

Abstract Submitted  
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**Numerical calculation of the neoclassical electron distribution function in an axisymmetric torus**<sup>1</sup> B.C. LYONS, S.C. JARDIN, PPPL, J.J. RAMOS, MIT PSFC — We solve for a stationary, axisymmetric electron distribution function ( $f_e$ ) in a torus using a drift-kinetic equation (DKE) with complete Landau collision operator. All terms are kept to gyroradius and collisionality orders relevant to high- temperature tokamaks (i.e., the neoclassical banana regime for electrons). A solubility condition on the DKE determines the non-Maxwellian pieces of  $f_e$  (called  $f_{NMe}$ ) to all relevant orders. We work in a 4D phase space  $(\psi, \theta, v, \lambda)$ , where  $\psi$  defines a flux surface,  $\theta$  is the poloidal angle,  $v$  is the total velocity, and  $\lambda$  is the pitch angle parameter. We expand  $f_{NMe}$  in finite elements in both  $v$  and  $\lambda$ . The Rosenbluth potentials,  $\Phi$  and  $\Psi$ , which define the collision operator, are expanded in Legendre series in  $\cos \chi$ , where  $\chi$  is the pitch angle, Fourier series in  $\cos \theta$ , and finite elements in  $v$ . At each  $\psi$ , we solve a block tridiagonal system for  $f_{NMe}$ ,  $\Phi$ , and  $\Psi$  simultaneously, resulting in a neoclassical  $f_e$  for the entire torus. Our goal is to demonstrate that such a formulation can be accurately and efficiently solved numerically. Results will be compared to other codes (e.g., NCLASS, NEO) and could be used as a kinetic closure for an MHD code (e.g., M3D-C1).

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