

モーメント・クローザーを用いた gyrokinetic 運動論的シミュレーション手法 Gyrokinetic simulation method with a moment closure

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Outline

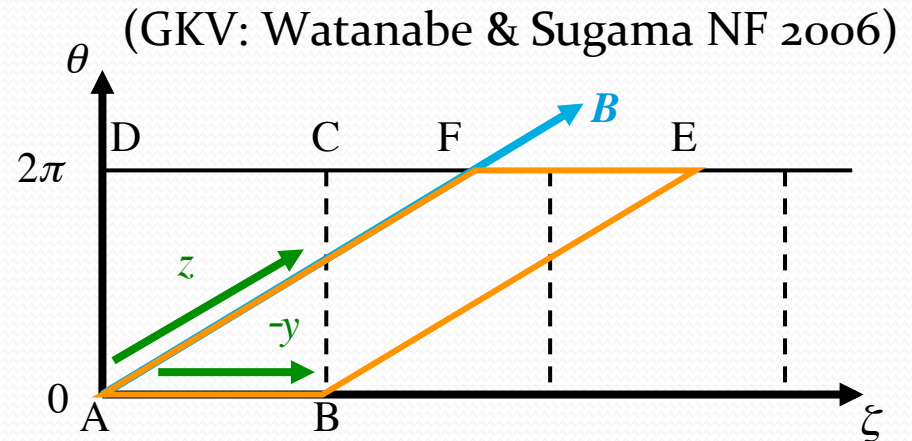
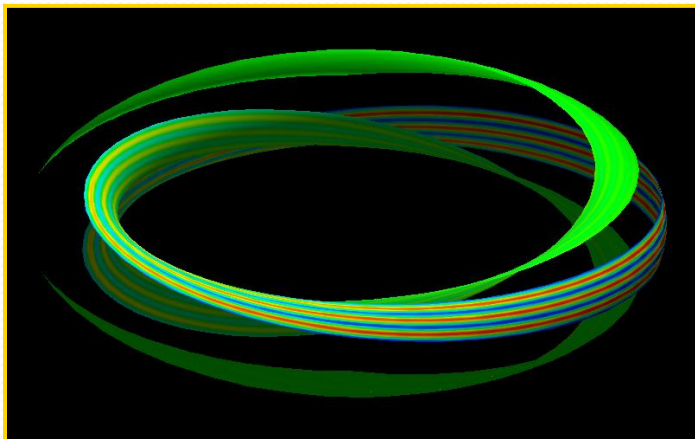
- Introduction
 - GKV code and its numerical difficulties
- Drift kinetic equation and moment closure
 - Moment extracted formulation
 - Application of the implicit scheme for the kinetic Alfvén wave propagation
- A numerical remark on the DK simulation
 - Maxwellian distribution on discretized grids in the finite velocity space
- Application of the semi-Lagrangian scheme for a finite β plasma
- Summary

Gyrokinetic Simulation Code: GKV

- Nonlinear gyrokinetic equation for perturbed gyrocenter distribution δf is numerically solved on the five-dimensional phase space, $(x, y, z, v_{\parallel}, \mu)$

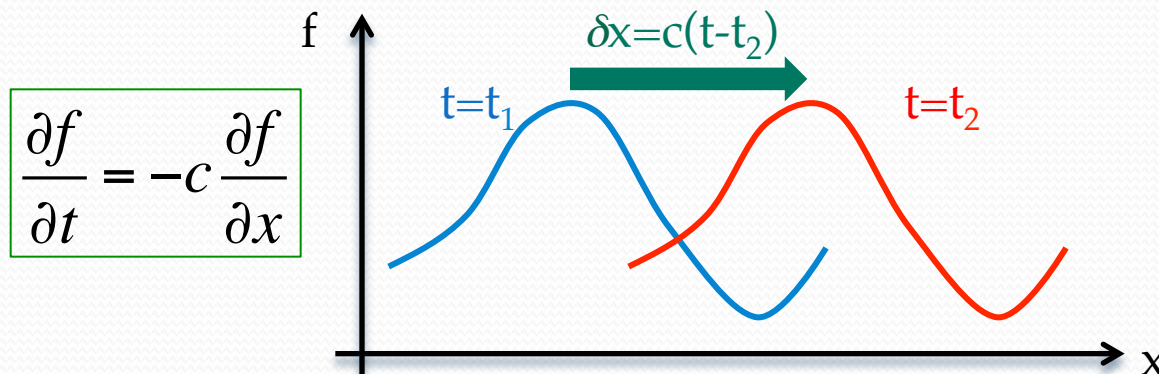
$$\left[\frac{\partial}{\partial t} + v_{\parallel} \hat{\mathbf{b}} \cdot \nabla + \mathbf{v}_d \cdot \nabla - \mu (\hat{\mathbf{b}} \cdot \nabla \Omega) \frac{\partial}{\partial v_{\parallel}} \right] \delta f + \frac{c}{B_0} \{ \psi, \delta f \} = (\mathbf{v}_* - \mathbf{v}_d - v_{\parallel} \hat{\mathbf{b}}) \cdot \frac{e \nabla \psi}{T_i} F_M + C(\delta f)$$

- Strong anisotropy of fluctuations is accurately resolved by using curvilinear coordinates along field lines.
- High resolution of 5-D phase space.



Numerical difficulties in multi-species gyrokinetic simulation

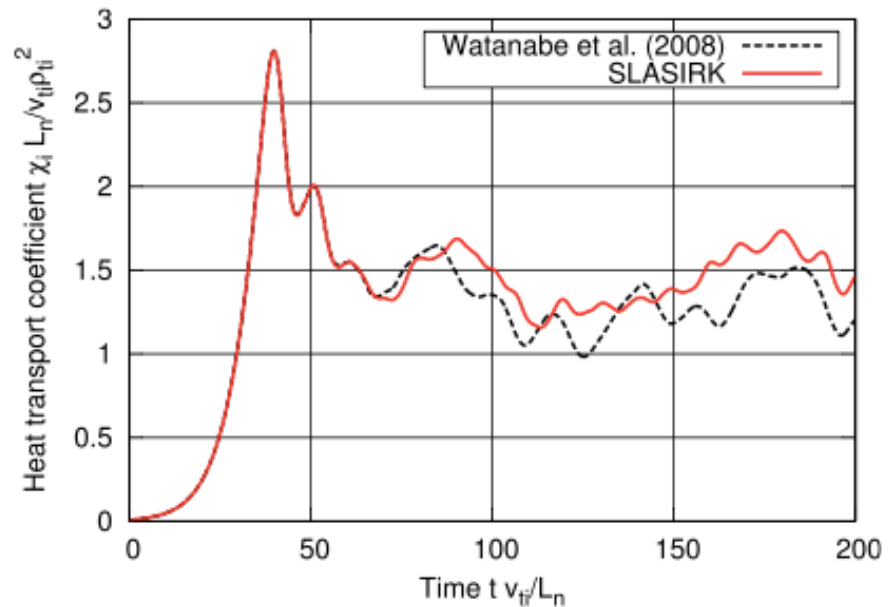
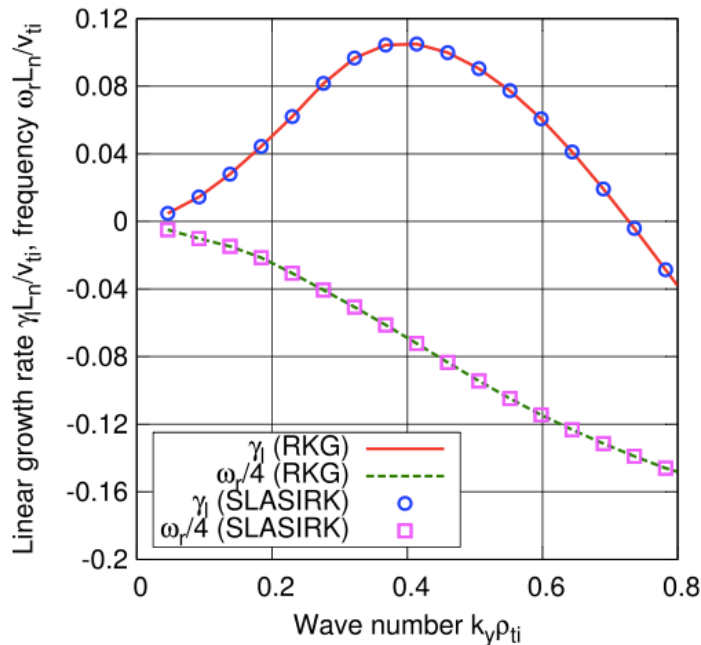
- Multi-species gyrokinetic (GK) simulation suffers from separation of typical time scales of **ions and electrons**
 - Fast electron motion along field lines restricts the time-step size of explicit schemes. (say, $\Delta t < 1.e-4 R_0 / C_s$ for LHD)
 - It also leads to slow convergence of recursive solvers in implicit time integrations.
 - Semi-Lagrangian (SL) method can trace drift motion of particles with time steps beyond the CFL condition.



Semi-Lagrangian scheme applied to GKV – Early work

(Maeyama et al CPC 2012)

- Semi-Lagrangian scheme applied to GKV code could successfully reduce the computational costs.



- Linear ITG benchmark with adiabatic electrons; $\Delta t=0.005$ for RKG but $\Delta t=0.1$ for Semi-Lagrangian + iRK
- Nonlinear benchmark for the ITG turbulence was also successful.

Numerical difficulties in multi-species gyrokinetic simulation

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 - It also leads to slow convergence of recursive solvers in implicit time integrations.
 - Semi-Lagrangian (SL) method can trace drift motion of particles with time steps beyond the CFL condition.
 - But, SL for GK equation is often numerically **unstable to electromagnetic fluctuations**, because the wave propagation direction may be different from those of particle motions.
- => We need new numerical techniques for multi-species GK.

Drift kinetic equation and moment closure

Revisiting the drift kinetic equation

- Let us consider the linearized DK equation for electrons

$$\frac{\partial f_e}{\partial t} + v_{\parallel} \frac{\partial f_e}{\partial z} = -\frac{q_e}{T_{e0}} v_{\parallel} F_M \frac{\partial}{\partial z} \left(\frac{\partial \phi}{\partial z} + \frac{1}{c} \frac{\partial A_{\parallel}}{\partial t} \right)$$

- 0th, 1st, and 2nd order moments are given by

$$n_e = \int f_e d^3v \quad n_0 U = \int v_{\parallel} f_e d^3v \quad n_e T_0 + n_0 T_{\parallel} = m_e \int v_{\parallel}^2 f_e d^3v$$

- Quasi-neutrality and the Ampere's law

$$\frac{q_i^2 n_0}{T_i} k_{\perp}^2 \rho_i^2 \phi = q_e n_e \quad k_{\perp}^2 A_{\parallel} = \frac{4\pi}{c} q_e n_0 U$$

- The above equations describe the parallel electron motion and the kinetic Alfvén waves

Equations of low-order moments and moment-extracted kinetic eq.

- The 0th and 1st order moment equations are given by

$$\frac{\partial n_e}{\partial t} = -n_0 \frac{\partial U}{\partial z} \quad n_0 m_e \frac{\partial U}{\partial t} = -q_e n_0 \left(\frac{\partial \phi}{\partial z} + \frac{1}{c} \frac{\partial A_{\parallel}}{\partial t} \right) - \frac{\partial}{\partial z} (n_e T_{e0} + n_0 T_{\parallel})$$

- Distribution function h_e where the low-order moments are extracted.

$$f_e = \frac{n_e}{n_0} F_M + \frac{m_e}{T_{e0}} U v_{\parallel} F_M + h_e \quad n_0 T_{\parallel} = m_e \int v_{\parallel}^2 h_e d^3 v$$

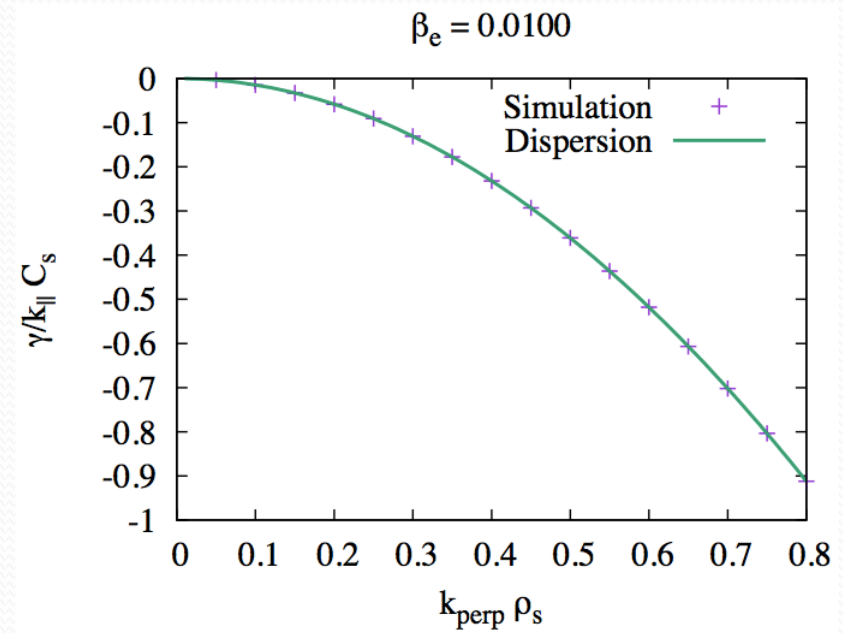
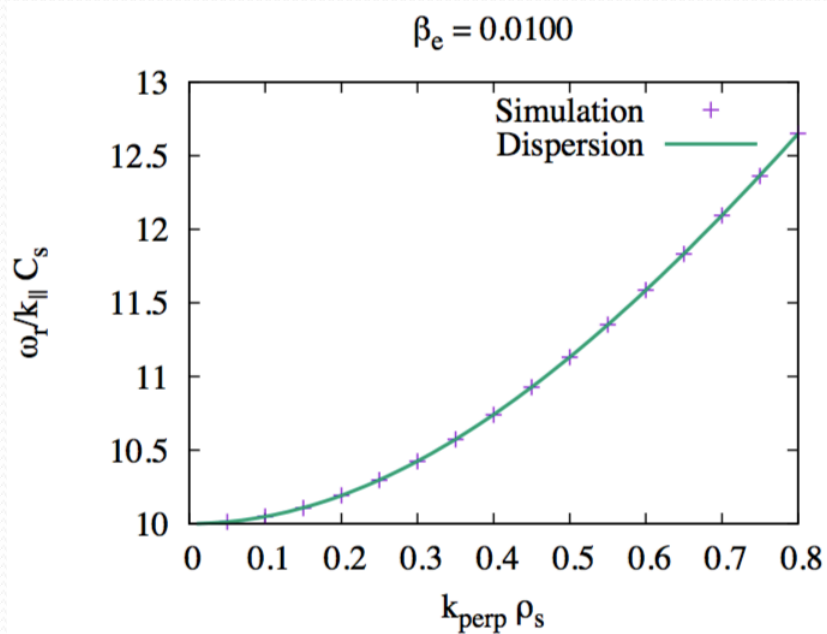
- The remnant drift kinetic equation for h_e

$$\frac{\partial h_e}{\partial t} + v_{\parallel} \frac{\partial h_e}{\partial z} = \left(1 - \frac{m_e v_{\parallel}^2}{T_{e0}} \right) F_M \frac{\partial U}{\partial z} + \frac{v_{\parallel}}{T_{e0}} F_M \frac{\partial T_{\parallel}}{\partial z}$$

with a constraint of $\int h_e d^3 v = \int v_{\parallel} h_e d^3 v = 0$

New formulation exactly describes the kinetic Alfvén wave

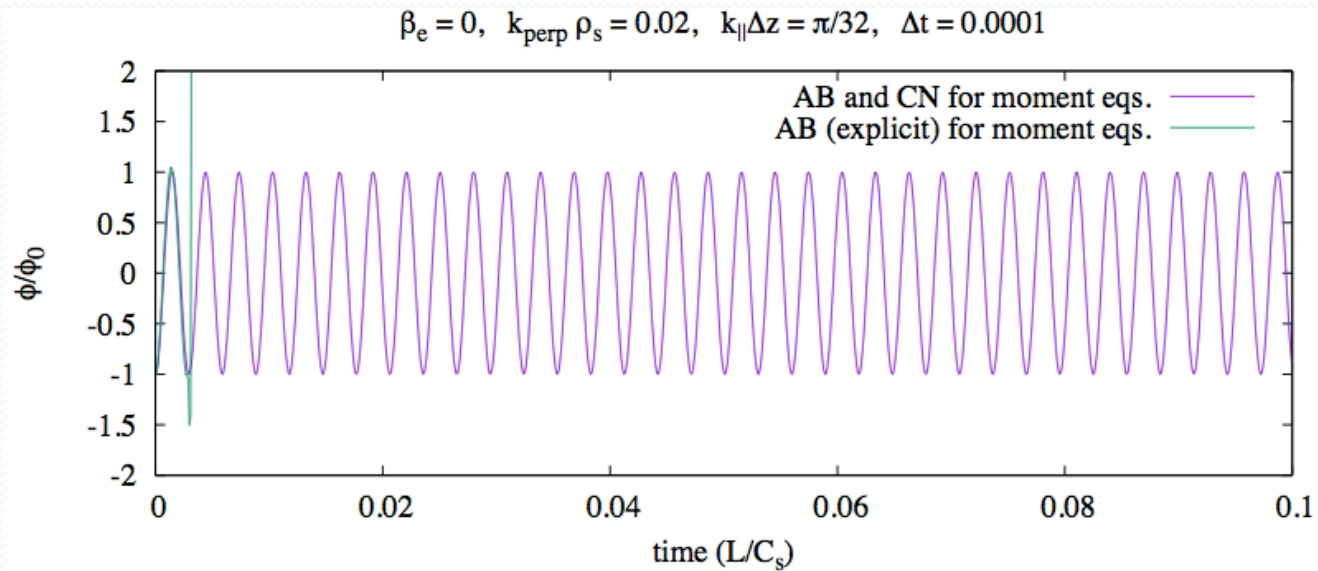
- Numerical solution of the DK equation with the moment extracted formulation successfully describes the KAWs.



- Ion polarization is included in the form of a long wave-length limit.

Application of an implicit solver to the fluid equations

- Adams-Bashforth + Crank-Nicolson
- An explicit integrator is used for the temperature gradient
- Stable solutions can be obtained even with a larger time step size than that given by the Courant number for ω_H .



- Explicit integrator is unstable for $k_{\text{perp}} \rho_s < 0.12$, while the implicit can be successfully applied to $k_{\text{perp}} \rho_s < 0.01$.

Summary

- We developed a numerical scheme for handling the fast electron motion in gyrokinetic simulation.
 - Solve a kinetic equation, of which 0th and 1st order moments are extracted, and the electron fluid equations.
 - Wave propagation can be solved implicitly, while an explicit scheme is applied to the kinetic equation.
 - Modified Maxwellian distribution of which 0th and 2nd order moments exactly satisfy the identities.
 - Application of the semi-Lagrangian scheme has also been tested, demonstrating its numerical stability and efficiency, where kinetic and fluid equations can be solved explicitly with a time step size beyond the CFL limit.