Differential formulation of the gyrokinetic Landau operator

Eero Hirvijoki^{1,†}, Alain J. Brizard² and David Pfefferlé¹

¹Princeton Plasma Physics Laboratory, Princeton, NJ 08543, USA ²Department of Physics, Saint Michael's College, Colchester, VT 05439, USA

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Subsequent to the recent rigorous derivation of an energetically consistent gyrokinetic collision operator in the so-called Landau representation, this paper investigates the possibility of finding a differential formulation of the gyrokinetic Landau collision operator. It is observed that, while a differential formulation is possible in the gyrokinetic phase space, reduction of the resulting system of partial differential equations to five dimensions via gyroaveraging poses a challenge. Based on the present work, it is likely that the gyrocentre analogues of the Rosenbluth–MacDonald–Judd potential functions must be kept gyroangle dependent.

Key words: fusion plasma, plasma nonlinear phenomena, plasma simulation

1. Introduction

Phase-space Lagrangian perturbation theory, formulated in terms of Lie transformations and first applied to the guiding-centre motion by Littlejohn (1982), provides a rigorous basis for gyrokinetics in that it yields exact energy and momentum invariants against which gyrokinetic codes can be validated. For a comprehensive overview of the topic, an exhaustive list of references could be pointed out. Instead, to maintain focus, we refer the reader to the review papers by Brizard & Hahm (2007) and Cary & Brizard (2009) and the references therein.

In spite of the success of Lie-transform perturbation theory, the current state-of-theart gyrokinetics tends to deviate from the path of strict rigorousness when discussion turns into the collision operator. It is our purpose to criticize neither the existing theory nor codes, but merely to point out how difficult consistency is to achieve if collisional effects are considered. And by rigour, we refer to the treatment of finite-Larmor-radius (FLR) effects that result in the collision operator from the coupling of configuration and velocity space coordinates in the gyrocentre transformation (see, e.g. Catto & Tsang 1977; Xu & Rosenbluth 1991).

This is not to say that significant effort would not have been made in order to develop collision operators for gyrokinetics. In contrast, several papers (see, e.g. Abel et al. 2008; Sugama, Watanabe & Nunami 2009; Li & Ernst 2011) provide expressions for both model and full linearized collision operators that are suitable for numerical implementation. These linear operators are typically presented in the form of

$$C^{GK}[h] = \langle e^{\rho \cdot \nabla} C[e^{-\rho \cdot \nabla} h] \rangle, \qquad (1.1)$$

†Email address for correspondence: ehirvijo@pppl.gov

where ρ is the Larmor radius and h is the non-adiabatic part of the perturbed distribution function. The prevalence of this expression, even during the current era of modern gyrokinetics, probably stems from history: in local flux tube simulations with periodic boundary conditions for the configuration space, Fourier transformation of the above expression admits straightforward and explicit gyroaveraging if ρ is taken to be of zeroth order with respect to the guiding-centre transformation. The caveat in these operators is, however, the lack of modern day push-forward and pull-back operators. While the push-forward-pull-back formalism is rigorously used on the left-hand side of the kinetic equation, equally accurate treatment of the collision operator has been all but neglected. After all, using push-forward and pull-back operators, the expression for the linearized collision operator should read

$$C^{GK}[h] = \langle T_{gy}^{-1} C[T_{gy}h] \rangle, \tag{1.2}$$

where T_{gy}^{-1} and T_{gy} denote the push-forward and pull-back operators between the particle and gyrocentre phase space. We believe the reason behind this discrepancy is the difficulty in explicitly and consistently applying the push forward to velocity derivatives present in the collision operator.

Some have derived collisional gyrokinetics even starting directly from many body gyrokinetics (see the elegant paper by Mishchenko & Könies 2007), but the pioneering step towards developing a collision operator, consistent with the modern day gyrokinetics, was put forward by Brizard (2004). He understood that the particle phase-space velocity derivative could be presented in terms of the non-canonical particle phase-space Poisson bracket. This discovery lead to the first consistent guiding-centre test particle operator and also to consistent bounce-averaged formalism (see Brizard et al. 2009). Recently, the development of consistent operators has accelerated (see Madsen 2013) and culminated in the results presented in Burby, Brizard & Qin (2015). By this date, the full nonlinear Landau collision operator has been transformed to gyrokinetic phase space and exact conservation laws together with H-theorem demonstrated up to arbitrary order with respect to the asymptotic gyrocentre transformation. For the first time, the equations behind gyrokinetics can thus be considered whole in the sense that both the Vlasov part and the collision operator can be transformed to the gyrocentre phase space using the same push-forward and pull-pack operators.

The rigorousness, however, comes with a heavy price tag. The gyrocentre Landau operator is an integro-differential operator, whose evaluation involves a rather complicated six-dimensional (6-D) integral over the gyrocentre phase space. In search of salvation, one remembers that, in particle phase space, the Landau operator can be converted into a system of coupled partial differential equations via the Rosenbluth-MacDonald-Judd (hereafter RMJ) potential functions (Rosenbluth, MacDonald & Judd 1957). One could thus expect the same to be true for the gyrocentre collision operator: after all, the gyrokinetic Landau operator results from the Lie transformation of the particle phase-space Landau operator. Furthermore, in particle phase space, the differential representation has allowed the use of fast elliptic solvers to speed up the evaluation of collisions, reducing the computational complexity from $O(n^2)$ for the integro-differential to $O(n \log n)$ for the differential formulation, as was demonstrated in Pataki & Greengard (2011). To be more precise, the RMJ potentials, in terms of which the velocity-space Fokker-Planck coefficients are often expressed, can be computed either by solving the corresponding Poisson equations, a problem that leads to sparse matrix inversion, or by evaluating the corresponding Green's function solutions, a problem which translates to full matrix-vector multiplication. Having a similar option for the gyrocentre collision operator could be especially important since the coupling of configuration and velocity-space coordinates in gyrokinetics actually increases the dimensionality of the collision operator, in contrast to what is often thought.

The purpose of this paper is, first of all, to present the differential formulation of the gyrocentre Landau collision operator and, secondly, to discuss the non-trivial issues that arise with it. We start by expressing the particle phase-space collisional kinetic theory in terms of the non-canonical Poisson bracket and illustrate the close analogy between using either the bracket or curvilinear velocity-space coordinates. Once this connection is demonstrated for the reader to feel less alienated with expressing everything in terms of brackets, the gyrocentre coordinate transformation is applied to particle phase-space collision operator and the differential form of the gyrokinetic collision operator presented and briefly compared to the Landau version. Finally, we discuss why the reduction of the 6-D differential version of the gyrocentre collision operator to a 5-D gyroaveraged version poses a challenge: the gyrocentre versions of the RMJ potentials cannot easily be reduced to five dimensions except in the limit of zero Larmor radius, in which case the operator naturally collapses to the axially symmetric particle phase-space collision operator.

2. Particle phase-space collision operator

Let us start by briefly recalling the collisional theory in particle phase space. In a multispecies plasma, the kinetic equation that describes the evolution of the particle distribution function f_s of species s due to both Hamiltonian motion and collisions with species \bar{s} is written as

$$\frac{\partial f_s}{\partial t} + \dot{z}^{\alpha} \frac{\partial f_s}{\partial z^{\alpha}} = C_{s\bar{s}}[f_s, f_{\bar{s}}], \tag{2.1}$$

where z = (x, v) are the phase-space coordinates, and the Fokker-Planck collision operator is defined as

$$C_{s\bar{s}}[f_s, f_{\bar{s}}] \equiv \frac{\partial}{\partial v^i} \left(D_{s\bar{s}}^{ij} \frac{\partial f_s}{\partial v^j} - K_{s\bar{s}}^i f_s \right). \tag{2.2}$$

Here the distribution function is normalized to the density according to $n_s(\mathbf{x}) = \int d\mathbf{v} f_s$ and the Latin indices i and j refer to Cartesian coordinates. The collisional velocity-space friction (or drag), $K_{s\bar{s}}^i$, and the collisional velocity-space diffusion, $D_{s\bar{s}}^{ij}$, describe the Coulomb interaction between particles of species s and \bar{s} .

The Coulomb friction and diffusion coefficients are compactly defined in terms of the RMJ potentials (Rosenbluth *et al.* 1957)

$$K_{s\bar{s}}^{i} = -\gamma_{s\bar{s}} \frac{m_{s}}{m_{\bar{s}}} \frac{\partial \phi_{\bar{s}}}{\partial v^{i}}, \quad D_{s\bar{s}}^{ij} = -\gamma_{s\bar{s}} \frac{\partial^{2} \psi_{\bar{s}}}{\partial v^{i} \partial v^{j}}, \tag{2.3a,b}$$

where $\gamma_{s\bar{s}} = e_s^2 e_{\bar{s}}^2 \ln \Lambda / (m_s \epsilon_0)^2$. The particle phase-space potentials are weighted integrals of the distribution function

$$\phi_{\bar{s}}(\boldsymbol{x},\,\boldsymbol{v}) = -\frac{1}{4\pi} \int d\bar{\boldsymbol{v}} f_{\bar{s}}(\boldsymbol{x},\,\bar{\boldsymbol{v}}) |\boldsymbol{v} - \bar{\boldsymbol{v}}|^{-1}, \tag{2.4}$$

$$\psi_{\bar{s}}(\boldsymbol{x},\,\boldsymbol{v}) = -\frac{1}{8\pi} \int d\bar{\boldsymbol{v}} f_{\bar{s}}(\boldsymbol{x},\,\bar{\boldsymbol{v}}) |\boldsymbol{v} - \bar{\boldsymbol{v}}|, \tag{2.5}$$

representing the free-space Green's function solutions to the following Poisson equations

$$\frac{\partial^2 \phi_{\bar{s}}}{\partial v^i \partial v^i} = f_{\bar{s}}, \quad \frac{\partial^2 \psi_{\bar{s}}}{\partial v^i \partial v^i} = \phi_{\bar{s}}. \tag{2.6a,b}$$

Furthermore, the diffusion and friction coefficients satisfy the relation

$$\frac{\partial}{\partial v^j} D_{s\bar{s}}^{ij} = \frac{m_{\bar{s}}}{m_s} K_{s\bar{s}}^i, \tag{2.7}$$

which can be used to show the equivalence between the Fokker-Planck operator and the Landau operator (Landau 1937)

$$C_{s\bar{s}}[f_s, f_{\bar{s}}] = \gamma_{s\bar{s}} \frac{m_s}{8\pi} \frac{\partial}{\partial v^i} \int d\bar{v} U_{ij} \left[\frac{f_{\bar{s}}}{m_s} \frac{\partial f_s}{\partial v^j} - \frac{f_s}{m_{\bar{s}}} \frac{\partial f_{\bar{s}}}{\partial \bar{v}^j} \right], \tag{2.8}$$

where the second-rank tensor U_{ij} is defined as the velocity-space Hessian of the relative speed $|u| = |v - \bar{v}|$ according to

$$U_{ij} = \frac{\partial^2 u}{\partial v^i \partial v^j} = \frac{1}{u} \left(\delta^{ij} - \frac{u^i u^j}{u^2} \right). \tag{2.9}$$

The tensor U_{ij} is often called the Coulomb kernel and is a projection operator with the null-space vector \mathbf{u} . Its trace $U_{ii} = 2/u$ is the Laplacian's Green function.

3. Poisson bracket formulation of kinetic theory

Before working out the expression for the gyrokinetic collision operator, let us prepare the stage by introducing the use of Poisson brackets in the particle phase space. In the paper by Brizard (2004) it was pointed out that, using the particle phase-space Poisson bracket, a momentum derivative of an arbitrary function could be written as $\{x, g\} \equiv \partial_p g$. With multiple species and different masses one can apply this idea into the collision operator by following

$$\frac{\partial f}{\partial v^i} = m_s \{ x^i, f \}_s,\tag{3.1}$$

where the bracket $\{\cdot,\cdot\}_s$, refers to the bracket that is derived from the Lagrangian one form of species s.

3.1. Friction-diffusion representation

As the Hamiltonian equations of motion are inherently expressed with the Poisson bracket, the kinetic equation can then be formulated as (Brizard 2004)

$$\frac{\partial f_s}{\partial t} + \{f_s, H_s\}_s = \{x^i, m_s^2 D_{s\bar{s}}^{ij} \{x^j, f_s\}_s - m_s K_{s\bar{s}}^i f_s\}_s, \tag{3.2}$$

where H_s is the Hamiltonian for the species s. While in his original work Brizard used this formalism to derive a guiding-centre test particle collision operator and, later, also a bounce-averaged test-particle operator (see Brizard $et\ al.\ 2009$), the use of brackets was extended in the paper by Madsen (2013) to express also the friction and diffusion coefficients according to

$$K_{s\bar{s}}^{i} = -\gamma_{s\bar{s}} \frac{m_{s}^{2}}{m_{\bar{s}}} \{x^{i}, \phi_{\bar{s}}\}_{s}, \tag{3.3}$$

$$D_{s\bar{s}}^{ij} = -\gamma_{s\bar{s}} m_s^2 \{ x^i, \{ x^j, \psi_{\bar{s}} \}_s \}_s. \tag{3.4}$$

Completing the picture for particle phase space, we use the brackets to express also the differential equations from which the RMJ potentials are solved

$$m_s^2 \{x^i, \{x^i, \phi_{\bar{s}}\}_s\}_s = f_{\bar{s}}, \quad m_s^2 \{x^i, \{x^i, \psi_{\bar{s}}\}_s\}_s = \phi_{\bar{s}}.$$
 (3.5*a*,*b*)

3.2. Landau representation

As an alternative to the RMJ potential formulation, one may start from the Landau form of the collision operator and similarly use the bracket notation to obtain (see Burby *et al.* 2015)

$$C_{s\bar{s}}[f_s, f_{\bar{s}}] = \gamma_{s\bar{s}} \frac{m_s^2}{8\pi} \{ x^i, \, \Gamma_{s\bar{s}}^i \}_s, \tag{3.6}$$

where the collisional flux $\Gamma_{s\bar{s}}^i$ is defined as the phase-space integral

$$\Gamma_{s\bar{s}}^{i} = \int d\bar{z} \, \delta(\bar{x} - x) U^{ij}(z, \bar{z}) (f_{\bar{s}}(\bar{z}) \{x^{i}, f_{s}\}_{s}(z) - f_{s}(z) \{\bar{x}^{i}, f_{\bar{s}}\}_{\bar{s}}(\bar{z})), \tag{3.7}$$

and the relative velocity needed for constructing the kernel $U^{ij}(z,\bar{z})$ is written as

$$u^{i}(z,\bar{z}) \equiv \{x^{i}, H_{s}\}_{s}(z) - \{\bar{x}^{i}, H_{\bar{s}}\}_{\bar{s}}(\bar{z}).$$
 (3.8)

Note that the brackets for both species s and \bar{s} are required. The delta function in the integrand of the collisional flux $\Gamma_{s\bar{s}}^i$ represents the local nature of the collision operator in the particle phase space but allows one to express the flux as phase-space integral rather than as a velocity-space integral. This detail is necessary for conducting the gyrocentre transformation of the Landau operator.

3.3. Equivalence to curvilinear coordinate system

At first sight, the idea of converting the partial derivatives to brackets might appear counterintuitive but it is in fact equivalent to expressing the derivatives in curvilinear coordinates u^{α} . Since for arbitrary functions f and g we have according to Liouville's theorem

$$\{f,g\} = \frac{1}{J} \frac{\partial}{\partial u^{\alpha}} [J\{f, u^{\alpha}\}g], \tag{3.9}$$

we can transform Cartesian velocity-space divergence into curvilinear coordinates according to

$$\frac{\partial A^{i}}{\partial v^{i}} = \frac{1}{J} \frac{\partial}{\partial u^{\alpha}} [J\{x^{i}, u^{\alpha}\}_{s} m_{s} A^{i}] = \frac{1}{J} \frac{\partial}{\partial u^{\alpha}} [JA^{\alpha}], \tag{3.10}$$

where *J* is the Jacobian of the transformation and $A^{\alpha} = A^{i} \{x^{i}, u^{\alpha}\}_{s} m_{s} = A^{i} \partial_{i} u^{\alpha}$ is the curvilinear component of the vector A^{i} .

This allows us to express the collision operator in a form

$$C[f_s, f_{\bar{s}}] = \frac{1}{J} \frac{\partial}{\partial u^{\alpha}} \left[J\{x^i, u^{\alpha}\}_s m_s^2 D_{s\bar{s}}^{ij} \{x^j, u^{\beta}\}_s \frac{\partial f}{\partial u^{\beta}} - J m_s K_{s\bar{s}}^i \{x^i, u^{\alpha}\}_s f \right]. \tag{3.11}$$

Similarly, we may express the Cartesian components of friction and diffusion coefficients according to

$$K_{s\bar{s}}^{i} = -\gamma_{s\bar{s}} \frac{m_{s}^{2}}{m_{\bar{s}}} \{x^{i}, u^{\beta}\}_{s} \frac{\partial \phi_{\bar{s}}}{\partial u^{\beta}}, \tag{3.12}$$

$$D_{s\bar{s}}^{ij} = -\gamma_{s\bar{s}}m_s^2 \left[\{x^i, u^\sigma\}_s \{x^j, u^\nu\}_s \frac{\partial^2 \psi_{\bar{s}}}{\partial u^\sigma \partial u^\nu} + \{x^i, u^\sigma\}_s \frac{\partial \{x^j, u^\nu\}_s}{\partial u^\sigma} \frac{\partial \psi_{\bar{s}}}{\partial u^\nu} \right]. \tag{3.13}$$

If we then define the transformed friction and diffusion coefficients

$$K_{s\bar{s}}^{\alpha} = m_s \{x^i, u^{\alpha}\}_s K_{s\bar{s}}^i, \tag{3.14}$$

$$D_{s\bar{s}}^{\alpha\beta} = m_s^2 \{ x^i, u^{\alpha} \}_s D_{s\bar{s}}^{ij} \{ x^j, u^{\beta} \}_s, \tag{3.15}$$

the collision operator can be compactly written as

$$C[f_s, f_{\bar{s}}] = \frac{1}{J} \frac{\partial}{\partial u^{\alpha}} \left[J D_{s\bar{s}}^{\alpha\beta} \frac{\partial f}{\partial u^{\beta}} - J K_{s\bar{s}}^{\alpha} f \right]. \tag{3.16}$$

The connection to curvilinear coordinates becomes transparent when we remind ourselves that the particle phase-space brackets were introduced to express velocityspace derivatives and that we have the identities

$$\{x^{i}, u^{\alpha}\}_{s} \{x^{i}, u^{\beta}\}_{s} = \frac{1}{m_{s}^{2}} g^{\alpha\beta},$$
 (3.17)

$$\{x^i, u^\alpha\}_s \frac{\partial \{x^i, u^\nu\}_s}{\partial u^\beta} = -\frac{1}{m_s^2} g^{\alpha\ell} \Gamma_{\beta\ell}^\nu, \tag{3.18}$$

where $g^{\alpha\beta}$ is the inverse of the metric tensor for the curvilinear velocity-space coordinates u^{α} and $\Gamma^{\nu}_{\alpha\beta}$ are the Christoffel symbols of the second kind. Our Poisson bracket expressions for the transformed friction and diffusion coefficients are thus equivalent to

$$K_{s\bar{s}}^{\alpha} = -\gamma_{s\bar{s}} \frac{m_s}{m_{\bar{s}}} g^{\alpha\beta} \frac{\partial \phi_{\bar{s}}}{\partial u^{\beta}}, \tag{3.19}$$

$$D_{s\bar{s}}^{\alpha\beta} = -\gamma_{s\bar{s}} \left[g^{\alpha\sigma} g^{\nu\beta} \frac{\partial^2 \psi_{\bar{s}}}{\partial u^{\sigma} \partial u^{\nu}} - g^{\alpha\sigma} g^{\beta\ell} \Gamma_{\sigma\ell}^{\nu} \frac{\partial \psi_{\bar{s}}}{\partial u^{\nu}} \right], \tag{3.20}$$

which are nothing more than the expressions one would find starting with curvilinear coordinates (Goncharov 2010).

4. Gyrocentre collision operator

We are finally ready to apply the gyrocentre transformation of the collision operator. It is obtained by replacing the particle phase-space Poisson bracket with the gyrocentre Poisson bracket, and by evaluating the particle phase-space quantities in terms of the

gyrocentre coordinates Z^{α} . This is the consequence of the transformation rules using the (species-wise) push-forward \mathcal{T}_{sgy}^{-1} and pull-back \mathcal{T}_{sgy} operators, which can be summarized in the following manner. Scalar functions transform according to

$$f_s(z) = (T_{sgy}F)(z) = (T_{sgy}F_s)(T_{sgy}^{-1}\mathbf{Z}) = F_s(\mathbf{Z}),$$
 (4.1)

$$F_s(\mathbf{Z}) = (T_{sgy}^{-1}f)(\mathbf{Z}) = (T_{sgy}^{-1}f_s)(T_{sgy}z) = f_s(z),$$
 (4.2)

where $f_s(z)$ is understood as an arbitrary function in the particle phase space of species s and $F_s(\mathbf{Z})$ as an image of f_s at the gyrocentre phase space. The Poisson bracket transforms under the chain rule for derivatives and the rule for scalar functions:

$$\{f(z), g(z)\} = \{F(Z), G(Z)\}_{gy}.$$
 (4.3)

Any phase-space integral transforms simply by the rule of transforming the differential volume element and the integrand giving

$$\int d\mathbf{z} f(\mathbf{z}) \equiv \int d\mathbf{Z} F(\mathbf{Z}) = \int d^6 Z \mathcal{J}_{gy} F(\mathbf{Z}), \tag{4.4}$$

where \mathcal{J}_{gy} is the gyrocentre phase-space Jacobian.

4.1. Friction-diffusion representation

Applying the transformation rules, we immediately find the gyrocentre Fokker-Planck collision operator

$$C_{s\bar{s}}[f_s, f_{\bar{s}}] \equiv C_{s\bar{s}}^{gy}[F_s, F_{\bar{s}}] = \{ T_{sgy}^{-1} X^i, \ m_s^2 T_{sgy}^{-1} D_{s\bar{s}}^{ij} \{ T_{sgy}^{-1} X^j, F_s \}_{sgy} \}_{sgy}$$

$$- \{ T_{sgy}^{-1} X^i, \ m_s T_{sgy}^{-1} K_{s\bar{s}}^i F_s \}_{sgy},$$

$$(4.5)$$

where the expressions for the pushed-forward particle phase-space friction and diffusion coefficients are

$$T_{sgy}^{-1}D_{s\bar{s}}^{ij} = -\gamma_{s\bar{s}}m_s^2 \{T_{sgy}^{-1}X^i, \{T_{sgy}^{-1}X^j, T_{sgy}^{-1}\psi_{\bar{s}}\}_{sgy}\}_{sgy}, \tag{4.6}$$

$$T_{sgy}^{-1}K_{s\bar{s}}^{i} = -\gamma_{s\bar{s}}\frac{m_{s}^{2}}{m_{\bar{s}}}\{T_{sgy}^{-1}X^{i}, T_{sgy}^{-1}\phi_{\bar{s}}\}_{sgy}.$$
(4.7)

Then, using the Liouville identity, we write the gyrocentre Fokker–Planck operator in the phase-space divergence form

$$C_{s\bar{s}}^{gy}[F_s, F_{\bar{s}}] = \frac{1}{\mathcal{J}_{sev}} \frac{\partial}{\partial Z^{\alpha}} \left[\mathcal{J}_{sgy} \mathcal{D}_{s\bar{s}}^{\alpha\beta} \frac{\partial F_s}{\partial Z^{\beta}} - \mathcal{J}_{sgy} \mathcal{K}_{s\bar{s}}^{\alpha} F_s \right], \tag{4.8}$$

where the gyrocentre friction and diffusion coefficients are defined as

$$\mathcal{D}_{s\bar{s}}^{\alpha\beta} = \Delta_s^{i\alpha} \Delta_s^{j\beta} T_{sev}^{-1} D_{s\bar{s}}^{ij}, \tag{4.9}$$

$$\mathcal{K}^{\alpha}_{s\bar{s}} = \Delta^{i\alpha}_{s} T^{-1}_{sgy} K^{i}_{s\bar{s}}, \tag{4.10}$$

with the so-called projection coefficients defined component-wise according to

$$\Delta_s^{i\alpha} = m_s \{ T_{sgy}^{-1} X^i, Z^\alpha \}_{sgy}, \tag{4.11}$$

and being analogous to Cartesian components of contravariant basis vectors.

Using the projection coefficients and the chain rule for the Poisson brackets, the expressions for the pushed-forward particle phase-space friction and diffusion coefficients become

$$T_{sgy}^{-1}D_{s\bar{s}}^{ij} = -\gamma_{s\bar{s}}\Delta_{s}^{i\sigma}\frac{\partial}{\partial Z^{\sigma}}\left(\Delta_{s}^{j\nu}\frac{\partial}{\partial Z^{\nu}}T_{sgy}^{-1}\psi_{\bar{s}}\right),\tag{4.12}$$

$$T_{sgy}^{-1}K_{s\bar{s}}^{i} = -\gamma_{s\bar{s}}\frac{m_{s}}{m_{\bar{s}}}\Delta_{s}^{i\sigma}\frac{\partial}{\partial Z^{\sigma}}T_{sgy}^{-1}\phi_{\bar{s}}.$$
(4.13)

After introducing the symmetric matrix

$$\Xi_s^{\alpha\beta} = \Delta_s^{i\alpha} \Delta_s^{i\beta},\tag{4.14}$$

we can finally express the gyrocentre friction and diffusion coefficients according to

$$\mathcal{K}^{\alpha}_{s\bar{s}} = -\gamma_{s\bar{s}} \frac{m_s}{m_{\bar{s}}} \Xi^{\alpha\beta}_s \frac{\partial T^{-1}_{sgy} \phi_{\bar{s}}}{\partial Z^{\beta}}, \tag{4.15}$$

$$\mathcal{D}_{s\bar{s}}^{\alpha\beta} = -\frac{\gamma_{s\bar{s}}}{2} \left[(\Xi_s^{\alpha\sigma} \Xi_s^{\beta\nu} + \Xi_s^{\beta\sigma} \Xi_s^{\alpha\nu}) \frac{\partial^2 T_{sgy}^{-1} \psi_{\bar{s}}}{\partial Z^{\sigma} \partial Z^{\nu}} \right]$$

$$+\left.\left(\Xi_{s}^{\alpha\sigma}\Delta_{s}^{k\beta}+\Xi_{s}^{\beta\sigma}\Delta_{s}^{k\alpha}\right)\frac{\partial\Delta_{s}^{k\nu}}{\partial Z^{\sigma}}\frac{\partial\mathcal{T}_{sgy}^{-1}\psi_{\bar{s}}}{\partial Z^{\nu}}\right],\tag{4.16}$$

where the symmetry in $\mathcal{D}_{s\bar{s}}^{\alpha\beta}$ with respect to α and β has been built in using the fact that the particle phase-space diffusion coefficient is symmetric with respect to i and j.

Referring to our earlier discussion of the particle phase-space operator in curvilinear coordinates, we see that the gyrocentre friction and diffusion coefficients appear no more exotic: they essentially have the same form as the particle phase-space friction and diffusion coefficients in curvilinear coordinates.

4.2. Gyrocentre Rosenbluth-MacDonald-Judd potential equations

To complete the gyrocentre transformation of the Fokker–Planck operator, expressions for the gyrocentre RMJ potentials are needed. Two different approaches are possible. One can obtain the potentials directly by transforming the particle phase-space integral definitions into the gyrocentre phase space according to

$$T_{sgy}^{-1}\phi_{\bar{s}} = -\frac{1}{4\pi} \int d\bar{\mathbf{Z}} \, \delta^{gy} \frac{F_{\bar{s}}}{u^{gy}},$$
 (4.17)

$$T_{sgy}^{-1}\psi_{\bar{s}} = -\frac{1}{8\pi} \int d\bar{Z} \, \delta^{gy} F_{\bar{s}} u^{gy},$$
 (4.18)

where the definitions for the gyrocentre delta function and relative velocity are

$$\delta^{gy}(\mathbf{Z}, \bar{\mathbf{Z}}) = \delta(T_{ggy}^{-1} \mathbf{X} - \bar{T}_{\bar{g}gy}^{-1} \bar{\mathbf{X}}), \tag{4.19}$$

$$\mathbf{u}^{gy} = \{ T_{ggy}^{-1} X, \mathcal{H}_s \}_s - \{ \bar{T}_{\bar{s}gy}^{-1} \bar{X}, \mathcal{H}_{\bar{s}} \}_{\bar{s}}, \tag{4.20}$$

and \mathcal{H} is the gyrocentre Hamiltonian. Alternatively, one may push the particle phase-space Poisson equations for the RMJ potentials into the gyrocentre phase space according to

$$m_s^2 \{ T_{sgy}^{-1} X^i, \{ T_{sgy}^{-1} X^i, T_{sgy}^{-1} \phi_{\bar{s}} \}_{sgy} \}_{sgy} = F_{\bar{s}},$$
 (4.21)

$$m_s^2 \{ T_{sov}^{-1} X^i, \{ T_{sov}^{-1} X^i, T_{sov}^{-1} \psi_{\bar{s}} \}_{sgy} \}_{sgy} = T_{sov}^{-1} \phi_{\bar{s}}.$$
 (4.22)

Converting the double brackets into double phase-space divergences similarly as was done for the collision operator, we find phase-space Poisson-like equations for the gyrocentre RMJ potentials

$$\frac{1}{\mathcal{J}_{sgy}} \frac{\partial}{\partial Z^{\alpha}} \left(\mathcal{J}_{sgy} \Xi_{s}^{\alpha\beta} \frac{\partial \mathcal{T}_{sgy}^{-1} \phi_{\bar{s}}}{\partial Z^{\beta}} \right) = F_{\bar{s}}, \tag{4.23}$$

$$\frac{1}{\mathcal{J}_{sgy}} \frac{\partial}{\partial Z^{\alpha}} \left(\mathcal{J}_{sgy} \Xi_{s}^{\alpha\beta} \frac{\partial \mathcal{T}_{sgy}^{-1} \psi_{\bar{s}}}{\partial Z^{\beta}} \right) = \mathcal{T}_{sgy}^{-1} \phi_{\bar{s}}, \tag{4.24}$$

where the matrix $\mathcal{Z}_s^{\alpha\beta}$ now clearly is analogous to the metric in curvilinear coordinates.

4.3. Landau representation

Similarly, using the transformation rules, one may compute the gyrocentre transformation of the particle phase-space Landau operator (for details, see Burby *et al.* 2015). When written in the friction–diffusion form, the coefficients in the gyrokinetic Landau operator become

$$\mathcal{D}_{s\bar{s}}^{\alpha\beta}(\mathbf{Z}) = \frac{\gamma_{s\bar{s}}}{8\pi} \Delta_{s}^{i\alpha} \Delta_{s}^{j\beta} \int d\bar{\mathbf{Z}} \, \delta^{gy} U^{ij} F_{\bar{s}}, \tag{4.25}$$

$$\mathcal{K}_{s\bar{s}}^{\alpha}(\mathbf{Z}) = \frac{\gamma_{s\bar{s}}}{8\pi} \frac{m_s}{m_{\bar{s}}} \Delta_s^{i\alpha} \int d\bar{\mathbf{Z}} \, \delta^{gy} U^{ij} \bar{\Delta}_{\bar{s}}^{j\beta} \frac{\partial F_{\bar{s}}}{\partial \bar{Z}^{\beta}}, \tag{4.26}$$

with U^{ij} now a function of the gyrocentre relative velocity \boldsymbol{u}^{gy} defined in (4.20). As such, these coefficients would seem to be simpler than the expressions we derived in (4.15) and (4.16) but they are difficult to evaluate due to the complicated expression for the gyrocentre delta function $\delta^{gy}(\mathbf{Z}, \bar{\mathbf{Z}})$ defined in (4.19).

5. Unresolved aspects of the gyrokinetic collision operator

As the ultimate goal of gyrokinetic theory is to eliminate the fast gyromotion time scale, the gyrocentre collision operator must be gyroaveraged. Assuming that the distribution functions are gyroangle independent (for the ordering, see Brizard 2004), the gyroaveraged 5-D collision operator is written as

$$\langle C_{s\bar{s}}^{gy}[\langle F_s \rangle, \langle F_{\bar{s}} \rangle] \rangle = \frac{1}{\mathcal{J}_{sgy}} \frac{\partial}{\partial Z^{\alpha}} \left[\mathcal{J}_{sgy} \langle \mathcal{D}_{s\bar{s}}^{\alpha\beta} \rangle \frac{\partial \langle F_s \rangle}{\partial Z^{\beta}} - \mathcal{J}_{sgy} \langle \mathcal{K}_{s\bar{s}}^{\alpha} \rangle \langle F_s \rangle \right]. \tag{5.1}$$

The important question is what happens for the gyrokinetic RMJ potentials and the corresponding differential equations, i.e. can the gyrocentre potential equations be reduced to five dimensions or not?

5.1. Existence of 5-D potential functions and Green's function operators?

In particle phase space, a distribution function with axial symmetry in the velocity space provides particle phase-space RMJ potentials that have the same axial symmetry. In gyrocentre phase space, only in the limit of zero Larmor radius have we been able to explicitly verify that the potentials indeed become gyroangle independent. At the zero-Larmor-radius limit, both the gyroaveraged gyrokinetic Landau collision integral and the differential equations defining the gyrokinetic RMJ potentials reduce to the corresponding axially symmetric particle phase-space expressions due to the gyrocentre transformation collapsing to an identity. If the FLR effects are included, the gyrokinetic RMJ potentials remain gyroangle dependent. This is

reflected in (4.23) and (4.24) in the sense that the tensor $\mathcal{E}_s^{\alpha\beta}$ is gyroangle dependent whereas, in the particle phase space, the velocity-space metric tensor, in either cylindrical or spherical coordinates, is axially symmetric. In short: given a gyroangle-independent gyrocentre distribution function, the gyrocentre potentials cannot be assumed gyroangle independent, and the differential equations (4.23) and (4.24) cannot be reduced to five dimensions.

One could, of course, assume the gyroangle dependency of the gyrokinetic potentials to be weak and approximate

$$T_{sgy}^{-1}\phi_{\bar{s}} \approx \langle T_{sgy}^{-1}\phi_{\bar{s}}\rangle,$$
 (5.2)

$$T_{ggy}^{-1}\psi_{\bar{s}} \approx \langle T_{ggy}^{-1}\psi_{\bar{s}}\rangle. \tag{5.3}$$

This would allow one to average the differential equations (4.23) and (4.24) over the gyroangle and to reduce them to five dimensions. This approximation will, however, most likely lead to violation of the important conservation properties of the gyrokinetic collision operator. Indeed, the latter can be proven only in the Landau formulation and the Landau and potential formulations are equivalent only if the gyroangle dependency of the potentials is preserved.

From a theoretical point of view, it would also be aesthetic to explicitly verify if the differential operator we have derived for the gyrocentre RMJ potentials in (4.23) and (4.24) satisfies

$$\frac{1}{\mathcal{J}_{sgy}} \frac{\partial}{\partial Z^{\alpha}} \left[\mathcal{J}_{sgy} \Xi_{s}^{\alpha\beta} \frac{\partial}{\partial Z^{\beta}} \left(\frac{\delta^{gy}}{u^{gy}} \right) \right] = -4\pi \delta(\mathbf{Z} - \bar{\mathbf{Z}}), \tag{5.4}$$

$$\frac{1}{\mathcal{J}_{sgy}} \frac{\partial}{\partial Z^{\alpha}} \left[\mathcal{J}_{sgy} \Xi_{s}^{\alpha\beta} \frac{\partial}{\partial Z^{\beta}} (\delta^{gy} u^{gy}) \right] = 2 \frac{\delta^{gy}}{u^{gy}}. \tag{5.5}$$

This is a property one would expect given the integral presentations (4.17) and (4.18).

5.2. Numerical considerations

Regardless of whether one opts for the Landau formulation or for the potential formulation, the computational effort in evaluating the gyrokinetic collision operator accurately will be significant. As an alternative, a stochastic approach could be chosen for emulating the Fokker–Planck operator using the corresponding stochastic differential equations (for a discussion on the guiding-centre test-particle operator, see Hirvijoki *et al.* 2013) which require only the evaluation of the friction and diffusion coefficients. The stochastic approach is, however, a completely different paradigm that is set aside in the present discussion.

Choosing the Landau formulation requires evaluation of a 6-D integral over phase space, regardless whether the distribution functions are assumed gyroangle independent. The reason is that the function $\delta^{gy}(\mathbf{Z}, \bar{\mathbf{Z}})$ is implicitly gyroangle dependent on both coordinate spaces and finding the many zeros of its argument is a difficult (if not intractable) nonlinear problem. Choosing the formulation in terms of the gyrokinetic RMJ potentials, 6-D Poisson-like differential equations need to be inverted instead. Evaluating weighted integrals for n points would result in a minimum effort of the order $O(n^2)$ whereas there is evidence that elliptic equations can be inverted with $O(n \log n)$ effort (see Pataki & Greengard 2011). As the number of points n at which the distribution function is presented can be expected to be large, the potential formulation could thus offer significant speed up for the evaluation of

the gyrokinetic collision operator, even if the boundary conditions for the potential equations must be computed from the integral expressions (there typically are far less boundary points than the total number of points).

As a final difficulty, the discrete implementation would have to satisfy the same conservation properties as the continuous operator does. In the case of axially symmetric particle phase-space operator or, equivalently, in the zero-Larmor-radius limit of the gyrokinetic operator, conservative numerical methods have been developed for both the Landau and the potential formulations (see e.g. Taitano *et al.* (2015), Hager *et al.* (2016)). On the other hand, it has been explicitly shown in Burby *et al.* (2015) that energetically consistent collisional gyrokinetics requires both the Vlasov and the collision operator to be treated equally at the same order with respect to the asymptotic gyrocentre transformation. Whether a conservative numerical method can be found for the gyrocentre operator including the FLR effects remains to be seen.

6. Summary

In this paper, a differential formulation for the Lie-transformed gyrocentre collision operator has been derived. This was achieved by transforming not only the particle phase-space Fokker–Planck friction and diffusion coefficients but also the particle phase-space Poisson equations that determine the Rosenbluth–MacDonald–Judd potential functions needed in evaluating the friction and diffusion coefficients. Our final results are summarized in the expressions for the gyroaveraged collision operator defined in (5.1), the gyrocentre friction and diffusion coefficients defined in (4.15) and (4.16) and the gyrocentre equivalents for the Rosenbluth–MacDonald–Judd potential functions defined either through the integrals (4.17) and (4.18) or the differential equations (4.23) and (4.24).

We also argued that the gyrocentre Rosenbluth–MacDonald–Judd potentials are not expected to be gyroangle independent, even if the gyrocentre distribution functions were to be. Thus, the differential equations for the gyrocentre Rosenbluth–MacDonald–Judd potentials cannot be averaged over the gyroangle but remain six-dimensional. However, with both integral and differential expressions available for the gyrocentre potential functions, we expect that fast elliptic solvers could be applied to solve the differential equations with boundary conditions evaluated from the integral expressions, similarly as is commonly done for the particle phase-space Rosenbluth–MacDonald–Judd potentials. As such, this recipe could offer significant speed up from $O(n^2)$ for the pure integral definitions to $O(n \log n)$ if the integral definitions are used only for evaluating boundary conditions for the differential equations.

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