

## Introduction

- Trustworthy gyrokinetic or two-fluid edge turbulence simulations require accurate representations of collisions
- Gyrokinetic simulations (largely) limited to model operators (e.g. Sugama operator) or drift-kinetic Landau operator
- Fluid simulations limited to analytical Braginskii approximation (Landau based, low order moments, infinite electron-ion mass ratio)
- Meaningful comparisons in fluid limit impossible
- Calculating (linearized) gyrokinetic complete Landau operator matrix elements nearly exactly can solve both problems

## Landau collision operator

• nonlinear:  $c^{ab}(f_a, f_b) = \frac{c^{ab}}{m_a} \nabla_a \cdot \int d^3 v_b K(\mathbf{u}) \cdot \left( \frac{f_b \nabla_a f_a}{m_a} - \frac{f_a \nabla_b f_b}{m_b} \right)$   $K(\mathbf{u}) := \frac{\mathbb{1} - \hat{\mathbf{u}}\hat{\mathbf{u}}}{u}$   $\mathbf{u} = \mathbf{v}_a - \mathbf{v}_b$

• linearized:  $C^{ab,T} \delta f_a + C^{ab,F} \delta f_b := c^{ab}(\delta f_a, f_{b0}) + c^{ab}(f_{a0}, \delta f_b)$   $\delta f_s := f_s - f_{s0}$

• matrix elements:  $\langle f_{a,i,l} | C^{ab} | f_{b,j,l} \rangle$   $f_{s,i,l}(\mathbf{v}) = L_l(v_{\parallel s}/v_s) P_l(v_s) \exp(-m_s v_s^2/(2T))$   
 $L_l$  Legendre polynomials  $P_l$  half-sided Hermite or similar polynomials

scalar products:  $\langle f_s | g_s \rangle := \int \frac{f_s(\mathbf{v}_s) g_s(\mathbf{v}_s)}{f_{s0}(\mathbf{v}_s)} d^3 v_s$  operator notation as in quantum mechanics

- regular (“unmagnetized”) matrix elements extremely efficient, hundreds of them in seconds on a laptop
- machine precision accuracy (14-15 digits) even for unrealistically high mass ratios ( $\sim 10^{10}$ )
- automatic conservation of all physical conservation laws, H-theorem
- can be used to compute standard unmagnetized transport coefficients with unprecedented precision

## Unmagnetized transport coefficients

- Spitzer transport coefficients illustrating convergence for increasing number n of radial polynomials (Z=1, infinite mass ratio):

n	$\lambda_{11}$	$\lambda_{12}$	$\lambda_{22}$	$\alpha_i$
4	1.9729113706547	1.3507439973966	3.5912394745740	3.5751695538083
10	1.9758136156134	1.3887606183611	4.1791659178893	3.9502079279755
20	1.9758224556276	1.3887475542516	4.1791860160905	3.9502080688609
40	1.9758225395211	1.3887474651190	4.1791861130853	3.9502080688623
60	1.9758225395566	1.3887474650753	4.1791861131411	3.9502080688623

[Belli 2012]	1.9757	1.3889	4.1789	
[Hinton 83]	1.975	1.389	4.174	3.91

with definitions [Hinton 83]

$$j_{\parallel,e} = \frac{enT\tau_{ei}}{m_e} (\lambda_{11}A_1 + \lambda_{12}A_2), \quad q_{\parallel,e} = -\frac{nT^2\tau_{ei}}{m_e} (\lambda_{11}A_1 + \lambda_{12}A_2), \quad q_{\parallel,i} = -\frac{nT^2\tau_{ii}}{m_i} \alpha_i A_3$$

$$A_1 = \partial_{\parallel} \ln p_e + \frac{e}{T_e} \partial_{\parallel} \phi, \quad A_2 = \partial_{\parallel} \ln T_e, \quad A_3 = \partial_{\parallel} \ln T_i$$

Braginskii style transport coefficients are related by [NRL 2019]

$$\sigma_{\parallel} = \lambda_{11} n e^2 \tau_{ei} / m_e, \quad R_{T,\parallel} = -\lambda_{12} / \lambda_{11} n \partial_{\parallel} T_e, \quad \kappa_{\parallel}^e = (\lambda_{22} - \lambda_{12}^2 / \lambda_{11}), \quad \kappa_{\parallel}^i = \alpha_i n T_i \tau_{ii} / m_i$$

- further Braginskii coefficients compared to best known values (Z=1, infinite mass ratio, all given digits significant):

	[NRL 2019]	accurate value
frictional heat flux coefficient	0.71	0.70287054493606
parallel electron thermal conductivity	3.2	3.203076425585
ion viscosity $\eta_0^i / (nT\tau_{ii})$	0.96	0.96529161180214
electron viscosity $\eta_0^e / (nT\tau_{ei})$	0.73	0.733488956541

- customary transport values not given in truly **independent** set of variables (e.g. parallel velocity depends on density gradient)
- in general: transport matrix in independent set of variables

## Conclusions:

- Inaccuracies of Landau operator larger than numerical ones
- Accuracy takes care of fundamental properties (self-adjointness, conservation laws, invariances, H-theorem)

## Gyrotransformation

- gyrokinetic matrix elements:  $\langle f_{gc,a,i_a,l_a} | C^{ab} | f_{gc,b,i_b,l_b} \rangle$   
 (i.e. gyrocenter coordinates)  $f_{gc,a,i_a,l_a}(\mathbf{v}_a) := \exp(-i\mathbf{k} \cdot \boldsymbol{\rho}_a) f_{a,i_a,l_a}(v_{\perp}, v_{\parallel})$   
 $\boldsymbol{\rho}_a = \frac{m_a}{q_a B} (\mathbf{B} \times \mathbf{v}_a)$
- gyrotransformation by expansion in regular matrix elements
- $k$  expansion in Chebyshev polynomials
- seconds on parallel machine for reasonable  $k\rho_i \lesssim 100$  to machine precision (10-14 digits)
- physical restrictions maintained due to **accuracy**
- can be used to calculate highly accurate perpendicular transport coefficients

## Magnetized transport coefficients

- perpendicular transport coefficients not corresponding to simple numbers in [NRL 2019] (all given digits are significant):

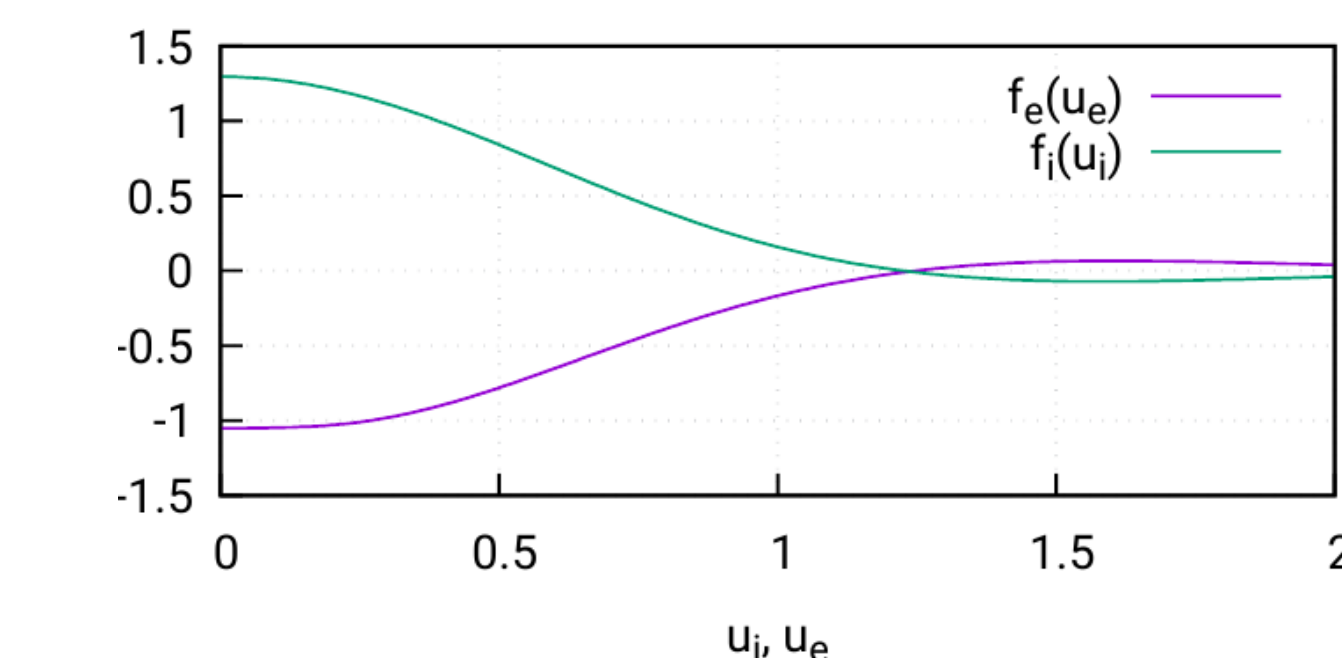
	[NRL 2019]	accurate value
perp. electron heat flux coefficient $m_e \omega_{ce}^2 \tau_{ei} \kappa_{\perp,e} / (nT_e)$	4.7	4.6642135623731056
perpendicular electron viscosity $\eta_1^e / (nT\tau_{ei})$	0.51	0.71213203435597
perp./parallel electron viscosity $\eta_2^e / (nT\tau_{ei})$	2.0	2.84853

- one value is so accurate that one can guess a (new) analytical expression:  $m_e \omega_{ce}^2 \tau_{ei} \kappa_{\perp,e} / (nT_e) = 13/4 + \sqrt{2}$
- the deviation by about  $\sqrt{2}$  for the other two values hints at a mix-up between  $\tau_{ei}$  and  $\tau_{ii}$  in the original publication
- all other values agree with [NRL 2019]

## Realistic ion/electron mass ratio

- several degrees of freedom to expand kinetic equation to obtain fluid transport coefficients, one possibility: replace stationary infinite mass ratio eigenstates corresponding to  $T_i, T_e, n_i, n_e, u$  by slowly changing eigenstates, e.g., one such eigenstate corresponds to  $\delta T_i = -\delta T_e, n_i = n_e = n_0, u = 0,$

heat exchanging distribution:



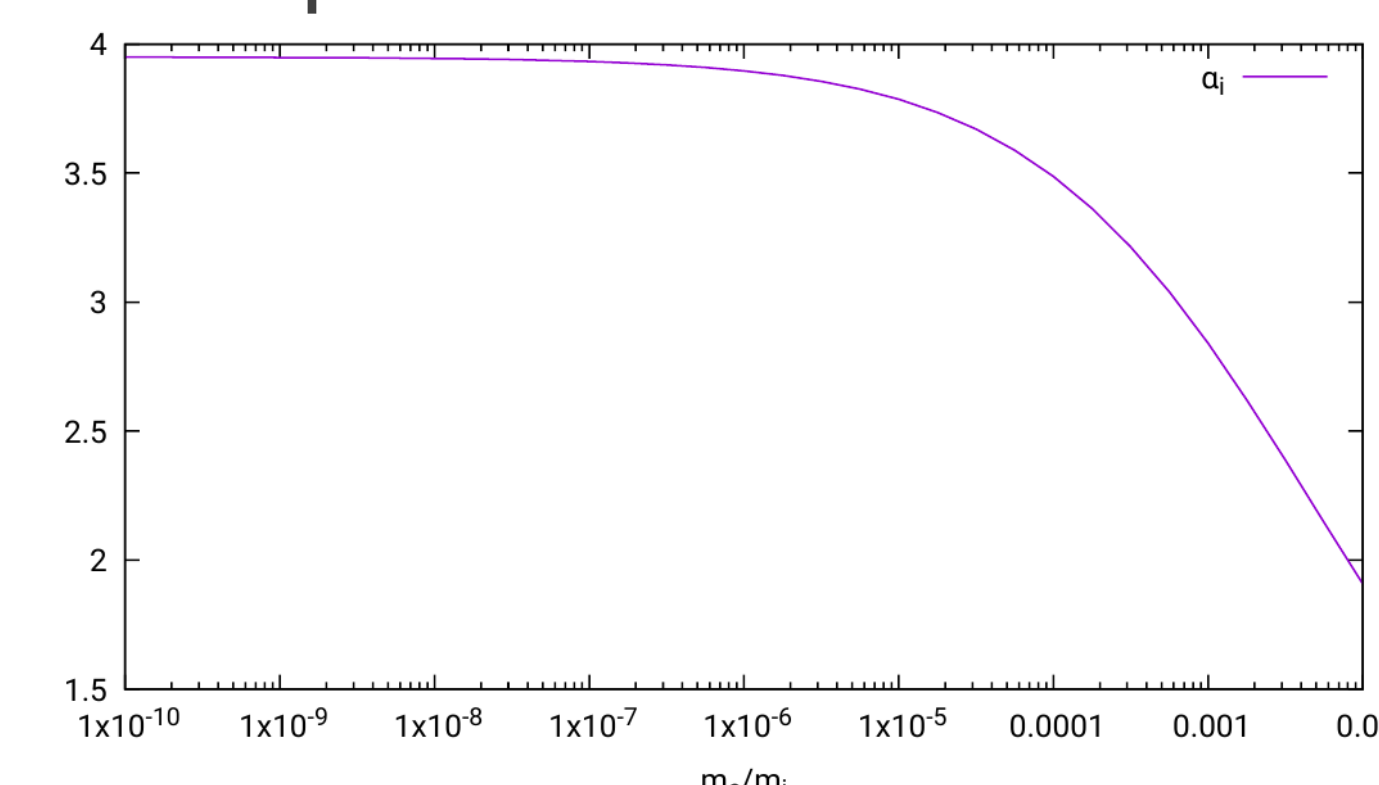
$$u_a := \sqrt{2T_a/m_a} \quad m_i/m_e = 100$$

$f_e(u_e), f_i(u_i)$  spherically symmetric perturbations of electron, ion distribution function

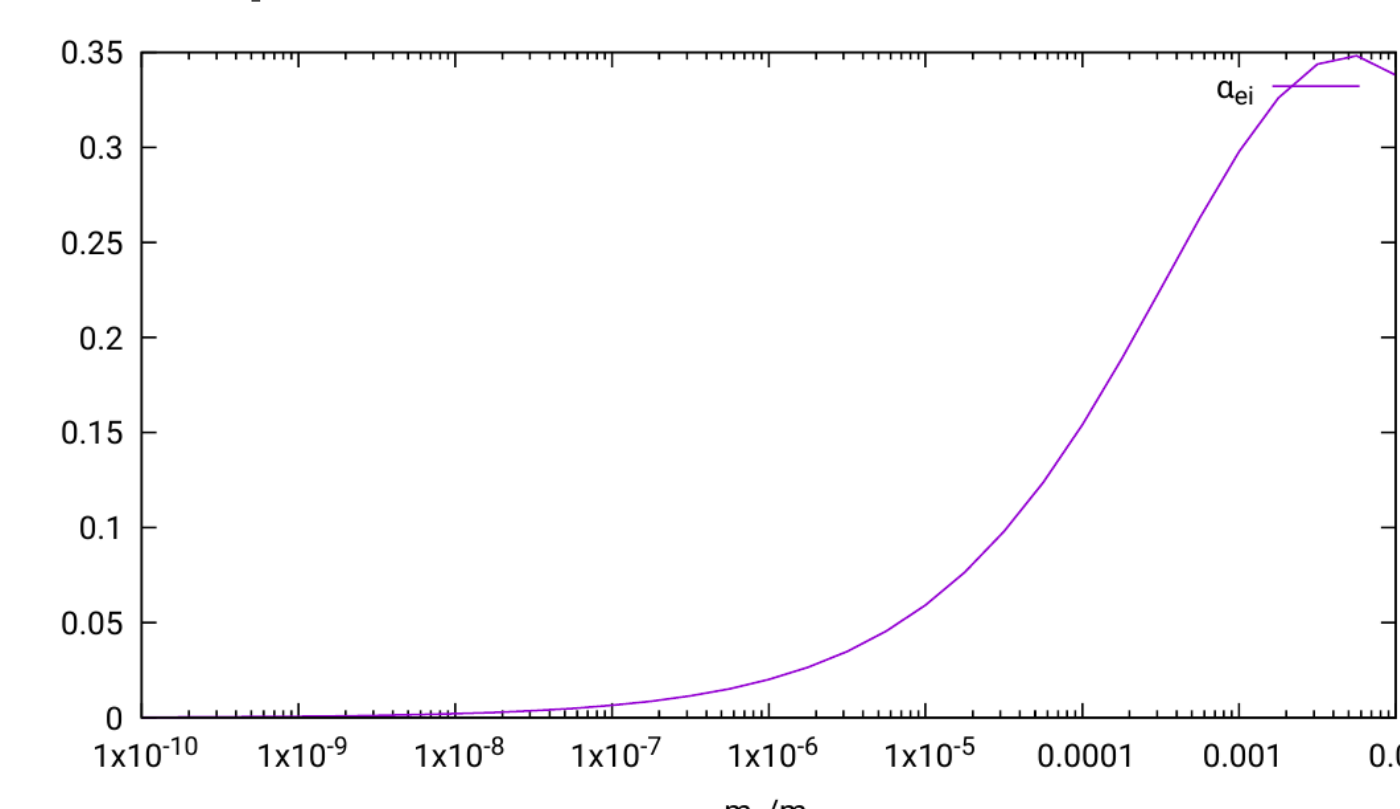
- significant changes from infinite mass ratio values

- e.g. for Deuterium:  $\alpha_i = 3.25485015$   $\eta_0^i = 0.913634368$

- dependence of parallel heat flux coefficient on mass ratio:



- mass ratio dependent electron-ion cross heat flux significant:



$$q_{i,e} = \alpha_{ei} n T_i \tau_{ii} / m_i \partial_{\parallel} T_e$$

- e.g. this term is completely absent in Braginskii.

- Sugama model operator in comparison:  $(\alpha_i, \eta_0^i) = (2.58, 0.760)$

- model operators/infinite mass ratio both partially far off
- Spin-off: Extremely accurate fluid transport coefficients
- Blueprint for more advanced operators (Balescu, Boltzmann)