

STOCHASTIC GALERKIN METHOD FOR CLOUD SIMULATION

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ABSTRACT. We develop a stochastic Galerkin method for a coupled Navier-Stokes-cloud system that models dynamics of warm clouds. Our goal is to explicitly describe the evolution of uncertainties that arise due to unknown input data, such as model parameters and initial or boundary conditions. The developed stochastic Galerkin method combines the space-time approximation obtained by a suitable finite volume method with a spectral-type approximation based on the generalized polynomial chaos expansion in the stochastic space. The resulting numerical scheme yields a second-order accurate approximation in both space and time and exponential convergence in the stochastic space. Our numerical results demonstrate the reliability and robustness of the stochastic Galerkin method. We also use the proposed method to study the behavior of clouds in certain perturbed scenarios, for examples, the ones leading to changes in macroscopic cloud pattern as a shift from hexagonal to rectangular structures.

1. INTRODUCTION

Clouds constitute one of the most important components in the Earth-atmosphere system. They influence the hydrological cycle and by interacting with radiation they control the energy budget of the system. However, clouds are one of the most uncertain components, which, unlike the atmospheric flows, cannot be modeled using first principles of physics.

Clouds are composed by myriads of water particles in different phases (liquid and solid), and thus they need to be described by a large ensemble in a statistical sense. A common way of obtaining such an ensemble is by using a mass or size distribution, which would lead to a Boltzmann-type evolution equation. Although there are some approaches available in literature to formulate cloud models in such a way [4, 18, 19], a complete and consistent description is missing. Since measurements of size distributions of cloud particles are difficult, we are often restricted to averaged quantities such as, for example, mass of water per dry air (mass concentrations). Therefore, models are often formulated in terms of so-called bulk quantities, that is, mass and number concentrations of the respective water species. Many cloud processes are necessary to describe the time evolution of the cloud as a statistical ensemble, that is, particle formation or annihilation, growth/evaporation of particles, collision processes, and sedimentation due to gravity. For each of the processes, we have to formulate a representative mathematical term in the sense of a rate equation. Although for some processes the physical mechanisms are quite understood, the formulation of the process rates usually contain uncertain parameters, thus cloud models come with inherent uncertainty. On the other hand, the initial conditions for atmospheric flows and the embedded clouds are also not perfectly constrained, leading to uncertainties in the environmental conditions. It is well-known from former studies that uncertainties in cloud processes and in environmental conditions can lead to drastic changes in simulations, thus these uncertainties influence predictability of moist atmospheric flows, clouds and precipitation in a crucial way; for instance, the distribution of latent heat is changed, which in turn can influence frontogenesis [16] or convection [15, 25].

For investigations of the impact of these uncertain cloud model parameters as well as the impact of variations in environmental conditions on atmospheric flows, sensitivity studies are usually carried out. Since one or more parameters are (randomly) varied, the Monte Carlo approach can be used. This, however, requires a large ensemble of simulations to be conducted, which makes Monte Carlo methods computationally expensive and requires a very fine sampling of the parameter space and possible environmental conditions. In most practical studies, a much smaller set of ensembles (with about 10 – 100 samples only) is used.

In order to improve both the efficiency and accuracy of a numerical method, we choose a different way of representing random variations by using spectral expansions in the stochastic space. This approach enables us to investigate the impact of variations in cloud model parameters and initial conditions on the evolution of moist flows with embedded clouds.

We consider a mathematical model of cloud physics that consists of the Navier-Stokes equations coupled with the cloud evolution equations for the water vapor, cloud water and rain. In this model developed in [23, 35] and presented in Section 2, the Navier-Stokes equations describe weakly compressible flows with viscous and heat conductivity effects, while microscale cloud physics is modeled by the system of advection-diffusion-reaction equations.

Meteorological applications typically inherit several sources of uncertainties, such as model parameters, initial and boundary conditions. Consequently, purely deterministic models are insufficient in such situations and more sophisticated methods need to be applied to analyze the influence of uncertainties on numerical solutions. In this paper, we study a stochastic version of the coupled Navier-Stokes-cloud model in order to account for uncertainties in input quantities. Our main goal is to design an efficient numerical method for quantifying uncertainties in solutions of the studied system. In recent years, a wide variety of uncertainty quantification methods has been proposed and investigated in the context of physical and engineering applications. These methods include stochastic Galerkin methods based on generalized polynomial chaos (gPC) [11, 34, 41, 42, 48], stochastic collocation methods [24, 46, 47], and multilevel Monte Carlo methods [29, 30, 38]. Each of these groups of methods has its own pros and cons. While results obtained by the Monte Carlo simulations are generally good, the approach is not very efficient due to a large number of realizations required. Stochastic collocation methods are typically more efficient than the Monte Carlo ones, since they only require solving the underlying deterministic system at the certain quadrature nodes in the stochastic space. These data are then used to reconstruct the gPC expansion using an appropriate set of orthogonal polynomials. The Monte Carlo as well as the stochastic collocation method fall into a class of the *non-intrusive methods*. Stochastic Galerkin methods offer an alternative approach for computing the gPC expansion. In general, they are more rigorous and efficient than the Monte Carlo and collocation ones; see, e.g., [13]. The stochastic Galerkin method is an *intrusive method* since it requires changes in the underlying code. In fact, one needs to solve a system of PDEs for the gPC expansion coefficients.

We develop a new stochastic Galerkin method for the coupled Navier-Stokes-cloud system. As it has already been mentioned above, the largest source of uncertainties is cloud physics. Therefore, we restrict our consideration to the case in which the uncertainties are only in the cloud physics representation; extension to full stochastic Navier-Stokes-cloud model is left to future studies. Thus, we need to solve the deterministic Navier-Stokes equations coupled with the PDE system for the gPC expansion coefficients for the cloud variables. Our numerical method is an extension of the approach proposed in [23] for the purely deterministic version of the coupled Navier-Stokes-cloud system. This method is based on the operator splitting approach, in which the system is split into the macroscopic Navier-Stokes equations and microscopic cloud model with random inputs. The Navier-Stokes equations are then solved by an implicit-explicit (IMEX) finite-volume method, while for the cloud equations we develop a stochastic Galerkin method based on the gPC. The resulting gPC coefficient system is numerically solved by a finite-volume method combined with an explicit Runge-Kutta method with an enlarged stability region [28].

The paper is organized as follows. We start in Section 2 with the description of the deterministic Navier-Stokes-cloud model. The numerical method for the deterministic model is presented in Sections 3. In Sections 4, we report on numerical experiments for well-known meteorological benchmarks—rising warm bubble and Rayleigh-Bénard convection—for the deterministic model. We then continue in Section 5 with the presentation of the stochastic model, which is followed by the description of the numerical method (Section 6) and presentation of the numerical experiments (Section 7) for the stochastic model. Our numerical results clearly demonstrate that the proposed stochastic Galerkin method is capable of quantifying the uncertainties of complex atmospheric flows.

2. DETERMINISTIC MATHEMATICAL MODEL

We study a mathematical model of cloud dynamics, which is based on the compressible nonhydrostatic Navier-Stokes equations for moist atmosphere (that is, mixture of ideal gases dry air and water vapor),

$$\begin{aligned} \rho_t + \nabla \cdot (\rho \mathbf{u}) &= 0, \\ (\rho \mathbf{u})_t + \nabla \cdot (\rho \mathbf{u} \otimes \mathbf{u} + p \text{Id} - \mu_m \rho (\nabla \mathbf{u} + (\nabla \mathbf{u})^\top)) &= -\rho g \mathbf{e}_3, \end{aligned} \tag{2.1}$$

$$(\rho\theta)_t + \nabla \cdot (\rho\theta\mathbf{u} - \mu_h\rho\nabla\theta) = S_\theta,$$

and evolution equations for cloud variables,

$$\begin{aligned} (\rho q_v)_t + \nabla \cdot (\rho q_v\mathbf{u} - \mu_q\rho\nabla q_v) &= \rho(-C + E), \\ (\rho q_c)_t + \nabla \cdot (\rho q_c\mathbf{u} - \mu_q\rho\nabla q_c) &= \rho(C - A_1 - A_2), \\ (\rho q_r)_t + \nabla \cdot (-v_q\rho q_r\mathbf{e}_3 + \rho q_r\mathbf{u} - \mu_q\rho\nabla q_r) &= \rho(A_1 + A_2 - E). \end{aligned} \quad (2.2)$$

Here, ρ is the density, $\mathbf{u} = (u_1, u_2, u_3)^\top$ is the velocity vector, θ is the moist potential temperature, p is the pressure, g is the acceleration due to gravity, μ_m is the dynamic viscosity, μ_h is the thermal conductivity, and μ_q is the cloud diffusivity. The cloud variables representing the mass concentration of water vapor, cloud droplets and rain drops, q_v , q_c , q_r , respectively, as well as the right-hand side (RHS) terms E , C , A_1 , A_2 will be defined below. We denote by t the time variable and by \mathbf{x} the space vector; $\mathbf{x} = (x_1, x_2, x_3)$ in the three-dimensional (3-D) and $\mathbf{x} = (x_1, x_3)$ in the two-dimensional (2-D) cases. Furthermore, $\mathbf{e}_3 = (0, 0, 1)^\top$ and $\mathbf{e}_3 = (0, 1)^\top$ in the 3-D and 2-D cases, respectively. We set $\mu_m = 10^{-3}$ and $\mu_h = 10^{-2} = \mu_q$. Note that the systems (2.1) and (2.2) are coupled through the source term S_θ , which represents the impact of phase changes and will be defined below, see (2.6). The temperature T can be obtained from the moist adiabatic ideal gas equation

$$T = \frac{R}{R_m} \theta \left(\frac{p}{p_0} \right)^{R_m/c_p}, \quad (2.3)$$

where $p_0 = 10^5$ Pa is the reference pressure at sea level. In addition to the usual definition of a potential temperature, we use $R_m = (1 - q_v - q_c - q_r)R + q_v R_v$ with the ideal gas constant of dry air $R = 287.05$ J/(kg·K), the gas constant of water vapor $R_v = 461.51$ J/(kg·K) and the specific heat capacity of dry air for constant pressure $c_p = 1005$ J/(kg·K). In order to close the system, we determine the pressure from the equation of state that includes moisture

$$p = p_0 \left(\frac{R\rho\theta}{p_0} \right)^{\gamma_m} \quad \text{with} \quad \gamma_m = \frac{c_p}{c_p - R_m}. \quad (2.4)$$

We note that in the dry case R_m reduces to R , $S_\theta = 0$ and the moist ideal gas equation as well as the moist equation of state become their dry analogon.

In this paper, we restrict our investigations to clouds in the lower part of the troposphere, that is, to clouds consisting of liquid droplets exclusively. All of the processes involving ice particles are left for future research. For the representation of liquid clouds in our model we use the so-called single moment scheme, that is, equations for the bulk quantities of mass concentrations of different water phases. For the representation of the relevant cloud processes we adapt a recently developed cloud model [35]. Note that for bulk models, the process rates cannot be derived completely from first principles. Consequently, some uncertain parameters show up naturally. This underlies the need of a rigorous sensitivity study which is the goal of the present paper.

Generally, we follow the standard approach in cloud physics modeling for separating hydrometeors of different sizes, as firstly introduced in [17]. This relies on the observations that small droplets have a negligible falling velocity. In addition, measurements indicate two different modes of droplets in the size distribution, which can be associated to small cloud droplets and large rain drops [43]. Thus, we use the cloud variables q_c and q_r indicating mass concentration of (spatially stationary) cloud droplets and (falling) rain drops, respectively, and the water vapor concentration q_v , that is,

$$q_\ell = \frac{\text{mass of the respective phase}}{\text{mass of dry air}} \quad \text{for} \quad \ell \in \{v, c, r\}.$$

The rest of this section is devoted to a description of the different terms on the RHS of (2.2), which represents the following relevant cloud processes, see [35].

2.1. Single particle properties.

- *General properties of a single water particle*

As we exclusively investigate water clouds, we can assume a spherical shape of water particles. For small cloud droplets this is a very good approximation, while for large rain drops drag effects change

their shape [39, 40]. However, for our investigations of ensembles of rain drops, the spherical shape approximation is appropriate. Thus, mass and radius of droplets are related by the usual equation

$$m = \frac{4}{3}\pi\rho_\ell r^3 \quad \Longleftrightarrow \quad r = \left(\frac{3}{4\pi\rho_\ell}\right)^{\frac{1}{3}} m^{\frac{1}{3}}$$

with the liquid water density $\rho_\ell = 10^3 \text{ kg m}^{-3}$. We make a general assumption that small cloud droplets are stationary, while large rain drops are accelerated by gravity. After balancing gravity by frictional forces, spherical rain drops fall with a terminal velocity, depending only on the drop mass and the density of air. According to [35], the terminal velocity for a droplet of mass m is given by

$$v_\tau(m) = \alpha m^\beta \left(\frac{m_\tau}{m_\tau + m}\right)^\beta \left(\frac{\rho_*}{\rho}\right)^{\frac{1}{2}}, \quad \alpha = 190.3 \text{ m s}^{-1} \text{ kg}^{-\beta}, \quad \beta = \frac{4}{15}, \quad m_\tau = 1.21 \cdot 10^{-5} \text{ kg},$$

with the reference density $\rho_* = 1.255 \text{ kg m}^{-3}$ at $T_* = 288 \text{ K}$ and $p_* = 101\,325 \text{ Pa}$. For masses $m \ll m_\tau$, we can approximate the terminal velocity by $v_\tau = \alpha m^\beta \sqrt{\frac{\rho_*}{\rho}}$; this approximation will be used in the description of the process accretion (collection of cloud droplets by rain drops).

- *Diffusion processes: Growth and evaporation*

Diffusion processes (transfer of water molecules to and from the liquid particle) can be described by the following growth equation:

$$\frac{dm}{dt} = -4\pi r D_v G \rho (q_* - q_v) f_v = -4\pi \underbrace{D_v G}_{=:d} \left(\frac{3}{4\pi\rho_\ell}\right)^{\frac{1}{3}} \rho (q_* - q_v) m^{\frac{1}{3}} f_v,$$

where D_v denotes the diffusion constant, G determines corrections due to the latent heat release for phase changes, and f_v is the ventilation correction for large particles taking into account the effect of flows around the falling spheres. A thermodynamics equilibrium is determined by the saturation mixing ratio $q_* = q_*(p, T) = \varepsilon p_s(T)/p$ with the saturation water vapor pressure over a liquid surface $p_s(T)$ given in [31]. By neglecting curvature effects, water particles grow for $q_v > q_*$ and evaporate for $q_v < q_*$, respectively. The diffusion constant is given according to [36]:

$$D_v = D_{v0} \left(\frac{T}{T_0}\right)^{1.94} \frac{p_0}{p}, \quad D_{v0} = 2.11 \cdot 10^{-5} \text{ m}^2 \text{ s}^{-1}, \quad T_0 = 273.15 \text{ K}, \quad p_0 = p_* = 101\,325 \text{ Pa}$$

and the impact of latent heat release is described by

$$G = \left[\left(\frac{L}{R_v T} - 1 \right) \frac{L p_s(T)}{R_v T^2} \frac{D_v}{K_T} + 1 \right]^{-1},$$

where the latent heat of vaporisation $L = 2.53 \cdot 10^6 \text{ J kg}^{-1}$ and the heat conduction of dry air is (see [12])

$$K_T = \frac{a_K T^{\frac{3}{2}}}{T + b_K 10^{\frac{c_K}{T}}}, \quad a_K = 0.002646 \text{ W m}^{-1} \text{ K}^{-\frac{5}{2}}, \quad b_K = 245.4 \text{ K}, \quad c_K = -12 \text{ K}.$$

Ventilation of large spherical particles of radius r can be taken into account using an empirical ventilation coefficient

$$f_v = a_v + b_v N_{\text{Sc}}^{\frac{1}{3}} N_{\text{Re}}^{\frac{1}{2}}, \quad a_v = 0.78, \quad b_v = 0.308,$$

where the Schmidt and Reynolds numbers are defined as

$$N_{\text{Sc}} = \frac{\mu}{\rho D_v} \quad \text{and} \quad N_{\text{Re}} = \frac{\rho}{\mu} v_\tau(2r), \quad (2.5)$$

respectively. In (2.5), μ is the dynamic viscosity of air, which is expressed according to [12] by

$$\mu = \frac{\mu_0 T^{\frac{3}{2}}}{T + T_\mu}, \quad \mu_0 = 1.458 \cdot 10^{-6} \text{ s Pa K}^{-\frac{1}{2}}, \quad T_\mu = 110.4 \text{ K}.$$

For cloud droplets, we neglect the ventilation correction, thus the mass rate of diffusion for a cloud droplet of mass m_c can be expressed as

$$\frac{dm_c}{dt} = d\rho(q_v - q_*)m_c^{\frac{1}{3}}.$$

For rain drops, growth due to the diffusion is negligible, and thus we obtain the mass rate for rain drops of mass m_r as

$$\frac{dm_r}{dt} = -d\rho(q_* - q_v)_+ \left[a_E m_r^{\frac{1}{3}} + b_E v_\tau(m_r)^{\frac{1}{2}} m_r^{\frac{1}{2}} \right], \quad a_E = a_v, \quad b_E = b_v \left(\frac{\mu}{\rho D_v} \right)^{\frac{1}{3}} \sqrt{\frac{2\rho}{\mu} \left(\frac{3}{4\pi\rho\ell} \right)^{\frac{1}{6}}}.$$

Here, $(\cdot)_+ := \max(\cdot, 0)$ denotes the positive part.

- *Collision of rain drops with cloud droplets: Accretion*

A spherical rain drop of mass m_r (radius r) falls with terminal velocity $v_\tau(m_r)$ through a volume $V = \pi r^2 v_\tau(m_r) \Delta t$ (during a time interval Δt) and collects cloud droplets of total mass $M_c = V \rho q_c$. Thus, the corresponding growth rate of the rain drop is given by

$$\frac{dm_r}{dt} = k'_2 \rho \pi q_c v_\tau(m_r) \left(\frac{3}{4\pi\rho\ell} \right)^{\frac{2}{3}} m_r^{\frac{2}{3}}$$

with an efficiency $k'_2 > 0$.

2.2. Ensemble/collective properties. For the description of clouds as an ensemble of water particles, we would have to introduce such averaged quantities as mass concentrations (as described above, that is, q_c and q_r) as well as number concentrations of cloud droplets, n_c , and rain drops, n_r . Since we do not extend the systems of equations for these two quantities, we introduce relations between mass and number concentrations in order to keep the main effects in a simplified way.

- *Formation of cloud droplets: Activation*

Cloud droplets can be formed by the activation of so-called cloud condensation nuclei (CCN). Liquid aerosol particles can grow by water vapor uptake to larger sizes; this effect can be described by the Köhler theory; see, e.g., [20, 33]. As described in detail in [35], we represent the cloud droplet number concentration n_c by a nonlinear relation

$$n_c = q_c \frac{N_\infty}{q_c + N_\infty m_0} \coth \left(\frac{q_c}{N_0 m_0} \right).$$

Here, N_∞ denotes the maximum number of CCN (depending on environmental conditions, e.g. clean or polluted air), m_0 can be interpreted as the activation mass of cloud droplets, and N_0 is the approximated number of activated droplets at $q_v = q_*$. In our investigations, we set these three parameters to the following values:

$$N_\infty = 8 \cdot 10^8 \text{ kg}^{-1}, \quad m_0 = 5.236 \cdot 10^{-16} \text{ kg}, \quad N_0 = 10^3 \text{ kg}^{-1}.$$

For the initialization of the cloud droplet production, we introduce an additional factor in case of supersaturation

$$C_{\text{act}} = N_0 d\rho(q_v - q_*)_+ m_0^{\frac{1}{3}}.$$

- *Relation between number and mass concentration for rain drops*

In contrast to the formulation in [35], we do not include another equation for the number concentration of rain drops. In a similar way as for cloud activation, we use a relation between n_r and q_r , that is, a closure of the form $n_r = f(q_r, c)$. Since we implicitly assume that the rain drops are distributed according to their size, this approach should be used for mimicking the shape of the distribution in a proper way. We propose the (non)linear relation

$$n_r = c_r q_r^\gamma, \quad 0 < \gamma \leq 1.$$

Assuming a constant mean mass of rain drops \bar{m}_r , we can determine the constants as $c_r = \bar{m}_r^{-1}$ and $\gamma = 1$. This approach would be meaningful for the case of a symmetric size distribution of rain

droplets, centered around the mean mass. However, it is well-known that size distributions of rain are usually skew to larger sizes, thus a linear relation is not appropriate. For sizes of rain drops often an exponential distribution is assumed, this leads to an exponent $\gamma = \frac{1}{4}$ and a coefficient $c_r = c_{r0}\rho^{-\frac{3}{4}}$ ($c_{r0} = 23752.6753 \text{ kg}^{-\frac{1}{4}} \text{ m}^{\frac{3}{4}}$, $[c_r] = \text{kg}^{-1}$) as derived in Appendix A.

- *Rates for diffusion processes*

For the mass mixing ratios q_c and q_r , the rates for the diffusion processes are given by multiplication of the single particle rates by the number concentration of the respective particles, namely:

$$\frac{dq_c}{dt} = n_c \frac{dm_c}{dt}, \quad \frac{dq_r}{dt} = n_r \frac{dm_r}{dt}.$$

Applying the relations between the mass and number concentrations as stated above, we obtain

$$C_1 = \frac{dq_c}{dt} = n_c d\rho(q_v - q_*) m_c^{\frac{1}{3}} \stackrel{m_c = \frac{q_c}{n_c}}{=} d\rho(q_v - q_*) \left(\frac{N_\infty}{q_c + N_\infty m_0} \coth \left(\frac{q_c}{N_0 m_0} \right) \right)^{\frac{2}{3}} q_c$$

and

$$E = -\frac{dq_r}{dt} = n_r d\rho(q_* - q_v)_+ \left[a_E m_r^{\frac{1}{3}} + b_E v_\tau(m_r) m_r^{\frac{1}{2}} \right] \stackrel{m_r = \frac{q_r}{n_r}}{=} d\rho(q_* - q_v)_+ \left[a_E c_r^{\frac{2}{3}} q_r^{\frac{1}{3} + \gamma \frac{2}{3}} + b_E c_r^{\frac{1}{2}} v_\tau(q_r)^{\frac{1}{2}} q_r^{\frac{1}{2} + \gamma \frac{1}{2}} \right]$$

using a reformulated terminal velocity

$$v_\tau(m_r) = v_\tau(q_r) = \alpha q_r^\beta \left(\frac{m_\tau}{q_r + m_\tau c_r q_r^\gamma} \right)^\beta \left(\frac{\rho_*}{\rho} \right)^{\frac{1}{2}}.$$

Note, that the rates for activation and diffusion growth of cloud droplets are combined in the model formulation, that is, $C = C_1 + C_{\text{act}}$.

- *Rate for accretion*

For the rate of accretion of rain water, we obtain

$$A_2 = \frac{dq_r}{dt} = n_r \frac{dm_r}{dt} = k_2 \rho \pi c_r^{\frac{1}{3}} \left(\frac{3}{4\pi\rho_\ell} \right)^{\frac{2}{3}} q_c v_\tau(q_r) q_r^{\frac{2+\gamma}{3}}.$$

Note that for compensating effects of the averaging the parameter k_2 can be adjusted (such that $k_2 = 0.8 \text{ kg} \neq k'_2$) and the impact of the uncertainty of this parameter is of high interest, since it cannot be measured or derived from the first principles.

- *Collision of cloud droplets, forming a rain drop: Autoconversion*

Beside the growth of an existing rain drop by collecting cloud droplets, a rain drop can be formed by the collision of two cloud droplets. According to [35], we formulate the growth rate similarly to population models, namely:

$$A_1 = \frac{dq_c}{dt} = k_1 \frac{\rho q_c^2}{\rho_\ell}.$$

Note that the coefficient k_1 cannot be measured or derived from the first principles. It is a free parameter, which must be fixed using parameter estimations. Thus, the impact of the uncertainty of this parameter is of high interest. In our deterministic experiments, we choose $k_1 = 0.0041 \text{ kg s}^{-1}$, as indicated in [35].

- *Sedimentation of rain mass mixing ratio*

We have introduced an additional convection term into the equation for the evolution of q_r in (2.2), that is, the term $\nabla \cdot (-v_q \rho q_r \mathbf{e}_3)$, where

$$v_q = v_q(q_r) = \alpha q_r^\beta \left(\frac{m_\tau}{q_r + m_\tau c_r q_r^\gamma} \right)^\beta \left(\frac{\rho_*}{\rho} \right)^{\frac{1}{2}}.$$

The parameter α can be derived empirically, but the influence of uncertainty in α is of high interest.

Note that activation and diffusion processes are formulated explicitly, in contrast to the usual approach of saturation adjustment (see, e.g., [22]), which is less accurate, but commonly used in operational weather forecast models. This explicit formulation introduces stiffness caused by modeling cloud processes on the RHS of the cloud equations with fractional exponents between -1 and 1 . In order to avoid the root evaluation for an argument that is close to zero, we introduce a cut-off function and replace ζ^ξ , $\xi \in (-1, 1)$, with

$$\begin{cases} \zeta^\xi, & \text{if } \zeta > 10^{-16}, \\ 0, & \text{otherwise.} \end{cases}$$

Due to phase changes (activation and growth/evaporation of water particles) latent heat is released or absorbed. These processes are modeled by the source term in (2.1):

$$S_\theta = \rho \frac{L\theta}{c_p T} (C - E), \quad (2.6)$$

whereas the terms $C = C_1 + C_{\text{act}}$ and E are explained in details in Section 2.2, *Rates for diffusion processes*.

Solving the Navier-Stokes equations (2.1) in a weakly compressible regime is known to cause numerical instabilities due to the multiscale effects. We follow the approach typically used in meteorological models, where the dynamics of interest is described by a perturbation of a background state, which is the hydrostatic equilibrium. The latter expresses a balance between the gravity and pressure forces. Denoting by \bar{p} , $\bar{\rho}$, $\bar{\mathbf{u}} = 0$, $\bar{\theta}$ and $\bar{\rho\theta}$ the respective background state, the hydrostatic equilibrium satisfies

$$\frac{\partial \bar{p}}{\partial x_3} = -\bar{\rho}g, \quad S_\theta = 0,$$

where \bar{p} is obtained from the equation of state (2.4)

$$\bar{p} = p(\bar{\rho\theta}) = p_0 \left(\frac{R\bar{\rho\theta}}{p_0} \right)^{\gamma_m}. \quad (2.7)$$

Let p' , ρ' , \mathbf{u}' , θ' and $(\rho\theta)'$ stand for the corresponding perturbations of the equilibrium state, then

$$p = \bar{p} + p', \quad \rho = \bar{\rho} + \rho', \quad \theta = \bar{\theta} + \theta', \quad \mathbf{u} = \mathbf{u}', \quad \rho\theta = \bar{\rho}\bar{\theta} + \bar{\rho}\theta' + \rho'\bar{\theta} + \rho'\theta' = \bar{\rho\theta} + (\rho\theta)'$$

The pressure perturbation p' is derived from (2.4) and (2.7) using the following Taylor expansion

$$p(\rho\theta) \approx p(\bar{\rho\theta}) + \frac{\partial p}{\partial(\rho\theta)} (\rho\theta - \bar{\rho\theta}) = \bar{p} + \gamma_m p_0 \left(\frac{R\bar{\rho\theta}}{p_0} \right)^{\gamma_m} \frac{(\rho\theta)'}{\bar{\rho\theta}},$$

which results in

$$p' \approx \gamma_m p_0 \left(\frac{R\bar{\rho\theta}}{p_0} \right)^{\gamma_m} \frac{(\rho\theta)'}{\bar{\rho\theta}}.$$

Consequently, a physical motivation from the hydrostatic balance state leads to a numerically preferable scaling and the perturbation formulation of the Navier-Stokes equations (2.1) then reads as

$$\begin{aligned} \rho'_t + \nabla \cdot (\rho\mathbf{u}) &= 0, \\ (\rho\mathbf{u})_t + \nabla \cdot (\rho\mathbf{u} \otimes \mathbf{u} + p' \text{Id} - \mu_m \rho (\nabla\mathbf{u} + (\nabla\mathbf{u})^\top)) &= -\rho' g \mathbf{e}_3, \\ (\rho\theta)'_t + \nabla \cdot (\rho\theta\mathbf{u} - \mu_h \rho \nabla\theta) &= S_\theta. \end{aligned} \quad (2.8)$$

For alternative representations of cloud dynamics and their numerical investigations, we refer the reader to [3, 37] and references therein.

3. NUMERICAL SCHEME FOR THE DETERMINISTIC MODEL

The numerical approximation of the coupled model (2.8), (2.2) is based on the second-order Strang operator splitting. Therefore, we split the whole system into the macroscopic Navier-Stokes flow equations and the microscopic cloud equations. The Navier-Stokes equations (2.8) are approximated by an IMEX finite-volume method and the cloud equations (2.2) are approximated by a finite-volume method in space and an explicit Runge-Kutta method with an enlarged stability region in time.

3.1. Operator form. Let $\mathbf{w} := (\rho', \rho\mathbf{u}, (\rho\theta)')^\top$ and $\mathbf{w}_q := (\rho q_v, \rho q_c, \rho q_r)^\top$ denote the solution vectors of (2.8) and (2.2), respectively. Then, the coupled system can be written as

$$\begin{aligned}\mathbf{w}_t &= -\nabla \cdot \mathbf{F}(\mathbf{w}) + \mathfrak{D}(\mathbf{w}) + \mathfrak{R}(\mathbf{w}), \\ (\mathbf{w}_q)_t &= -\nabla \cdot \mathbf{F}_q(\mathbf{w}_q) + \mathfrak{D}_q(\mathbf{w}_q) + \mathfrak{R}_q(\mathbf{w}_q),\end{aligned}$$

where \mathbf{F} and \mathbf{F}_q are advection fluxes and \mathfrak{D} , \mathfrak{R} and \mathfrak{D}_q , \mathfrak{R}_q denote the diffusion and reaction operators of the respective systems. They are given by

$$\begin{aligned}\mathbf{F}(\mathbf{w}) &:= (\rho\mathbf{u}, \rho\mathbf{u} \otimes \mathbf{u} + p' \text{Id}, \rho\theta\mathbf{u})^\top, \\ \mathfrak{D}(\mathbf{w}) &:= (0, \nabla \cdot (\mu_m \rho (\nabla \mathbf{u} + (\nabla \mathbf{u})^\top)), \nabla \cdot (\mu_h \rho \nabla \theta))^\top, \\ \mathfrak{R}(\mathbf{w}) &:= (0, -\rho' g \mathbf{e}_3, S_\theta)^\top, \\ \mathbf{F}_q(\mathbf{w}_q) &:= (\rho q_v \mathbf{u}, \rho q_c \mathbf{u}, \rho q_r \mathbf{u} - v_q \rho q_r \mathbf{e}_3)^\top, \\ \mathfrak{D}_q(\mathbf{w}_q) &:= (\nabla \cdot (\mu_q \rho \nabla q_v), \nabla \cdot (\mu_q \rho \nabla q_c), \nabla \cdot (\mu_q \rho \nabla q_r))^\top, \\ \mathfrak{R}_q(\mathbf{w}_q) &:= (-C + E, C - A_1 - A_2, A_1 + A_2 - E)^\top.\end{aligned}\tag{3.1}$$

In order to derive an asymptotically stable, accurate and computationally efficient scheme for the Navier-Stokes equations, we first split the equations into linear and nonlinear parts; see [6, 23] and references therein. Consequently, we introduce

- $\mathbf{F}(\mathbf{w}) = \mathbf{F}_L(\mathbf{w}) + \mathbf{F}_N(\mathbf{w})$ with $\mathbf{F}_L(\mathbf{w}) := (\rho\mathbf{u}, p' \text{Id}, \bar{\theta}\rho\mathbf{u})^\top$ and $\mathbf{F}_N(\mathbf{w}) := (0, \rho\mathbf{u} \otimes \mathbf{u}, \theta' \rho\mathbf{u})^\top$;
- $\mathfrak{D}(\mathbf{w}) = \mathfrak{D}_L(\mathbf{w}) + \mathfrak{D}_N(\mathbf{w})$ with

$$\begin{aligned}\mathfrak{D}_L(\mathbf{w}) &:= (0, \mu_m (\Delta(\rho\mathbf{u}) + \nabla(\nabla \cdot (\rho\mathbf{u}))), \mu_h \Delta(\rho\theta)')^\top \text{ and} \\ \mathfrak{D}_N(\mathbf{w}) &:= (0, -\mu_m ((\Delta\rho)\mathbf{u} + (D^2\rho)\mathbf{u} + \nabla\mathbf{u}\nabla\rho + \nabla\rho\nabla \cdot \mathbf{u}), \mu_h (\Delta(\bar{\rho}\theta) - \theta\Delta\rho - \nabla\rho \cdot \nabla\theta))^\top;\end{aligned}$$
- $\mathfrak{R}(\mathbf{w}) = \mathfrak{R}_L(\mathbf{w}) + \mathfrak{R}_N(\mathbf{w})$ with $\mathfrak{R}_L(\mathbf{w}) := (0, -\rho' g \mathbf{e}_3, 0)^\top$ and $\mathfrak{R}_N(\mathbf{w}) := (0, 0, S_\theta)^\top$.

We would like to point out that the choice of the linear and nonlinear operators is crucial. We choose the linear part to model linear acoustic and gravitational waves as well as linear viscous fluxes. The nonlinear part describes nonlinear advective effects together with the remaining nonlinear viscous fluxes and the influence of the latent heat. We will use the following notation:

$$\mathcal{L} := -\nabla \cdot \mathbf{F}_L(\mathbf{w}) + \mathfrak{D}_L(\mathbf{w}) + \mathfrak{R}_L(\mathbf{w}) \quad \text{and} \quad \mathcal{N} := -\nabla \cdot \mathbf{F}_N(\mathbf{w}) + \mathfrak{D}_N(\mathbf{w}) + \mathfrak{R}_N(\mathbf{w}).$$

3.2. Discretization in space. The spatial discretization is realized by a finite-volume method. We take a cuboid computational domain $\Omega \subset \mathbb{R}^d$, which is divided into N uniform Cartesian cells. The cells are labeled in a certain order using a single-index notation. For simplicity of notation, we assume that the cells are cubes with the sides of size h so that $|C_i| = h^d$. We also introduce the notation $S(i)$ for the set of all neighboring cells of cell C_i , $i = 1, \dots, N$.

We assume that at a certain time t the approximate solution is realized in terms of its cell averages

$$\mathbf{w}_i(t) \approx \frac{1}{h^d} \int_{C_i} \mathbf{w}(\mathbf{x}, t) \, d\mathbf{x} \quad \text{and} \quad (\mathbf{w}_q)_i(t) \approx \frac{1}{h^d} \int_{C_i} \mathbf{w}_q(\mathbf{x}, t) \, d\mathbf{x}, \quad i = 1, \dots, N.$$

In order to simplify the notation, we will now omit the time dependence of $\mathbf{w}_i(t)$ and $(\mathbf{w}_q)_i(t)$. Next, we introduce the notation $\mathbf{w}_h := \{\mathbf{w}_i\}_{i=1}^N$ and $(\mathbf{w}_q)_h := \{(\mathbf{w}_q)_i\}_{i=1}^N$ and consider the following approximation of the advection, diffusion and reaction operators:

$$\begin{aligned}\mathcal{A}_i(\mathbf{w}_h) &= (\mathcal{A}_L)_i(\mathbf{w}_h) + (\mathcal{A}_N)_i(\mathbf{w}_h) \approx \frac{1}{h^d} \int_{C_i} \nabla \cdot \mathbf{F}_L(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x} + \frac{1}{h^d} \int_{C_i} \nabla \cdot \mathbf{F}_N(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x}, \\ \mathcal{D}_i(\mathbf{w}_h) &= (\mathcal{D}_L)_i(\mathbf{w}_h) + (\mathcal{D}_N)_i(\mathbf{w}_h) \approx \frac{1}{h^d} \int_{C_i} \mathfrak{D}_L(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x} + \frac{1}{h^d} \int_{C_i} \mathfrak{D}_N(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x}, \\ \mathcal{R}_i(\mathbf{w}_h) &= (\mathcal{R}_L)_i(\mathbf{w}_h) + (\mathcal{R}_N)_i(\mathbf{w}_h) \approx \frac{1}{h^d} \int_{C_i} \mathfrak{R}_L(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x} + \frac{1}{h^d} \int_{C_i} \mathfrak{R}_N(\mathbf{w}(\mathbf{x}, t)) \, d\mathbf{x}.\end{aligned}$$

Analogous notation will be used for the approximations $(\mathcal{A}_q)_i(\mathbf{w}_h)$, $(\mathcal{D}_q)_i(\mathbf{w}_h)$ and $(\mathcal{R}_q)_i(\mathbf{w}_h)$ of the cloud operators.

3.2.1. *Advection.* The advection terms are discretized using flux functions as follows:

$$\begin{aligned}(\mathcal{A}_L)_i(\mathbf{w}_h) &= \frac{1}{h} \sum_{j \in S(i)} H_{ij}^L(\mathbf{w}_h) \sum_{k=1}^d n_{ij}^{(k)}, \\ (\mathcal{A}_N)_i(\mathbf{w}_h) &= \frac{1}{h} \sum_{j \in S(i)} H_{ij}^N(\mathbf{w}_h) \sum_{k=1}^d n_{ij}^{(k)}, \\ (\mathcal{A}_q)_i((\mathbf{w}_q)_h) &= \frac{1}{h} \sum_{j \in S(i)} (H_q)_{ij}((\mathbf{w}_q)_h) \sum_{k=1}^d n_{ij}^{(k)},\end{aligned}$$

where the numerical fluxes H_{ij}^L , H_{ij}^N and $(H_q)_{ij}$ approximate the corresponding fluxes between the computational cells C_i and C_j , and $n_{ij}^{(k)}$ denotes the k -th component of the outer normal unit vector of cell C_i in the direction of cell C_j . We use the Rusanov numerical flux for H_{ij}^N and $(H_q)_{ij}$ and the central flux for H_{ij}^L . For $(\mathcal{A}_N)_i(\mathbf{w}_h)$ and $(\mathcal{A}_q)_i((\mathbf{w}_q)_h)$ a discretization is obtained via a MUSCL-type approach using piecewise linear reconstructions with the minmod limiter. It is well-known this approach yields an approximation, which is almost second-order accurate as its accuracy deteriorates at sudden changes, i.e. jumps, discontinuities or large curvatures. The numerical fluxes are given by

$$\begin{aligned}H_{ij}^L(\mathbf{w}_h) &= \frac{1}{2} (\mathbf{F}_L(\mathbf{w}_j) + \mathbf{F}_L(\mathbf{w}_i)), \\ H_{ij}^N(\mathbf{w}_h) &= \frac{1}{2} (\mathbf{F}_N(\mathbf{w}_{ij}^+) + \mathbf{F}_N(\mathbf{w}_{ij}^-)) - \frac{\lambda_{ij}}{2} (\mathbf{w}_{ij}^+ - \mathbf{w}_{ij}^-), \\ (H_q)_{ij}((\mathbf{w}_q)_h) &= \frac{1}{2} (\mathbf{F}_q((\mathbf{w}_q)_{ij}^+) + \mathbf{F}_q((\mathbf{w}_q)_{ij}^-)) - \frac{(\lambda_q)_{ij}}{2} ((\mathbf{w}_q)_{ij}^+ - (\mathbf{w}_q)_{ij}^-).\end{aligned}\tag{3.2}$$

Here, \mathbf{w}_{ij}^- , \mathbf{w}_{ij}^+ and $(\mathbf{w}_q)_{ij}^-$, $(\mathbf{w}_q)_{ij}^+$ denote the corresponding interface values, which are computed using a piecewise linear reconstruction so that

$$\mathbf{w}_{ij}^- = \mathbf{w}_i + \mathbf{s}_{ij} \frac{h}{2} \sum_{k=1}^d n_{ij}^{(k)}, \quad \mathbf{w}_{ij}^+ = \mathbf{w}_j - \mathbf{s}_{ij} \frac{h}{2} \sum_{k=1}^d n_{ij}^{(k)},$$

where the slopes \mathbf{s}_{ij} are computed by the minmod limiter,

$$\mathbf{s}_{ij} = \frac{1}{h} \text{minmod}(\mathbf{w}_j - \mathbf{w}_i, \mathbf{w}_i - \mathbf{w}_{j^*}) \sum_{k=1}^d n_{ij}^{(k)},$$

applied in a component-wise manner. Here,

$$\text{minmod}(a, b) = \begin{cases} a, & \text{if } |a| < |b| \text{ and } ab > 0, \\ b, & \text{if } |b| < |a| \text{ and } ab > 0, \\ 0, & \text{if } ab \leq 0, \end{cases}$$

and $(\mathbf{w}_q)_{ij}^-$ and $(\mathbf{w}_q)_{ij}^+$ are obtained similarly. Thereby C_{j^*} is the other neighboring cell of C_i in the opposite direction from C_j . Finally, the values λ_{ij} and $(\lambda_q)_{ij}$ are given by

$$\lambda_{ij} = \max \left\{ \sigma \left(\frac{\partial \mathbf{F}_N(\mathbf{w}_{ij}^-)}{\partial \mathbf{w}} \right), \sigma \left(\frac{\partial \mathbf{F}_N(\mathbf{w}_{ij}^+)}{\partial \mathbf{w}} \right) \right\}, \quad (\lambda_q)_{ij} = \max \left\{ \sigma \left(\frac{\partial \mathbf{F}_q((\mathbf{w}_q)_{ij}^-)}{\partial \mathbf{w}_q} \right), \sigma \left(\frac{\partial \mathbf{F}_q((\mathbf{w}_q)_{ij}^+)}{\partial \mathbf{w}_q} \right) \right\},$$

where σ denotes the spectral radius of the corresponding Jacobians.

Remark 3.1. Note that in the computation of H_{ij}^L in (3.2), we use the cell averages rather than the point values at the cell interfaces for the following two reasons. First, the flux is second-order accurate. Second, in Section 3.3, we will treat the linear part of the flux implicitly and this is much easier to do when the numerical flux is linear as well.

3.2.2. Diffusion. The components of the discrete diffusion operators are discretized in a straightforward manner using second-order central differences.

3.2.3. Reaction. The reaction terms are discretized by a direct evaluation of the reaction operators at the cell centers:

$$\mathcal{R}_i(\mathbf{w}_h) = \mathfrak{R}_L(\mathbf{w}_i) + \mathfrak{R}_N(\mathbf{w}_i), \quad (\mathcal{R}_q)_i((\mathbf{w}_q)_h) = \mathfrak{R}_q((\mathbf{w}_q)_i).$$

After the spatial discretization, we obtain the following system of time-dependent ODEs:

$$\frac{d}{dt} \mathbf{w}_i = -\mathcal{A}_i(\mathbf{w}_h) + \mathcal{D}_i(\mathbf{w}_h) + \mathcal{R}_i(\mathbf{w}_h), \quad (3.3)$$

$$\frac{d}{dt} (\mathbf{w}_q)_i = -(\mathcal{A}_q)_i((\mathbf{w}_q)_h) + (\mathcal{D}_q)_i((\mathbf{w}_q)_h) + (\mathcal{R}_q)_i((\mathbf{w}_q)_h). \quad (3.4)$$

This system has to be solved using an appropriate ODE solver as discussed in Section 3.3.

3.3. Discretization in time. Let \mathbf{w}_h^n and $(\mathbf{w}_q)_h^n$ denote the numerical approximation of the solutions $\mathbf{w}_h(t)$ and $(\mathbf{w}_q)_h(t)$ at the discrete time level t^n . We evolve the solution to the next time level $t^{n+1} = t^n + \Delta t^n$, where Δt^n is the size of the Strang operator splitting time step. In the operator splitting approach, we first numerically solve the ODE system (3.3) with $\Delta t_{\text{NS}}^n = \Delta t^n/2$, we then numerically integrate the ODE system (3.4) with Δt^n and finally we solve the system (3.3) again with Δt_{NS}^n .

Notice that the system (3.3) may be very stiff as the Navier-Stokes equations are in the weakly compressible regime. We therefore follow the approach in [6] (see also [5]), and employ the second-order ARS(2,2,2) IMEX method from [2]:

$$\begin{aligned} \mathbf{w}_h^{n+\frac{1}{4}} &= \mathbf{w}_h^n + \beta \Delta t_{\text{NS}}^n \left(\mathcal{L} \left(\mathbf{w}_h^{n+\frac{1}{4}} \right) + \mathcal{N} \left(\mathbf{w}_h^n \right) \right), \\ \mathbf{w}_h^{n+\frac{1}{2}} &= \mathbf{w}_h^n + \Delta t_{\text{NS}}^n \left(\delta \mathcal{N} \left(\mathbf{w}_h^n \right) + (1 - \delta) \mathcal{N} \left(\mathbf{w}_h^{n+\frac{1}{4}} \right) \right) + \Delta t_{\text{NS}}^n \left(\beta \mathcal{L} \left(\mathbf{w}_h^{n+\frac{1}{2}} \right) + (1 - \beta) \mathcal{L} \left(\mathbf{w}_h^{n+\frac{1}{4}} \right) \right), \end{aligned} \quad (3.5)$$

where $\alpha = 1 - 1/\sqrt{2}$, $\delta = 1 - 1/2\beta$, $t^{n+\frac{1}{2}} = t^n + \Delta t_{\text{NS}}^n$, $t^{n+\frac{1}{4}} = t^n + \Delta t_{\text{NS}}^n/2$, and Δt_{NS}^n satisfies the following CFL condition:

$$\max_{s=1,2,3} \max_{i=1,\dots,N} (|(u_s)_i|) \frac{\Delta t_{\text{NS}}^n}{h} < 0.5.$$

For solving the linear systems arising in (3.5), we use the generalized minimal residual (GMRES) method combined with a preconditioner, the incomplete LU factorization (ILU). As it was shown in [6] (see also [5]), the resulting method is both accurate and efficient in the weakly compressible regime.

The ODE system (3.4) is also stiff, but its stiffness only comes from the diffusion and power-law-type source terms. We therefore efficiently solve it using the large stability domain third-order Runge-Kutta method

from [28]. We have utilized the ODE solver DUMKA3, which is a free software that can be found in [27]. We note that DUMKA3 selects time steps automatically, but in order to improve its efficiency, one needs to provide the code with a time step stability restriction for the forward Euler method; see [27, 28]. This bound is obtained by $\min\{\Delta t^n, \Delta t_{\text{cloud}}^n\}$, where $\Delta t_{\text{cloud}}^n$ satisfies the following CFL condition for the cloud system:

$$\max_{s=1,2} \max_{i=1,\dots,N} (|(u_s)_i|, |(u_3)_i + v_q|) \frac{\Delta t_{\text{cloud}}^n}{h} < 0.5.$$

4. DETERMINISTIC NUMERICAL EXPERIMENTS

In this section, we test the numerical method described in Section 3. The experimental order of convergence is computed for the so-called free convection of a moist warm air bubble and the structure formation in cloud dynamics is shown in the Rayleigh-Bénard convection. The latter will be simulated in both the 2-D and 3-D cases.

4.1. Free convection of a moist warm air bubble in 2-D. We start with the well-known meteorological benchmark describing the free convection of a smooth warm air bubble; see, e.g., [8, 10].

Example 1. In this experiment, the warm bubble rises and deforms axisymmetrically due to the shear friction with the surrounding air at the warm/cold air interface, gradually forming a mushroom-like shape. The warm air bubble is placed at (3500 m, 2000 m) with the initial perturbation:

$$\begin{aligned} \rho'(\mathbf{x}, 0) &= -\bar{\rho}(\mathbf{x}) \frac{\theta'(\mathbf{x}, 0)}{\theta(\mathbf{x}) + \theta'(\mathbf{x}, 0)}, \quad \bar{\rho}(\mathbf{x}) = \frac{p_0}{R\bar{\theta}(\mathbf{x})} \pi_e(\mathbf{x})^{\frac{1}{\gamma-1}}, \quad \pi_e(\mathbf{x}) = 1 - \frac{g x_3}{c_p \bar{\theta}}, \\ \mathbf{u}(\mathbf{x}, 0) &= 0, \\ \theta'(\mathbf{x}, 0) &= \begin{cases} 2 \cos^2\left(\frac{\pi r}{2}\right), & r := \sqrt{(x_1 - 3500)^2 + (x_3 - 2000)^2} \leq 2000, \\ 0, & \text{otherwise,} \end{cases} \end{aligned}$$

where $\bar{\theta} = 300$ K and $p_0 = \bar{p} = 10^5$ Pa. The experiment was simulated in a domain $\Omega = [0, 7000] \times [0, 5000]$ m². For the cloud variables we choose the following initial conditions:

$$q_v(\mathbf{x}, 0) = 0.08 \theta'(\mathbf{x}, 0), \quad q_c = 10^{-3} \theta'(\mathbf{x}, 0), \quad q_r = 10^{-5} \theta'(\mathbf{x}, 0).$$

Furthermore, we apply the no-flux boundary conditions $\mathbf{u} \cdot \mathbf{n} = 0$, $\nabla \rho' \cdot \mathbf{n} = 0$, $\nabla(\rho\theta)' \cdot \mathbf{n} = 0$, $\nabla(\rho q_\ell) \cdot \mathbf{n} = 0$, $\ell \in \{v, c, r\}$.

In Figure 1, we show the potential temperature θ and cloud variables q_v , q_c and q_r , computed on a 160×160 mesh at $t = 150$ and 200 s. One can clearly observe condensation taking place on the interface between cold and warm air and leading to cloud formation in this region. In consequence, rain is formed in the clouds and falls towards the surface. Note that the order of magnitude of the different water mass concentrations is very different, that is, $q_v \gg q_c \gg q_r$, as expected. The experimental convergence study for the cloud and flow variables is presented in Tables 1 and 2, respectively. The experimental order of convergence (EOC) has been computed in the following way:

$$EOC = \log_2 \left(\frac{\|v_{N,\Delta t} - v_{2N,\Delta t/2}\|_{L^2(\Omega)}}{\|v_{2N,\Delta t/2} - v_{4N,\Delta t/4}\|_{L^2(\Omega)}} \right),$$

where $v_{N,\Delta t}$ is the numerical solution computed on a grid with $N \times N$ grid cells and using a fixed time step Δt . As one can clearly see, the expected second order of accuracy has been achieved. For comparison, we present in Figures 2 and 3 the errors measured in the L^1 -, L^2 - and L^∞ -norms. They all give similar results.

4.2. Rayleigh-Bénard convection. In this experiment, we study a natural convection that is used to model structure formation. It occurs in a planar flow between two horizontal plates, where the lower one is heated from below and the upper one is cooled from above. Due to the presence of buoyancy, and hence gravity, the fluid develops a regular pattern of convection roles, known as the Bénard cells. In 3-D, these convection roles form additionally hexagonal structures; see, e.g., [1, 14, 32].

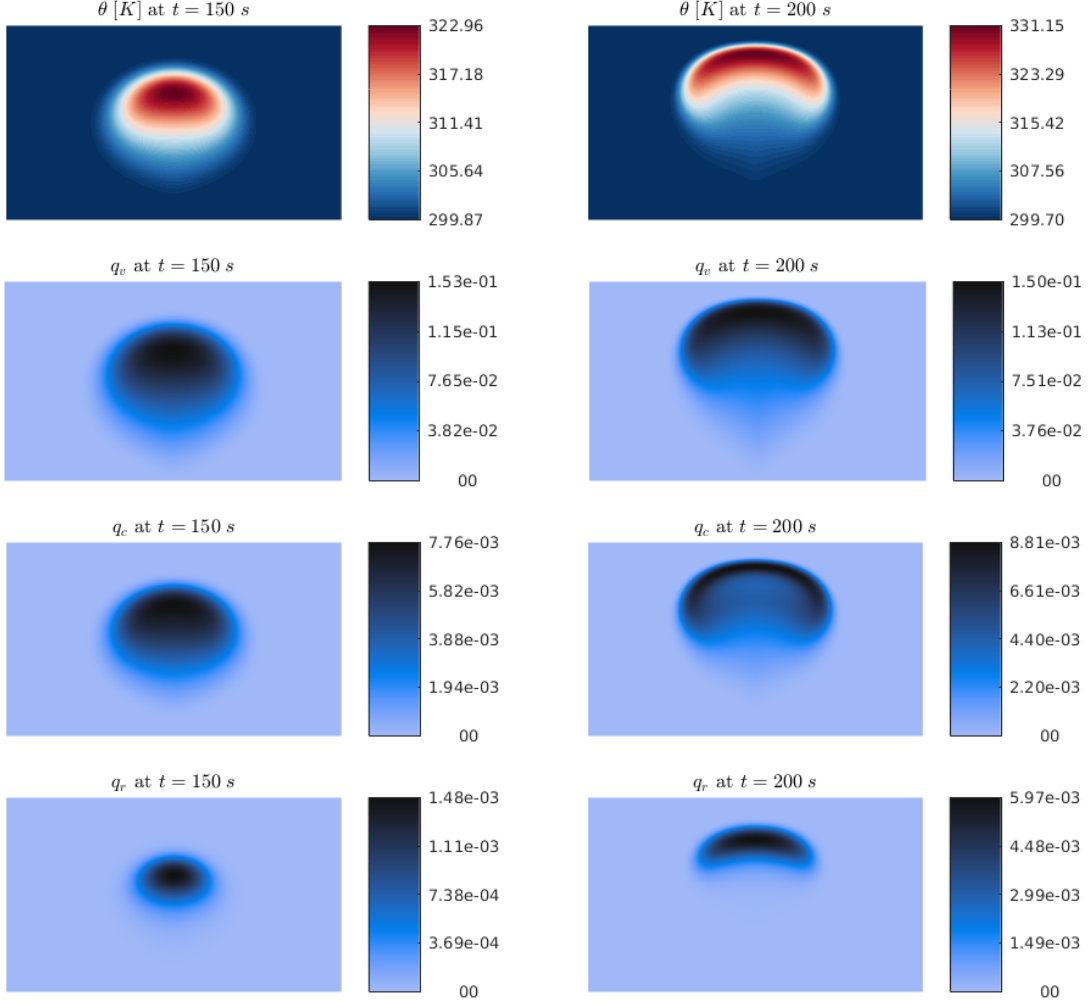


FIGURE 1. *Example 1: Potential temperature θ , water vapor concentration q_v , cloud drops concentration q_c and rain concentration q_r at times $t = 150$ (left column) and 200 s (right column) simulated on a 160×160 mesh.*

N	L^2 -error in ρq_v	EOC	L^2 -error in ρq_c	EOC	L^2 -error in ρq_r	EOC
10	6.870e+00	–	1.019e-01	–	1.159e-03	–
20	1.711e+00	2.01	2.544e-02	2.00	3.152e-04	1.88
40	4.271e-01	2.00	6.380e-03	2.00	1.240e-04	1.35
80	1.080e-01	1.98	1.611e-03	1.99	4.952e-05	1.32
160	2.703e-02	2.00	4.042e-04	1.99	1.571e-05	1.67
320	6.765e-03	2.00	1.016e-04	1.99	5.666e-06	1.47

TABLE 1. *Example 1: L^2 -errors and EOC for the cloud variables computed at time $t = 10$ s using $\Delta t = 20/N$.*

For our numerical simulations, we prescribe the following initial conditions:

$$\rho'(\mathbf{x}, 0) = -\bar{\rho}(\mathbf{x}) \frac{\theta'(\mathbf{x}, 0)}{\bar{\theta}(\mathbf{x}) + \theta'(\mathbf{x}, 0)}, \quad \bar{\rho}(\mathbf{x}) = \frac{p_0}{R\bar{\theta}(\mathbf{x})} \pi_e(\mathbf{x})^{\frac{1}{\gamma-1}}, \quad \pi_e(\mathbf{x}) = 1 - \frac{g x_3}{c_p \bar{\theta}},$$

N	L^2 -error in ρ'	EOC	L^2 -error in ρu_1	EOC	L^2 -error in ρu_2	EOC	L^2 -error in $(\rho\theta)'$	EOC
10	7.522e-01	–	1.134e+02	–	5.805e+01	–	1.213e+02	–
20	1.757e-01	2.10	2.607e+01	2.12	1.744e+01	1.74	3.494e+01	1.80
40	4.418e-02	2.00	5.604e+00	2.22	5.462e+00	1.67	9.641e+00	1.86
80	1.147e-02	1.95	1.436e+00	1.96	1.658e+00	1.72	2.584e+00	1.90
160	3.170e-03	1.85	3.972e-01	1.85	5.875e-01	1.50	7.577e-01	1.77
320	9.810e-04	1.70	1.159e-01	1.78	2.420e-01	1.28	2.556e-01	1.57

TABLE 2. Example 1: L^2 -errors and EOC for the flow variables computed at time $t = 10s$ using $\Delta t = 20/N$.

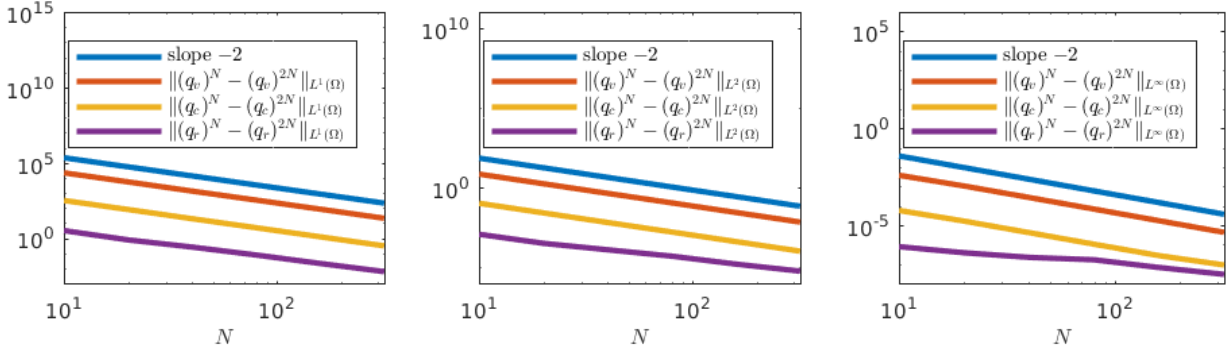


FIGURE 2. Example 1: Comparison of different error norms for the cloud variables q_v , q_c and q_r computed at time $t = 10s$ using $\Delta t = 20/N$.

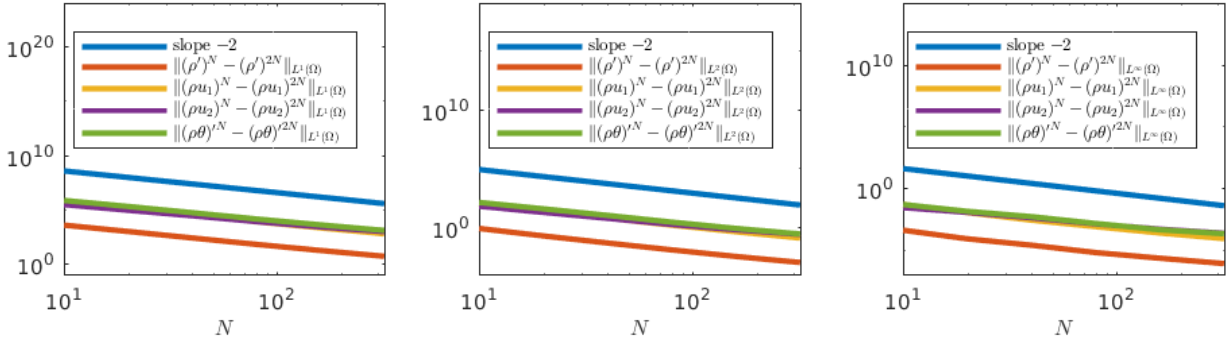


FIGURE 3. Example 1: Comparison of different error norms for the flow variables ρ' , ρu_1 , ρu_2 and $(\rho\theta)'$ at time $t = 10s$ using $\Delta t = 20/N$.

$$\mathbf{u}(\mathbf{x}, 0) = 0, \quad \theta'(\mathbf{x}, 0) = \eta(\mathbf{x}), \quad \bar{\theta}(\mathbf{x}) = 284 - \frac{1}{1000}x_3,$$

where $p_0 = \bar{p} = 10^5$ Pa and $\eta(\mathbf{x})$ is a random perturbation uniformly distributed in $[-0.0021, 0.0021]$. For the cloud equations, the following initial data are used:

$$q_v(\mathbf{x}, 0) = 2 \cdot 10^{-5} \bar{\theta}(\mathbf{x}), \quad q_c = 0, \quad q_r = 0.$$

We apply periodic boundary conditions in horizontal direction and the following conditions vertically: $\mathbf{u} \cdot \mathbf{n} = 0$, $\nabla \rho' \cdot \mathbf{n} = 0$, $\nabla(\rho q_\ell) \cdot \mathbf{n} = 0$, $\ell \in \{v, c, r\}$ with the Dirichlet boundary conditions for the potential

temperature,

$$\theta(x_3 = 0) = 284 \text{ K} \quad \text{and} \quad \theta(x_3 = 1000) = 283 \text{ K}.$$

Example 2: 2-D case. In Figures 4–7, we present time snapshots of the numerical solution computed in a domain $\Omega = [0, 5000] \times [0, 1000] \text{ m}^2$ that has been discretized using 320×320 mesh cells. The potential temperature, water vapor mixing ratio, cloud mass and rain mass concentration are presented at two distinct times ($t = 800$ and 1400 s) in Figures 4–7, where we can clearly see the formation and evolution of a pattern. At an earlier time $t = 800 \text{ s}$, one can observe the formation of small convective cells, visible as narrow finger-like structures reaching towards the top of the domain. Inside these cells, the potential temperature is enhanced, partly due to the upward transport of higher values from below and partly due to phase changes and thus latent heat release. Also the mass concentrations of water vapor and cloud water follow the small scale structure and show enhanced values inside the fingers. Even at this early stage, rain can be formed at the top of the domain, since there the cloud water concentration is high enough for autoconversion and accretion. Nevertheless, the structure of rain water is very different since after the formation of rain, it is vertically transported due to sedimentation leading to vertically smeared structures. By a later time $t = 1400 \text{ s}$, much larger structures, which are similar to classical structures for dry thermal convection, have been formed. In the variables θ , q_v and q_c , the spatially extended convective cells can be clearly seen. In contrast, rain water is not following the convective structure although some larger features can be seen. In general, smearing due to sedimentation is again a major feature of the rain mass concentration.

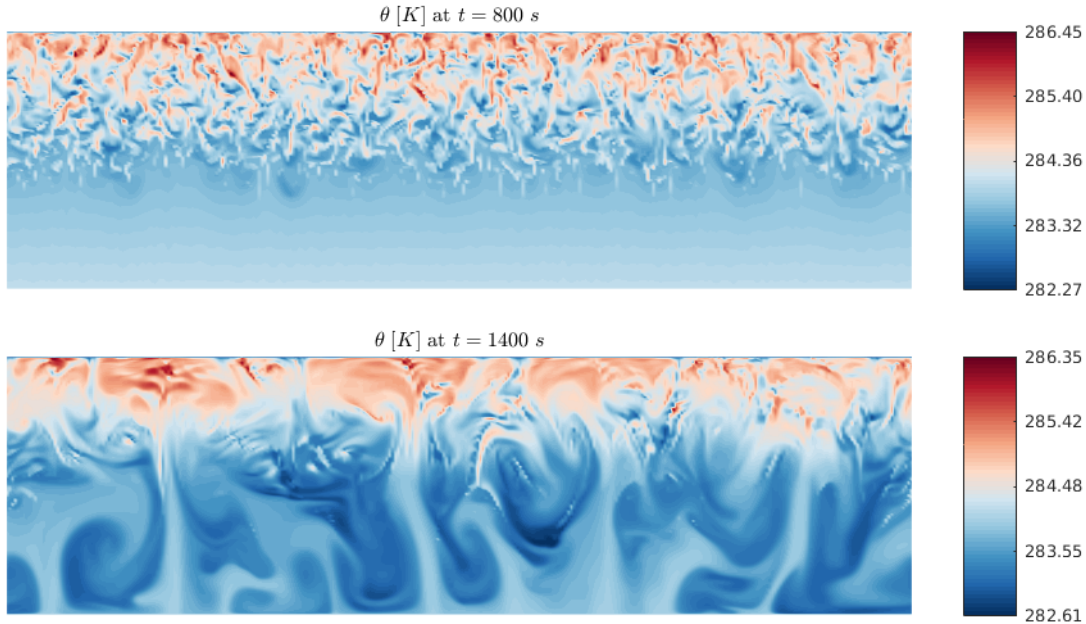


FIGURE 4. *Example 2: Time evolution of the potential temperature θ .*

Example 3: 3-D case. In this example, we compute the numerical solution in a domain $\Omega = [0, 5000] \times [0, 5000] \times [0, 1000] \text{ m}^3$ that has been discretized using $50 \times 50 \times 50$ mesh cells. Without moisture, this would be the classical setup for dry Rayleigh-Bénard convection. While for the dry case phase diagrams for resulting patterns, as e.g. rolls and hexagons, are available, see, e.g., [7], only little is known about patterns in moist Rayleigh-Bénard convection. Few studies with reduced order models indicate the possibility of roll-like structures or the formation of hexagons, e.g. [44], but no results for compressible Navier-Stokes equations coupled with full cloud microphysics are available, at least to our knowledge. Thus, we would expect similar patterns as evolving in dry convection, although the latent heat release might change these patterns partially. Figures 8–11 show θ , q_v , q_c and q_r computed at times $t = 1200, 1600, 1800$ and 2000 s . In order to better visualize the computed structures, we have plotted the solution in a slightly smaller domain

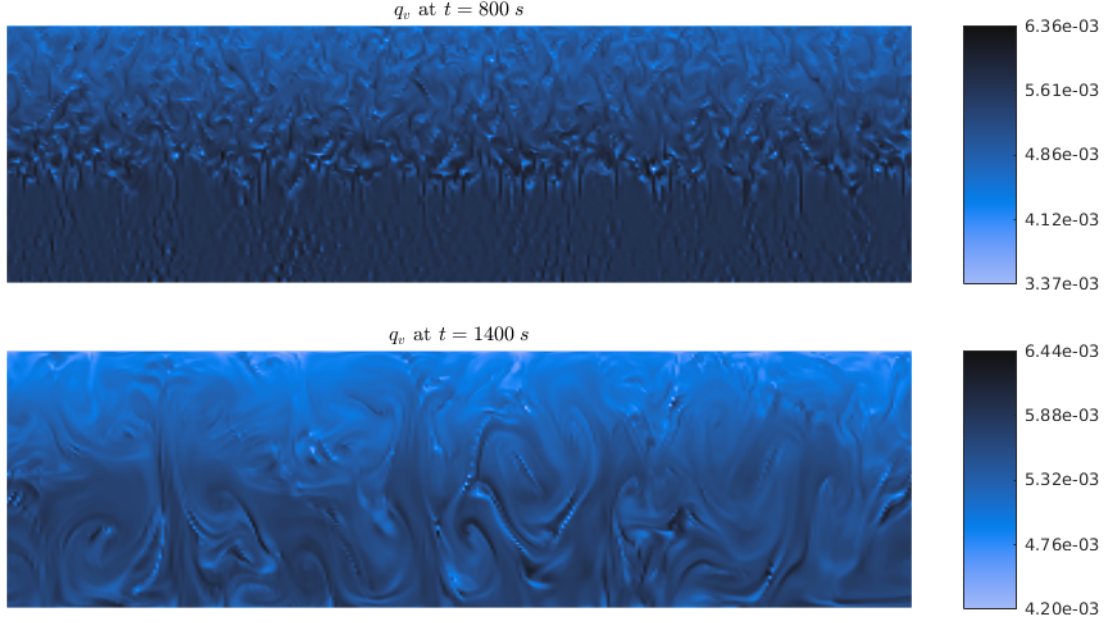


FIGURE 5. *Example 2: Time evolution of the water vapor concentration q_v .*

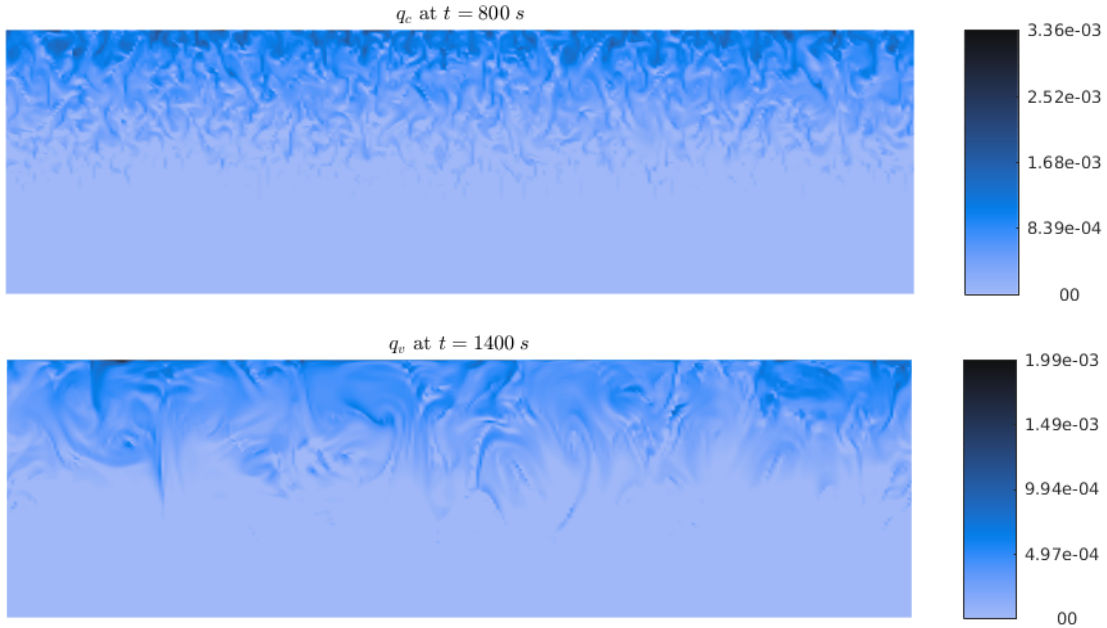


FIGURE 6. *Example 2: Time evolution of the cloud drops concentration q_c .*

$[0, 5000] \times [0, 5000] \times [0, 980] \text{ m}^3$. As in the 2-D case, different structures are formed in the different variables. At an earlier time $t = 1200 \text{ s}$, small scale structures can be seen at the top layer of the domain, especially in the potential temperature (Figure 8) and cloud water (Figure 10). As time progresses, these structures aggregate and reorganize to quasi-hexagonal structures, which would be typical for classical dry thermal convection—Rayleigh-Bénard convection. These structures in potential temperature seem to be quite robust as the overall scales and pattern do not change from $t = 1800 \text{ s}$ to $t = 2000 \text{ s}$. As one can see in Figure

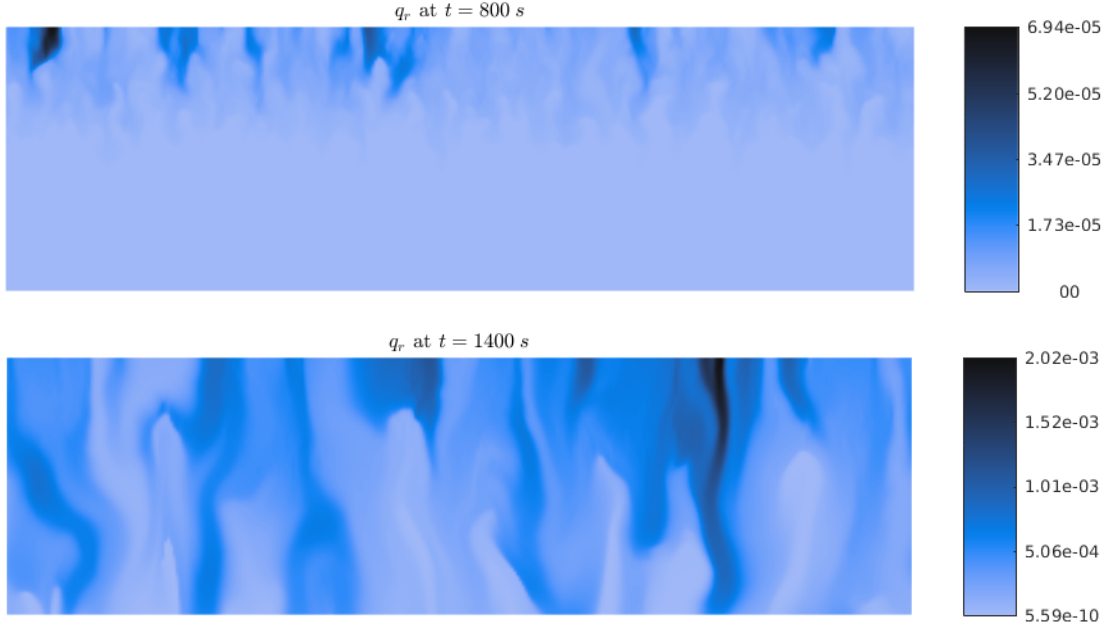


FIGURE 7. *Example 2: Time evolution of the rain concentration q_r .*

10, the structures in the cloud water are very similar since the variables θ and q_c are closely connected; the structures are mainly visible in horizontal planes. For the rain distribution, the evolved structures are quite different since rain is formed at regions with high cloud water (that is, at the top layers) and then transported by sedimentation leading to more pronounced pattern in the vertical direction, since sedimentation is the dominant process after the rain has formed.

Remark 4.1. Let us note that the Rayleigh-Bénard convection can be understood as a very simplified model for atmospheric convection in the turbulent planetary boundary layer. In [32, 45], numerical simulations for moist Rayleigh-Bénard convection have been realized using the Boussinesq approximation, a simplified equation of state, and the rigid-lid boundary conditions at the top and bottom of the computational domain. Our mathematical model is more general and takes weakly compressible effects into account. Numerical experiments presented in Examples 2 and 3 are in good agreement with the results presented in the literature, but the focus of those studies differs from ours.

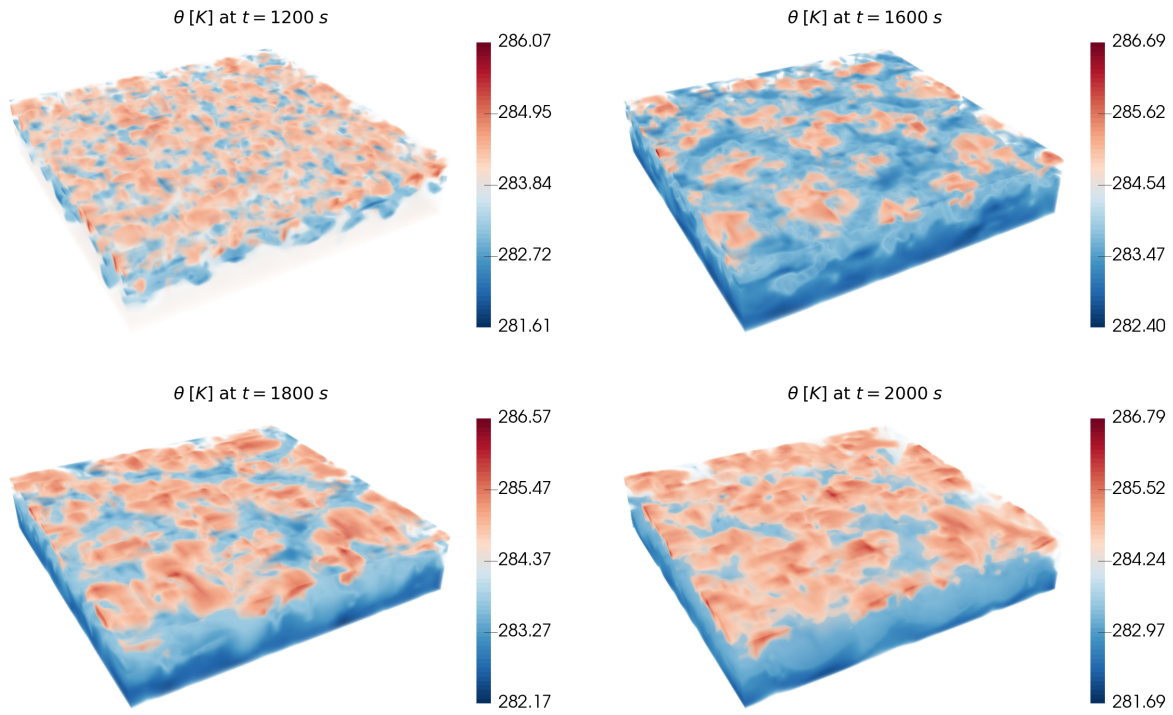


FIGURE 8. *Example 3: Time evolution of the potential temperature θ .*

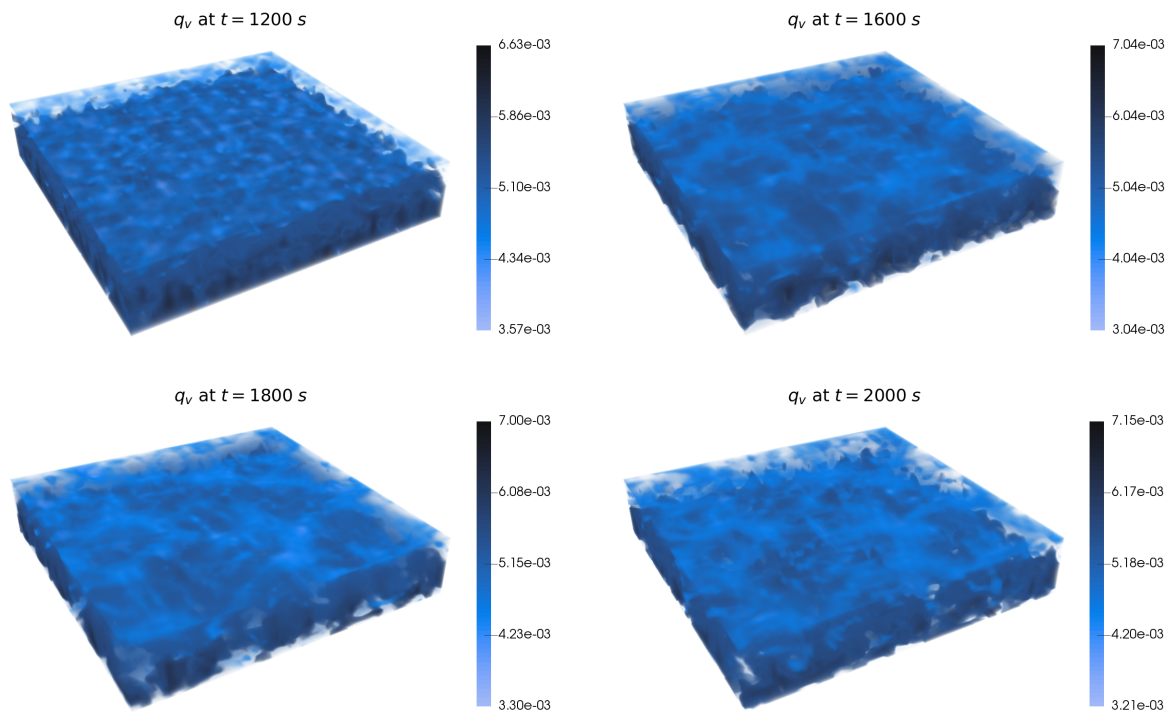


FIGURE 9. *Example 3: Time evolution of the water vapor concentration q_v .*

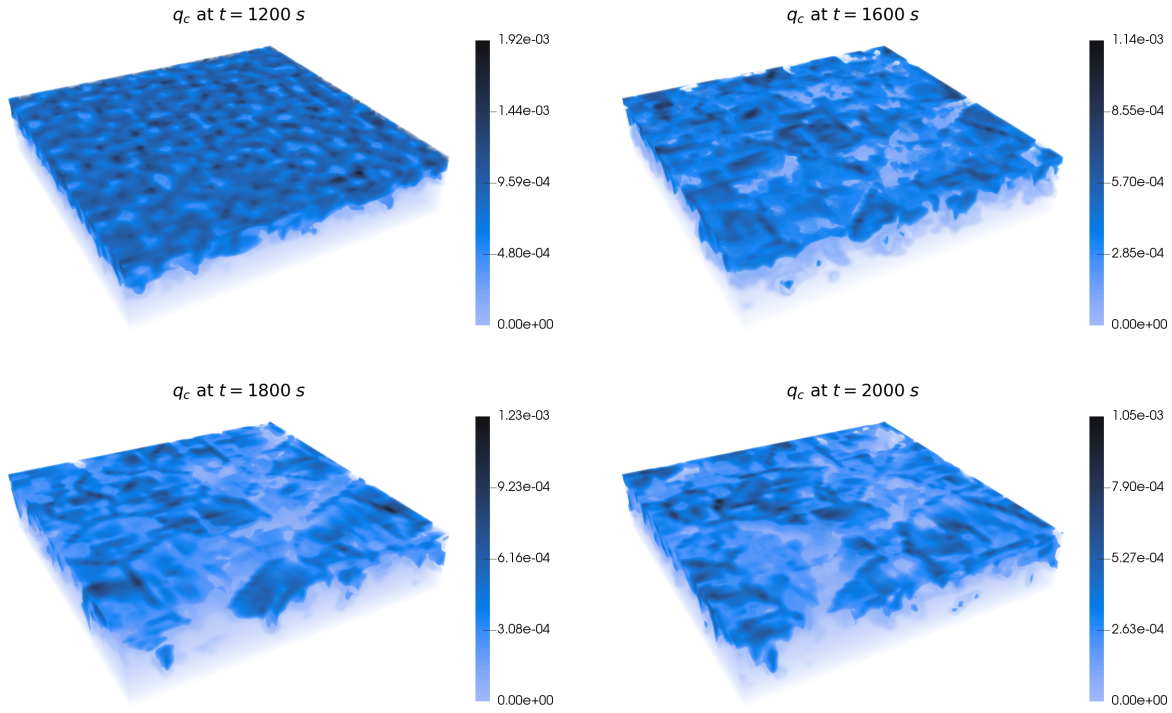


FIGURE 10. *Example 3: Time evolution of the cloud drops concentration q_c .*

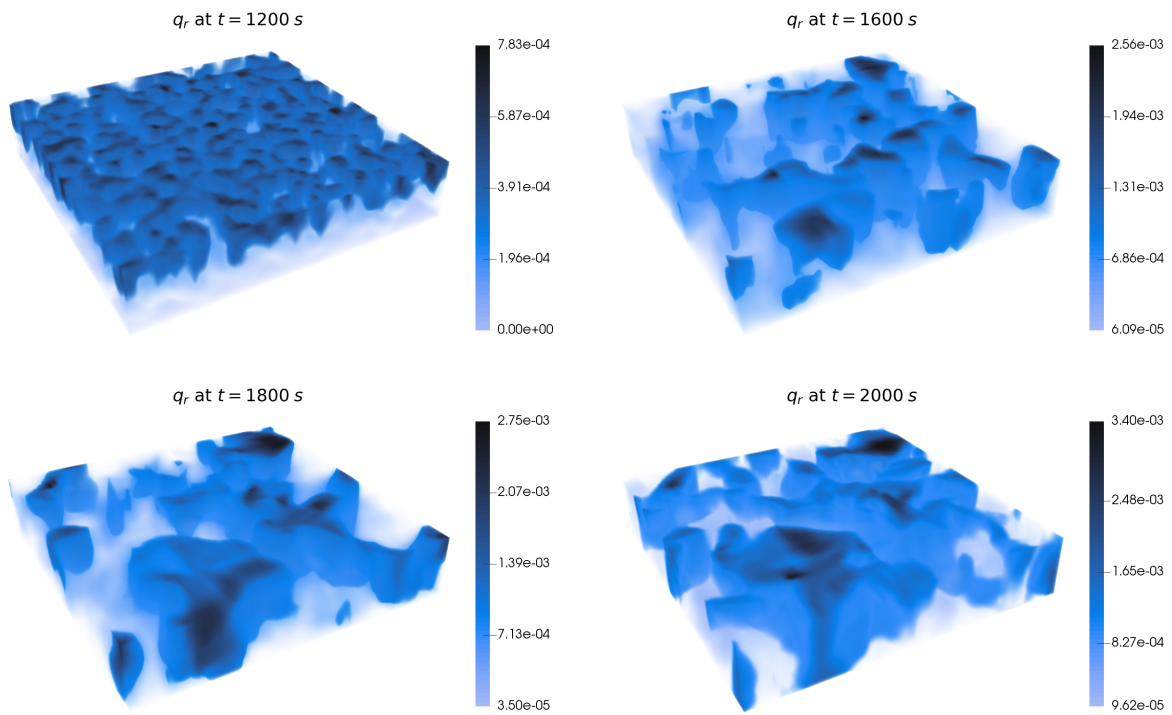


FIGURE 11. *Example 3: Time evolution of the rain concentration q_r .*

5. STOCHASTIC MATHEMATICAL MODEL

In meteorological applications, it is typical that initial or boundary data as well as parameters are uncertain. In order to take this into account and analyze the influence of such uncertainties on the output of numerical simulations, we apply the stochastic Galerkin method that will be described below. In this paper, we consider the case where the uncertainty arises from the initial data or some coefficients in the microphysical cloud parametrizations. In order to mathematically describe the uncertainty, we introduce a random variable ω . We assume that either the initial data or some well-chosen model parameters depend on ω , that is,

$$(\rho q_\ell)|_{t=0} = (\rho q_\ell)(\mathbf{x}, t = 0, \omega) \quad \text{with } \ell \in \{v, c, r\}$$

or

$$k_1 = k_1(\omega), \quad k_2 = k_2(\omega), \quad \alpha = \alpha(\omega).$$

Consequently, the solution at later time will also depend on ω , that is, $(\rho q_\ell)(\mathbf{x}, t, \omega)$ for $\ell \in \{v, c, r\}$, and the system (2.2) for cloud variables will read as

$$\begin{aligned} ((\rho q_v)(\omega))_t + \nabla \cdot ((\rho q_v)(\omega) \mathbf{u} - \mu_q \rho \nabla q_v(\omega)) &= \rho(-C(\omega) + E(\omega)), \\ ((\rho q_c)(\omega))_t + \nabla \cdot ((\rho q_c)(\omega) \mathbf{u} - \mu_q \rho \nabla q_c(\omega)) &= \rho(C(\omega) - A_1(\omega) - A_2(\omega)), \\ ((\rho q_r)(\omega))_t + \nabla \cdot ((\rho q_r)(\omega)(-v_q(\omega) \mathbf{e}_3 + \mathbf{u}) - \mu_q \rho \nabla q_r(\omega)) &= \rho(A_1(\omega) + A_2(\omega) - E(\omega)). \end{aligned} \quad (5.1)$$

From now on we will stress the dependence on ω , but we will omit the dependence on \mathbf{x} and t to simplify the notation. We would like to point out that the solution of the Navier-Stokes equations (2.8) will also depend on ω , because of the source term S_θ . In this paper, we will consider a simplified situation by replacing

$$S_\theta(\omega) = \rho \frac{L\theta}{c_p T} \{C((\rho q_v)(\omega), (\rho q_c)(\omega)) - E((\rho q_v)(\omega), (\rho q_r)(\omega))\}$$

in (2.8) by \bar{S}_θ which only depends on the expected values of the cloud variables

$$\bar{S}_\theta := \rho \frac{L\theta}{c_p T} \{C(\mathbb{E}[\rho q_v], \mathbb{E}[\rho q_c]) - E(\mathbb{E}[\rho q_v], \mathbb{E}[\rho q_r])\}.$$

This ensures that all of the fluid variables, ρ' , $\rho \mathbf{u}$ and $(\rho \theta)'$, remain deterministic.

6. NUMERICAL SCHEME FOR THE STOCHASTIC MODEL

In this section, we describe a generalized polynomial chaos stochastic Galerkin (gPC-SG) method for the system of cloud equations (5.1). Such method belongs to the class of intrusive methods and the use of the Galerkin expansion leads to a system of deterministic equations for the expansion coefficients. In the gPC-SG method, the solution is sought in the form of a polynomial expansion

$$\rho q_\ell(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{\rho q_\ell})_k(\mathbf{x}, t) \Phi_k(\omega) \quad \text{with } \ell \in \{v, c, r\}, \quad M \geq 0, \quad (6.1)$$

where $\Phi_k(\omega)$, $k = 0, \dots, M$, are polynomials of k -th degree that are orthogonal with respect to the probability density function $\mu(\omega)$. The choice of the orthogonal polynomials $\{\Phi_k(\omega)\}_{k=0}^M$ depends on the distribution of ω . In our case, we use a uniformly distributed $\omega \in \Gamma = (-1, 1)$, which defines the Legendre polynomials, which satisfy

$$\int_{\Gamma} \Phi_k(\omega) \Phi_{k'}(\omega) \mu(\omega) d\omega = \frac{1}{2k+1} \delta_{kk'} \quad \text{for } 0 \leq k, k' \leq M, \quad (6.2)$$

where $\delta_{kk'}$ is the Kronecker symbol and Γ is the sample space.

We use the same expansion for the uncertain coefficients,

$$k_1(\omega) = \sum_{k=0}^M (\widehat{k_1})_k \Phi_k(\omega), \quad k_2(\omega) = \sum_{k=0}^M (\widehat{k_2})_k \Phi_k(\omega), \quad \alpha(\omega) = \sum_{k=0}^M \widehat{\alpha}_k \Phi_k(\omega), \quad (6.3)$$

for the source terms on the RHS of (5.1),

$$\begin{aligned}
\rho(-C(\mathbf{x}, t, \omega) + E(\mathbf{x}, t, \omega)) &=: R_1(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{r}_1)_k(\mathbf{x}, t) \Phi_k(\omega), \\
\rho(C(\mathbf{x}, t, \omega) - A_1(\mathbf{x}, t, \omega) - A_2(\mathbf{x}, t, \omega)) &=: R_2(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{r}_2)_k(\mathbf{x}, t) \Phi_k(\omega), \\
\rho(A_1(\mathbf{x}, t, \omega) + A_2(\mathbf{x}, t, \omega) - E(\mathbf{x}, t, \omega)) &=: R_3(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{r}_3)_k(\mathbf{x}, t) \Phi_k(\omega),
\end{aligned} \tag{6.4}$$

as well as for the raindrop fall velocity,

$$v_q(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{v}_q)_k(\mathbf{x}, t) \Phi_k(\omega). \tag{6.5}$$

Since $\rho(\mathbf{x}, t) = \widehat{\rho}_0(\mathbf{x}, t)$, we also obtain

$$q_\ell(\mathbf{x}, t, \omega) = \sum_{k=0}^M (\widehat{q}_\ell)_k(\mathbf{x}, t) \Phi_k(\omega) \quad \text{with} \quad (\widehat{q}_\ell)_k(\mathbf{x}, t) = \frac{(\widehat{\rho q}_\ell)_k(\mathbf{x}, t)}{\rho(\mathbf{x}, t)} \quad \text{for } \ell \in \{v, c, r\}, \quad k = 1, \dots, M. \tag{6.6}$$

We note that if $\rho(\mathbf{x}, t)$ is very small, the computation of the coefficients $(\widehat{q}_\ell)_k(\mathbf{x}, t)$ should be desingularized; see [21, formulae (5.16)–(5.18)].

Applying the Galerkin projection to (5.1) yields

$$\begin{aligned}
\langle (\rho q_v)_t + \nabla \cdot (\rho q_v \mathbf{u} - \mu_q \rho \nabla q_v), \Phi_k \rangle &= \langle \rho(-C + E), \Phi_k \rangle, \\
\langle (\rho q_c)_t + \nabla \cdot (\rho q_c \mathbf{u} - \mu_q \rho \nabla q_c), \Phi_k \rangle &= \langle \rho(C - A_1 - A_2), \Phi_k \rangle, \\
\langle (\rho q_r)_t + \nabla \cdot (\rho q_r(-v_q \mathbf{e}_3 + \mathbf{u}) - \mu_q \rho \nabla q_r), \Phi_k \rangle &= \langle \rho(A_1 + A_2 - E), \Phi_k \rangle,
\end{aligned} \tag{6.7}$$

for $k = 0, \dots, M$, where $\langle \cdot, \cdot \rangle$ is the scalar product in our probability space which is given through

$$\langle u, v \rangle = \int_{-1}^1 u(\omega) v(\omega) \mu(\omega) d\omega.$$

We now substitute (6.1), (6.4)–(6.6) into (6.7) and use the orthogonality property (6.2) to obtain the following $3(M+1) \times 3(M+1)$ deterministic system for the gPC coefficients:

$$\begin{aligned}
\frac{\partial}{\partial t} (\widehat{\rho q}_v)_k + \sum_{s=1}^d \frac{\partial}{\partial x_s} ((\widehat{\rho q}_v)_k u_s) - \mu_q \sum_{s=1}^d \left(\frac{\partial \rho}{\partial x_s} \frac{\partial}{\partial x_s} (\widehat{q}_v)_k + \rho \frac{\partial^2}{\partial x_s^2} (\widehat{q}_v)_k \right) &= (\widehat{r}_1)_k, \\
\frac{\partial}{\partial t} (\widehat{\rho q}_c)_k + \sum_{s=1}^d \frac{\partial}{\partial x_s} ((\widehat{\rho q}_c)_k u_s) - \mu_q \sum_{s=1}^d \left(\frac{\partial \rho}{\partial x_s} \frac{\partial}{\partial x_s} (\widehat{q}_c)_k + \rho \frac{\partial^2}{\partial x_s^2} (\widehat{q}_c)_k \right) &= (\widehat{r}_2)_k, \\
\frac{\partial}{\partial t} (\widehat{\rho q}_r)_k - \frac{\partial}{\partial x_d} \widehat{\alpha}_k + \sum_{s=1}^d \frac{\partial}{\partial x_s} ((\widehat{\rho q}_r)_k u_s) - \mu_q \sum_{s=1}^d \left(\frac{\partial \rho}{\partial x_s} \frac{\partial}{\partial x_s} (\widehat{q}_r)_k + \rho \frac{\partial^2}{\partial x_s^2} (\widehat{q}_r)_k \right) &= (\widehat{r}_3)_k,
\end{aligned} \tag{6.8}$$

for $k = 0, \dots, M$. Here, the coefficients $\{\widehat{\alpha}_k\}_{k=0}^M$ are obtained using the following expansion:

$$v_q(\mathbf{x}, t, \omega) (\rho q_r)(\mathbf{x}, t, \omega) = \sum_{j=0}^M (\widehat{v}_q)_j(\mathbf{x}, t) \Phi_j(\omega) \sum_{m=0}^M (\widehat{\rho q}_r)_m(\mathbf{x}, t) \Phi_m(\omega) =: \sum_{k=0}^M \widehat{\alpha}_k(\mathbf{x}, t) \Phi_k(\omega).$$

The coefficients $\{(\widehat{r}_1)_k, (\widehat{r}_2)_k, (\widehat{r}_3)_k\}_{k=0}^M$, as well as $\{\widehat{\alpha}_k\}_{k=0}^M$ are calculated via discrete Legendre transform (DLT) and inverse discrete Legendre transform (IDLTL), which can be briefly described as follows.

- DLT: First, the Galerkin projection applied to the expansion $f(\mathbf{x}, t, \omega) = \sum_{k=0}^M \widehat{f}_k(\mathbf{x}, t) \Phi_k(\omega)$ yields

$$\widehat{f}_k(\mathbf{x}, t) = \frac{2k+1}{2} \int_{-1}^1 f(\mathbf{x}, t, \omega) \Phi_k(\omega) d\omega \quad \text{for } 0 \leq k \leq M. \quad (6.9)$$

Then, approximating the integral in (6.9) using the Gauss-Legendre quadrature leads to

$$\text{DLT} \left[\{f(\mathbf{x}, t, \omega_\ell)\}_{\ell=0}^M \right] = \left\{ \widehat{f}_k(\mathbf{x}, t) \right\}_{k=0}^M = \left\{ \frac{2k+1}{2} \sum_{l=0}^M \beta_l f(\mathbf{x}, t, \omega_\ell) \Phi_k(\omega_\ell) \right\}_{k=0}^M,$$

where β_ℓ are the Gauss-Legendre quadrature weights and ω_ℓ is the ℓ -th root of Φ_{M+1} .

- IDLT: Given the coefficients $\{\widehat{f}_k\}_{k=0}^M$, we compute the function f through the gPC expansion

$$\text{IDLT} \left[\left\{ \widehat{f}_k(\mathbf{x}, t) \right\}_{k=0}^M \right] = \{f(\mathbf{x}, t, \omega_\ell)\}_{\ell=0}^M = \left\{ \sum_{k=0}^M \widehat{f}_k(\mathbf{x}, t) \Phi_k(\omega_\ell) \right\}_{\ell=0}^M.$$

Consequently, we obtain

$$\{(\widehat{r}_1)_k\}_{k=0}^M = \text{DLT} \left[R_1 \left(\text{IDLT} \left[\{(\widehat{\rho q}_v)_k\}_{k=0}^M \right], \text{IDLT} \left[\{(\widehat{\rho q}_c)_k\}_{k=0}^M \right], \text{IDLT} \left[\{(\widehat{\rho q}_r)_k\}_{k=0}^M \right] \right) \right],$$

and analogously for $\{(\widehat{r}_2)_k\}_{k=0}^M$, $\{(\widehat{r}_3)_k\}_{k=0}^M$ and $\{\widehat{\alpha}_k\}_{k=0}^M$.

Remark 6.1. We stress that since the values $\Phi_k(\omega_\ell)$, $0 \leq k, \ell \leq M$, are needed every time either DLT or IDLT is applied, one can pre-compute them for the code efficiency.

For the spatial and temporal discretizations of the system (6.8), we apply the same finite volume method as described in Section 3.2 and the same large stability domain explicit time integration method mentioned in Section 3.3. As in the deterministic case, we implement the ODE solver DUMKA3, which we provide with the following time step stability restriction for the forward Euler method:

$$\max_{s=1,2} \max_{i=1,\dots,N} (|(u_s)_i|, |(u_3)_i + v_q(\omega_l)|) \frac{\Delta t_{\text{cloud}}^n}{h} < 0.5,$$

which should be satisfied for all of the Legendre roots ω_ℓ , $\ell = 0, \dots, M$.

7. STOCHASTIC NUMERICAL EXPERIMENTS

In this section, we conduct numerical experiments with the stochastic Galerkin method described in Section 6 for the free convection of a moist warm air bubble and the Rayleigh-Bénard convection. We demonstrate the influence of uncertainty in initial data as well as in cloud parameters on the solution of the coupled Navier-Stokes-cloud model (2.8), (5.1). In all of our numerical examples below, we take $M = 3$. Our extensive tests, from which we present here only a selected part, showed that this was sufficient. Indeed, as documented in Example 6, high-order stochastic coefficients typically have very small influence on a solution (see Figure 15), and thus can be neglected. Similar behavior was observed in other experiments.

7.1. Free convection of a smooth warm air bubble. In this test, we modify Example 1 by randomly perturbing either the initial data or selected model parameters.

Example 4: 2-D case with stochastic initial data. We begin by considering the following experiment with a 10% perturbation of the initial water vapor concentration:

$$\begin{aligned} (\widehat{q}_v)_0(\mathbf{x}, 0) &= 0.02\theta'(\mathbf{x}, 0), \quad (\widehat{q}_v)_1(\mathbf{x}, 0) = 0.1(\widehat{q}_v)_0(\mathbf{x}, 0), \quad (\widehat{q}_v)_k(\mathbf{x}, 0) = 0 \quad \text{for } 2 \leq k \leq M, \\ (\widehat{q}_c)_k(\mathbf{x}, 0) &= (\widehat{q}_r)_k(\mathbf{x}, 0) = 0 \quad \text{for } 0 \leq k \leq M. \end{aligned}$$

We compute the solution using different meshes until the final time $t = 10s$.

The experimental convergence study for the cloud and flow variables is presented in Figure 12. Similarly to the deterministic case, one can observe second-order convergence in space and time. In order to test the convergence in the stochastic space, we obtain a reference solution computed by the stochastic Galerkin

method with 20 stochastic modes. The convergence study is presented in Figure 13, where we plot the difference between the approximate and reference solutions, both computed using a mesh with 160×160 cells and $\Delta t = 0.01$ at time $t = 10s$. One can clearly see a spectral convergence with the rate $e^{-0.3M}$.

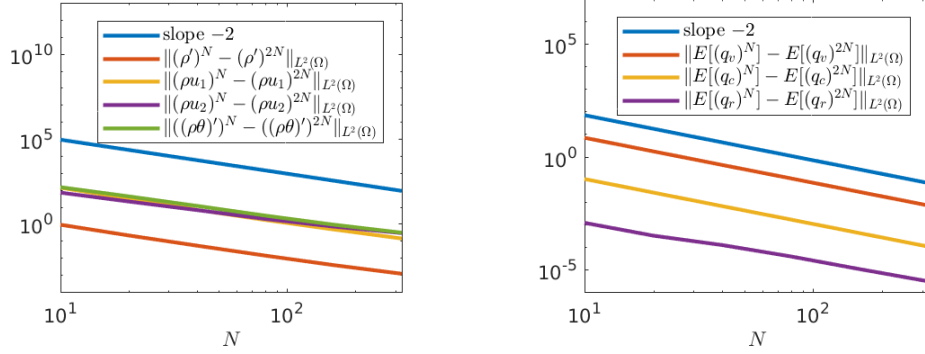


FIGURE 12. *Example 4: Convergence study for the flow variables (left) and the expected values of $(\rho q_\ell)_k$, $\ell \in \{v, c, r\}$ (right) in space and time.*

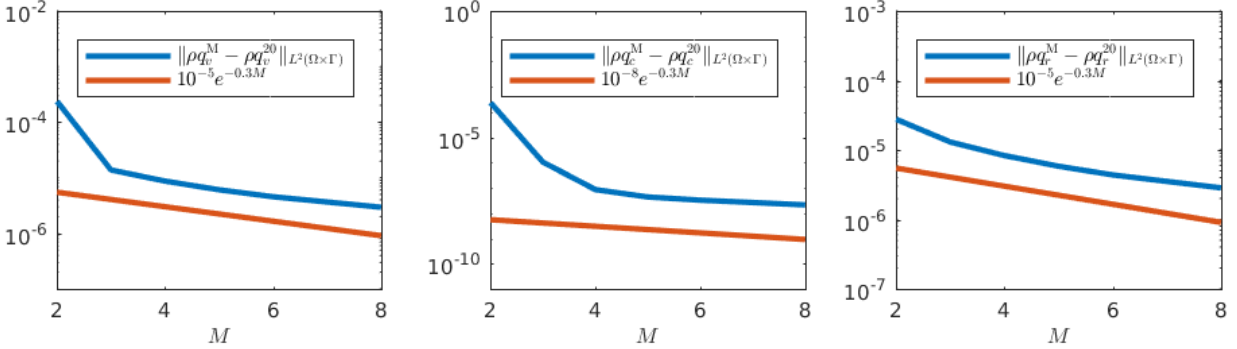


FIGURE 13. *Example 4: Convergence study for the cloud variables (ρq_ℓ) , $\ell \in \{v, c, r\}$ in the stochastic space.*

Example 5: 2-D case with stochastic parameters. In this experiment, we perturb the three selected model parameters (k_1 , k_2 and α) by 50% each:

$$\begin{aligned} (\hat{k}_1)_0 &= 0.0041, & (\hat{k}_1)_1 &= 0.5(\hat{k}_1)_0, & (\hat{k}_1)_k &= 0 \quad \text{for } 2 \leq k \leq M, \\ (\hat{k}_2)_0 &= 0.8, & (\hat{k}_2)_1 &= 0.5(\hat{k}_2)_0, & (\hat{k}_2)_k &= 0 \quad \text{for } 2 \leq k \leq M, \\ \hat{\alpha}_0 &= 190.3, & \hat{\alpha}_1 &= 0.5\hat{\alpha}_0, & \hat{\alpha}_k &= 0 \quad \text{for } 2 \leq k \leq M. \end{aligned}$$

These parameters were proposed in [35] as the most sensitive model parameters. We study the convergence in the stochastic space. To this end, we plot in Figure 14 the difference between the approximate and reference (obtained with 20 stochastic modes) solutions, both computed using a mesh with 160×160 cells and $\Delta t = 0.01$ at time $t = 10s$. As in Example 4, one can observe a spectral convergence with the rate $e^{-0.3M}$.

Example 6: Comparison of the stochastic Galerkin and stochastic collocation methods in the 3-D case. In this comparison test, we consider free convection of a moist smooth warm air bubble with perturbed initial data as in Example 4. Numerical solutions are computed on $50 \times 50 \times 50$ mesh at time $t = 1s$. We compare the performance of the stochastic Galerkin and stochastic collocation methods. The collocation method (see, e.g., [9]) is an interpolation method in the stochastic space, which uses a deterministic model with the values of the stochastic variable taken at collocation points suitably chosen on the interval $(-1, 1)$; here, we use the Gauss-Legendre points.

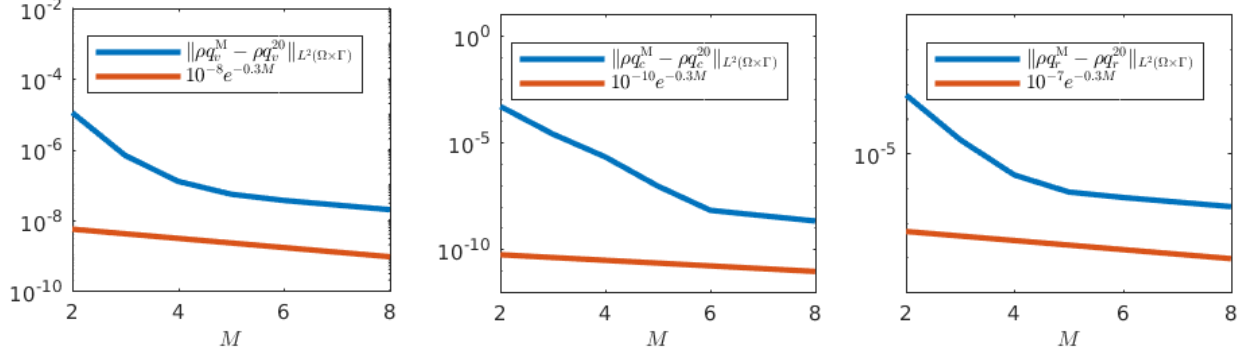


FIGURE 14. Example 5: Convergence study for the cloud variables (ρq_ℓ) , $\ell \in \{v, c, r\}$ in the stochastic space.

In Figure 15, the $L^2(\Omega)$ -norms of the stochastic coefficients of the stochastic Galerkin ($\|(\widehat{\rho q_\ell})_m\|_{L^2(\Omega)}$) and stochastic collocation ($\|(\widetilde{\rho q_\ell})_m\|_{L^2(\Omega)}$) methods for $\ell \in \{v, c, r\}$ and $m = 0, \dots, 19$ are shown. One can observe an exponential decay with respect to m and good agreement between both methods that demonstrates the reliability of the stochastic Galerkin method. We note, however, that the norms of the solutions computed by these two different methods are not equal since the stochastic Galerkin method uses the expected values of the cloud variables in the Navier-Stokes equations.

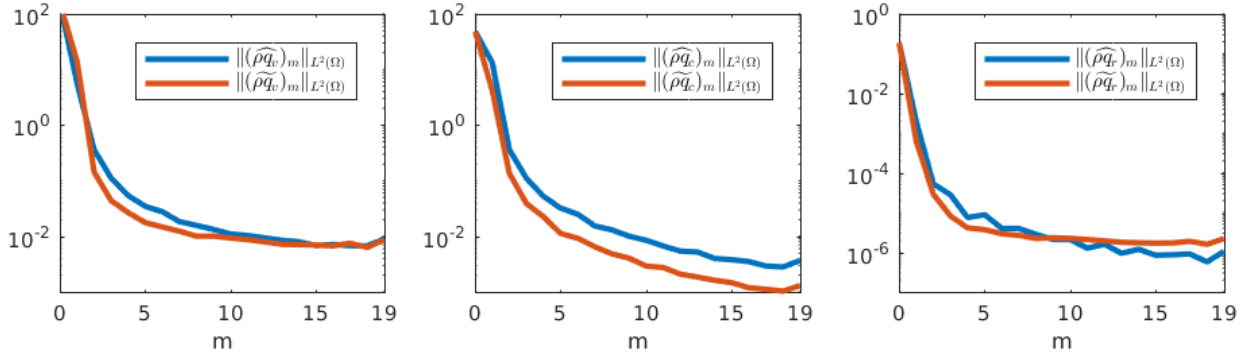


FIGURE 15. Example 6: $L^2(\Omega)$ -norms of the stochastic coefficients computed by the stochastic Galerkin ($(\widehat{\rho q_\ell})_m$) and stochastic collocation ($(\widetilde{\rho q_\ell})_m$) methods, $\ell \in \{v, c, r\}$.

In Figure 16, we compare the CPU times consumed by the stochastic Galerkin and stochastic collocation methods with the same number M of stochastic modes/collocation points. Since the stochastic Galerkin method solves the Navier-Stokes equations just once instead of M times, as needed by the stochastic collocation method does, it is expected to outperform the stochastic collocation method. This has been confirmed by our simulations.

7.2. Rayleigh-Bénard convection. In this section, we present results of uncertainty study for the Rayleigh-Bénard convection in both 2-D and 3-D. We investigate uncertainty propagation, which is triggered either by the initial data or cloud parameters.

Example 7: 2-D case with stochastic initial data. In this experiment, we choose the same initial data for the flow variables as in Section 4.2 and perturb the initial data in q_v by 5%, 10%, 20% and 50%:

$$\begin{aligned} (\widehat{q}_v)_0(\mathbf{x}, 0) &= 2 \cdot 10^{-5} \bar{\theta}, \quad (\widehat{q}_v)_1(\mathbf{x}, 0) = \nu (\widehat{q}_v)_0(\mathbf{x}, 0), \quad (\widehat{q}_v)_k = 0 \quad \text{for } 2 \leq k \leq M, \\ (\widehat{q}_c)_k(\mathbf{x}, 0) &= (\widehat{q}_r)_k(\mathbf{x}, 0) = 0 \quad \text{for } 0 \leq k \leq M, \end{aligned} \quad (7.1)$$

where $\nu = 0.05, 0.1, 0.2$ and 0.5 , respectively. It should be observed that uniform perturbations in the initial conditions for q_v may lead to either reduced or enhanced water vapor concentrations as compared to

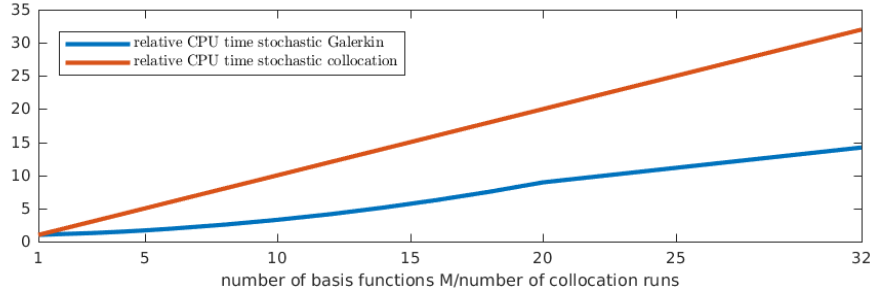


FIGURE 16. *Example 6: Relative CPU times consumed by the stochastic Galerkin and stochastic collocation methods.*

the deterministic simulations. Since the temperature gradient is quite small, a change in q_v translates to an (almost) linear change in saturation ratio, which directly controls cloud formation. Thus, in the case of positive perturbations, a higher water vapor concentration leads to earlier cloud formation and, in addition, a higher potential temperature change since more water is available in the system. On the other hand, lower values of q_v lead to a time delay in the formation of clouds, even if small convective cells are driven by the dry unstable situation. In a feedback cycle, a reduced or even delayed formation of cloud water propagates further to a weaker rain formation. Finally, the evaporation of rain water leads to a strong cooling effect of the lower layers of the domain, which also crucially depends on the amount of sedimenting rain water. These effects have to be taken into account for the evaluation of the different perturbation scenarios.

Numerical solutions for both the potential temperature θ (Figure 17) and the expected value of the cloud drops concentration q_c (Figure 18) are computed at time $t = 1400s$ using 320×320 mesh cells and presented for $\nu = 0, 0.05, 0.1, 0.2$ and 0.5 . For a better comparison, we have used the same range of values for different perturbations in all of the plots.

From the results shown in Figure 17 for the potential temperature θ , one can observe two major features that increase in strength with increasing strength of perturbations. First, larger perturbations lead to more concise fine structures, that is, for the deterministic simulation ($\nu = 0$), the variable is quite smooth, whereas for large perturbations, large gradients on a very small scale appear. This feature can also be recognized for the cloud water, that is, for the expected values $\mathbb{E}[q_c]$ of cloud water concentrations in Figure 18. This is probably due to the fact that even small variations in water vapor have a strong impact on cloud formation, since the activation of cloud droplet is basically a threshold process. For values closer to saturation, even small variations in vertical upward motions can trigger cloud formation. Thus, the small scale variations are more prominent in the massively perturbed scenarios. The second feature is even more striking. Increasing the perturbation in the initial water vapor distribution leads to a stronger vertical gradient in potential temperature, that is, at low levels the temperature is much smaller than in the deterministic case. This feature is due to cooling of sedimenting rain water. For simulations with a high water vapor loading, more cloud and thus more rain is formed, which is subsequently falling down into lower levels and cools the environment by evaporation. Since this process is very effective, the temperature can be reduced drastically. Note that this feature is well-known for the real atmosphere: Falling rain can cool lower levels efficiently, so that a transition from melting rain droplets to snow can be possible for winter seasons. The efficient formation of rain also leads to a strong reduction in cloud water, since accretion can eat up cloud droplets in the lower level; see Figure 18. For reduced values of initial water vapor, processes of cloud and precipitation formation is strongly reduced. However, on average the positive perturbations dominate the statistical picture.

In Figure 19, we plot the standard deviation of the cloud drops concentration. Here, we compute the standard deviation for a function $f(\mathbf{x}, t, \omega) = \sum_{k=0}^M \widehat{f}_k(\mathbf{x}, t) \Phi_k(\omega)$ as

$$\sigma(f(\mathbf{x}, t, \omega)) = \sqrt{\sum_{k=1}^M \frac{1}{2k+1} \widehat{f}_k(\mathbf{x}, t)^2}, \quad (7.2)$$

which follows from the orthogonal property (6.2).

In order to investigate the influence of perturbations, the time evolution of the mean expected value per m^2 as well as the mean standard deviation per m^2 for the cloud variables are presented in Figure 20. In d -space dimensions these quantities can be computed in the following way:

$$\begin{aligned} \mathbb{E} \left[\frac{h^d}{|\Omega|} \sum_{i=1}^N (q_\ell)_i \right] &= \frac{h^d}{|\Omega|} \sum_{i=1}^N \mathbb{E} [(q_\ell)_i] = \frac{h^d}{|\Omega|} \sum_{i=1}^N \overline{((q_\ell)_i)_0}, \\ \sigma \left(\frac{h^d}{|\Omega|} \sum_{i=1}^N (q_\ell)_i \right) &= \frac{h^d}{|\Omega|} \sqrt{\sum_{k=1}^M \left(\sum_{i=1}^N \overline{((q_\ell)_i)_k} \right)^2 \frac{1}{2k+1}}, \end{aligned}$$

where N is the number of mesh cells and $\ell \in \{v, c, r\}$.

These averaged quantities show the qualitative difference in the different perturbation scenarios, as already described above. For low perturbations, the difference of the expectation value is quite small and also the standard deviation remains small. For increasing perturbations, the spread is increased. As noted before, the averaged quantities are dominated by the positive perturbations, leading to (i) earlier cloud formation, (ii) thicker clouds due to more available water vapor, and (iii) to enhanced rain formation. These three features can be seen very nicely in the strongest perturbations ($\nu = 0.5$), with a large drop in water vapor concentration accompanied by a strong increase in cloud water and earlier onset of precipitation. We would also like to note that the spread is only given by the standard deviation, whereas the actual minima (for instance, almost no cloud formation) cannot be seen directly, although these scenarios are possible.

Example 8: 2-D case with stochastic parameters. In the following experiment we study uncertainty propagation due to incomplete information about the model parameters which is a very typical problem arising in atmospheric science. We chose the same initial data for the flow and cloud variables as in Section 4.2. More precisely, we take the following initial cloud variables:

$$(\widehat{q}_v)_0(\mathbf{x}, 0) = 0.02\theta'(\mathbf{x}, 0), \quad (\widehat{q}_v)_k(\mathbf{x}, 0) = (\widehat{q}_c)_k(\mathbf{x}, 0) = (\widehat{q}_r)_k(\mathbf{x}, 0) = 0 \quad \text{for } 1 \leq k \leq M.$$

Consequently, in order to investigate uncertainty propagation in the numerical solution we choose 10%, 20% and 50% perturbation of these coefficients, namely, we take

$$\begin{aligned} (\widehat{k}_1)_0 &= 0.0041, \quad (\widehat{k}_1)_1 = \nu(\widehat{k}_1)_0, \quad (\widehat{k}_1)_k = 0 \quad \text{for } 2 \leq k \leq M, \\ (\widehat{k}_2)_0 &= 0.8, \quad (\widehat{k}_2)_1 = \nu(\widehat{k}_2)_0, \quad (\widehat{k}_2)_k = 0 \quad \text{for } 2 \leq k \leq M, \\ \widehat{\alpha}_0 &= 190.3, \quad \widehat{\alpha}_1 = \nu\widehat{\alpha}_0, \quad \widehat{\alpha}_k = 0 \quad \text{for } 2 \leq k \leq M, \end{aligned}$$

where $\nu = 0.1, 0.2$ and 0.5 , respectively. The numerical solution is computed at time $t = 1400s$ on a 320×320 mesh. Figures 21 and 22 present the potential temperature θ and the expected values of the cloud drops concentration q_c , respectively, for $\nu = 0, 0.1, 0.2$ and 0.5 .

In these scenarios with perturbed cloud model parameters, the overall structures are more stable and changes are less pronounced than in Example 7. Nevertheless, the main features of the variations are obviously driven by precipitation processes since the only perturbed parameters are those determining rain processes. Again, one key feature of the perturbed scenarios is the signature of evaporating rain in lower levels of the 2-D domain. For positive perturbations, that is, larger parameters k_1 , k_2 and α , rain formation is enhanced (more rain is formed from cloud water, due to larger k_1 and k_2) and sedimentation is enhanced (more rain is falling downwards due to larger α). Thus, more rain water is transported downwards into subsaturated regions, which is then evaporated inducing cooling due to latent heat consumption. These positive variations again dominate the potential temperature field; see Figure 21. For the cloud water field (Figure 22), one can

observe higher expected values for stronger perturbations. This is probably due to the fact that more rain is evaporated in lower levels, thus more water vapor is then available for cloud formation in upward motions of the convective cells. This redistribution of water vapor as well as the reduction of rain for negative perturbation lead to larger variations of cloud water in the 2-D domain, as can be seen in the standard deviation for the cloud drops concentration q_c , as depicted in Figure 23.

The expected values as well as the standard deviation for the cloud variables vary considerably with respect to the perturbation of model parameters; see Figure 24. Our numerical experiments indicate that the standard deviation increases in time and also depends on the size of the parameter perturbation. Indeed, the larger the parameter perturbation, the higher is the standard deviation of the cloud variables. The size of the corresponding standard deviations is depicted in the right column of Figure 24. As expected, rain formation sets on earlier for large perturbations, leading also to a strong decrease in the overall water vapor in comparison to less perturbed scenarios. However, the variations in all water variables increase a lot from the onset of precipitation to later times, and even the spread in cloud water increases in contrast to the time evolution in Example 7. Generally, the spread in the mean values is smaller than in Example 7; changes in initial data produce a larger variation, that means the model physics is quite stable with respect to perturbations in rain process formulations.

Example 9: 3-D case with stochastic initial data. Similarly to Example 7, we now investigate the uncertainty quantification in the 3-D Rayleigh-Bénard convection for stochastically perturbed initial data of the cloud variables given by (7.1). The numerical solution is computed in a domain $\Omega = [0, 5000] \times [0, 5000] \times [0, 1000] \text{ m}^3$, which is discretized using $50 \times 50 \times 50$ mesh cells. In Figure 25, we show the influence of 0%, 10%, 20% and 50% perturbation on the potential temperature θ as well as the expected value of the cloud drop concentration q_c . For the perturbation scenarios in the case of the 3-D moist Rayleigh-Bénard convection, one can see overall a qualitatively similar picture as for the corresponding 2-D simulations. However, due to an additional spatial direction, the 3-D structures can further change their patterns. For positive perturbations, clouds can be formed even at low vertical upward motions. The latent heat release increases the vertical motions in the convective cells, which leads to additional feedback, such as stronger cloud formation, which in turn leads to formation of larger amount of rain water. As in the 2-D case, the potential temperature distribution changes considerably at lower levels of the 3-D domain since evaporative cooling of precipitation is a dominant process. Similarly, the mean cloud water distribution is crucially changed. For positive perturbations, which are dominant on average in the beginning, more cloud water is formed and is later removed. Thus, less cloud water is available in the domain at later times. This can also be seen in the time evolution of spatially averaged variables, as shown in Figure 26, where the time evolution of the mean expected value per m^3 and the mean standard deviation per m^3 of the cloud variables are plotted. The time evolution of these variables is very similar to the 2-D case, but the 3-D scenarios show a very interesting feature, which is not available in 2-D. For larger perturbations, the pattern of the convective cells is changed. While convective cells have a quasi hexagonal shape for the deterministic simulation, they are ordered in a different way for larger perturbations. At upper levels, a more roll-like structure or even rectangular shape is formed. Thus, there is a transition of structures due to perturbations in the initial conditions. This feature has not been documented until now. A more detailed analysis of these structures is left for future studies. The change in structures in these experiments is very likely a result from different initial conditions. Of course, our results are preliminary and the observed parameter dependences and possible grid alignment effects of the solution structures deserve further attention in future work.

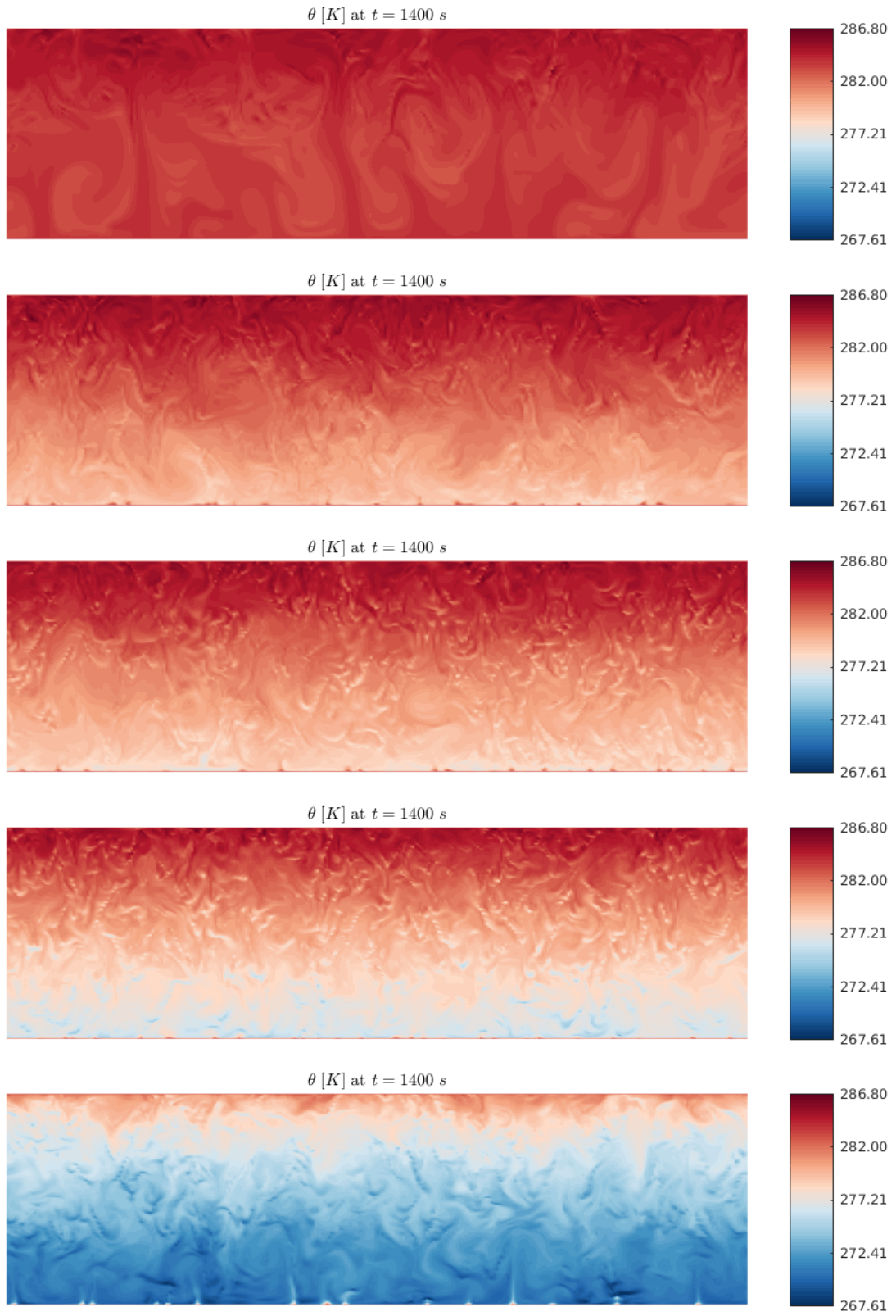


FIGURE 17. *Example 7: Potential temperature θ for $\nu = 0, 0.05, 0.1, 0.2$ and 0.5 (from top to down).*

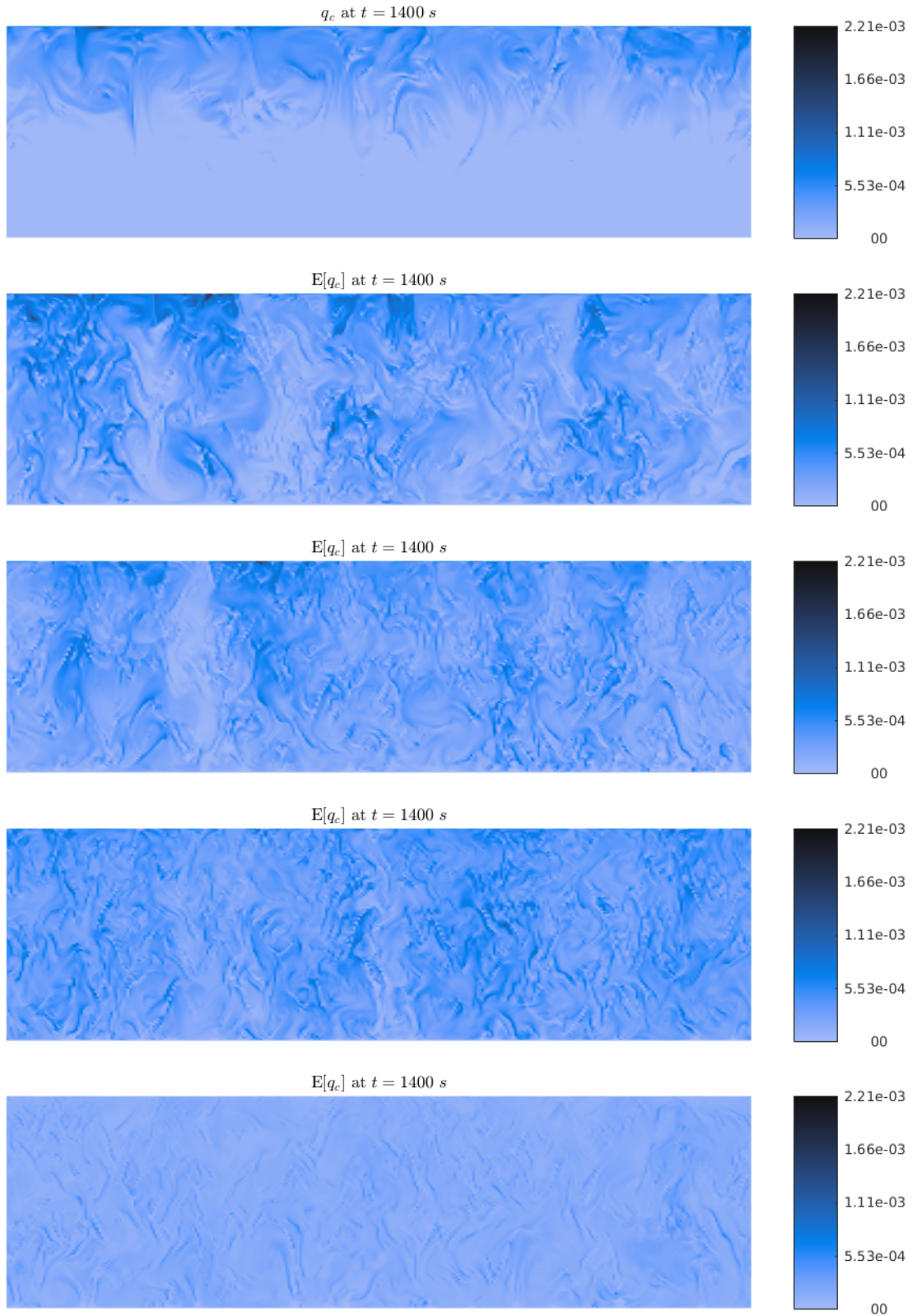


FIGURE 18. Example 7: Cloud drops concentration q_c for $\nu = 0, 0.05, 0.1, 0.2$ and 0.5 (from top to down).

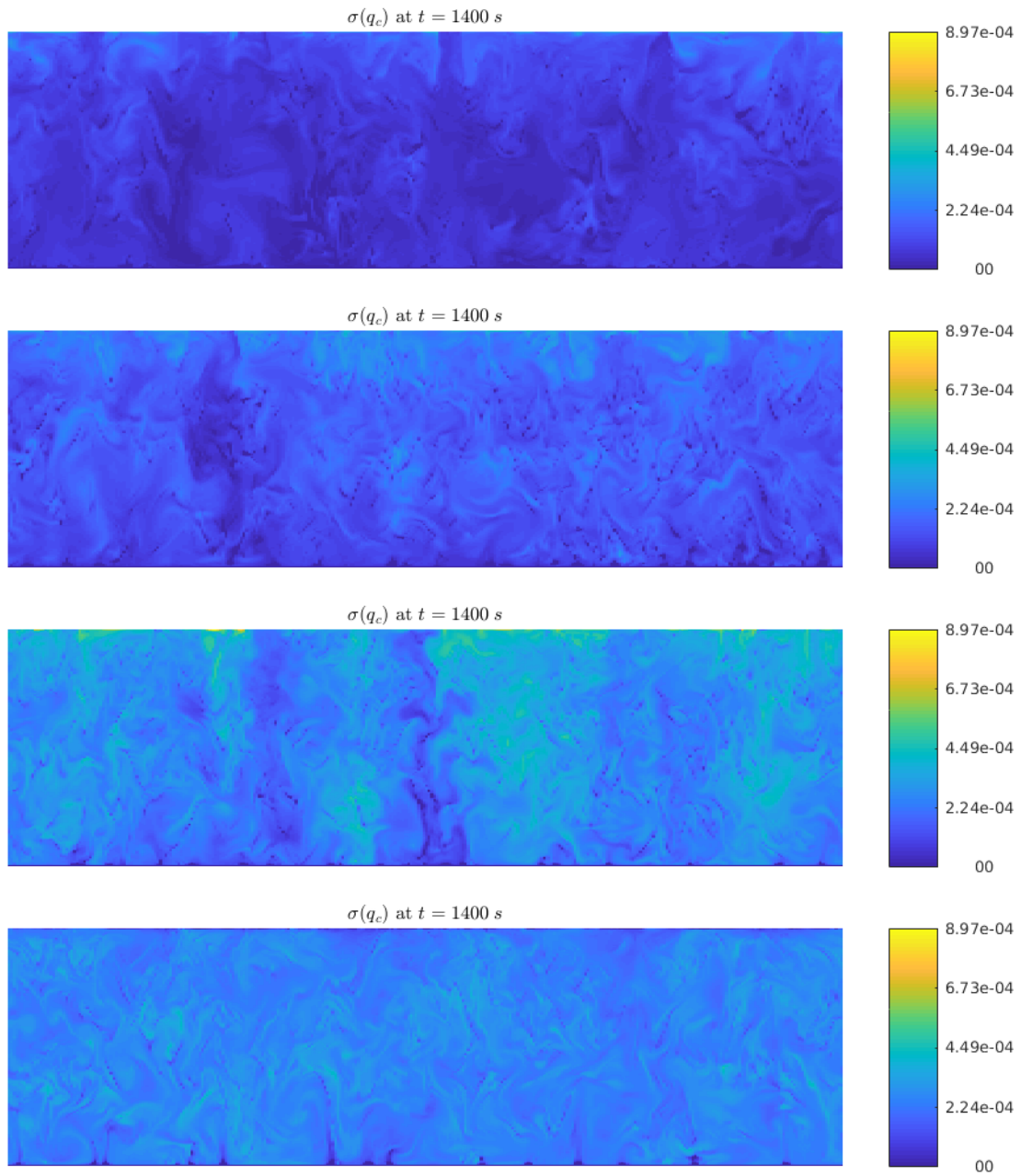


FIGURE 19. *Example 7: Standard deviation of the cloud drops concentration q_c for $\nu = 0.05, 0.1, 0.2$ and 0.5 (from top to down).*

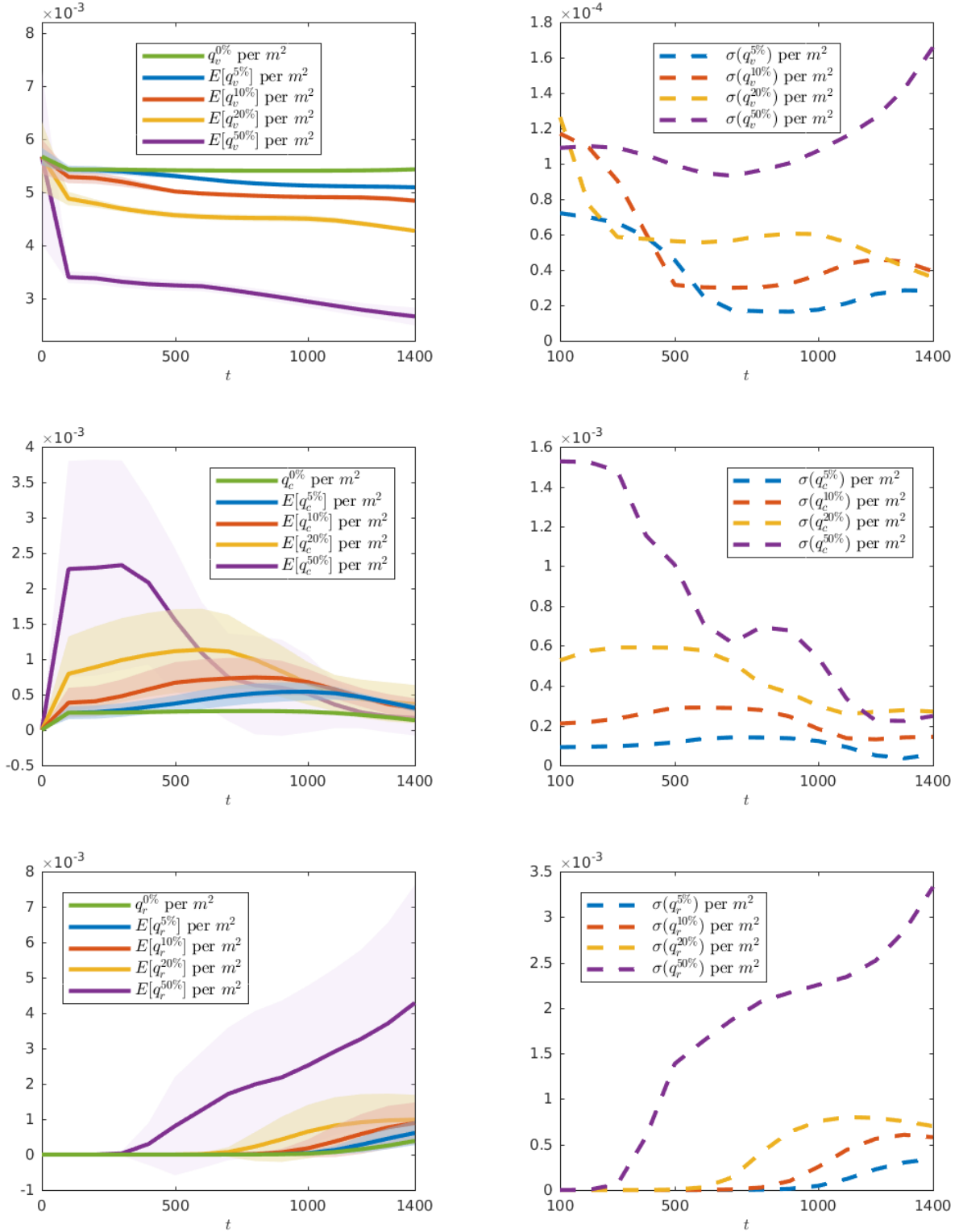


FIGURE 20. Example 7: Time evolution of the expected values with their standard deviations for the cloud variables per m^2 (shaded region, left column) and standard deviation (right column) for $\nu = 0.05, 0.1, 0.2$ and 0.5 (from top to down).

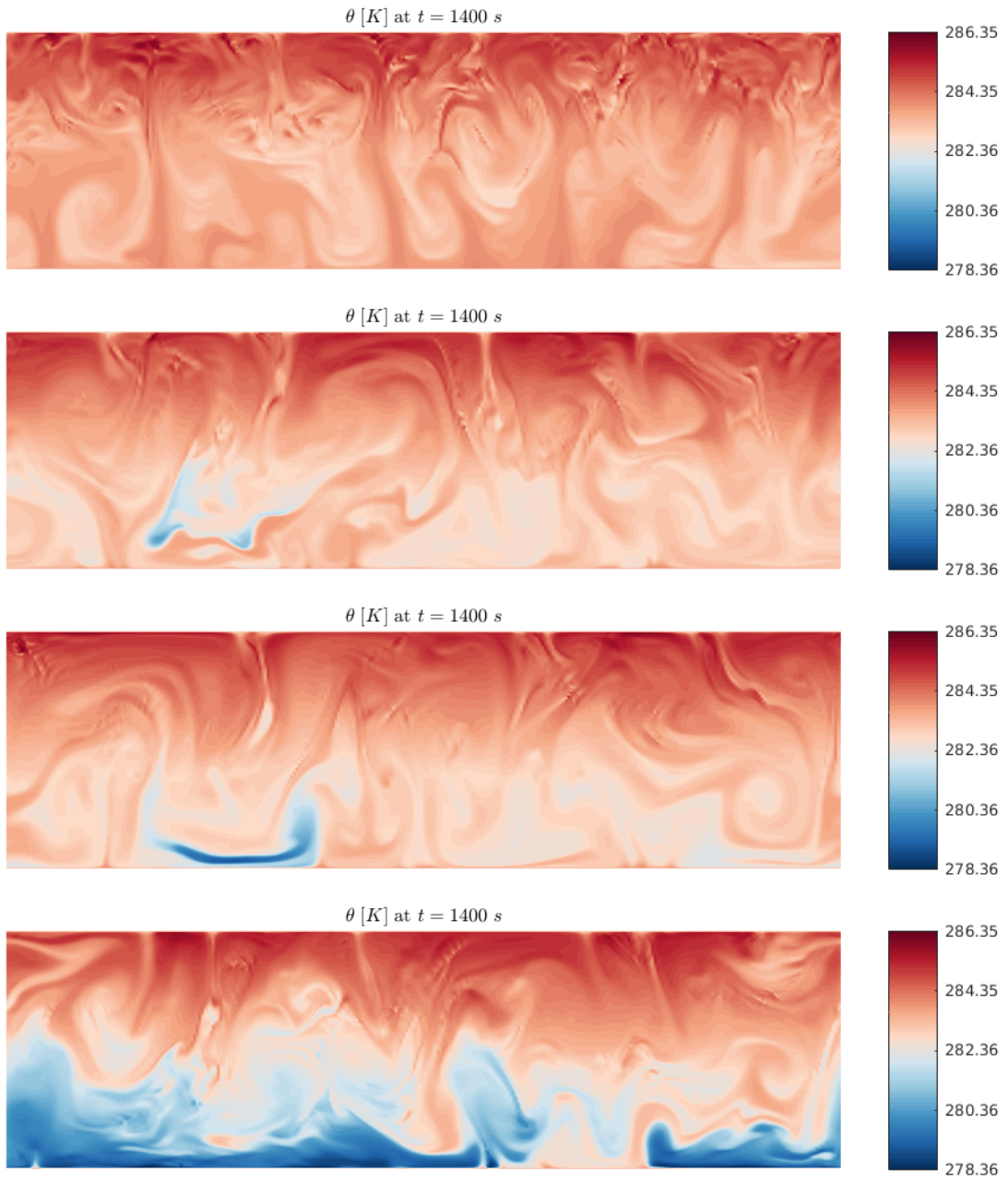


FIGURE 21. *Example 8: Potential temperature θ for $\nu = 0, 0.1, 0.2$ and 0.5 (from top to down).*

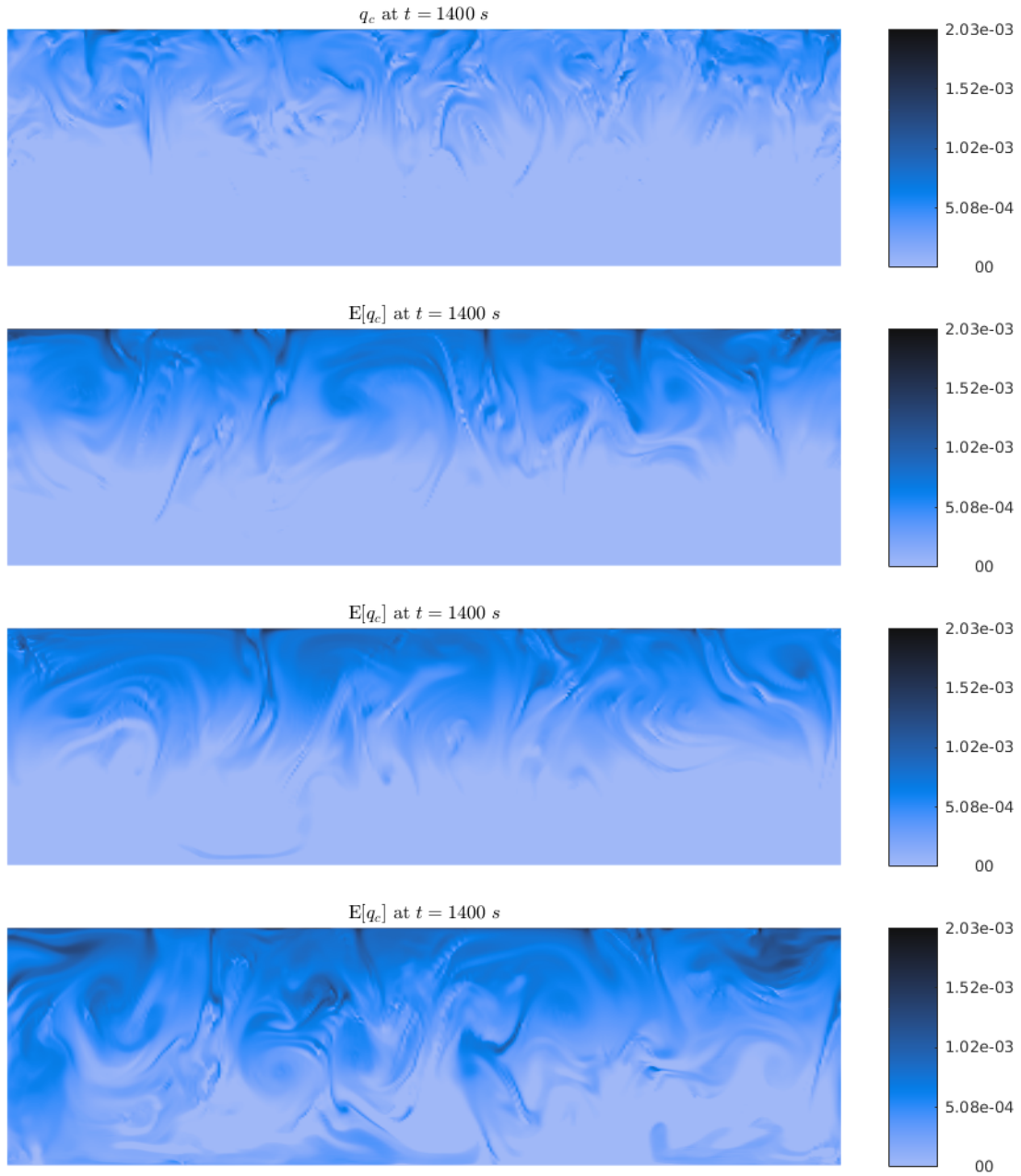


FIGURE 22. Example 8: Cloud drops concentration q_c for $\nu = 0, 0.1, 0.2$ and 0.5 (from top to down).

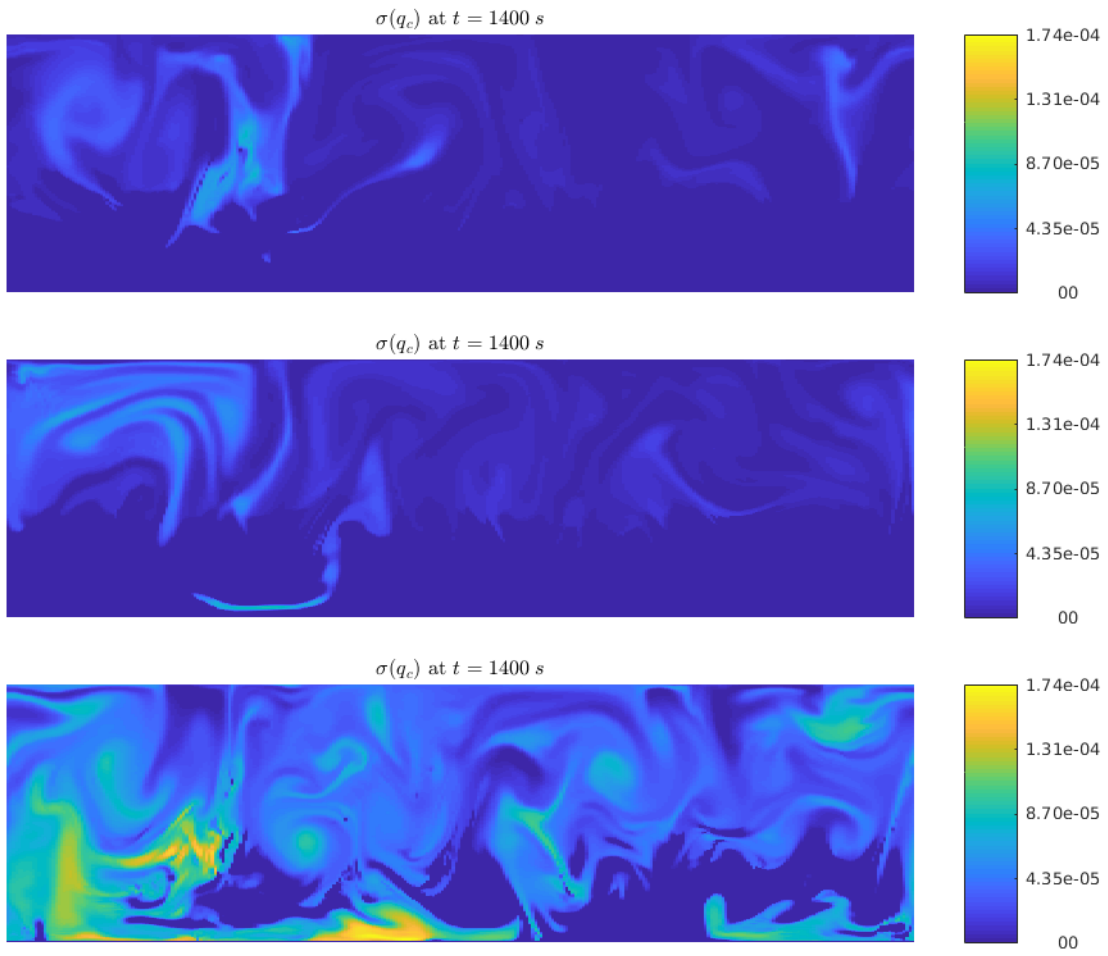


FIGURE 23. *Example 8: Standard deviation of the cloud drops concentration q_c for $\nu = 0.1, 0.2$ and 0.5 (from top to down).*

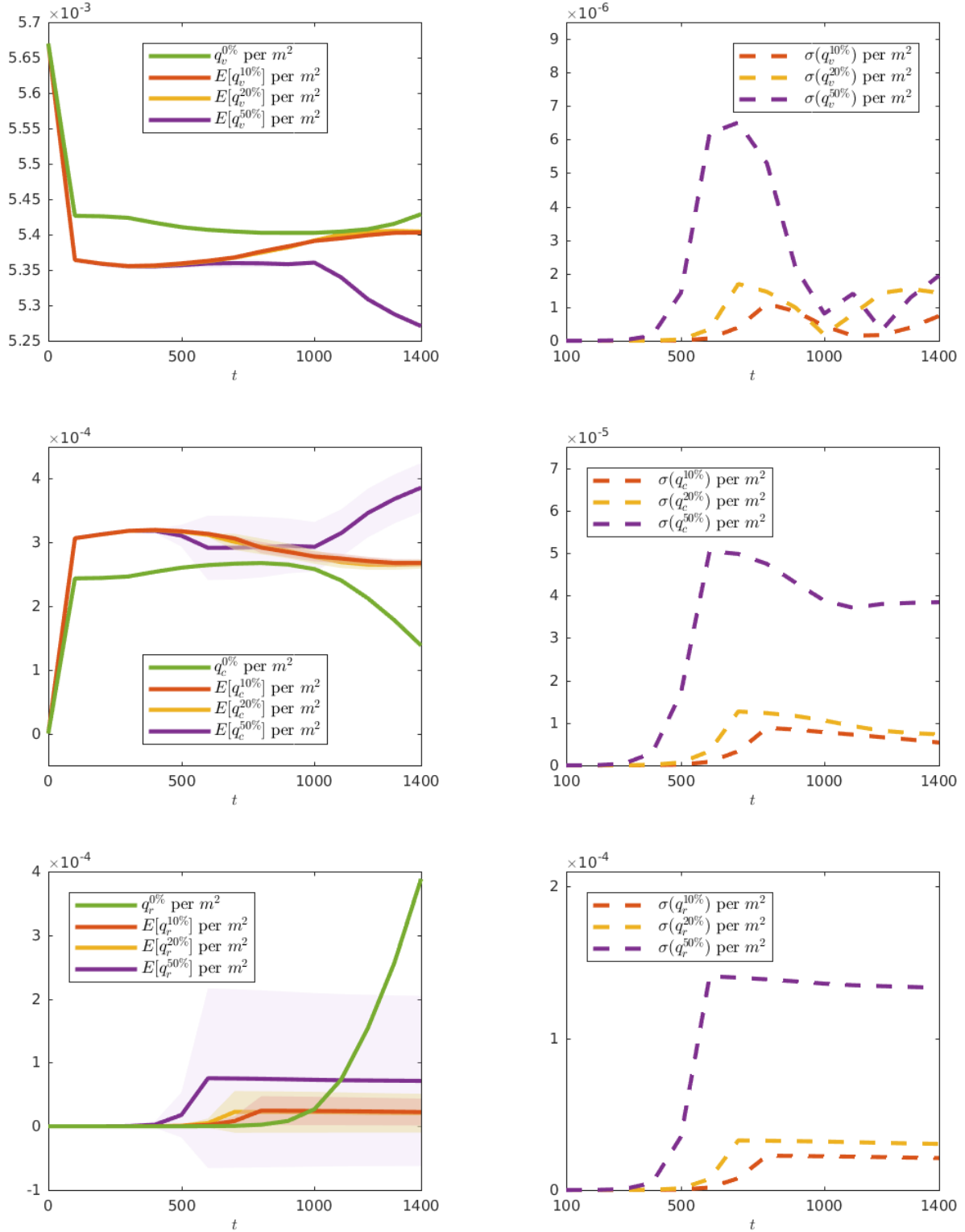


FIGURE 24. Example 8: Time evolution of the expected values with their standard deviations for the cloud variables per m^2 (shaded region, left column) and standard deviation (right column) for $\nu = 0.1, 0.2$ and 0.5 (from top to down).

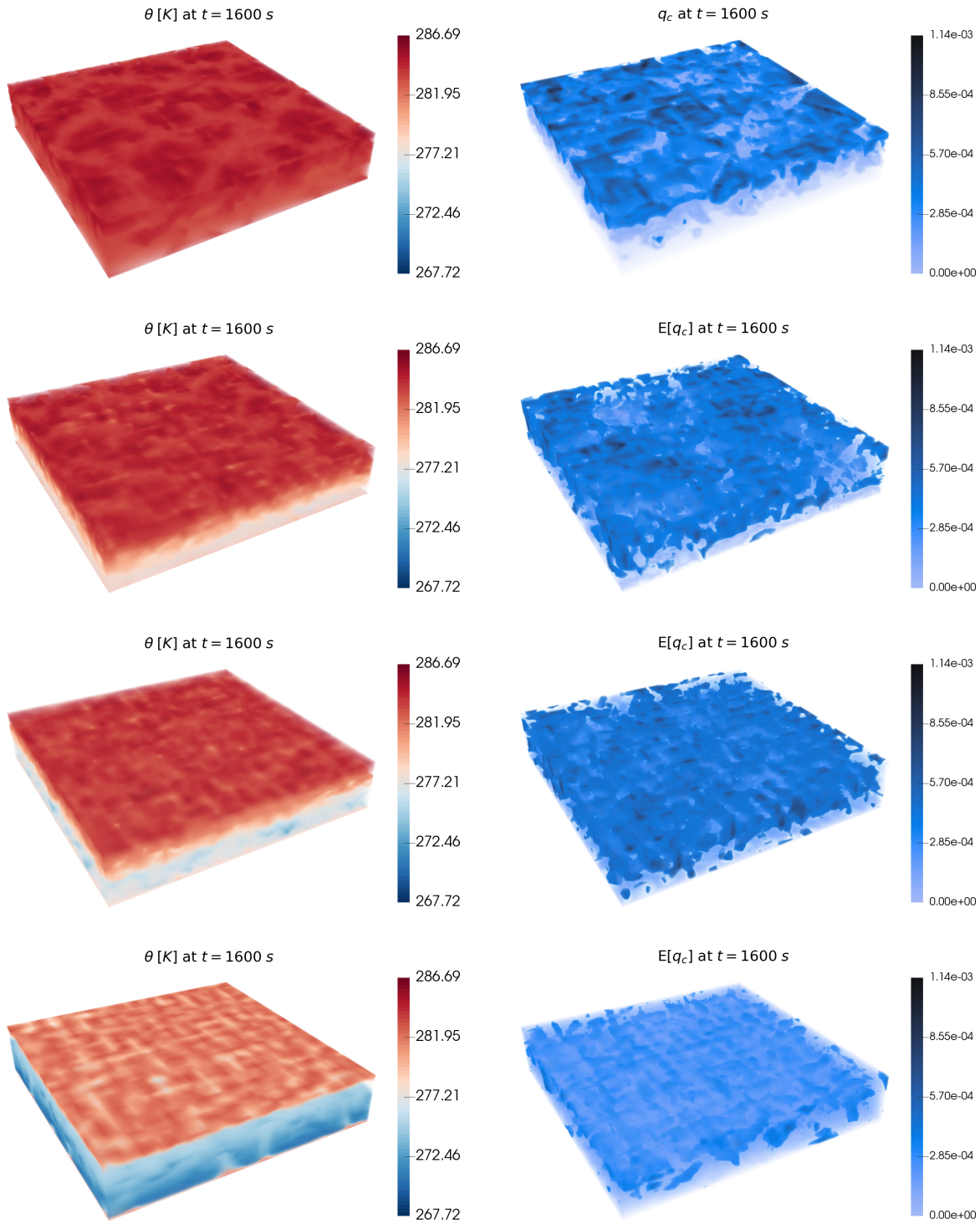


FIGURE 25. Example 9: Potential temperature θ (left column) and cloud drops concentration q_c (right column) using 0%, 10%, 20% and 50% (from top to down) perturbation of the initial data in q_v .

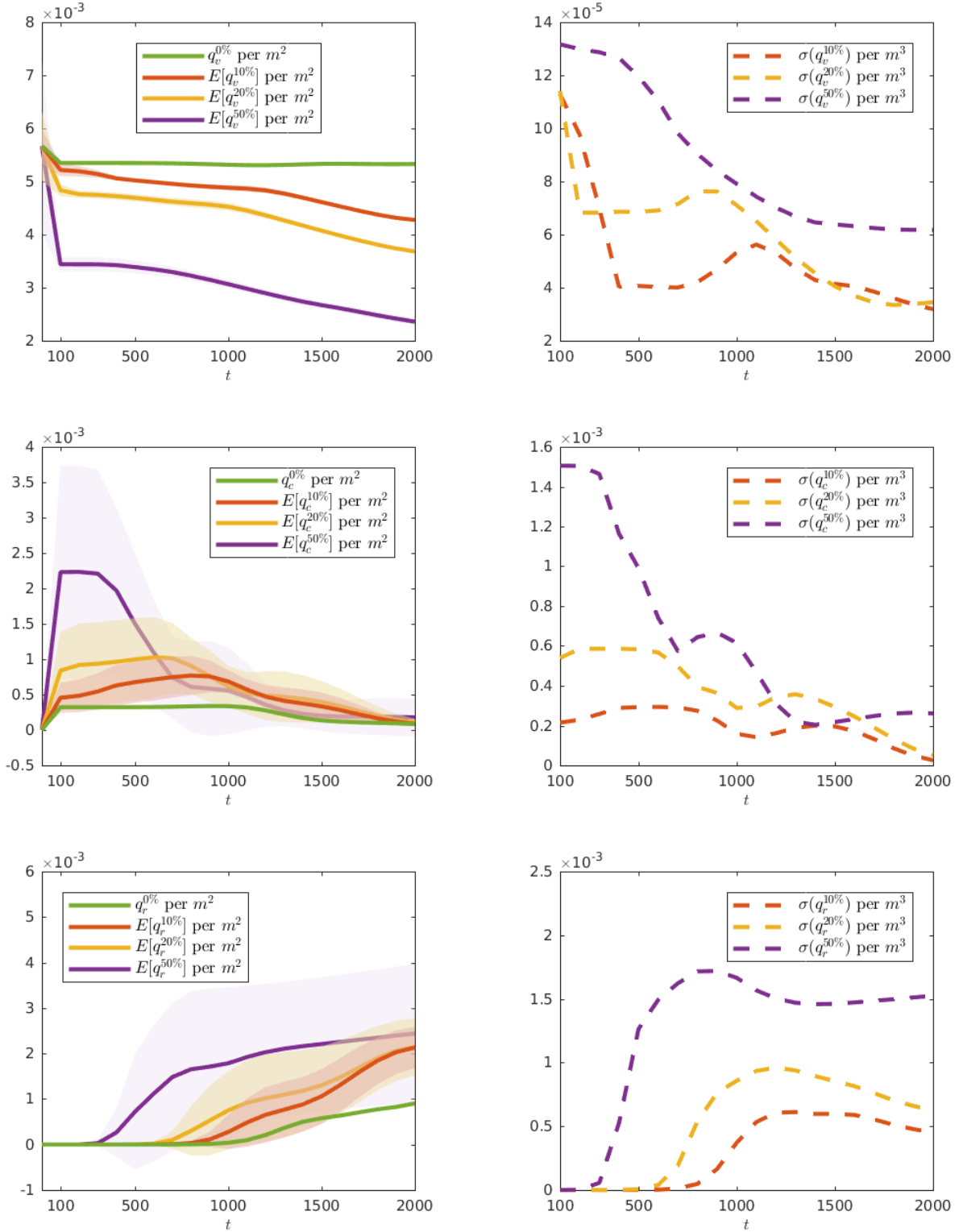


FIGURE 26. Example 9: Time evolution of the expected values with their standard deviations for the cloud variables per m^3 (shaded region, left column) and standard deviation (right column) using 0%, 10%, 20% and 50% (from top to down) perturbation of the initial data in q_v .

8. CONCLUSION

In the present paper, we have studied uncertainty propagation in an atmospheric model that combines the Navier-Stokes equations for weakly compressible fluids (2.1) with the cloud equations (2.2). The latter has been recently proposed in [35] and is based on the so-called single moment approach considering the evolution equations for the mass concentrations of the water vapor, cloud drops and rain. Our numerical strategy is based on the stochastic Galerkin method that combines a finite-volume method for space-time discretization with a spectral approximation in the stochastic space. We point out that atmospheric flows are weakly compressible which leads to the low Mach number problem. One therefore needs to use a finite-volume method, which is accurate and efficient in the low Mach number regime; see [5, 6]. To this end, we have chosen a suitable linear-nonlinear splitting between the fast and slow flow variables and the second-order IMEX discretization in time (the ARS (2,2,2) scheme) as described in Section 3. Coupling between the cloud model (2.2) and the Navier-Stokes system (2.8) is realized numerically by the second-order Strang splitting. The cloud equations are approximated in space by the finite-volume method and in time using the explicit third-order Runge-Kutta method with an enlarged stability region as explained in Section 3. Note that microscopic cloud dynamics requires a smaller time step than the flow dynamics and thus several microscopic cloud subiterations are realized within one macroscopic splitting time step, whose size is dictated by the flow dynamics. To the best of our knowledge, this is the first contribution that combines an accurate and efficient method for the weakly compressible Navier-Stokes equations with the stochastic Galerkin method for the uncertainty quantification of time evolution of the mass densities of water vapor, cloud drops and rain.

We have conducted extensive numerical benchmarking for both the deterministic and stochastic models and present the obtained numerical results in Sections 4 and 7. In the latter, we took into account the uncertainties in both initial data and cloud model parameters. Our numerical study clearly demonstrates applicability of the stochastic Galerkin method for the uncertainty quantification in complex atmospheric models. We have obtained interesting results illustrating the behavior of clouds in different perturbed scenarios and demonstrated that perturbations in the initial conditions can crucially change the time evolution of the moist Rayleigh-Bénard convection. In particular, it has been shown that for larger perturbation, on average the positive perturbations dominate the expectation values, although the standard deviation can be quite large. The main feature is a strong evaporative cooling in lower levels of the (2-D/3-D) domain due to enhance rain formation and sedimentation into low humidity levels. In the 3-D case, a change in the formed pattern can be seen, changing from hexagons to rolls/rectangles, which is quite surprising. For perturbations in parameters for rain processes, the results are also dominated by the positive part of the perturbations. Since rain processes are affected, the spread in the time evolution is increasing since the changes depend crucially on the interaction of formed rain drops with other variables. Overall, it seems that for cloud physics, the expectation values are dominated by the positive perturbations leading to a change in the distributions. This is an interesting topic for detailed studies in future. Our further goal is to extend the developed numerical method to the fully random Navier-Stokes-cloud system by considering random weakly compressible Navier-Stokes equations. We are also interested in considering different random effects, such as initial data, boundary data and model parameters simultaneously, which would require a multivariate stochastic Galerkin method. This will allow one to quantify more precisely the propagation of small scale stochastic errors initiated at cloud scales to macroscopic scales of flow dynamics.

Acknowledgement

The research leading to these results has been done within the sub-project A2 of the Transregional Collaborative Research Center SFB/TRR 165 “Waves to Weather” funded by the German Science Foundation (DFG). The work of A. Chertock was supported in part by NSF grants DMS-1521051 and DMS-1818684. The work of A. Kurganov was supported in part by NSFC grant 11771201.

The authors gratefully acknowledge the support of the Data Center (ZDV) in Mainz for providing computation time on MOGON and MOGON II clusters and of NSF RNMS Grant DMS-1107444 (KI-Net). M. Lukáčová-Medviďová and B. Wiebe would like to thank K. Werth and L. Yelash (University of Mainz) for fruitful discussions. P. Spichtinger would like to thank P. Reutter (University of Mainz) for fruitful discussions on the results of the 2-D/3-D simulations.

9. APPENDIX A: CLOSURE FOR SINGLE MOMENT SCHEMES

The number concentration of rain drops can be approximated by a function of the respective mass concentration $n_r = f(q_r, c_r)$. Since we implicitly assume that rain drops are distributed according to a size distribution, this approach should be used for mimicking the shape of the distribution in a proper way. If we use a constant mean mass of rain drops, the function will be a simple linear relation $n_r = \frac{1}{\bar{m}_r} q_r$. We extend this approach and propose the following nonlinear relation:

$$n_r = c_r \cdot q_r^\gamma, \quad 0 < \gamma \leq 1.$$

Using this approach, one can replace the quantity n_r in the processes related to rain drop number concentration. For the simple case of a constant mean mass \bar{m}_r , we can determine the constants as $c_r = \bar{m}_r^{-1}$ and $\gamma = 1$. This approach would be meaningful for the case of a symmetric size distribution of rain droplets centered around the mean mass. However, it is well-known that size distributions of rain are usually skew to larger sizes and thus a linear relation is inappropriate. For sizes of rain drops, an exponential distribution is often assumed (see [26]), namely:

$$f(r) = B_r e^{-\lambda r}$$

with a constant parameter $B_r = 2 \cdot 10^7 \text{ m}^{-4}$ and the drop radius r . Using the general moments of the distribution,

$$\mu_k[r] = \frac{\Gamma(k+1)}{\lambda^{k+1}} B_r$$

with the gamma function $\Gamma(x) := \int_0^\infty t^{x-1} \exp(-t) dt$, we obtain

$$\rho n_r = \mu_0[r] = \frac{B_r}{\lambda} \quad \text{and} \quad \rho q_r = \frac{4}{3} \pi \rho \ell \frac{\Gamma(4)}{\lambda^4} B_r.$$

Using these relations, one can derive the following function for the number concentration n_r :

$$n_r = \underbrace{B_r^{\frac{3}{4}} \rho^{-\frac{3}{4}} (8\pi\rho\ell)^{-\frac{1}{4}} q_r^{\frac{1}{4}}}_{=c_r} = c_r q_r^\gamma, \quad \gamma = \frac{1}{4}.$$

We stress that c_r is, in fact, a function of the air density ρ , that is, $c_r = c_{r0} \cdot \rho^{-\frac{3}{4}}$.

10. APPENDIX B: EXPLICIT FORMULATION OF THE CLOUD EQUATIONS

We present the equations of microphysical processes in an explicit way as they are used in our numerical experiments. In Tables 3 and 4, we present physical constants and model parameters with their values used in our numerical simulations.

$$\begin{aligned} n_c &= q_c \frac{8 \cdot 10^8}{q_c + 4.1888 \cdot 10^{-7}} \coth\left(\frac{q_c}{5.236 \cdot 10^{-13}}\right), \quad C_{\text{act}} = 6.2832 \cdot 10^{-3} D_v G \rho (q_v - q_*)_+, \\ C_1 &= 0.7796 D_v G (q_v - q_*) \left(\frac{8 \cdot 10^8}{q_c + 4.1888 \cdot 10^{-7}} \coth\left(\frac{q_c}{5.236 \cdot 10^{-13}}\right)\right)^{\frac{2}{3}} \rho q_c, \\ p_s(T) &= \exp\left\{54.842763 - 6763.22/T - 4.21 \ln T + 0.000367T + \tanh(0.0415(T - 218.8))\right. \\ &\quad \left. \cdot (53.878 - 1331.22/T - 9.44523 \ln T + 0.014025T)\right\}, \\ n_r &= 23752.6753 \rho^{-\frac{3}{4}} q_r^{\frac{1}{4}}, \quad r = \left(\frac{1.21 \cdot 10^{-5}}{q_r + 0.2874 \rho^{-\frac{3}{4}} q_r^{\frac{1}{4}}}\right)^{\frac{4}{15}}, \\ E &= -0.7796 D_v G (q_* - q_v)_+ \left(644.5198 \sqrt{\rho q_r} + 17.5904 \mu^{-\frac{1}{6}} D_v^{-\frac{1}{3}} \sqrt{\alpha r} \rho^{\frac{13}{24}} q_r^{\frac{91}{120}}\right), \\ A_1 &= 10^{-3} k_1 \rho q_c^2, \quad A_2 = 0.3846 \alpha k_2 \rho^{\frac{1}{4}} q_c r q_r^{\frac{61}{60}}, \quad v_q = 1.1068 \alpha q_r^{\frac{4}{15}} r \rho^{-\frac{1}{2}}. \end{aligned}$$

Constant	Description
$p_* = 101\,325$ Pa	reference pressure
$T_* = 288$ K	reference temperature
$T_0 = 273.15$ K	melting temperature
$\rho_* = 1.225$ kg m ⁻³	reference air density
$\rho_l = 1000$ kg m ⁻³	density of liquid water
$R_v = 461.52$ J kg ⁻¹ K ⁻¹	specific gas constant, water vapor
$c_p = 1005$ J kg ⁻¹ K ⁻¹	specific heat capacity, dry air
$g = 9.81$ m s ⁻²	acceleration due to gravity
$L = 2.53 \times 10^6$ J kg ⁻¹	latent heat of vaporization
$\varepsilon = \frac{M_{\text{mol},v}}{M_{\text{mol},a}} = 0.622$	ratio of molar masses of water and dry air
$D_0 = 2.11 \times 10^{-5}$ m ² s ⁻¹	diffusivity constant

TABLE 3. *Physical constants and reference quantities, [35].*

Parameter	Description
$\alpha = 190.3 \pm 0.5 \cdot 190.3$ m s ⁻¹ kg ^{-β}	parameter for terminal velocity
$k_1 = 0.0041 \pm 0.5 \cdot 0.0041$ kg s ⁻¹	parameter for autoconversion
$k_2 = 0.8 \pm 0.5 \cdot 0.8$ kg	parameter for accretion
$\beta = \frac{4}{15}$	parameter for terminal velocity
$m_t = 1.21 \times 10^{-5}$ kg	parameter for terminal velocity
$N_0 = 1000$ m ⁻³	parameter for activation
$N_\infty = 8 \times 10^8$ kg ⁻¹	parameter for activation
$m_0 = 5.236 \times 10^{-16}$ kg	parameter for activation
$a_E = 0.78$	parameter for evaporation
$a_v = 0.78$	parameter for ventilation
$b_v = 0.308$	parameter for ventilation

TABLE 4. *Model parameters, [35].*

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