DNS for studying entrainment and mixing processes

ZHENG GAO

STONY BROOK UNIVERSITY

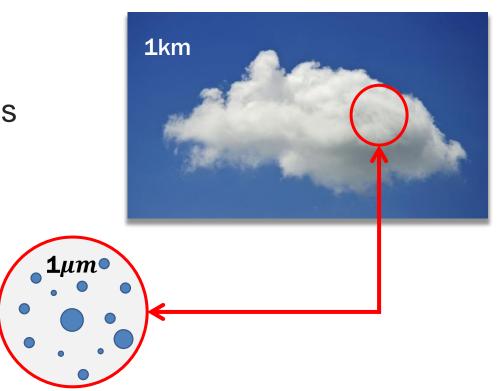


Outline

- Background
- Models
- Numerical method
- Simulation
- Results

Background

- Cloud structure in microscale
- Turbulence and interactions
- Studying with DNS
- Entrainment and mixing



Mathematics Models

$$\frac{\partial \mathbf{u}}{\partial t} + (\mathbf{u} \cdot \nabla \mathbf{u}) = -\frac{1}{\rho_0} \nabla \mathbf{p} + \mu \Delta \mathbf{u} + f(\mathbf{q}, T)$$

$$\nabla \cdot \mathbf{u} = 0$$

$$\frac{\partial \mathbf{q}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{q} = -C_d + \kappa \Delta \mathbf{q}$$

$$\frac{\partial \mathbf{T}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{T} = \frac{L}{c_p} C_d + \mu_T \Delta \mathbf{T}$$

Vapor mixing ratio

Temperature

Numerical Methods

- Projection method to decouple velocity and pressure (HYPRE and PETSc)
- WENO scheme to evaluate advection (no oscillation, high order)
- Crank-Nicolson for diffusion (stable)
- Totally second order of accuracy

Mathematics Models

$$S(X,t) = \frac{q_v(X,t)}{q_{v,s}} - 1$$

$$\frac{dR_i(t)}{dt} = A_3 \frac{S(x,t)}{R_i(t)}$$

Condensational growth

$$\frac{dX(t)}{dt} = V(t)$$

$$\frac{dV(t)}{dt} = \frac{1}{\tau_p} [u(X, t) - V(t)] + g$$
Droplets motion

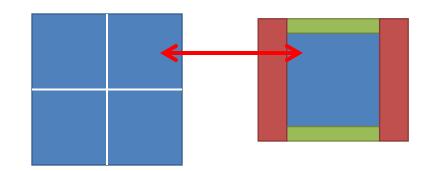
Numerical Methods (cont.)

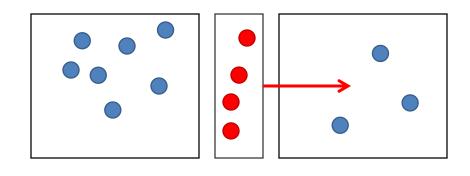
- Implicit Euler scheme for particle motion (stable)
- Explicit Euler scheme for condensation (efficient)
- Two way interaction, water mass conserved.
- S = 0, equilibrium state, no water exchange

$$\frac{dR_i(t)}{dt} = A_3 \frac{S(x,t)}{R_i(t)}$$

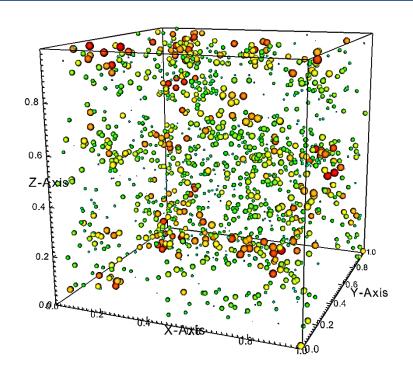
Parallel computing

- MPI (MPICH2)
- Parallelization of field (add buffer)
- Parallelization of particles (send and receive)
- Statistics analysis

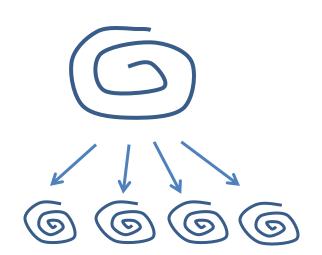




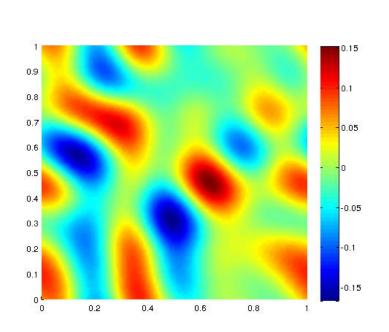
- Simulation box
- √1m³ domain
- ✓ Periodic boundary condition
- √64³ or 128³ mesh grid
- ✓ Particles are uniformly placed on supersatured region



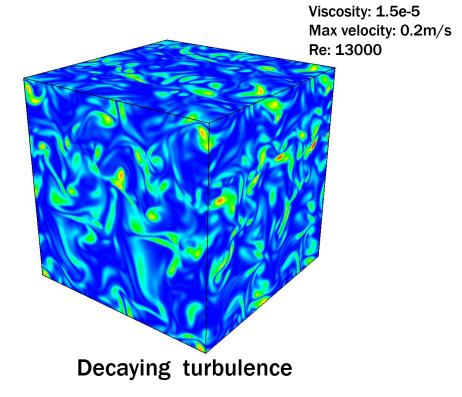
- Turbulence
- ✓ Energy input only in large wave length
- ✓ Energy cascades to small length automatically
- ✓ Energy dissipate in Kolmogorov length scale



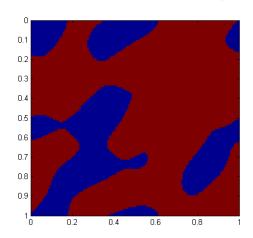
Turbulence initialization



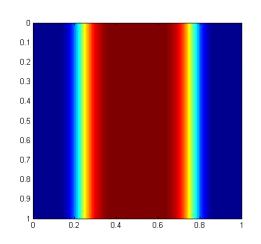
Energy input from large scale Isotropic



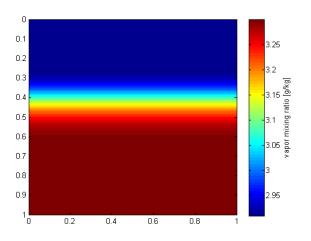
Vapor mixing ratio



Interior
[1] Andrejczuk (04 – 09)
Case 1



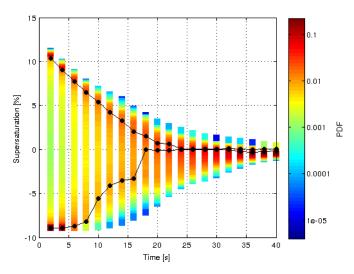
Boundary
[2] Kumar (12)
Case 2



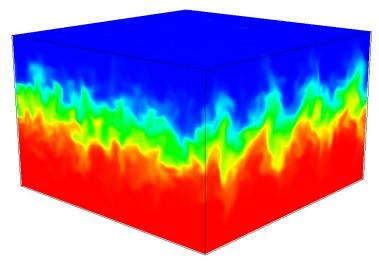
Top Ours Case 3

Vapor mixing ratio

Vapor mixing ratio changes with time



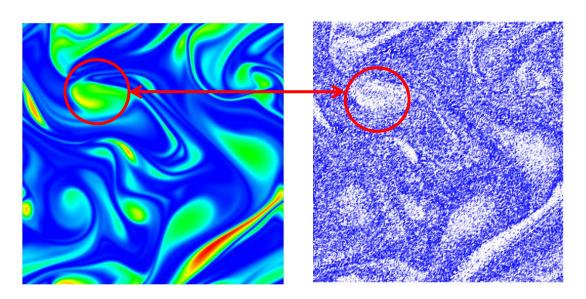
Supersaturation with time



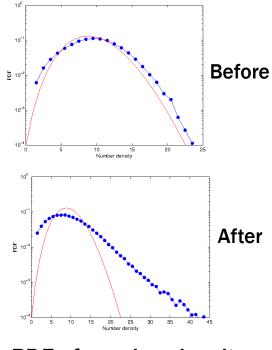
Vapor mixing ratio at t = 2s

- Particles
- ✓ Initial position: collocated with s > 0
- ✓Initial velocity: 0m/s
- ✓ Initial size: uniform size (10um)
- √ Consider sedimentation and inertial

Preferential Concentration

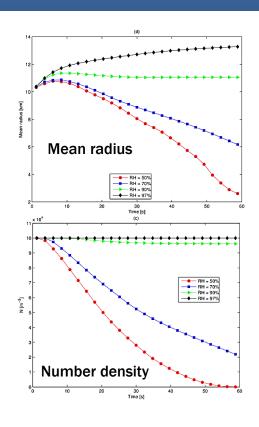


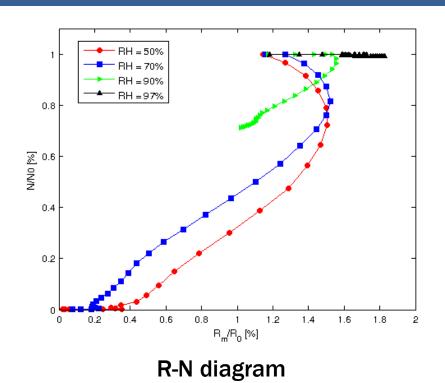
Enstrophy and number density Enhance collision rates



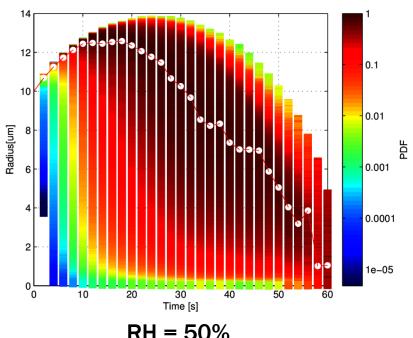
PDF of number density

Radius and Number density

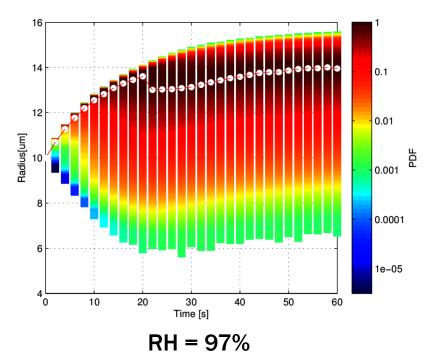




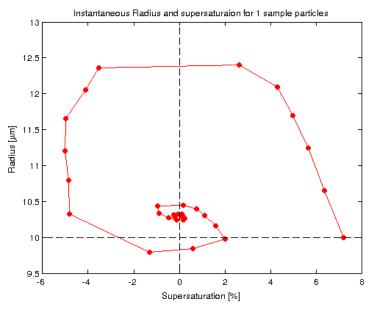
Radius spectrum



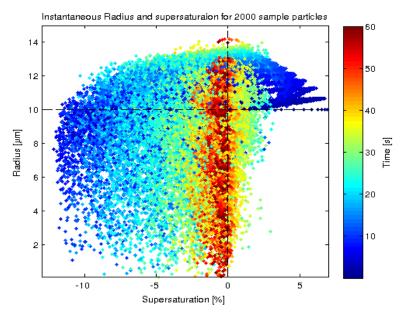
RH = 50%



Radius and supersaturation

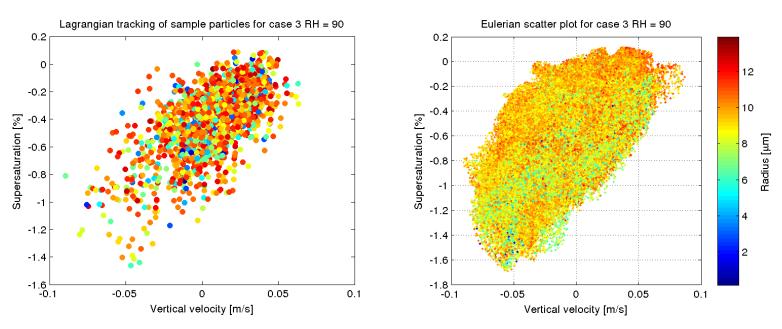


Lagrangian tracking of sample particle



Lagrangian tracking of 2000 particles

Supersaturation and vertical velocity



Supersaturation and vertical velocity at final state in Lagrangian (left) and Eulerian view (right)

Future work

- Larger domain and mesh refinement
- Adding external force from larger scale
- Collision and coalescence
- Particle point vs. particle resolved
- Thank you!