2011). 225

226 A blended compressible/pseudo-incompressible model

In Klein (2009) the task of incorporating the time derivative term in (1c) and modelling the fully-compressible dynamics was left for future work. Here we aim to merge the compressible, pseudo-incompressible, and thermodynamically consistent discretizations in the "p- ρ -formulation" for the momentum equation in a single numerical model featuring

- a conservative discretization with respect to $\rho, \rho \mathbf{v}, \rho \theta \equiv P$,
- second-order accuracy,

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- time steps independent of the sound speed,
- a continuous transition between pseudo-incompressible and compressible forms,
- a well-balanced discretization that does not rely on subtraction of a background state.

For $\alpha = 0$ the two pseudo-incompressible models with the "p- ρ -formulation" of the pressure

gradient term are retrieved. Then, setting $\beta = 1$ selects the thermodynamically consistent

The blended equations are given as follows, for $\alpha \in \{0, 1\}$:

$$\rho_t + \nabla \cdot (\rho \mathbf{v}) = 0, \tag{18a}$$

$$(\rho \mathbf{v})_t + \nabla \cdot (\rho \mathbf{v} \circ \mathbf{v}) + \nabla p = -g \mathbf{k} \left(\rho + (1 - \alpha) \beta \frac{\rho_0}{\gamma p_0} p' \right), \tag{18b}$$

$$\alpha P_t + \nabla \cdot (P\mathbf{v}) = 0. \tag{18c}$$

 $(\mathrm{PI}_{\rho,p}^{\mathrm{tc}})$ model whereas setting $\beta=0$ retrieves the "naive" pseudo-incompressible $(\mathrm{PI}_{\rho,p})$ 239 model. We note that in $\text{PI}_{\rho,p}$ and $\text{PI}_{\rho,p}^{\text{tc}}$ the density ρ takes the role of the pseudo-density, which was denoted by ρ^* in (17b), and necessitates the additional term for thermodynamic consistency in the momentum equation (18b) for $(\alpha, \beta) = (0, 1)$. As the model parameter α is adjusted from 0 to 1, the effect of pressure perturbations on density is retrieved in 243 a continuous fashion. This formulation recovers the fully-compressible (FC) dynamics for $\alpha = 1$. A summary of the model configurations is given in Table 1. 245 System (18) features unapproximated mass and momentum equations for $\alpha \in \{0,1\}$ 246 when $\beta = 1$. The reason is that the $\text{PI}_{\rho,p}^{\text{tc}}$ model is equivalent to Durran's original pseudo-247 incompressible model with the " π - θ -formulation" of the pressure gradient term. Klein et al. 248 (2013) observe that the model satisfies an energy conservation law with a definition of the 249 total energy that is an interpolation between those of the fully-compressible and the pseudo-250 incompressible models. The model's internal wave dispersion properties for realistic stratifi-251 cations are close to those of the limiting models. This follows from related analyses for the 252

limiting models by Klein (2010) and the fact that the underlying Sturm-Liouville problems
depend smoothly on the defining data. We also refer to Vasil et al. (2013) for related analysis
and relegate further discussion to a future publication.

In (18) the α and β parameters are introduced to formulate the FC, PI^{tc}_{ρ,p}, and PI_{ρ,p} models conveniently in one and the same set of equations. Only discrete values $\alpha, \beta \in \{0, 1\}$ make sense to begin with. Yet, let us consider the resulting model equations for any $\alpha \in [0, 1]$.

A seamless discretization that allows integration of (18) for any of these values can be used to our advantage in some meteorologically interesting situation.

Suppose we are to initialize one of the well-known test cases of a rising warm-air bubble or flow over a mountain. As in "real meteorology", we are not interested in acoustic perturbations and would like to simulate acoustically balanced flows. Yet, we have no analytical way to determine the balanced pressure distributions that would be associated with given initial data for potential temperature and velocity.

However, knowing that the pseudo-incompressible models provide good approximations 266 to compressible flows free of sound waves, we can attempt to generate reasonable approxi-267 mations to the missing pressure fields by starting a simulation pseudo-incompressibly with 268 $\alpha = 0$ for, say, S_1 time steps. Within the next S_2 time steps we increase α continuously 269 from 0 to 1, and after time step $S_1 + S_2$ we maintain $\alpha = 1$ to operate the model in fully-270 compressible mode. This procedure should generate a compressible flow simulation that is 271 balanced with respect to acoustic modes essentially from the start. Promising related results 272 for the rising bubble test are discussed in section 4 below. 273

We conjecture that such a smooth blending of balanced and unbalanced model equations within a common discretization framework could substantially contribute to resolving similar balancing issues in the context of data assimilation.

77 3. Numerical Framework

A semi-implicit finite volume method is used to approximate the dynamics of the blended model. The scheme is a variant and extension of the soundproof solver described in Klein (2009). An outline is presented here, for more details see Appendix. The discrete solution of (18) is obtained by the following time stepping procedure, say from t^n to t^{n+1} :

- An explicit predictor solves an auxiliary hyperbolic system obtained by replacing the pressure gradient in the momentum equation (18b) with its value at time level t^n . This step yields second-order accurate ρ , θ and P;
- A first elliptic corrector solves for the cell-centered pressure time increment $\delta p = p^{n+1} p^n$ by enforcing consistency with the pressure equation (18c). This step also corrects the advecting fluxes in (18a) and (18b);
- The solution of a second elliptic problem is used to correct the pressure-related momentum flux for fully second-order accurate updates of the cell-centered momenta.

For the time discretization we divide the simulation time interval [0,T] into N subinter-290 vals, with $t_0 = 0$, $t^{n+1} = t^n + (\Delta t)^n$ for $n = 0, 1, \dots N - 1$. For any variable X, we denote 291 $X^n = X(t^n)$. $(\Delta t)^n = O(T/N)$ denote the time steps. In the implementation, a dynamically 292 adaptive choice of the time step based on fixing the Courant number is implemented, see 293 Appendix for details. The spatial domain is divided into primary computational cells $C_{i,j}$ 294 (finite volumes) with $i=1,\ldots,\mathcal{N}_x,\ j=1,\ldots,\mathcal{N}_z$, in two dimensions according to a carte-295 sian grid arrangement. The cells $C_{i,j}$ are separated by interfaces $I_{i+1/2,j}$, $I_{i,j+1/2}$ as shown in 296 Fig. 1. The extension to three dimensions is straightforward. The primary variables $\rho, \rho \mathbf{v}, P$ 297 are stored at the centers of the primary cells $C_{i,j}$. Pressures are computed at centers of the 298 primary cells $C_{i,j}$ in the first correction step and at the centers of the dual cells $\overline{C}_{i+1/2,j+1/2}$ 299 shown in Fig. 1 in the second correction step.

In the first sub-step for a full time step $t^n \to t^{n+1}$, the following auxiliary hyperbolic system, obtained from (18) by freezing p and p' at time level t^n , is solved (Klein 2009):

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) = 0, \tag{19a}$$

$$\frac{\partial \rho \mathbf{v}}{\partial t} + \nabla \cdot (\rho \mathbf{v} \circ \mathbf{v} + p^n \mathbf{I}) = -g \mathbf{k} \left(\rho + (1 - \alpha) \beta \frac{\rho_0}{\gamma p_0} (p')^n \right), \tag{19b}$$

$$\frac{\partial P}{\partial t} + \nabla \cdot (P\mathbf{v}) = 0. \tag{19c}$$

A two-stage strong stability-preserving Runge-Kutta method (Gottlieb et al. 2001) is used for time integration here (Klein (2009) instead used a MUSCL technique and directional operator splitting). The spatial discretization at any stage of the Runge-Kutta time integrator is performed with a finite volume approach. That is, discrete variables X_C , $X = \rho$, $\rho \mathbf{v}$, P, are defined as approximations of the cell averages set at the cell centers:

$$X_C = \frac{1}{|C|} \int_C X \, dx + O\left(\Delta x^2\right),\tag{20}$$

where |C| is the cell volume. To achieve second-order accuracy in space, piecewise linear reconstruction of P, \mathbf{v} , and the advected quantities $(1/\theta, \mathbf{v}/\theta)$ is applied within the grid cells. The reconstructed values are used to determine any data required at grid cell interfaces and to evaluate the numerical flux functions. The pressure variables p^n , $(p')^n$ are set at the grid nodes.

New values of X_C are obtained from the old ones subtracting the net outflow fluxes at the boundaries and adding the contributions from the source terms:

$$\rho_C^{n+1,*} = \rho_C^n - \Delta t \left(\widetilde{\nabla} \cdot \left(P \mathbf{v} \, \theta^{-1} \right) \right)_C^{n+\frac{1}{2},*}, \tag{21a}$$

$$(\rho \mathbf{v})_C^{n+1,*} = (\rho \mathbf{v})_C^n - \Delta t \left(\widetilde{\nabla} \cdot \left(P \mathbf{v} \circ \mathbf{v} \theta^{-1} + p^n \mathbf{I} \right) \right)_C^{n+\frac{1}{2},*} - \Delta t \, g \mathbf{k} \left(P/\theta + (\rho')^n \right)_C^{n+\frac{1}{2},*}, \quad (21b)$$

$$P_C^{n+1,*} = P_C^n - \Delta t \left(\widetilde{\nabla} \cdot (P\mathbf{v}) \right)_C^{n+\frac{1}{2},*}, \tag{21c}$$

where $\rho' = (1 - \alpha)\beta(\rho_0/\gamma p_0)p'$. The superscripts $(\cdot)^{n+1/2,*}$ in (21) indicate effective time

averaged terms as they emerge from the chosen time integrator, and the asterisk indicates
quantities evaluated in the course of the predictor step.

Note, we have rewritten the ρg term in the momentum equation (21b) in terms of P and θ using the equation of state (given by (2) for the FC model and (15) for the $PI_{\rho,p}$ and $PI_{\rho,p}^{tc}$ models) where in the pseudo-incompressible cases $P^{n+\frac{1}{2},*} \equiv P_0$. In the compressible case, in agreement with second order accuracy we use $P^{n+\frac{1}{2},*} = P^n + \frac{1}{2}\delta p \left(\partial P/\partial p\right)$, where δp here is the pressure increment computed in the correction step of the previous time loop. The derivative of P with respect to p is computed using the equation of state.

By writing ρg in this way we were able to decouple the buoyancy term from the small advective flux divergence errors that arise in the predictor step. Potential temperature effects can fully be accounted for in the predictor, because potential temperature is accurately advected and not affected by the divergence errors. However, the pressure does react to divergence errors. By relying on accurate pressure information computed during the previous time steps, the buoyancy term is shielded from this effect. As a result, this formulation was found to give models increased stability for larger time steps.

We have used the following symbolic notation to abbreviate the balance of a numerical flux, say **q**, across grid cell boundaries,

$$\widetilde{\nabla} \cdot \mathbf{q}_C = \frac{1}{|C|} \sum_{I \in \mathcal{I}_C} \mathbf{q}_I \cdot \mathbf{n} = \frac{1}{|C|} \oint_{\partial C} \mathbf{q} \cdot \mathbf{n} \, d\ell + O(\Delta x^2) \,. \tag{22}$$

Here ∂C is the boundary of cell C. See Appendix for further details on the numerical scheme used in the predictor.

Note that we discretize advection by considering $P\mathbf{v}$ as the carrier flux that transports (upwind) values of the advected quantities $(1/\theta, \mathbf{v}/\theta, 1)$. This has turned out to be advantageous in many respects, e.g., in the construction of a positivity preserving advection scheme in Klein (2009) (see also Smolarkiewicz et al. (2014) and references therein).

We consciously refrain from going into more detail here because many different combinations of second-order accurate finite volume space discretizations and time integrators can more or less interchangeably be employed for the predictor step, provided they are used in conjunction with a well-balanced discretization of the pressure-gradient and gravity terms, see, e.g., Botta et al. (2004); Klein (2009). The details of the scheme used to generate the results of section 4 are given in the Appendix.

346 At the end of the predictor step,

- the scalar variables ρ , θ and P are second-order accurate (Klein 2009),
- the advecting fluxes $(P\mathbf{v})^{n+1/2}$ do not comply with the divergence constraint for $\alpha = 0$, and they do not provide a *stable* update of P for $\alpha > 0$, and
 - using the old time level pressure in the momentum equation (21b) prevents the scheme from being fully second-order accurate.

Crucially, for all values of α the time step used is limited by a CFL stability condition
(Courant et al. 1928) independent of sound speed (see Appendix), so that we sidestep the
stiffness induced by sound waves.

55 Step 2: First Correction

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The first correction step, which is the first of two linearly implicit substeps, corresponds to the MAC-projection in projection methods for incompressible flows (Bell et al. 1991). The advecting fluxes $P\mathbf{v}$ used in the predictor step do not abide by a semi-implicit discretization of the P equation for the FC model and by the divergence constraint for the $PI_{\rho,p}$ and $PI_{\rho,p}^{tc}$ models. In the first correction, an elliptic equation for a cell-centered pressure update $\delta p = p^{n+1} - p^n$ is derived by approximating (18c) at the half time level $t^{n+1/2}$, i.e., by reconsidering

$$\left[\alpha \left(\frac{\partial P}{\partial t}\right) + \nabla \cdot (P\mathbf{v})\right]^{n + \frac{1}{2}} = 0.$$
 (23)

The predictor step is discretized with second-order accuracy in time. As a consequence, the advecting fluxes $(P\mathbf{v})^{n+1/2,*}$ already include a first-order accurate update to the half time

level according to the auxiliary equation system (19), and this is sufficient to maintain secondorder accuracy for advection. Yet, for stability reasons an implicit correction is added that accounts for the influence of the new time level pressure gradient in the momentum equation in the following form (Klein 2009):

$$(P\mathbf{v})^{n+\frac{1}{2}} = (P\mathbf{v})^{n+\frac{1}{2},*} - \frac{\Delta t}{2} \theta^{n+\frac{1}{2},*} \nabla \delta p.$$
 (24)

Again, the asterisk denotes predicted values. Since $\Delta t \, \delta p = \Delta t \, (p^{n+1} - p^n) = O((\Delta t)^2)$,
this correction does not affect the second-order accuracy of advection. For $\alpha \neq 0$, the time
derivative term is transformed as:

$$\left(\frac{\partial P}{\partial t}\right)^{n+1/2} = \left(\frac{\partial P}{\partial p}\frac{\partial p}{\partial t}\right)^{n+1/2} = \left(\frac{\partial P}{\partial p}\right)^{n+1/2,*} \frac{\delta p}{\Delta t} + O\left((\Delta t)^2\right).$$
(25)

Using (24) and (25) in (23) we obtain the elliptic problem for any $\alpha \in [0, 1]$,

$$-\alpha \left(\frac{\mathcal{C}_{H}^{n+\frac{1}{2},*}}{\Delta t} \delta p \right)_{C} + \widetilde{\nabla} \cdot \left(\frac{\Delta t}{2} \, \theta^{n+\frac{1}{2},*} \, \nabla \delta p \right)_{C} = \widetilde{\nabla} \cdot \left((P\mathbf{v})^{n+\frac{1}{2},*} \right)_{C}, \tag{26}$$

373 where

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$$C_H^{n+1/2,*} = \left(\frac{\partial P}{\partial p}\right)^{n+1/2,*}.$$
 (27)

Expression (26) is responsible for determining stable time increments of P in the compressible model ($\alpha = 1$), whereas it enforces the divergence constraint for $\alpha = 0$.

With the solution of (26) δp at hand, the advecting flux corrections read

$$\delta P \mathbf{v} \cdot \mathbf{n} = -\frac{\Delta t}{2} \theta \nabla \delta p \cdot \mathbf{n}, \tag{28}$$

and the predicted values are corrected by,

$$\rho_C^{n+1} = \rho_C^{n+1,*} - \Delta t \, \widetilde{\nabla} \cdot (\delta P \mathbf{v} \, \theta^{-1})_C,
(\rho \mathbf{v})_C^{n+1,**} = (\rho \mathbf{v})_C^{n+1,*} - \Delta t \, \widetilde{\nabla} \cdot (\delta P \mathbf{v} \circ \mathbf{v} \theta^{-1})_C,
P_C^{n+1} = P_C^{n+1,*} - \Delta t \, \widetilde{\nabla} \cdot (\delta P \mathbf{v})_C.$$
(29)

where the advected variables θ^{-1} and $\mathbf{v}\theta^{-1}$ are evaluated at $(\cdot)^{n+1/2,*}$. The second asterisk indicates that the obtained value of the momentum is due to receive a second correction as described below.

Note that (26) turns into a standard Poisson pressure projection equation for the pseudoincompressible cases when $\alpha = 0$. In these cases, the correction of P in (29) automatically
yields $P^{n+1} \equiv P_0$ up to the tolerance in the divergence term with which the Poisson equation
was solved. Thus, in the pseudo-incompressible cases, the pressure variable P is restored to
its background value as a result of the first correction as it should be.

Thus far we have stabilized the advecting fluxes by incorporating an implicit pressure gradient contribution. We have not yet corrected the first-order error committed in the predictor step for the momentum equation by using the old time level pressure. This task is left to the second correction.

Step 3: Second Correction

The use of the old time level pressure in the momentum equation (21b) makes the predictor step first order accurate w.r.t. momentum. In a second correction step, the pressure and the momentum flux are corrected to achieve second-order accuracy and stability. Suppose we have already calculated an appropriate pressure update $\delta p = p^{n+1} - p^n$, then the correction of momentum reads

$$(\rho \mathbf{v})_C^{n+1} = (\rho \mathbf{v})_C^{n+1,**} - \frac{\Delta t}{2} \left(\widetilde{\nabla} \cdot (\delta p \mathbf{I})_C + \mathbf{k} \sigma \, \delta p \right) , \tag{30}$$

396 where

$$\sigma = (1 - \alpha) \beta \frac{g\rho_0}{\gamma p_0}. \tag{31}$$

Interpolating δp as computed in the first correction from the cell centers to the cell interfaces and using these data to evaluate (30) turns out to generate an unstable update. We avoid this by solving a second elliptic problem for a node-centered pressure variable (see similar procedures in Almgren et al. (1998); Schneider et al. (1999); Klein (2009); Vater and Klein (2009)). To derive the second elliptic equation, we multiply (30) by θ^{n+1} taking into account that the scalars ρ , P, θ have already attained their final values after the first correction and are unchanged in the second one. This yields

$$(P\mathbf{v})_C^{n+1} = (P\mathbf{v})_C^{n+1,**} - \frac{\Delta t}{2} \theta_C^{n+1} \left(\widetilde{\nabla} \cdot (\delta p \, \mathbf{I})_C + \mathbf{k} \sigma \, \delta p \right) . \tag{32}$$

404 As in the first correction we insert (32) into

$$\alpha \left(\frac{\partial P}{\partial t}\right)^{n+1/2} + \nabla \cdot \left(\frac{2-\alpha}{2} \left(P\mathbf{v}\right)^{n+1} + \frac{\alpha}{2} \left(P\mathbf{v}\right)^{n}\right) = 0, \tag{33}$$

where, for $\alpha=1$, a second-order accurate midpoint discretization with no off-centering is considered. After node-centered space discretization of the divergence, we obtain the elliptic problem:

$$-\alpha \left(\frac{\mathcal{C}_{H}^{n+1}}{\Delta t} \delta p\right)_{\overline{C}} + \widetilde{\nabla} \cdot \left(\frac{(2-\alpha)\Delta t}{4} \theta^{n+1} \left(\nabla \delta p + \mathbf{k} \sigma \delta p\right)\right)_{\overline{C}} = \widetilde{\nabla} \cdot \left(\frac{2-\alpha}{2} (P\mathbf{v})^{n+1,**} + \frac{\alpha}{2} (P\mathbf{v})^{n}\right)_{\overline{C}}, \quad (34)$$

where \mathcal{C}_H^{n+1} is defined by (27) using the corrected value of P.

As in the first correction, we obtain a Helmholtz equation for $\alpha = 1$ where the zero-order term accounts for compressibility. The difference between FC ($\alpha = 1$) and PI^{tc}_{ρ,p} ($\alpha = 0$) is a modified structure of the system matrix.

We note that in the fully-compressible case a backward difference (BDF2) discretization can be used, as done in Vater (2013). In that case, and for $\alpha = 1$, (34) is replaced with

$$-\left(\frac{3C_H^{n+1}}{2\Delta t}\delta p\right)_{\overline{C}} + \frac{2}{3}\Delta t \,\widetilde{\nabla} \cdot \left(\theta^{n+1}\widetilde{\nabla}\delta p\right)_{\overline{C}} = \widetilde{\nabla} \cdot (P\mathbf{v})_{\overline{C}}^{n+1} - \left(\frac{C_H^{n+1}}{2\Delta t}\delta p^{\text{old}}\right)_{\overline{C}}, \quad (35)$$

where $\delta p^{\text{old}} = p^n - p^{n-1}$ denotes the old time level pressure increment.

A nine-point stencil is used for the discretization of the laplacian (34) or (35), which is obtained as follows: the nodal values define continuous piecewise bilinear pressure distributions on the primary control volumes. We integrate their gradients analytically over the boundaries of the dual cells that are centered on the grid nodes. The solution δp is accordingly defined in the centers of the dual cells, \overline{C} . Straightforward numerical integration of pressures over the primary cell interfaces can thus be employed in evaluating the second

momentum correction in (30). After the nodal pressures have been updated to the new time 421 level as well, all variables are now second-order accurate and ready for the next time step. 422 See details of the discretization in the Appendix. 423

Numerical Results 4.

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In this section, we present the results of the simulations performed with our semi-implicit method. The aim is to show that the model numerics produces results in agreement with its 426 theoretical properties in different configurations. First, a convergence study in the FC config-427 uration is presented. Then, results with fully-compressible (FC) and pseudo-incompressible 428 $(PI_{\rho,p})$ models are compared on simulations of thermal perturbations. The impact of the 429 thermodynamic consistency $(PI_{\rho,p}^{tc})$ term is also evaluated. 430 The numerical model is implemented in an object oriented C++ environment based on 431 the SAMRAI framework for mesh refinement (Hornung et al. 2006). Krylov-type methods 432 with algebraic multi-grid preconditioners as included in the Hypre library (Falgout et al. 433 2006) are used to solve the linear systems in the correction step. Our coding framework is 434 fully parallelized and 3d-ready. However, an extensive analysis of its parallel efficiency lies 435 outside the scope of the present work.

Convergence study 437

First, we assess the accuracy properties of the FC model on a case of pure transport 438 in a highly idealized setting with g = 0. The case (Kadioglu et al. 2008) consists of a 439 travelling rotating vortex in the doubly periodic unit-square-shaped domain $\Omega = [0, 1]^2$ m². The vortex is axisymmetric and rotates counterclockwise with unitary velocity. Density is modelled by a smooth, non-constant function and a constant and a unitary transport velocity $\mathbf{v} = (1,1)^T \,\mathrm{m\,s^{-1}}$ is superimposed. The vortex is an exact solution for the zero 443 Mach number incompressible equations, to which $\mathrm{PI}_{\rho,p}^{\mathrm{tc}}$ and $\mathrm{PI}_{\rho,p}$ reduce in the absence of gravity (Klein 2009). With the pressure field correctly initialized, it is an exact solution for the fully-compressible equations as well. We refer to Kadioglu et al. (2008) for the initial data not reported here for brevity. Note that some of the coefficient in the expression for initial pressure were incorrectly reported in Kadioglu et al. (2008), the correct expression is available upon request.

In the compressible case, the initial distribution for P is derived via the equation of state (3). Reference physical quantities are set as follows:

$$\rho_{\text{ref}} = 0.5 \text{ kg m}^{-3}, \ p_{\text{ref}} = 101625 \text{ Pa}, \ T_{\text{ref}} = 706.098 \text{ K},$$
(36)

corresponding to a maximum Mach number $M_{\text{max}} = \max(\|\mathbf{v}\|_{\text{RMS}}/\sqrt{\gamma p/\rho}) = 4.96\text{E-}03$. The high value of T_{ref} is computed from p_{ref} and ρ_{ref} considered in Kadioglu et al. (2008) and enables an easier comparison with their results for the density.

The flow is simulated by running the FC semi-implicit model ($\alpha \equiv 1$) on a grid with 192 cells in both directions at CFL = 0.45, that is, constant $\Delta t = \Delta t_{\rm A} = 9.7 \text{E-}04 \text{ s}$ and $\Delta x = 5.21 \text{E-}03 \text{ m}$. These data correspond to a sound-speed based CFL_S = CFL/ $M_{max} \approx 90.72$.

The vortex is transported by the background unitary velocity. Due to the doubly periodic boundary, the initial configuration is reproduced unchanged at time T=1 s (figure 2). Similar results (not shown) are obtained for momentum and P in FC runs and for all variables except for P (which is constant) in $PI_{\rho,p}^{tc}$ runs.

Furthermore, the numerical solution converges quadratically in the maximum norm (Figure 3). The experimental order of accuracy is in agreement with the theoretical accuracy of the scheme presented in Section 3. Similar results are obtained with $PI_{\rho,p}^{tc}$ runs (not shown). The FC results shown above validate the use of the fully-compressible flow solver that

extends the pseudo-incompressible framework of Klein (2009).

467 Rising bubble

Next, we consider a warm air bubble test case in the domain $\Omega = (x, z) \in [-10, 10] \times [0, 10]$ km². We set the following initial data for a homentropic atmosphere (Botta et al. 2004):

$$p(z) = p_{\text{ref}} \left(1 - \Gamma \frac{g\rho_{\text{ref}}}{p_{\text{ref}}} z \right)^{\frac{1}{\Gamma}}, \quad \rho(z) = \rho_{\text{ref}} \left(\frac{p(z)}{p_{\text{ref}}} \right)^{\frac{1}{\gamma}}, \quad \rho_{\text{ref}} = \frac{p_{\text{ref}}}{RT_{\text{ref}}}, \tag{37}$$

where, in agreement with Klein (2009), $\rho_{\rm ref}$, $p_{\rm ref}$, g, and $T_{\rm ref}$ have the values 1 kg m⁻³, 8.61E04 N m⁻², 10 m s⁻², and 300 K, respectively, and $\Gamma = (\gamma - 1)/\gamma$. The background potential temperature θ is constant. The homentropic setting (37) is perturbed with a smoothed cone-shaped thermal perturbation θ' , given by (Klein 2009):

$$\theta'(x,z) = \begin{cases} \delta\theta \cos^2(\frac{\pi}{2}r) & (r \le 1) \\ 0 & \text{otherwise} \end{cases}, \qquad \begin{cases} \delta\theta = 2 \text{ K} \\ r = 5\sqrt{(\frac{x}{L})^2 + (\frac{z}{L} - \frac{1}{5})^2} \\ L = 10 \text{ km} \end{cases}$$
(38)

The initial velocity is zero. Lateral boundary conditions are periodic, with solid walls on top and bottom boundaries.

We run our semi-implicit trapezoidal scheme on a grid with $\Delta x = \Delta z = 125$ m, i.e. 160×80 cells, and CFL = 0.5. In the first five steps a buoyancy-driven time step ($\Delta t = \Delta t_{\rm B} \approx 21.69$ s) is used. Due to growing velocities, the advection-driven time step is used for the remainder of the simulation. Towards the end of the simulation, values of $\Delta t \approx 4.6$ s are attained.

Driven by buoyancy, the warm bubble rises and rolls up on the sides (figure 4). The
amplitude of the thermal perturbation at final time T = 1000 s is in agreement with the
results in Klein (2009), as shown in table 2. However, the $PI_{\rho,p}$ bubble rises faster, is not as
wide and exhibits a phase shift with respect to both the $PI_{\rho,p}^{tc}$ and the FC models (figure 5).

The discrepancies in the $PI_{\rho,p}$ model come from neglecting the effect of pressure perturbations on the buoyancy. The extra buoyancy term present in the $PI_{\rho,p}^{tc}$ model reduces
buoyancy near the top of the bubble due to an increase in pressure near the bubble top and

increases buoyancy at the two tails due to a pressure decrease near the tails. Furthermore, the overall buoyancy of the bubble decreases causing a decrease in the phase speed. Therefore the $PI_{\rho,p}^{tc}$ bubble is both lower and wider than the $PI_{\rho,p}$ model and, as a result, resembles the FC model more closely.

Results with $PI_{\rho,p}^{tc}$ as measured in a one-dimensional cut of θ' at height z=7500 m match the FC results within a 2 per cent error (table 3).

Results with the PI_{ho,p} model do not differ substantially from FC results at the end of the simulation at T=1000 s. The different dynamics of the FC case can be detected in the onset of sound waves in the initial stages of the simulation. With the FC model ($\alpha=1$) the initial potential temperature perturbation triggers acoustic waves. These are visible in the upper left panel of Figure 6, which displays pressure increments at time t=26.6 s in a run of the FC model with $\Delta t = \Delta t_{\rm I} = 1.9$ s. The oscillations are due to the initial hydrostatic pressure distribution from (37) not being acoustically balanced.

The presence of associated pressure oscillations is confirmed by a time series over the first 350 s of the pressure time increment values recorded at the point (x, z) = (-7.5, 5) km marked with a cross in the upper left panel of Figure 6. The time series are shown in the upper right, lower left and lower right panels of Figure 6. The upper right and lower left plots are relative to simulations at constant $\Delta t = \Delta t_{\rm I} = 1.9$ s. The simulation relative to the lower right panel is at CFL ≈ 0.5 as in Figure 4.

FC model results (solid lines in all plots) display oscillations triggered by the initial pressure imbalance. The amplitude of the acoustic oscillations in the small time step case (upper right panel) is ninefold the amplitude of the large time step runs (lower right panel). The effect is suppressed in the $PI_{\rho,p}$ runs (dashed lines) except for an initial transient. Note that in the large-time step run the initial transient masks the amplitude of the acoustics. Therefore, the data of the first time step was removed in the lower right panel of Figure 6. In the case of the $PI_{\rho,p}$ model, pressure is determined by the solution of a time-independent Poisson problem, which describes the pressure field in the absence of sound waves. $PI_{\rho,p}$ is

considered here because the extra $PI_{\rho,p}^{tc}$ term does not modify the results as far as acoustics are concerned. On the one hand, the reduction in the amplitude of the large time step acoustic oscillations shows that the semi-implicit method is able to handle acoustic oscillations at CFL numbers independent of the sound speed. On the other hand, the effect of acoustics is not completely suppressed in the large-time step, either.

However, thanks to the blending feature, the code is able to continuously transition from the PI_{ρ ,p} configuration to the FC configuration. The lower left panel of Figure 6 shows the time series of pressure increments for blended runs. We set the transition parameter α from section 2 to zero for S_1 time steps. Then, α increases linearly to $\alpha = 1$ over S_2 time steps. Starting at the time step number $S_1 + S_2$, the code runs compressibly with $\alpha = 1$.

In the lower left panel of Figure 6, the thin solid line in the background denotes the 526 fully-compressible run. The dashed-dotted curve and thick solid curves were obtained with 527 $S_2 = 20$ and $S_2 = 40$, respectively. There are no disturbances for the first $S_1 = 10$ pseudo-528 incompressible steps in these two pressure graphs, and the results coincide with those from 529 the run of the $PI_{\rho,p}$ model (dashed line in the right panels). Perturbations arise in the 530 transitional period and fully develop after $S_1 + S_2$ time steps. The oscillations' amplitudes 531 in the blended runs are considerably lower than those of the FC run and they are lower for 532 the larger S_2 value, i.e. the longer transitional period. 533

Results in the lower left panel of figure 6 demonstrate the capabilities of the blended model. Acoustic perturbations are absent when the model runs in pseudo-incompressible mode with $\alpha = 0$ and they emerge significantly damped after the transition to $\alpha = 1$ in fully-compressible mode. Therefore, when blended continuously with the compressible discretization, the soundproof limit discretization can be used to actively control imbalances in the initial data. The oscillation amplitudes are substantially reduced also when larger time steps are employed as seen in the lower right panel of figure 6.

Finally, as in Almgren et al. (2006a), which presents a pseudo-incompressible code for stellar hydrodynamics, we compare plots of the Mach number in the initial stages of FC, PI_{ρ ,p} and blended runs. Results at time t=21.66 s, that is, time step number 57 at $\Delta t = \Delta t_{\rm I} = 0.38$ s, are displayed in Figure 7. The mushroom-shaped FC result (left panel) reveals the initial onset of sound waves due to pressure imbalances already inspected in Figure 6, while the PI_{ρ ,p} plot (middle panel) and blended plot (right panel) show no perturbation away from the bubble. A very small time step was considered in this case following Almgren et al. (2006a) in order to track more closely the dynamics in the initial stages.

Density current

This test (Straka et al. 1993) consists of a negative potential temperature perturbation in a $[-25.6, 25.6] \times [0, 6.4]$ km² homentropic atmosphere (37),

$$T' = \begin{cases} 0 & \text{K} & \text{if } r > 1 \\ -15 \left[1 + \cos(\pi r) \right] / 2 & \text{K} & \text{if } r < 1 \end{cases}$$
(39)

where $r=\{[(x-x_c)/x_r]^2+[(z-z_c)/z_r]^2\}^{0.5},\ x_c=0$ km, $x_r=4$ km, $z_c=3$ km and $z_r=2$ km. From $\theta=T(p/p_{\rm ref})^{-\Gamma}$ we derive the potential temperature perturbation and density distribution,

$$\theta'(x,z) = \frac{T'}{1 - \Gamma \frac{g\rho_{\text{ref}}}{p_{\text{ref}}}z}, \qquad \rho(z) = \rho_{\text{ref}} \left(\frac{p(z)}{p_{\text{ref}}}\right)^{\frac{1}{\gamma}} \frac{\theta_{\text{ref}}}{\theta_{\text{ref}} + \theta'}, \tag{40}$$

where $\theta_{\rm ref} = T_{\rm ref}$. The boundary conditions are periodic on the left and right boundary, solid walls on the top and bottom boundary. Furthermore, we add an artificial diffusion term $\rho\mu\nabla^2\mathbf{v}$ to the right hand side of the momentum equation ($\rho\mu\nabla^2\theta$ in the P equation), with $\mu = 75 \text{ m}^2 \text{ s}^{-1}$ as in Straka et al. (1993). The initial velocity is set to zero, and the reference quantities are $T_{\rm ref} = 300 \text{ K}$, $p_{\rm ref} = 10^5 \text{ Pa}$, $\rho_{\rm ref} = p_{\rm ref}/(RT_{\rm ref})$.

The models are run with $\Delta x = 50$ m and CFL = 0.5. Thus, the time step is $\Delta t = \Delta t_{\rm B} \approx$ 4.65 s for the first three steps and then the advective time step is used. For the FC model, a backward difference approach in the second projection is used, see equation (35). Due to the symmetrical nature of the test case, only the plots for the subdomain $[0, 19.2] \times [0, 4.8]$ km² are shown.

Obtained values of the final thermal perturbation and the front positions as calculated by the FC and PI^{tc}_{ρ,p} models (Figure 8 and table 4) are in line with results in the literature (Straka et al. 1993; Restelli and Giraldo 2009). In contrast to the rising bubble case, the extra buoyancy term in the PI^{tc}_{ρ,p} model results in an overall increase in the buoyancy of the bubble. This increase in buoyancy causes the bubble to fall slower and reduces the phase speed when compared with the PI_{ρ,p} model. This can be seen in the farther front position and in the horizontal cut at height z = 1200 m (Figure 9) of the PI_{ρ,p} model when compared to both the FC and PI^{tc}_{ρ,p} models. As a result, the PI_{ρ,p} model displays considerable deviations (higher than 40 per cent) relative to FC runs (Table 5). For the PI^{tc}_{ρ,p} model, the deviation from FC is lower than 5 per cent.

575 Inertia-gravity waves

Next, we consider a thermally stratified atmosphere with stable stratification of potential temperature $\partial\theta/\partial z>0$. In particular, as in Restelli and Giraldo (2009); Skamarock and Klemp (1994), we take:

$$\theta(z) = T_{\text{ref}} \exp\left(\frac{N^2}{g}z\right),$$
(41)

where N denotes the buoyancy frequency. With $N=0.01~\rm s^{-1},~g=9.81~\rm m\,s^{-2},~and~T_{ref}=300~\rm K,$ we have $\theta\in[300,~332.19]~\rm K$ for $z\in[0,~10]~\rm km$. The other variables are defined as:

$$p(z) = p_{\text{ref}} \left\{ 1 - \frac{g}{N^2} \Gamma \frac{g\rho_{\text{ref}}}{p_{\text{ref}}} \left[1 - \exp\left(-\frac{N^2 z}{g}\right) \right] \right\}^{\frac{1}{\Gamma}}, \tag{42}$$

$$\rho(z) = \rho_{\text{ref}} \left(\frac{p(z)}{p_{\text{ref}}}\right)^{\frac{1}{\gamma}} \exp\left(-\frac{N^2 z}{g}\right), \qquad \rho_{\text{ref}} = \frac{p_{\text{ref}}}{R T_{\text{ref}}}, \tag{43}$$

with $p_{\text{ref}} = 10^5$ Pa. On top of the background stratification (41)–(42), in a $[0, 300] \times [0, 10]$ km² domain we consider the perturbation (Skamarock and Klemp (1994) and Figure 10 left panel):

$$\theta'(x, z, 0) = 0.01 \text{ K} * \frac{\sin(\pi z/H)}{1 + [(x - x_c)/a]^2}$$
(44)

with H=10 km, $x_c=100$ km, a=5 km. In addition, there is a background horizontal 584 flow $u = 20 \text{ m s}^{-1}$. The simulations are performed with at advective CFL = 0.3, that is 585 $\Delta t = \Delta t_A \approx 3.75$ s. The grid spacing is $\Delta x = \Delta z = 250$ m and the trapezoidal time 586 integrator is employed for the FC model. In agreement with published work (Restelli and 587 Giraldo 2009), the Coriolis term is neglected here because of the small length of the channel. 588 Unlike the previous test cases, here the dynamics is chiefly wavelike rather than vertically 589 buoyancy-driven. Inertia-gravity waves develop in the horizontal direction (Figure 10). As 590 in the previous test case, only the FC contour plots are presented in Figure 10 as the $PI_{\rho,p}^{tc}$ 591 and $PI_{\rho,p}$ plots are visually indistinguishable. 592

A quantitative comparison between the FC, $PI_{\rho,p}^{tc}$ and $PI_{\rho,p}$ results and the results of 593 Restelli and Giraldo (2009) is reported in table 6. Maxima and minima of perturbations of 594 velocity components, potential temperature and Exner pressure at final time T=3000 s are 595 in line with published work. 596

The left panel of Figure 11 shows a one-dimensional cut of the potential temperature 597 perturbation at z = 5000 m. As in the previous cases, the $\text{PI}_{\rho,p}$ model displays a higher 598 phase speed than the $PI_{\rho,p}^{tc}$ and FC models due to the neglect of pressure perturbations in 599 the buoyancy term. The region of the leftmost crest is magnified in Figure 11 to highlight 600 the difference in the phase speed of the $PI_{\rho,p}$ model (dashed-dotted line) with respect to the 601 $PI_{\rho,p}^{tc}$ model (starred markers) and the FC model (solid line). 602

The right panel of Figure 11 shows the differences between the FC cut and the $PI_{\rho,p}^{tc}$ cut (dashed line) and between the FC cut and the $\text{PI}_{\rho,p}$ cut (solid line). The amplitude of the 604 difference is larger in the latter case due to the phase shift highlighted on the left panel. The 605 result is quantified in Table 7 which shows relative RMS and max errors of the FC cut with 606 respect to the $PI_{\rho,p}^{tc}$ and $PI_{\rho,p}$ cuts. Relative $PI_{\rho,p}$ -FC errors are threefold the $PI_{\rho,p}^{tc}$ -FC ones 607 608

Finally, as in Restelli and Giraldo (2009) we define conservation errors as:

$$C_{\phi} = \frac{|\left(\phi_{\text{tot}}\right)_{T} - \left(\phi_{\text{tot}}\right)_{0}|}{\left(\phi_{\text{tot}}\right)_{0}},\tag{45}$$

where $\phi_{\text{tot}} = \int_{\Omega} \phi \, d\mathbf{x}$ denotes the volumetric integral of ϕ in the domain Ω . Subscripts 0

and T denote initial and final time, respectively. We expect our scheme to conserve density ρ and horizontal momentum density ρu . Though our model does not conserve total energy ρE , we report conservation scores for that variable, too. For the FC model, results for P are also reported. Values of the conservation error for ρ , ρu , P, and ρE are fairly low for the three model configurations (table 8). Note, in table 8 we define the total energy variable as

$$E = \frac{1}{\rho} \frac{p}{\gamma - 1} + \frac{\mathbf{v}^2}{2} + gz. \tag{46}$$

where $p = p_0$ in (46) for the $PI_{\rho,p}^{tc}$ and $PI_{\rho,p}$ cases as shown in Klein and Pauluis (2011). Numerical analysis of the P-conservation is only meaningful for the FC model, since in the incompressible cases $P = P_0(z)$ holds.

We have presented a blended weakly compressible computational model with seamless

access to thermodynamically consistent pseudo-incompressible dynamics, these two repre-

senting the limiting cases of a family of models depending on one parameter. For each

5. Discussion and conclusions

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member of the model family, the numerical discretization is the same up to certain weights in the stencil of the implicit corrector invoked to enable advection-based time steps in simulations of small to mesoscale systems. 624 This seamless and straightforward compressible-to-soundproof model transition can be 625 realized in any flow solver that features the density and the mass-weighted potential temper-626 ature as prognostic variables for the thermodynamics, together with flux-based formulations 627 of their determining equations. Weak checkerboard modes were observed in the runs of 628 gravity-driven flows for very small time steps. We attribute them to the fact that the diver-629 gence of the cell-centered velocity is controlled in the second correction through a discrete 630 elliptic problem derived from the linearized acoustic equations on the Arakawa B-Grid with 631 a standard stencil. This grid arrangement allows for oscillatory modes with phase vectors 632 pointing roughly along the grid diagonals (see Figure 8 of Arakawa and Lamb (1977)). These 633

modes might be controllable by adopting a staggered grid arrangement (Arakawa C-grid) or by adopting an inf-sup stable discretization of the elliptic operator on the B-Grid as in Vater and Klein (2009).

The key observation enabling the blending is that, at least for an ideal gas with constant specific heat capacities, $\rho\theta$ is a function of pressure alone. Thus the transport equation for $\rho\theta$ is equivalent to the pressure evolution equation and lends itself naturally for implicit pressure formulations. Once available, such a seamless framework can be used, e.g., for a clean comparison of compressible and soundproof models that is not affected by sizeable differences between the respective model discretizations (see Smolarkiewicz and Dörnbrack (2008); Smolarkiewicz et al. (2014) for comparable arguments).

As a further potentially attractive application of such a modeling tool we suggest the 644 filtering of unbalanced initial data. For given initial data, a matching pressure field and a 645 related divergence correction that would guarantee a nearly sound-free subsequent evolution 646 are generally not available. With a blended soundproof-compressible framework, one can 647 generate accurate balanced pressure and velocity fields by running the model in soundproof 648 mode for a few time steps and then making the transition to fully-compressible over another 649 few steps. This idea may also be transferred to other nearly balanced situations, such as 650 hydrostatic and geostrophic, but exploring this is left for future work. In the framework 651 of techniques for atmospheric data assimilation (Rabier 2005), the resulting ability of a 652 computational model to manage and regularly embed new, unbalanced input in a balanced fashion and without invoking additional filtering procedures appears quite attractive. This 654 capability can also be more generally useful when one has to map externally obtained data 655 into a multi-dimensional finite volume scheme as analyzed in Zingale et al. (2002). 656

Besides the aforementioned blending features, there are other noteworthy aspects of the scheme. First, we discretize the equations in full form without subtraction of a background state, maintaining accuracy by adopting a well-balanced discretization of the pressure gradient and gravity terms as discussed in Botta et al. (2004); Klein (2009). Second, we cast the momentum equation in terms of pressure and density instead of the more common Exner pressure and potential temperature. The former choice guarantees conservation of momentum in the absence of external forces and increases flexibility with a view to implementing more general equations of state (Klein and Pauluis 2011).

Code performance was assessed in a number of configurations. The second-order accuracy
of the scheme was verified on a smooth benchmark. Then, standard test cases consisting
of buoyant thermal perturbations were considered, where our data confirmed no substantial difference between the compressible and pseudo-incompressible results. For the latter,
including the linearized effect of pressure on density in the gravity term results not only in
thermodynamic consistency (Klein and Pauluis 2011) but also in improved accuracy. Our
findings are consistent with Davies et al. (2003); Klein et al. (2010), thus confirming the
validity of the pseudo-incompressible model at small to mesoscales and for realistic stratifications.

As mentioned, we are planning to extend the present general stategy to include addi-674 tional dominant balances relevant for larger scale flows, specifically to the hydrostatic and 675 geostrophic limits. This goal appears feasible in view of recent related work. For example, 676 successful results have been achieved by EULAG model users (Prusa and Gutowski 2011; 677 Szmelter and Smolarkiewicz 2011; Smolarkiewicz et al. 2014) with compressible, anelastic, 678 and pseudo-incompressible models on the synoptic and planetary scales. Furthermore, al-679 ternatives have been explored to merge hydrostatic models with fully-compressible (Janjic et al. 2001) or soundproof ones. Careful consideration will be needed to identify the correct 681 large-scale limiting model in the light of recent suggestions of unified multiscale reduced 682 models by Durran (2008) and Arakawa and Konor (2009); Konor (2014). 683

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Details of the numerical scheme

Here we elaborate on the aspects of the numerical scheme omitted in the main text.

701 Predictor

We use a second-order accurate, explicit two-stage strong stability-preserving Runge-Kutta method for time integration (Gottlieb et al. 2001). For the Ordinary Differential Equation:

$$\frac{du}{dt} = L(u),\tag{A1}$$

where L denotes a differential operator, the method reads:

$$u^{(1)} = u^n + \Delta t L(u^n), \tag{A2}$$

$$u^{n+1} = \frac{1}{2}u^n + \frac{1}{2}u^{(1)} + \frac{1}{2}\Delta t L(u^{(1)}), \tag{A3}$$

where $u^{(1)}$ denotes the first stage solution.

The spatial discretization is performed with a finite volume approach, see, e.g., LeVeque (2002). Discrete variables are defined as approximations of cell averages set at cell centers, with the exception of dynamic pressure, set at cell nodes. The new cell-centered values are obtained from the old ones subtracting the net outflow flux at the boundaries and adding the contribution from the source term, expressions (21a)–(21b)–(21c) in the main text.

The discretization of the fluxes is performed according to the following steps:

712 i. The velocity at the interfaces is determined by averaging the neighbouring leftmost and rightmost cell-centered values \mathbf{v}_L and \mathbf{v}_R :

$$\mathbf{v} = \frac{1}{2}(\mathbf{v}_L + \mathbf{v}_R),\tag{A4}$$

where, for a second-order method, \mathbf{v}_L and \mathbf{v}_R have to be linearly reconstructed/limited.

Considering the interface $(x_{i+1/2}, y_j)$, and omitting the subscript j for simplicity, the reconstructed values of the horizontal velocity u are:

$$u_L = u_i + \frac{1}{2}\psi \left(u_i - u_{i-1}, u_{i+1} - u_i\right), \tag{A5}$$

$$u_R = u_{i+1} - \frac{1}{2}\psi \left(u_{i+1} - u_i, u_{i+2} - u_{i+1}\right), \tag{A6}$$

717 where:

$$\psi(a,b) = \frac{a+b}{2} \tag{A7}$$

for centered slopes. Our implementation features also an option for slope limiters, for which ψ would have a different functional form. Upwind fluxes F_P for the P variable are computed by means of the obtained velocity:

$$F_P = F_P^+ + F_P^-,$$
 (A8)

where:

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$$F_P^+ = P_L \max(\mathbf{v}, 0), \qquad F_P^- = P_R \min(\mathbf{v}, 0),$$
 (A9)

and the subscripts L and R denote cell-centered leftmost and rightmost values of the variable.

ii. Fluxes for the remaining quantities are referred to the carrier flux $P\mathbf{v}$ and derived using (A9) as

$$F_{\phi} = F_P^+ \, \phi_L + F_P^- \, \phi_R \tag{A10}$$

where $\phi \in \{1/\theta, \mathbf{v}/\theta\}$. The contribution from the pressure term is incorporated in the momentum flux adding the pressure value at the center of the cell interface, obtained via average of the adjacent nodal values.

729 Pressure update

730 The nodal pressure update at the end of the time step proceeds as follows:

i. An auxiliary cell-centered pressure p_c is computed from P using the inverse of the equation of state (2). The result is then interpolated to the nodes:

$$p_c^{n+1} = \left(\frac{P^{n+1,**}}{\rho_{ref}T_{ref}}\right)^{\gamma} p_{ref} - p_{ref}, \qquad p_c^{n+1} \longrightarrow p_{EOS}^{n+1}.$$
(A11)

733 ii. The obtained value is weighted with the old time level pressure update with the solution of (34) or (35), δp :

$$p^{n+1} = \alpha p_{EOS}^{n+1} + (1 - \alpha) (p^n + \delta p).$$
 (A12)

When the model runs in pseudo-incompressible mode with $\alpha = 0$, the node-centered pressure 735 increment δp is summed to the old time level value. In compressible mode, with $\alpha = 1$, the 736 new nodal pressure is locked to P imposing the equation of state at a discrete level. Other solutions are possible and were tested. For example, as a pseudo-incompressible 738 update, an interpolated value of the solution δp_c of the first correction equation (26) can be 739 summed to the old time level pressure value. This was used in the thermal perturbations 740 simulated with the fully-compressible model initially run in pseudo-incompressible mode. 741 In that case the solution of the second Poisson problem only serves as a correction to the 742 momentum flux, expression (30), not as an update for the nodal pressure value. 743

744 Time step choice

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The explicit time integration method adopted in the predictor step must be consistent with
the CFL stability condition for advection (Courant et al. 1928), and a similar constraint
for internal wave dynamics since these processes are handled explicitly in our scheme. In
particular, we dynamically compute the time step size at each time loop according to:

$$\Delta t = \min\left(\Delta t_{\rm I}, \Delta t_{\rm A}, \Delta t_{\rm B}\right) \tag{A13}$$

where $\Delta t_{\rm I}$ is an externally imposed value of the time step. $\Delta t_{\rm A}$ is the advective time step:

$$\Delta t_{\rm A} = \frac{\rm CFL} \Delta x}{\max_{\Omega} (\|\mathbf{v}\|_2)},\tag{A14}$$

where CFL ≤ 1 and $\|\cdot\|_2$ is the discrete L^2 norm. $\Delta t_{\rm B}$ is a buoyancy-dependent time step:

$$\Delta t_{\rm B} = \text{CFL} \sqrt{\frac{\Delta x \, \min_{\Omega} \theta}{g \, \max_{\Omega} \Delta \theta}},\tag{A15}$$

where $\max_{\Omega} \Delta \theta = \max_{\Omega} \theta - \min_{\Omega} \theta$ is the maximum potential temperature perturbation.

Dynamically adaptive time stepping is standard on computational fluid dynamics and for two time level schemes it's implementation is quite straightforward (LeVeque 2002).

Well-balanced treatment of vertical pressure gradient and gravity term

In the envisaged atmospheric applications, flow patterns arise as perturbations around a hydrostatically balanced state, where the vertical pressure gradient offsets the gravitational force

$$\frac{\partial p}{\partial z} = -\rho g. \tag{A16}$$

Therefore, an essential characteristic of a numerical method in this context is the capabil-758 ity of mimicking the hydrostatic balance at the discrete level. This means, for instance, that 759 the numerical discretization should introduce no perturbations on an initially motionless at-760 mospheric setting. The feature is especially nontrivial for models as the ones presented here 761 whose analytical formulation relies on full variables, unlike other non-hydrostatic fully com-762 pressible models (e.g., Skamarock and Klemp (2008); Restelli and Giraldo (2009)) wherein 763 the unknowns are themselves perturbations around a background hydrostatically balanced 764 reference state. 765

Here we adopt the approach of Botta et al. (2004), who describe the implementation of a discrete Archimedes' principle, and in the following we present the parts of our implementation tuned to take into account the hydrostatic balance. 769 Initialization

Since the problem is inherently one-dimensional, we focus on the vertical coordinate for the moment. First, let the initial data for pressure p(z) and density $\rho(z)$ be given in the form of a homentropic or stably stratified atmosphere as in expressions (37) or (42) above. Next:

- p(z) is initialized in cell centres z_j , $j=1,\ldots,\mathcal{N}_z$ and nodes $z_{j-1/2}$, $j=1,\ldots,\mathcal{N}_z+1$ according to its analytical expression (37) or (42);
- $\rho(z)$ is initialized at z_j using a discretized form of (A16), i.e.

$$\rho(z_j) = -\frac{1}{q\Delta z} [p(z_{j+1/2}) - p(z_{j-1/2})], \quad j = 1, \dots, \mathcal{N}_z.$$
(A17)

where Δz is the vertical grid spacing.

778 Predictor step

The value of the pressure at the center of the cell face needed for the momentum flux computation in expression (19b) is computed as follows:

$$p(z_j) = \frac{1}{2} \left\{ p(z_{j+1/2}) + p(z_{j-1/2}) - g \left[2f(z_j) - f(z_{j+1/2}) - f(z_{j-1/2}) \right] \right\}$$
(A18)

781 for $j = 1, ..., N_z$, where:

$$f(z) = \int_0^z \rho(z') dz' \tag{A19}$$

and the square bracket in (A18) represents a hydrostatic modification of the simple average.

 $Boundary\ conditions$

The so-called "solid wall" boundary conditions are adjusted to take into account hydrostatic balance. As customary in finite differences and finite volume codes (LeVeque 2002),

we implement fully reflecting boundaries using "ghost cells". The strategy involves attaching two dummy cells to the boundary in which the value of all the variables except for the
normal velocity is mirrored from the two innermost cells, whereas the normal velocity value
is taken with opposite sign.

We modify the process for the mirrored variables in that we retrieve in the ghost cells
the hydrostatically-consistent values. For instance, for the pressure in the first lower ghost
cell (cell 0) we have:

$$p_{z_0} = p(z_1) + g \int_{z_0}^{z_1} \rho(z) dz$$
 (A20)

and similar expressions hold for the upper values.

Final locking of pressure and P variables

The third modification involves the interpolation from nodes to cell centers or *vice versa*,
which in the case without gravity is a standard linear interpolation. Here, a correction taking
into account hydrostaticity is introduced. In particular, for the cell-to-node interpolation
used in the pressure update (A11) after the second correction step:

• For the lower boundary nodes:

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$$p(x_{i+1/2}, z_{1/2}) = 0.5(p_{NW} + p_{NE}), \quad \forall i = 1, \dots, \mathcal{N}_x$$
 (A21)

where p_{NW} and p_{NE} denote the pressure values obtained with analytical integration downwards from the hydrostatic pressure values in the adjacent upper left and upper right cell, respectively.

• For the upper boundary nodes:

$$p(x_{i+1/2}, z_{\mathcal{N}_z+1/2}) = 0.5(p_{SW} + p_{SE}), \quad \forall i = 1, \dots, \mathcal{N}_x$$
 (A22)

where p_{SW} and p_{SE} denote the pressure values obtained with analytical integration upwards from the hydrostatic pressure values in the adjacent lower left and lower right cell, respectively.

• For the internal nodes:

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$$p(x_{i+1/2}, z_{j+1/2}) = 0.25(p_{SW} + p_{SE} + p_{NW} + p_{NE}), \quad \forall i = 1, \dots, \mathcal{N}_x, \ j = 1, \dots, \mathcal{N}_z - 1$$
(A23)

Finally, we remark that issues due to neglect of hydrostatic balance at the discrete level manifest less in the incompressible than in the fully-compressible version of our method. In the former, small spurious perturbations due to inexact balancing, for instance, at the boundary are projected away in the correction step, while in the latter P and pressure are locked through the equation of state, thus requiring a careful adjustment.

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Model name	Abbreviation	(α, β)
Fully-compressible Thermodynamically Consistent Pseudo-incompressible	$\operatorname{FC}_{\operatorname{PI}_{ ho,p}^{\operatorname{tc}}}$	(1,0) $(0,1)$
Non-thermodynamically Consistent Pseudo-incompressible	$\operatorname{PI}_{ ho,p}$	(0,1) $(0,0)$

Table 1. Model configurations used in the numerical scheme.

	$\theta'_{ m max}$	$z_{ m max}$	$x_{\max} - x_{\min}$
FC	1.64 K	8183 m	$6637~\mathrm{m}$
$\mathrm{PI}^{\mathrm{tc}}_{ ho,p}$	$1.64~\mathrm{K}$	$8187~\mathrm{m}$	6648 m
$\mathrm{PI}_{ ho,p}$	$1.65~\mathrm{K}$	$8469~\mathrm{m}$	$6278 \mathrm{\ m}$

Table 2. Rising bubble results: maximum temperature perturbation θ'_{\max} , attained height z_{\max} , and horizontal extension $x_{\max} - x_{\min}$ at final time T = 1000 s for FC, $\text{PI}^{\text{tc}}_{\rho,p}$, and $\text{PI}_{\rho,p}$ models. The values refer to the external contour at 0.25 K.

	$E_{\mathrm{rel}}^{\mathrm{rms}}(\theta')$	$E_{\mathrm{rel}}^{\mathrm{max}}(\theta')$	$E_{\mathrm{rel}}^{\mathrm{max}}(\theta_{\mathrm{max}}')$
$PI_{\rho,p}^{tc}$ -FC	0.017	0.018	1.07E-03
$\mathrm{PI}_{\rho,p}^{n} ext{-FC}$	0.57	0.57	3.61E-02

Table 3. Rising bubble results: relative root-mean square error $E_{\rm rel}^{\rm rms}$ and maximum error $E_{\rm rel}^{\rm max}$ on potential temperature perturbation profile θ' and maximum error $E_{\rm rel}^{\rm max}$ on the maximum perturbation amplitude $\theta'_{\rm max}$ for the ${\rm PI}_{\rho,p}^{\rm tc}$ and ${\rm PI}_{\rho,p}$ cuts at z=7500 m with respect to the FC cut as in figure 5.

	$\theta_{ m max}'$	x_{max}
FC	-10.14 K	$15476~\mathrm{m}$
$\mathrm{PI}^{\mathrm{tc}}_{ ho,p}$	-10.17 K	$15456~\mathrm{m}$
$\mathrm{PI}_{ ho,p}$	-9.96 K	$15676~\mathrm{m}$

Table 4. Density current results: maximum temperature perturbation θ'_{max} and front position x_{max} at final time T=900 s. x_{max} is the rightmost intersection of the 1 K contour with the bottom boundary.

	$E_{\mathrm{rel}}^{\mathrm{rms}}(\theta')$	$E_{\mathrm{rel}}^{\mathrm{max}}(\theta')$	$E_{\mathrm{rel}}^{\mathrm{max}}(\theta_{\mathrm{max}}')$
$\mathrm{PI}^{\mathrm{tc}}_{ ho,p} ext{-FC}$	0.046	0.090	1.93E-03
$\mathrm{PI}_{\rho,p} ext{-}\mathrm{FC}$	0.441	0.584	0.026

TABLE 5. Density current results: relative root-mean square error $E_{\rm rel}^{\rm rms}$ and maximum error $E_{\rm rel}^{\rm max}$ on potential temperature perturbation profile θ' and maximum error $E_{\rm rel}^{\rm max}$ on the maximum perturbation amplitude $\theta'_{\rm max}$ for the ${\rm PI}_{\rho,p}^{\rm tc}$ and ${\rm PI}_{\rho,p}$ cuts at z=1200 m with respect to the FC cut as in figure 9.

	u'_{\max}	u'_{\min}	$w'_{\rm max}$	w'_{\min}	$\theta_{ m max}'$	$ heta'_{\min}$	π'_{\max}	π'_{\min}
FC	1.054E-2	-1.060E-2	2.739E-3	-2.262E-3	2.808E-3	-1.526E-3	7.75E-7	-5.27E-7
$\mathrm{PI}^{\mathrm{tc}}_{\rho,p}$	1.063E-2	-1.063E-2	2.645E-3	-2.424E-3	2.808E-3	-1.526E-3	1.18E-5	-6.56E-7
		-1.362E-2						
REF	1.064E-2	-1.061E-2	2.877E-3	-2.400E-3	2.808E-3	-1.511E-3	9.11E-7	-7.13E-7

Table 6. Inertia-gravity wave results: maxima and minima of horizontal velocity u, vertical velocity w, potential temperature θ and Exner pressure $\pi = T\theta^{-1}$ perturbations at final time T = 3000 s in the present study and Restelli and Giraldo (2009) (denoted with REF).

	$E_{\mathrm{rel}}^{\mathrm{rms}}(\theta')$	$E_{\mathrm{rel}}^{\mathrm{max}}(\theta')$
$\mathrm{PI}^{\mathrm{tc}}_{ ho,p} ext{-FC}$	0.039	0.055
$\mathrm{PI}_{ ho,p}^{\prime} ext{-FC}$	0.132	0.16

Table 7. Inertia-gravity wave results: relative root-mean square error $E_{\rm rel}^{\rm rms}$ and maximum error $E_{\rm rel}^{\rm max}$ on potential temperature perturbation profile θ' for the ${\rm PI}_{\rho,p}^{\rm tc}$ and ${\rm PI}_{\rho,p}$ cuts at z=5000 m with respect to the FC cut as in figure 11.

	$C_{ ho}$	$C_{\rho u}$	C_P	$C_{\rho E}$
FC	1.15E-09	8.05E-11	5.68E-09	1.98E-09
$\mathrm{PI}^{\mathrm{tc}}_{ ho,p}$	6.77E-10	9.66E-10	\	3.99E-09
$\operatorname{PI}_{\rho,p}^{n}$	8.90E-10	8.55E-10	\	4.21E-09
REF	1.67E-08	2.60E-07	\	1.64E-08

Table 8. Inertia-gravity wave results: conservation errors for density, horizontal momentum density, P and total energy density (see text for definitions) in the present study and in Restelli and Giraldo (2009), denoted with REF.

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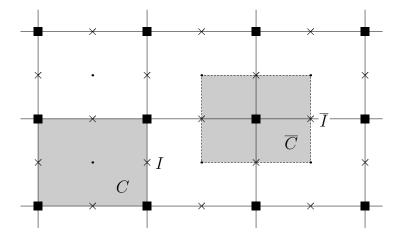


Fig. 1. Computational grid for the numerical scheme. Solid lines define cells; dashed lines define dual cells, used for the second correction. Dots, squares and crosses denote cell centers, nodes, and interface centers, respectively.

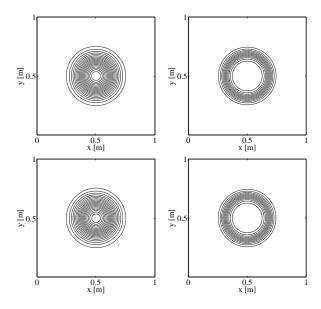


FIG. 2. Smoothed rotating vortex results: density (left) and pressure (right). The upper row shows initial data. The lower row shows computed values at T=1 s with the FC model. Contours are plotted every 0.025 kg m^{-3} in $[0.525, 0.975] \text{ kg m}^{-3}$ for density, every 0.025 Pa in the interval [-0.025, -0.3] Pa for pressure. The domain is discretized with 192 cells in each direction with CFL = 0.45.

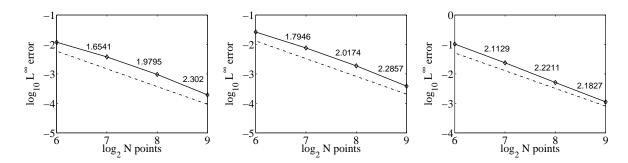


Fig. 3. Smoothed rotating vortex results: density (left), momentum norm (middle) and pressure (right) convergence story. Errors are shown in the maximum norm of computed solutions at T=1 s on grids with 64^2 , 128^2 , 256^2 , and 512^2 cells with respect to computed solutions on a reference grid with 1024^2 cells. The numbers inside the graphs are the experimental rates of convergence between subsequent grid refinements. The dashed-dotted line represents quadratic slope.

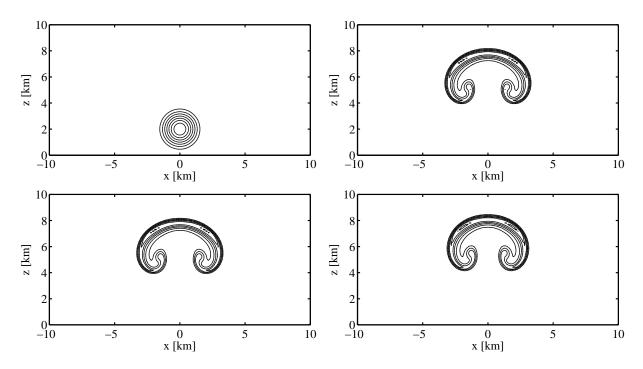


FIG. 4. Rising bubble results. Panels show potential temperature initial data (upper left) and computed value at T=1000 s with the FC (upper right), $\mathrm{PI}_{\rho,p}^{\mathrm{tc}}$ (lower left) and $\mathrm{PI}_{\rho,p}$ models (lower right). Contours are plotted every 0.25 K starting at 300.25 K.

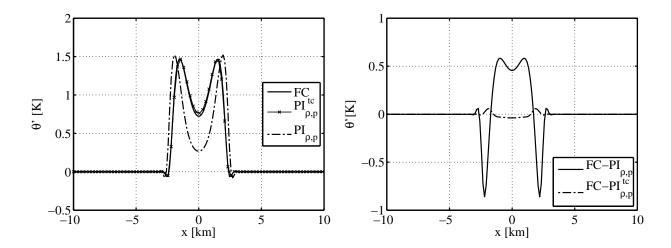


FIG. 5. Rising bubble results: potential temperature perturbation at final time T=1000 s. The left panel shows a horizontal cut of the final θ' at height z=7500 m of the FC (solid line), $\operatorname{PI}_{\rho,p}^{\operatorname{tc}}$ (cross-marked line) and $\operatorname{PI}_{\rho,p}$ (dashed-dotted line). The right panel shows the difference from the FC cut of the $\operatorname{PI}_{\rho,p}^{\operatorname{tc}}$ cut(solid line) and the $\operatorname{PI}_{\rho,p}$ cut(dashed-dotted line).

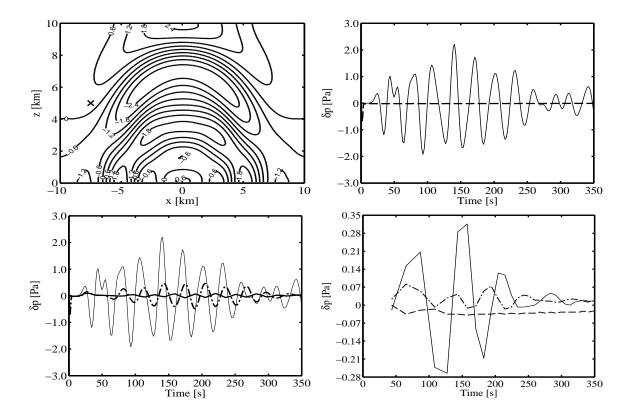


FIG. 6. Rising bubble results, nodal pressure time increment δp . The upper left panel shows contours of δp every .6 Pa starting at -3 Pa, time step 14 (t=26.6 s), FC model. The right panels shows the value of δp over the first 350 s measured at (x,z) = (-7.5,5) km for FC (solid line) and PI_{ρ,p} (dashed line) configurations. In the upper right panel the time step is constant and $\Delta t=1.9$ s. The lower left panel displays the value of δp over the first 350 s measured at the same location. Blended runs at constant $\Delta t=1.9$ s with $S_1=10$ initial pseudo-incompressible steps and $S_2=20$ (dashed-dotted line) and $S_2=40$ (thick solid line) transitional steps are compared with the fully-compressible run, $S_1=S_2=0$ (thin sold line). The dashed-dotted line in the lower right panel refers to a blended run with $S_1=0$, $S_2=3$. In the lower right panel the time step is determined by CFL = 0.5 (initial $\Delta t \approx 21.69$ s) and the data for the first time step is removed.

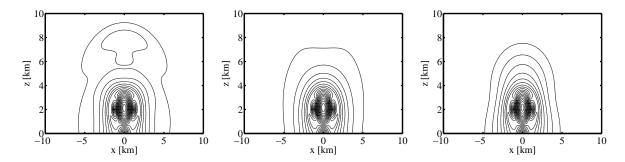


FIG. 7. Rising bubble results: Mach number M at time step 56 ($T \approx 21.66 \,\mathrm{s}$ for $\Delta t = 0.38 \,\mathrm{s}$); left: FC model, $S_1 = S_2 = 0$; middle: $\mathrm{PI}_{\rho,p}$ model; right: $\mathrm{PI}_{\rho,p}$ -then-FC model, $S_1 = 10$, $S_2 = 40$. Contours are plotted every 10^{-4} in the range [0.0001, 0.002].

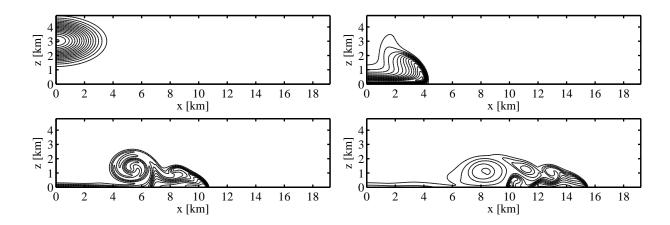


FIG. 8. Density current results: potential temperature perturbation. Panels shows initial data (upper left), FC results at $t=300\,\mathrm{s}$ (upper right), $t=600\,\mathrm{s}$ (lower left) and at $t=900\,\mathrm{s}$ (lower right). Contours are plotted every 1 K from -16.5 K to -0.5 K.

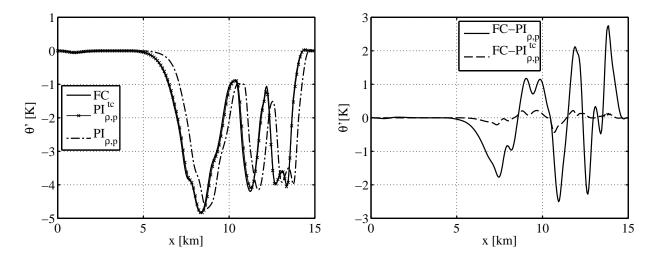


FIG. 9. Density current results: potential temperature perturbation at final time T=900 s. The left panel shows a horizontal cut at height z=1200 m. The right panel shows the difference from the FC profile of the $\mathrm{PI}_{\rho,p}^{\mathrm{tc}}$ profile (solid line) and of the $\mathrm{PI}_{\rho,p}$ profile(dashed line).

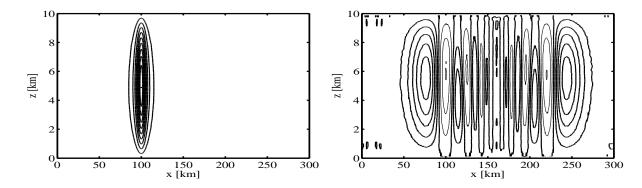
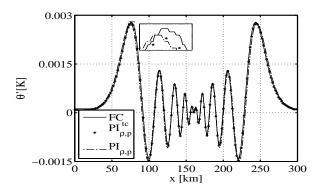


FIG. 10. Inertia-gravity wave results: potential temperature perturbation. The left panel shows initial data, contours every 10^{-3} K; the right panel shows FC result at T=3000 s, contours every $5\cdot 10^{-4}$ K in the range [-0.0015,0.003] K. Thin lines denote negative contours.



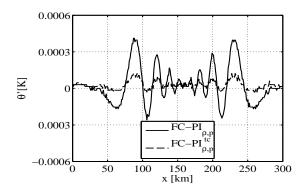


FIG. 11. Inertia-gravity wave results: potential temperature perturbation at final time. The left panel shows a horizontal cut at height z=5000 m for the FC model (solid line), the $\mathrm{PI}_{\rho,p}^{\mathrm{tc}}$ model (stars), and the $\mathrm{PI}_{\rho,p}$ model (dashed-dotted line). The region of the leftmost crest is magnified to highlight the higher phase speed of the $\mathrm{PI}_{\rho,p}$ model. The right panel shows the difference from the FC cut for the $\mathrm{PI}_{\rho,p}^{\mathrm{tc}}$ cut (solid line) and the $\mathrm{PI}_{\rho,p}$ cut (dashed line).