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On the two approaches to incorporate wave-particle resonant effects into global test particle simulations

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Key Points:

- We discuss two approaches to incorporate resonant effects into test particle simulation models
- Detailed approaches have been shown for continuous stochastic differential equations (SDE) and the mapping technique, respectively
- In contrast to continuous SDE, the mapping technique allows one to simulate non-linear resonant effects

Abstract

Energetic electron dynamics in the Earth’s radiation belts and near-Earth plasma sheet are controlled by multiple processes operating on very different time scales: from storm-time magnetic field reconfiguration on a timescale of hours to individual resonant wave-particle interactions on a timescale of milliseconds. The most advanced models for such dynamics either include test particle simulations in electromagnetic fields from global magnetospheric models, or those that solve the Fokker-Plank equation for long-term effects of wave-particle resonant interactions. The most prospective method, however, would be to combine these two classes of models, to allow the inclusion of resonant electron scattering into simulations of electron motion in global magnetospheric fields. However, there are still significant outstanding challenges that remain regarding how to incorporate the long term effects of wave-particle interactions in test-particle simulations. In this paper, we describe in details two approaches that incorporate electron scattering in test particle simulations: stochastic differential equation approach and the mapping technique. Both approaches assume that wave-particle interactions can be described as a probabilistic process that changes electron energy, pitch-angle, and thus modifies the test particle dynamics. To compare these approaches, we model electron resonant interactions with field-aligned whistler-mode waves in dipole magnetic fields. This comparison shows advantages of the mapping technique in simulating the nonlinear resonant effects, but also underlines that more significant computational resources are needed for this technique in comparison with the stochastic differential equation approach. We further discuss applications of both approaches in improving existing models of energetic electron dynamics.

1 Introduction

One key element in substorm magnetosphere dynamics is plasma sheet injections into the inner magnetosphere (Baker et al., 1996; Birn et al., 1997; Nakamura et al., 2002; Angelopoulos et al., 2008; Gabrielse et al., 2012). Simulations of the energetic particle transport during such injections require modeling of large-scale magnetic field reconfiguration and particle responses to a wide variety of kinetic processes, such as wave-particle resonant interactions and scattering by the magnetic field gradients. The most advanced approach here is the test-particle modeling in electromagnetic fields of global (magnetosphere) MHD or hybrid simulations (e.g., Perroomian & El-Alaoui, 2008; Birn et al., 2004; Ashour-Abdalla et al., 2005). This approach can well resolve meso-scale electromagnetic field structures, like plasma injection fronts (e.g., Wiltberger et al., 2015), and can reproduce main details of energetic electron (Ashour-Abdalla et al., 2011; Liang et al., 2014; Pan et al., 2014; Birn et al., 2014, 2022; Zhou et al., 2018; Sorathia et al., 2018) and ion (Perroomian & Zelenyi, 2001; Birn et al., 2015, 2017; Ukhorskiy et al., 2018) transport and energization. MHD simulations with sufficiently high spatial resolution can reproduce magnetic field gradients around injection (dipolarization) fronts and magnetotail current sheets, and thus may adequately describe electron scattering by magnetic field-line curvatures (Eshetu et al., 2018, 2019; Desai et al., 2021). Moreover, global hybrid simulations resolving ion kinetics (Lin et al., 2014; Lu et al., 2016, 2017) can reproduce kinetic Alfvén wave dynamics (Lin et al., 2017; Cheng et al., 2020), making it possible to simulate the plasma sheet electron and ion acceleration by field-aligned transient electric fields (i.e., the main ion-kinetic feature of plasma injections (Chaston et al., 2012, 2015; Ergun et al., 2015; Hull et al., 2020)). Although schemes to include electron kinetics into global simulations are under development and verification (e.g., Chen et al., 2017; Walker et al., 2018; Alho et al., 2022), neither existing global MHD nor global hybrid simulations can resolve electron-scale waves, and thus cannot describe the wide range of electron resonant phenomena associated with plasma injections (see discussion in Mozer et al., 2015; Malaspina et al., 2018; Ukhorskiy et al., 2022; Artemyev, Neishtadt, & Angelopoulos, 2022).

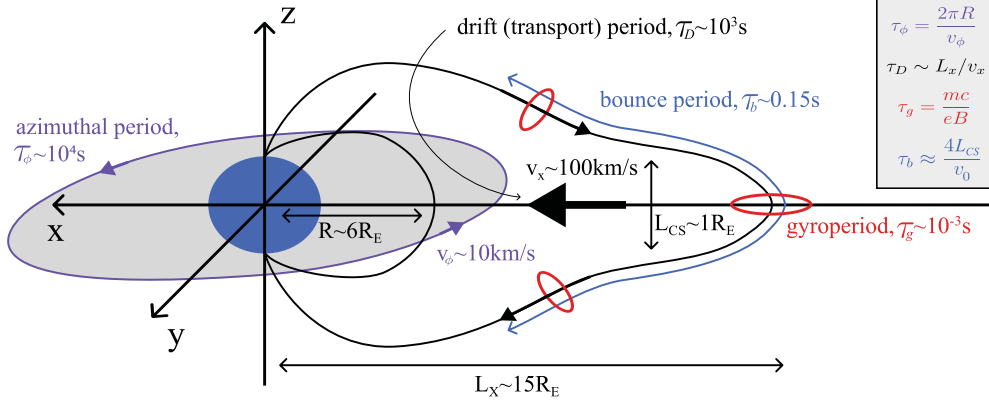


Figure 1. Schematic of four main time-scales of electron motion during plasma sheet injections into the inner magnetosphere. Four main time-scales: electron gyroperiod $\tau_g = mc/eB \sim 10^{-3}$ s (with a typical magnetic field magnitude of $B = 10$ nT), electron bounce period $\tau_b \sim 4L_{CS}/v_0 \sim 10^{-1}$ s (with the current sheet thickness being $L_{CS} = 1R_E$ and v_0 being the thermal velocity of 100 keV electrons), $\tau_D \sim L_x/v_x \sim 10^3$ s is electron transport time by plasma flows ($v_x \sim 100$ km/s), $\tau_\phi = 2\pi R/v_\phi \sim 10^4$ s is the electron azimuthal drift period (the radial distance of the injection region is assumed to be $R = 6R_E$ and the electron energy is 100 keV).

76 It is actually rather challenging to incorporate wave-particle resonant interactions
 77 into test particle simulations of plasma injections, because of their vastly different timescales.
 78 Figure 1 shows a schematic of the four main types of electron motions: cyclotron rota-
 79 tion with the timescale of $\tau_g \sim 10^{-3}$ s, bounce motion between magnetic mirrors with
 80 the time scale of $\tau_b \sim 10^{-1}$ s (for 100keV electrons), earthward transport with the time-
 81 scale of $\tau_D \sim 10^3$ s, and azimuthal motion around the Earth with the time-scale of $\tau_\phi \sim$
 82 10^4 s (for 100keV electrons). The ratio of the timescales of the fastest and the slowest
 83 motions can be $\tau_\phi/\tau_g \sim 10^7$ or $\tau_D/\tau_g \sim 10^6$. Individual wave-particle interaction, e.g.,
 84 via cyclotron resonance, occurs over $\sim \tau_g$, and thus such interactions may not be directly
 85 incorporated into test particle simulations of the plasma injections, which occur over \sim
 86 τ_D .

87 A possible solution of this problem was proposed by Elkington et al. (2018, 2019),
 88 who suggest that wave-particle interactions can be incorporated as stochastic perturba-
 89 tions of electron test orbits (see also Michael et al., 2023; Chan et al., 2023). The sim-
 90 plified version of this approach combines test particle equations of motion and contin-
 91 uous stochastic differential equations (SDEs), which are characteristics for the Fokker-
 92 Plank diffusion equation (Tao et al., 2008; Zheng et al., 2014). The Fokker-Plank equa-
 93 tion is used in the quasi-linear theory (Vedenov et al., 1962; Andronov & Trakhtengerts,
 94 1964; Kennel & Engelmann, 1966), which describes electron scattering by low intensity
 95 waves (see Karpman, 1974; Albert, 2001, 2010; Frantsuzov et al., 2023, for discussions
 96 on the wave intensity limitations in inhomogeneous background magnetic field, like in
 97 the Earth’s magnetosphere). Applications of the quasi-linear diffusion approximation has
 98 been well developed for the Earth’s magnetosphere (Lyons & Williams, 1984; Schulz &
 99 Lanzerotti, 1974), and the main parameters of this approximation, the diffusion rates,
 100 are widely evaluated and used for the observed wave characteristics in the magnetotail
 101 (e.g., Panov et al., 2013; Zhang et al., 2015; Ni et al., 2011, 2012, 2016) and inner mag-
 102 netosphere (see reviews by Shprits et al., 2008; Artemyev, Agapitov, et al., 2016; Li &
 103 Hudson, 2019; Thorne et al., 2021, and references therein).

104 The main limitation of the quasi-linear models is the requirement of a small wave
 105 intensity. A significant portion of most intense electromagnetic whistler-mode waves (Wilson
 106 et al., 2011; Tyler et al., 2019; Zhang et al., 2018, 2019), electrostatic whistler-mode waves
 107 (Cully et al., 2008; C. Cattell et al., 2008; C. A. Cattell et al., 2015; Agapitov et al., 2014,
 108 2015), and electromagnetic ion cyclotron waves (e.g., Wang et al., 2017; Tonoian et al.,
 109 2022) can exceed their threshold amplitudes and likely resonate with electrons nonlin-
 110 early. Such nonlinear resonant interactions include effects of phase bunching and phase
 111 trapping (see Nunn, 1971; Karpman et al., 1975; Inan & Bell, 1977; Solovév & Shklyar,
 112 1986; Albert, 1993; Itin et al., 2000), which can significantly modify the characteristics
 113 of wave-particle interactions (see reviews by Shklyar & Matsumoto, 2009; Albert et al.,
 114 2013; Artemyev, Neishtadt, Vainchtein, et al., 2018, and references therein). Nonlinear
 115 effects can be incorporated into the kinetic equation (i.e., modified Fokker-Plank equa-
 116 tion) for electron distribution functions (e.g., Omura et al., 2015; Hsieh & Omura, 2017;
 117 Artemyev, Neishtadt, et al., 2016; Artemyev, Neishtadt, Vasiliev, & Mourenas, 2018),
 118 but the corresponding characteristic equations will be different from those in the SDE
 119 approach. The main difference is the probability distribution function of pitch-angle/energy
 120 jumps due to resonant interactions: nonlinear effects cannot be described by Gaussian
 121 probability distributions that are often adopted to model the diffusive scattering within
 122 SDE approach. Recently, an alternative to SDE approach was proposed in (Artemyev,
 123 Neishtadt, Vainchtein, et al., 2018; Zheng et al., 2019), where a non-Gaussian probabili-
 124 ty distribution of pitch-angle/energy jumps has been incorporated into equations of mo-
 125 tion for resonant electrons. This approach resembles the generalization of the classical
 126 mapping technique (e.g., Chirikov, 1979; Zaslavskii et al., 1989; Khazanov et al., 2014)
 127 to systems with a finite probability of very large pitch-angle/energy jumps due to phase
 128 trapping (see Artemyev et al., 2020). The mapping technique can reproduce many ob-
 129 served effects of nonlinear wave-particle interactions (e.g., Zhang et al., 2022; Artemyev,
 130 Zhang, et al., 2022) and in principle can be incorporated into test-particle codes (Artemyev,
 131 Neishtadt, & Angelopoulos, 2022). Therefore, both SDE and mapping technique can be
 132 used to include wave-particle resonant interactions into models of energetic electron dy-
 133 namics in global electromagnetic fields provided by MHD/hybrid simulations. The map-
 134 ping technique should generalize the SDE approach, but it is yet to be investigated whether
 135 the mapping equations can describe diffusive scattering and nonlinear resonant effects
 136 with the same accuracy level. The mapping technique usually adopts analytical equa-
 137 tions to model pitch-angle/energy jumps, which do not include diffusive scattering (Artemyev
 138 et al., 2020). Recently Lukin et al. (2021) has generalized the mapping technique for the
 139 entire probability distribution function of pitch-angle/energy jumps, but it remains to
 140 be verified for electron cyclotron resonances with whistler-mode waves.

141 In this paper, we combine two approaches from (Artemyev et al., 2020; Lukin et
 142 al., 2021) to construct the mapping technique that operates with the probability distri-
 143 bution function of pitch-angle/energy jumps for electron cyclotron resonances with whistler-
 144 mode waves. We also compare results from this newly developed mapping technique to
 145 those from the SDE approach, which operates by a single characteristic of such proba-
 146 bility distribution functions of pitch-angle/energy jumps – distribution variance, which
 147 dictates the diffusion rate. We examine two wave intensities: small intensity for the dif-
 148 fusive interaction, when SDE and mapping are expected to provide the same results, and
 149 large intensity with nonlinear wave-particle interactions, when the mapping technique
 150 is validated by full test particle simulations. Therefore, this paper demonstrates the ap-
 151 proaches to incorporate the long-term effects of quasi-linear and nonlinear wave-particle
 152 interactions into global scale test-particle models.

153 The rest of the paper starts with introducing basic characteristics of electron reso-
 154 nant interactions with whistler-mode waves in the dipole magnetic field (see Sect. 2).
 155 Then in Sect. 3, we introduce the main property of electron ensemble dynamics – the
 156 probability distribution function of energy/pitch-angle jumps in a single resonance. This
 157 distribution is used to introduce SDE approach (in Sect. 4) and mapping technique (in

158 Sect. 5). Results obtained from SDE and mapping approaches are compared with test
 159 particle simulations in Sect. 6. Finally, we discuss the applicability of both methods and
 160 briefly summarize our conclusions in Sect. 7.

161 2 Resonant electron interactions with whistler-mode waves

We examine the interaction of relativistic electrons (rest mass m_e , charge $-e$, speed of light c) with field-aligned whistler-mode waves moving in the dipole magnetic field $B_0(\lambda)$, where λ is the magnetic latitude. Electron dynamics in such a system can be described by the following Hamiltonian (Albert, 1993; Vainchtein et al., 2018):

$$H = \sqrt{m_e^2 c^4 + c^2 p_z^2 + 2I_x \Omega_0(\lambda) m_e c^2} + \sqrt{\frac{2I_x \Omega_0(\lambda)}{m_e c^2} \frac{eB_w(\lambda)}{k(\lambda)}} \cos(\phi + \psi) \quad (1)$$

where (z, p_z) and (ψ, I_x) are two pairs of conjugate variables, the field-aligned coordinate and moment, and electron gyrophase and magnetic moment. The electron cyclotron frequency $\Omega_0(\lambda)$ is defined as

$$\Omega_0(\lambda) = \Omega_{eq} \frac{\sqrt{1 + 3 \sin^2(\lambda)}}{\cos^6(\lambda)} \quad (2)$$

where $\Omega_{eq} = eB_0(0)/m_e c$ is the equatorial cyclotron frequency determined by the radial distance from the Earth to the equatorial crossing of the magnetic field line, i.e., L -shell. In the dipole field, the magnetic latitude, λ , is related to the field-aligned coordinate z as:

$$\frac{dz}{d\lambda} = R_E L \sqrt{1 + 3 \sin^2(\lambda)} \cos(\lambda) \quad (3)$$

162 where R_E is the Earth radius. Note that in Hamiltonian (1), the magnetic moment, I_x ,
 163 conjugate to the gyrophase, ψ , should be normalized in such a way that $I_x \Omega_0$ has a di-
 164 mension of energy.

The second term in Hamiltonian (1) describes the wave contribution to the electron dynamics. The wave amplitude B_w is modeled as:

$$B_w(\lambda) = \varepsilon B_0(0) f(\lambda), \quad f(\lambda) = \begin{cases} \tanh(c_\lambda(\lambda - \lambda_0)), & \lambda \geq \lambda_0 \\ 0, & \lambda < \lambda_0 \end{cases} \quad (4)$$

165 where ε parameter controls $B_w/B_0(0)$, function $f(\lambda)$ determines the wave latitudinal pro-
 166 file that agrees with the wave generation around the equator, i.e., wave amplitude growth
 167 within the generation/amplification region $\Delta\lambda$, which is controlled by the value of c_λ :
 168 $\Delta\lambda \sim 1/c_\lambda$, followed by saturation (see typical $B_w(\lambda)$ profiles from statistical models
 169 in Agapitov et al., 2013, 2018). Note that we assume the waves only exist in one hemi-
 170 sphere, $z > 0$, because the wave field and wave propagation equations are symmetric
 171 in two hemispheres ($z \rightarrow -z$ does not change the system equation). This assumption
 172 allows us to simplify the calculations, because in this case the wave-particle interaction
 173 occurs only during a quarter of the bounce period when the particles move northward
 174 from the equator.

The wave phase is given by equation $\dot{\phi} = k(\lambda)\dot{z} - \omega$, where the wave frequency ω is constant and the wave number $k(\lambda)$ is determined by the cold plasma dispersion relation (Stix, 1962):

$$ck(\lambda) = \Omega_{pe}(\lambda) \left(\frac{\Omega_0(\lambda)}{\omega} - 1 \right)^{-1/2} \quad (5)$$

The plasma frequency Ω_{pe} is given by the empirical function (Denton et al., 2006)

$$\Omega_{pe}(\lambda) = \Omega_{pe,eq} \cos^{-5/2}(\lambda) \quad (6)$$

175 with the equatorial values $\Omega_{pe,eq}$ given by the empirical function (Sheeley et al., 2001)
 176 $\Omega_{pe,eq}/\Omega_0(0) \approx L$.

Wave phase ϕ linearly depends on time, $\partial\phi/\partial t = \omega = const$, and thus Hamiltonian (1) has an integral of motion h (see, e.g., review by Shklyar & Matsumoto, 2009):

$$h = \sqrt{m_e^2 c^4 + c^2 p_z^2 + 2I_x \Omega_0(\lambda) m_e c^2} - \omega I_x \quad (7)$$

In the absence of wave perturbation, $B_w = 0$, particle energy $m_e c^2(\gamma - 1)$ and equatorial pitch-angle $\alpha_{eq} = \arcsin(2I_x \Omega_0(0)/m_e c^2(\gamma^2 - 1))$ are conserved; here

$$\gamma = \sqrt{1 + \left(\frac{p_z}{m_e c}\right)^2 + \frac{2I_x \Omega_0(\lambda)}{m_e c^2}}$$

177 is the gamma factor. Wave-particle resonant interactions can change particle's energy
 178 and equatorial pitch-angle (change I_x), but due to the conservation of $h(\gamma, I_x)$ waves move
 179 electrons along a specific curve in the energy, pitch-angle space. Therefore, for a fixed
 180 value of h (i.e., when all particles have the same initial h), the wave-particle resonant
 181 interaction becomes a 1D problem, and we may just examine energy changes, whereas
 182 changes of equatorial pitch-angle (or I_x) can be determined from $h = const$.

To obtain basic characteristics of wave-particle resonant interactions, we integrate electron equations of motion (Hamiltonian equations) with 4th order Runge-Kutta scheme with an adaptive time step (1/20 of the local electron gyroperiod). Throughout the rest of the paper, we use the following dimensionless variables:

$$H \rightarrow H m_e c^2, \quad p_z = p m_e c, \quad I_x \rightarrow I_x \frac{m_e c^2}{\Omega_{eq}}, \quad t \rightarrow t/\Omega_0(0), \quad z \rightarrow z \frac{c}{\Omega_{eq}} \quad (8)$$

Thus, the dimensionless Hamiltonian and integral of motion h take the following forms

$$H = \sqrt{1 + p_z^2 + 2I_x \Omega_0(\lambda)} + \sqrt{2I_x \Omega_0(\lambda)} \frac{\varepsilon f(\lambda)}{k(\lambda)} \cos(\phi + \psi) \quad (9)$$

$$h = \sqrt{1 + p_z^2 + 2I_x \Omega_0(\lambda)} - \omega I_x \quad (10)$$

183 where $\Omega(\lambda) \rightarrow \Omega(\lambda)\Omega_0(0)$. To demonstrate the result, we use the following parameters
 184 throughout: $h = 3/2$, $\lambda_0 = 5^\circ$, and $c_\lambda = 180/\pi$. Each integration starts from the
 185 equatorial plane ($\lambda = z = 0$) and stops there after N_{res} resonant wave-particle inter-
 186 actions. Since waves only exist in the northern hemisphere at a fixed frequency, and prop-
 187 agate along magnetic field lines (i.e., resonate with electrons through the first-order cy-
 188 clotron resonance only), each electron bounce period corresponds to one wave-particle
 189 interaction on a time scale of a quarter bounce period. Effects of multiple resonances within
 190 one bounce period (e.g., due to oblique wave propagation or wave frequency variation
 191 with time, see Shklyar & Matsumoto, 2009; Artemyev et al., 2021; Hsieh & Omura, 2023)
 192 can be incorporated into this approach by including additional wave terms into Eq. (9).

193 Depending on the wave magnitude, there are two possible regimes of wave-particle
 194 resonant interactions: diffusive scattering and nonlinear resonant interactions. Figure
 195 2 shows the profiles of electron energy evolution with time for both regimes (all electrons
 196 have the same initial energy and pitch-angle, but random initial gyrophases). For small
 197 wave amplitudes: (a) the interaction is linear and energy evolution with time is stochas-
 198 tic. After each wave-particle interaction, the electron energy undergoes a small (com-
 199 pared to electron initial energy) positive or negative jump with almost equal probabili-
 200 ties to increase or decrease the energy. This process (with normalized diffusion coeffi-
 201 cients) can be approximated by the Wiener stochastic process, i.e., the evolution of elec-
 202 tron distribution function can be described by the Fokker-Planck equation or, equivalently,
 203 by stochastic differential equations. For large wave amplitudes: (b) electron dyn-
 204 amics cannot be described with the diffusive approach. Most of the time, particle en-
 205 ergy undergoes small negative jumps (but positive jumps are also possible, see Albert
 206 et al., 2022) and there is a nonzero probability of large positive jumps, caused by par-
 207 ticle phase trapping.

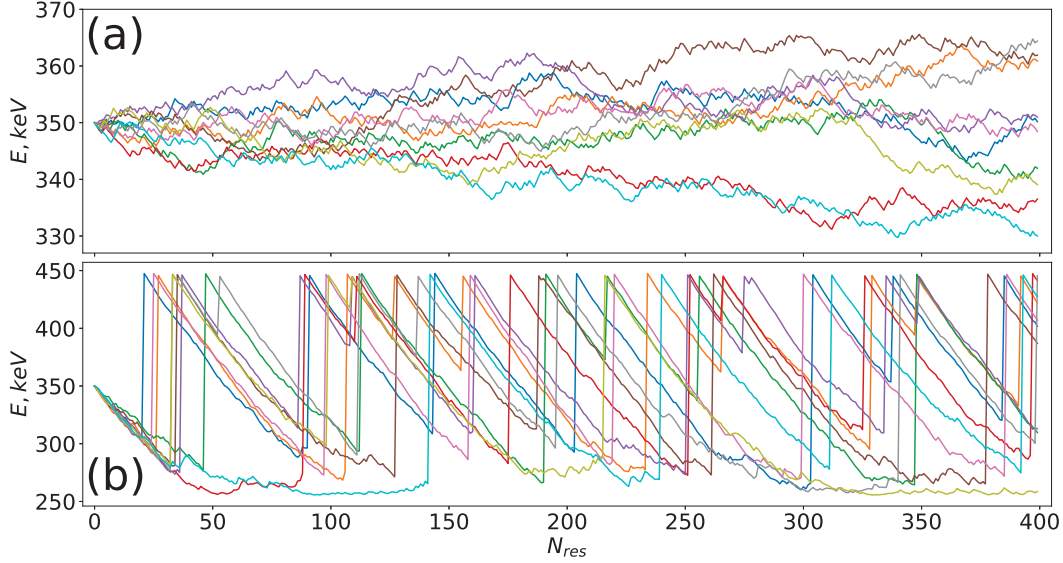


Figure 2. Examples of electron energy dynamics for the system with diffusive scattering (a) and nonlinear resonant effects (b). All electrons have the same initial energy and pitch-angle, but random initial gyrophases. Energy is plotted versus number of resonant interactions (i.e., number of electron bounce periods).

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3 Probability distributions of energy jumps

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To quantify variations of electron energy due to wave-particle resonant interactions, we use test particle simulations and evaluate the distributions of energy jumps as a function of initial energy E_0 , $\Delta E(E_0)$. Figure 3 shows such distributions for several initial electron energies and two different magnitudes of the wave field. For each histogram, we integrated trajectories of $N_p = 32768$ test particles with random initial gyrophases and the same initial energy and h values. Each integration includes one bounce period (i.e., a single resonant interaction). For small wave amplitudes (panels (a-d)), electron energy jumps are small and randomly (but symmetrically) distributed around zero. For sufficiently high wave amplitudes (panels (e-h)), there appear nonlinear resonant effects: most of the electrons lose their energy due to the phase bunching with $\Delta E < 0$, while a small population of phase trapped particles (panels (g-h)) gain a significant portion of energy with $\Delta E > 0$ comparable to their initial energy, E_0 . Changes of the electron pitch-angle (and I_x) are directly determined by the energy changes due to the conservation of the integral of motion (7). Therefore, the probability distribution of $\Delta E(E_0)$ fully characterizes the evolution of electron distribution function and can be used as a basic input for the SDE approach or mapping technique. Note that for intense, but very low-coherent waves, the probability distribution function of $\Delta E(E_0)$ will be symmetric relative to $\Delta E(E_0) = 0$, and thus will largely resemble distributions from panels (a-d) (see Zhang et al., 2020; An et al., 2022; Gan et al., 2022; Frantsuzov et al., 2023).

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4 Stochastic differential equations

In the limit of small wave field amplitudes, the wave-particle resonant interaction is diffusive (e.g., Kennel & Engelmann, 1966; Lyons & Williams, 1984; Schulz & Lanzetta, 1974; Allanson et al., 2022, and references therein). Note that in inhomogeneous magnetic field, these interactions are diffusive even for monochromatic waves (see Albert, 2010; Shklyar, 2021). Therefore, for such systems one can use the Fokker-Planck

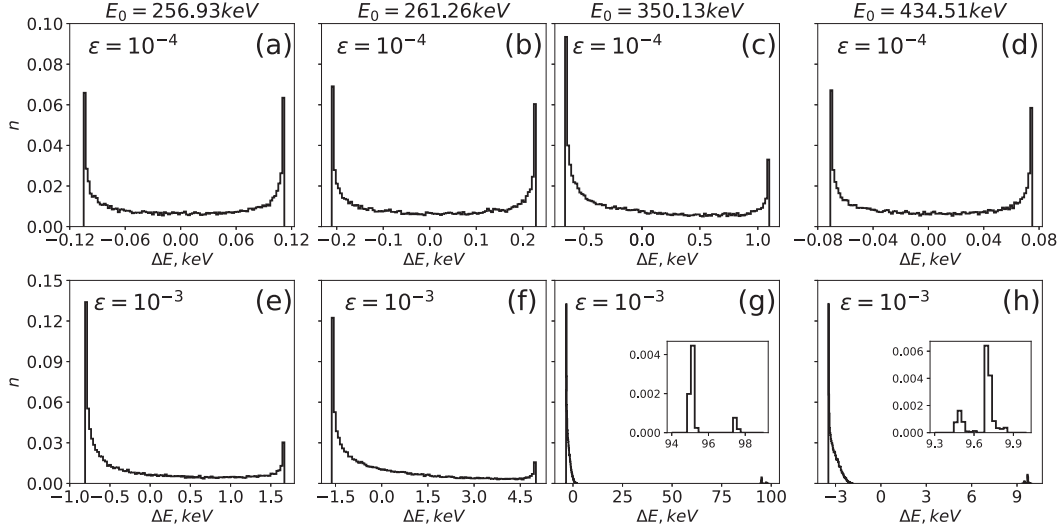


Figure 3. Examples of ΔE probability distributions for systems without nonlinear resonant effects (a-d) and with nonlinear resonant effects (e-h). In panels (e,f), we show simulation results for E_0 without trapping: a small population of $\Delta E > 0$ is due to the positive phase bunching effect (see Albert et al., 2022). In panels (g,h), we show simulation results with phase trapping; the inserted panels show the expanded view of the population with large $\Delta E > 0$.

equation to describe the evolution of the electron distribution function $f(\boldsymbol{\chi})$ of the velocity vector $\boldsymbol{\chi}$:

$$\frac{\partial f(\boldsymbol{\chi}, t)}{\partial t} = - \sum_{i=1}^N \frac{\partial}{\partial \chi_i} (\mu_i(\boldsymbol{\chi}, t) f(\boldsymbol{\chi}, t)) + \sum_{i=1}^N \sum_{j=1}^N \frac{\partial^2}{\partial \chi_i \partial \chi_j} (D_{ij}(\boldsymbol{\chi}, t) f(\boldsymbol{\chi}, t)) \quad (11)$$

where t is time, $\boldsymbol{\mu}(\boldsymbol{\chi}(t), t)$ is an N -dimension vector of the drift coefficient, \mathbf{D} is an $N \times N$ matrix of diffusion coefficients (Lyons & Williams, 1984; Schulz & Lanzerotti, 1974; Albert, 2018). Instead of solving the Fokker-Planck equation, we can use the corresponding Ito stochastic differential equations to integrate trajectories of *quasi-particles* (Tao et al., 2008; Zheng et al., 2014):

$$\boldsymbol{\chi}(t + \Delta t) = \boldsymbol{\chi}(t) + \boldsymbol{\mu}(\boldsymbol{\chi}(t), t) \Delta t + \boldsymbol{\sigma}(\boldsymbol{\chi}(t), t) d\mathbf{W}_t \quad (12)$$

where Δt is the time step over which we calculate the change of $\boldsymbol{\chi}$, $\boldsymbol{\sigma}(\boldsymbol{\chi}(t), t)$ is an $N \times N$ matrix related to the diffusion coefficients written in such a way that $\mathbf{D} = \frac{1}{2} \boldsymbol{\sigma} \boldsymbol{\sigma}^T$, \mathbf{W}_t is an N -dimension standard Wiener process; $d\mathbf{W}_t = \sqrt{\Delta t} \mathbf{N}$, where \mathbf{N} is a vector of standard normal random values, $N_i \sim N(0, 1)$.

The term *quasi-particles* means that we do not directly integrate the equations of motion, but treat the change of $\boldsymbol{\chi}$ as a stochastic process and approximate it by the equation (12). As a result, two *quasi-particles* having equal initial conditions $\boldsymbol{\chi}_0$ may have different trajectories $\boldsymbol{\chi}(t)$ (in the numerical integration of the equation (12), one can fix the seed of the pseudo random generator to preserve the sequence of random numbers and make the results repeatable). We will examine electron distributions in the energy space, and thus the Ito equation (12) can be rewritten in the following form

$$E(t + \Delta t) = E(t) + \mu_E(E(t)) \Delta t + \sqrt{2D_E(E(t))} dW_t \quad (13)$$

This equation describes the energy evolution for fixed h given by Eq. (10), i.e., for a monochromatic wave we may reduce the energy, pitch-angle evolution to the energy-only evolution and calculate the associated pitch-angle changes from h conservation. For generic

236 wave spectra, there are three diffusion rates (energy, pitch-angle, and mixed energy-pitch-
 237 angle), and thus we would need to solve a system of equations for the energy and pitch-
 238 angle evolution (see detail in Tao et al., 2008). Note that if we rewrite the Fokker-Planck
 239 equation in terms of energy, additional coefficients of variable transformation from ve-
 240 locity (momentum) to energy and pitch-angle (Lamé coefficients) should be added (e.g.,
 241 Glauert & Horne, 2005), but the Ito equation will still have the same form.

The main limitation of this approach is the requirement on small energy changes ΔE for each wave-particle interaction, i.e., it is only applicable for systems without non-linear effects of phase trapping. Equation (13) includes the drift $\mu_E(E_0)$ and bounce averaged diffusion $D_E(E_0)$ coefficients, which can be estimated using the distribution of energy jumps $\Delta E(E_0)$:

$$D_E(E_0) = \frac{1}{2\tau_b(E_0)} \left(\frac{1}{N_p} \sum_{i=1}^{N_p} \Delta(E_i)^2(E_0) - \left(\frac{1}{N_p} \sum_{i=1}^{N_p} \Delta E_i(E_0) \right)^2 \right) \quad (14)$$

For the drift term, we use

$$\mu_E(E_0) = \frac{dD_E(E_0)}{dE_0} \quad (15)$$

242 which keeps the divergence-free form of the Fokker-Planck equation (Lichtenberg & Lieber-
 243 man, 1983; Sinitzyn et al., 2011; Lemons, 2012; Zheng et al., 2019; Allanson et al., 2022).
 244 Note that using such estimates of drift and diffusion coefficients, we implicitly assume
 245 that the $\Delta E(E_0)$ distributions are symmetric relative to the mean ΔE value. Figures
 246 4(a-b) show the numerically obtained ΔE distributions (black) and their fittings to sym-
 247 metric Gaussian distributions (this fitting outputs the variance $\sim D_E(E_0)$). However,
 248 this assumption may be violated even in the case of small wave amplitudes (see Figure
 249 3(c)), when we can over- or underestimate the drift and diffusion coefficients. We will
 250 discuss the corresponding uncertainties in Section 6. Figure 4(c) shows the drift and dif-
 251 fusion coefficients as a function of the electron initial energy E_0 .

Each electron bounce period corresponds to a single wave-particle interaction (as commented above), and thus the time step of integration Δt in the equation 13 should be set equal to $\tau_b(E_0)$. The main advantage of the SDE approach (and also of the mapping technique, which will be discussed in the Section 5), in comparison with the test particle approach, is the significant reduction of computational time for long-term simulations (see also discussions in Lukin et al., 2021): we integrate trajectories of quasi-particles with a time step equal to the bounce period, considering only the effects of wave-particle interactions and do not trace particles during adiabatic paths of their motion. So at each step of integration, we recalculate the electron energy as

$$E_{i+1} = E_i + \mu_E(E_i)\tau_b(E_i) + \sqrt{2D_E(E_i)\tau_b(E_i)}N_i$$

252 where $N_i \sim N(0, 1)$ is the standard normal random number.

253 Panels (d,e) in Figure 4 show examples of the electron energy profiles calculated
 254 by direct integration of the Hamiltonian equations or using the SDE approach. Particles
 255 having equal initial conditions will have different trajectories from these two approaches,
 256 due to randomness of resonant interactions, but statistical properties of the evolution
 257 of the electron ensemble should be the same. We have verified this property by a set of
 258 simulated evolution of the electron ensembles (not shown).

259 5 Mapping technique

260 For high wave amplitudes, the SDE approach is no longer applicable, because of
 261 the nonlinear effects of wave-particle interactions. In this case, the Wiener stochastic pro-
 262 cess cannot describe the evolution of electron energy as there is a finite probability of

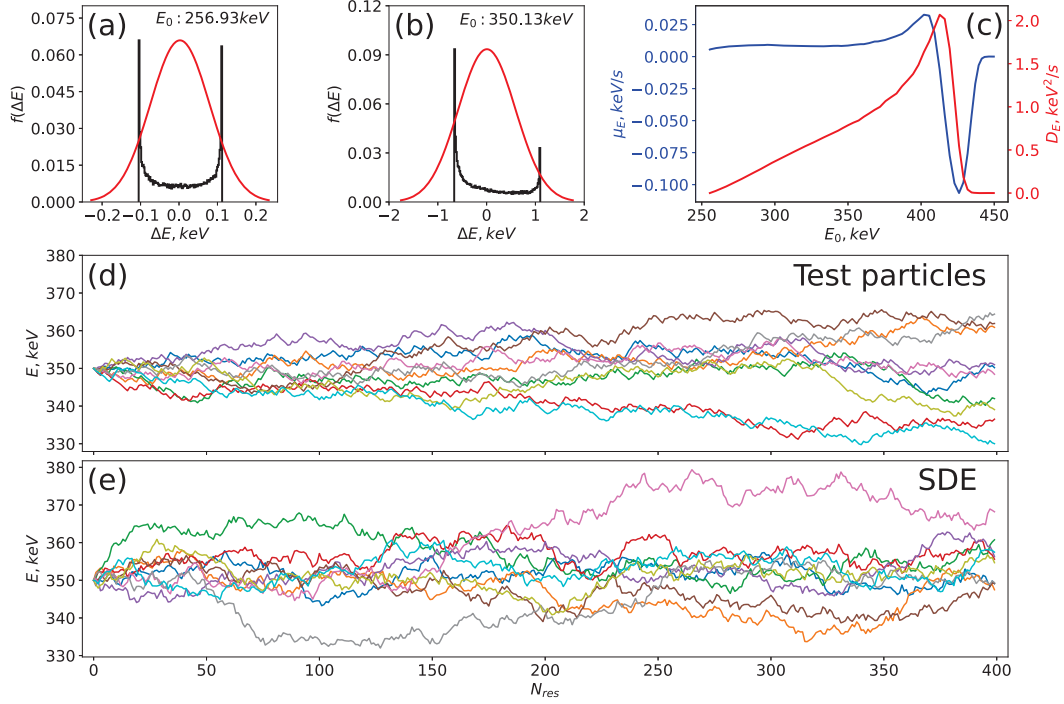


Figure 4. Panels (a,b) show ΔE distributions (in black), obtained after a single wave-particle interaction, and Gaussian distributions with the same variance (in red). Panel (c) shows the energy diffusion rate and drift term versus initial energy. Panels (d,e) show examples of ten trajectories of long-term electron energy dynamics (~ 300 resonances) for a system without nonlinear resonant effects.

263 large energy jumps caused by the phase trapping. For such systems, the Ito SDE can
 264 be generalized by introducing a new stochastic process accounting for the phase trap-
 265 ping and phase bunching. This process can be represented by a series of mapping func-
 266 tions depending on electron energy (the changes of pitch-angle are related to energy changes
 267 through the integral of motion (7)).

268 Figure 5(a) shows the distribution of energy jumps for the wave amplitude $\varepsilon = 10^{-3}$
 269 and initial electron energy $E_0 = 257$ keV. For this particular energy, there is no elec-
 270 tron phase trapping: the resonant latitude for this energy and h (which determines the
 271 equatorial pitch-angle) corresponds to zero probability of electron trapping, because wave
 272 intensity increases slower along the resonant trajectory than the background magnetic
 273 field inhomogeneity (see detailed equations determining the trapping probability in the
 274 Appendix of Vainchtein et al., 2018, and in references therein). Hence the procedure to
 275 construct the mapping function is straightforward: using this distribution, one can calcu-
 276 late the cumulative function of energy jumps $F(\Delta E, E_0) \in [0, 1]$ (right vertical axis
 277 in Figure 5(a)) and then use the inverse function $\Delta E(U) = F^{-1}(U, E_0), U \in [0, 1]$ as
 278 the generator of electron energy jumps for each wave-particle interaction.

Figure 5(b) shows the distribution of energy jumps for an initial electron energy
 of $E_0 = 350$ keV. There is a small probability of electron phase trapping in this case,
 thus we can divide the ΔE distribution into *bunching* ($|\Delta E|/E_0 \ll 1$) and *trapping* ($\Delta E/E_0 \sim$
 1) parts, and then calculate corresponding cumulative functions: $F_b(\Delta E, E_0)$ and $F_t(\Delta E, E_0)$.
 Note that this separation of ΔE distribution into two parts is not necessary, but it sim-
 plifies the calculations by allowing a linear interpolation for the cumulative function. There
 is a gap of ΔE between two parts of ΔE distribution, i.e., for the case shown in Figure
 5(b) the electron energy cannot change by the value between ~ 4 and ~ 95 keV (pro-
 hibited values of ΔE in this case). If we would not separate the ΔE distribution into
 two parts, this ΔE gap will require a separate treatment. Using *trapping* ΔE distribu-
 tions, we may estimate the probability of electron trapping $p_{trap}(E_0)$ (shown in Figure
 5(c)) as a function of the initial electron energy:

$$p_{trap}(E_0) = \frac{N(\Delta E(E_0)/E_0 \sim 1)}{N_p} \quad (16)$$

279 where N_p is the total number of particles used to construct the $\Delta E(E_0)$ distribution and
 280 $N(\Delta E(E_0)/E_0 \sim 1)$ is the number of trapped particles.

We use the subscript *b* for the bunching part of the $\Delta E(E_0)$ distribution and also
 for the entire $\Delta E(E_0)$ distribution if the probability of trapping is equal to zero. Then
 the generalization of the Ito SDE is straightforward: we can replace the Wiener process
 in equation (13) with the constructed mapping functions. At each integration step, we
 generate two uniform random numbers $u_1, u_2 \in U(0, 1)$ and recalculate electron energy
 as

$$E_{i+1} = E_i + \begin{cases} F_t^{-1}(u_2, E_i), & u_1 \leq p_{trap}(E_i) \\ F_b^{-1}(u_2, E_i), & u_1 > p_{trap}(E_i) \end{cases} \quad (17)$$

281 Note that the gap between ΔE distributions due to phase trapping and bunching
 282 (see Figure 5(b)) depends on system parameters (mostly on the wave field model) and,
 283 e.g., for the system with small wave-packets, the energy of phase-trapped electrons could
 284 not be far away from the initial energy (e.g., Omura et al., 2015). In this case, there is
 285 no need to divide the distributions into *bunching* and *trapping* parts and thus only one
 286 cumulative function, $F(\Delta E, E_0)$, needs to be constructed. In this case, the computation
 287 scheme slightly simplifies: at each iteration one needs to calculate a single, uniform ran-
 288 dom number $u \in U(0, 1)$ and the energy change can be calculated in the same man-
 289 ner: $E_{i+1} = E_i + F^{-1}(u, E_i)$.

290 Figures 5(d,e) show several profiles of electron energy as a function of the number
 291 of wave-particle resonances, calculated by integration of the test particle trajectories and

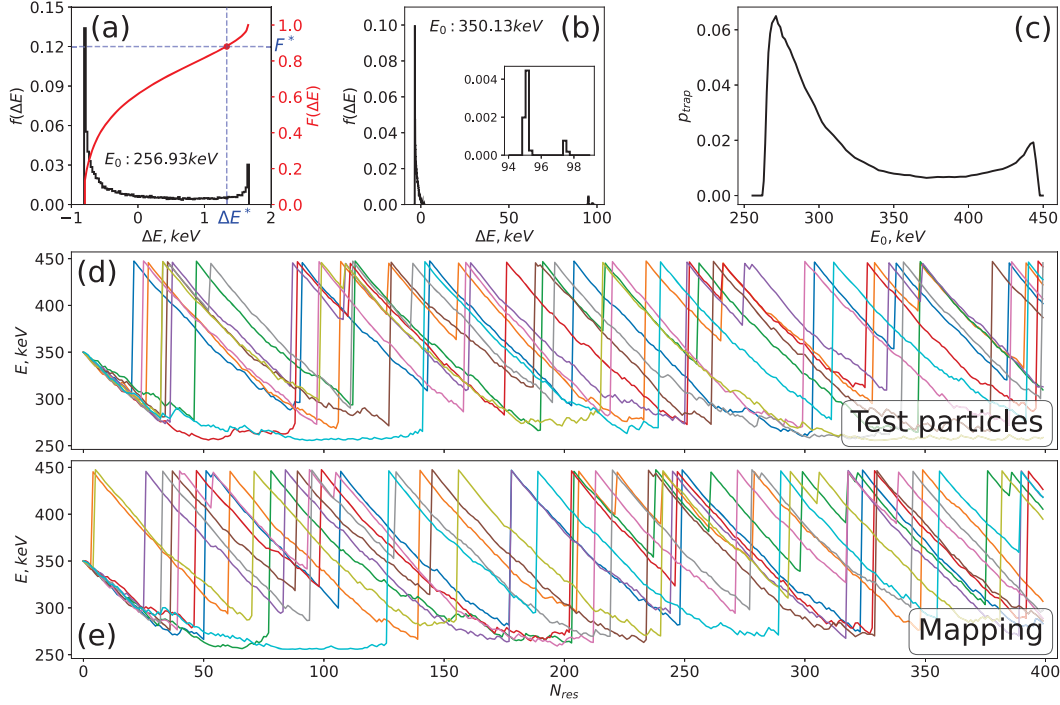


Figure 5. Panels (a,b) show ΔE distributions and the corresponding cumulative probability distribution function (only on panel (a)). Panel (c) shows the probability of electron phase trapping versus the initial energy. Panels (d,e) show examples of ten trajectories of long-term electron energy dynamics (~ 300 resonances) for a system with nonlinear resonant effects.

292 using the mapping technique, respectively. Both approaches show very similar electron
 293 energy dynamics: long-duration electron drift to smaller energy due to the phase bunch-
 294 ing and rare large jumps to higher energy due to the phase trapping. Note that for both
 295 SDE and mapping techniques, we still need to use test particles to calculate the initial
 296 distributions of $\Delta E(E_0)$, but having these distributions (requiring statistics of single wave-
 297 particle interactions), we can use the simplified integration scheme with a time step equal
 298 to the bounce period.

299 6 Method validation

300 For small wave amplitudes, we can apply all three methods (test particles, SDE,
 301 and mapping) to simulate the evolution of the electron distribution function, and these
 302 three methods are expected to give the same results. The test particles approach is ex-
 303 pected to be more precise, because it does not rely on the constructed ΔE distribution
 304 and is based on the full set of equations of motion. The main advantage of SDE and map-
 305 ping techniques is their computational efficiency in long-term simulations, so they can
 306 be less accurate, but should still describe the main features of the evolution of electron
 307 distributions and their results should statistically repeat those from the test particles ap-
 308 proach. Figure 6 (left panels) shows the evolution of the electron distribution function
 309 for a small wave amplitude ($\varepsilon = 10^{-4}$) at four time instants. Hereinafter $N_p = 32768$
 310 test particles are used to compute the changes of the distribution functions for all men-
 311 tioned approaches. Without nonlinear resonant effects, both SDE and mapping technique
 312 show results that are consistent with the test particle simulation. In this case, the evo-
 313 lution is diffusive (as expected for quasi-linear theory) and shows a spread of the initially

314 localized electron phase space density peak. The difference between SDE and the test
 315 particle simulation, most clearly seen around $E \sim 420$ keV, is due to the overestima-
 316 tion of the diffusion coefficients. We evaluate diffusion coefficients as a half of the vari-
 317 ance of ΔE distributions, and thus we assume that ΔE distributions are symmetric rela-
 318 tive to the mean value. However, even in the case of low-amplitude waves (see, e.g., Fig-
 319 ure 3(c)) this assumption may not work, which will result in an overestimation of the
 320 diffusion rate. The mapping technique does not require any assumptions about ΔE dis-
 321 tributions, and thus it performs better even in the case of low-amplitude waves.

322 Figure 6 (right panels) shows the evolution of the electron distribution function for
 323 a large wave amplitude ($\varepsilon = 10^{-3}$) at four time instants. For such intense waves, the
 324 SDE approach becomes inapplicable, but we can still compare results of test particle sim-
 325 ulations and the mapping technique. The mapping technique accounts for nonlinear reso-
 326 nant effects (e.g., phase trapping) and describes well the evolution of electron distri-
 327 butions. After several wave-particle resonant interactions (see top right panel of Figure
 328 6), the main electron population propagates to lower energies due to the phase bunch-
 329 ing, while a small population becomes trapped by waves and gains energy. During the
 330 drift of the main population to the smaller energies, the probability of particle trapping
 331 increases (see Figure 5(c)) and more particles become trapped and accelerated. Accel-
 332 erated particles appear in the resonant latitudes where no more phase trapping is pos-
 333 sible, and thus these particles start losing their energy due to the phase bunching. Around
 334 the time when the main population (at the initial peak of electron phase space density)
 335 reaches the left boundary of the allowed energies, the processes of phase bunching and
 336 phase trapping statistically compensate each other. This results in formation of a plateau
 337 in the distribution function (see the bottom right panel in Figure 6). Such an evolution
 338 of the electron distribution is consistent with theoretical predictions for the system with
 339 multiple nonlinear resonances (Artemyev et al., 2019).

340 The main uncertainty of the mapping technique arises at the edges of the simula-
 341 tion domain due to the finite grid size (discretization) in the ΔE distribution. In this
 342 approach, we treat the electron energy change as a probabilistic process and in Eqs. (13)
 343 and (17), electrons can reach the energy $E < E_{min}$ outside of the simulation domain.
 344 In the subsequent calculations, therefore, we will use the mapping functions correspond-
 345 ing to the energy E_{min} to calculate electron energy change. This underestimates the prob-
 346 ability of positive and overestimate probability of negative energy jumps, so that par-
 347 ticles can stay at the edges of the energy grid for a long time. The SDE approach also
 348 suffers from this boundary effect. This problem can be eliminated by introducing bound-
 349 ary conditions, e.g., reflective boundaries, or by normalizing the ΔE distribution around
 350 boundaries (see discussions in the Appendix of Artemyev et al., 2021).

351 7 Discussion and conclusions

352 In this study we compare two approaches that incorporate the wave-particle reso-
 353 nant effects into test particle simulations: SDE approach and mapping technique. Both
 354 approaches simplify the characterization of wave-particle interactions and reduce reso-
 355 nant effects to a ΔE distribution of energy jumps during a single interaction; in partic-
 356 ular, SDE further simplifies this description and operates only by the second moment
 357 (variance) of this distribution. The comparison shows that these two approaches pro-
 358 vide the same results for the system with low-amplitude waves, whereas electron non-
 359 linear resonant interactions with intense waves may be only described well by the map-
 360 ping technique. Note that the diffusion approximation is not only applicable to low in-
 361 tensity waves, but also to systems with very intense, low-coherent waves where the res-
 362 onance overlapping results in destruction of nonlinear effects (Tao et al., 2013; Zhang
 363 et al., 2020; An et al., 2022; Gan et al., 2022; Frantsuzov et al., 2023). Let us discuss ad-
 364 vantages and limitations of this technique.

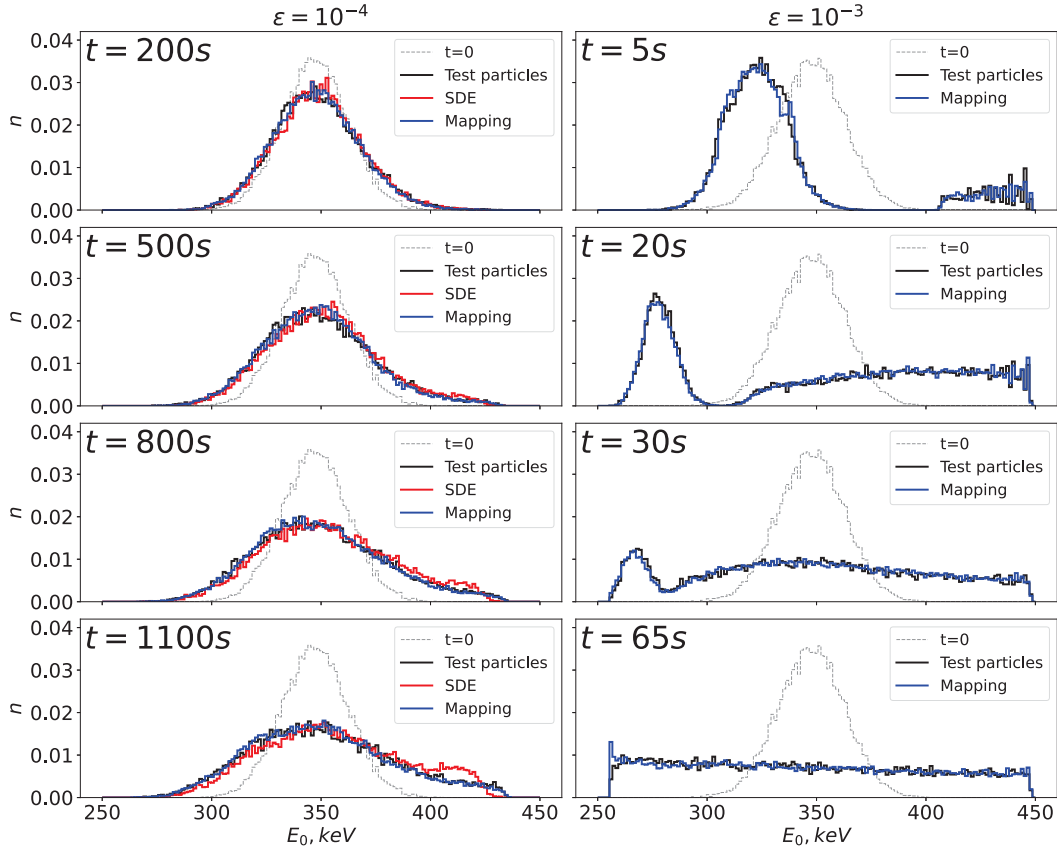


Figure 6. Evolution of the electron energy distribution for a systems without nonlinear resonant effects (left panels) and with nonlinear resonant effects (right panels). Grey color shows the initial distribution, black color shows test particle simulation results, red color shows results obtained with SDE, and blue color shows results for the mapping technique. Time (in seconds) is calculated under the assumption of a single resonant interaction per bounce period, $\tau_b(E)$, where τ_b is evaluated at L -shell = 6 and for equatorial pitch-angles derived from Eq. (7).

365 The mapping technique is based on ΔE distributions, which for fixed system pa-
 366 rameters depends only on the initial electron energy E_0 , i.e., we deal with 2D distribu-
 367 tions of energy (or pitch-angle) jumps, $F(\Delta E, E_0)$, which are used as described in Sec-
 368 tion 5. These distributions can be determined with any required accuracy from a set of
 369 short-term test particle simulations. However, in realistic space plasma systems, we deal
 370 with some ensemble of waves that can be described by the distribution of wave ampli-
 371 tudes and frequencies, $\mathcal{P}_w(B_w, \omega)$; for reasonable discretization levels and available sta-
 372 tistical wave datasets, this wave distribution usually consists of $\sim 10^2 - 10^3$ different
 373 pairs of (B_w, ω) (see, e.g., Artemyev, Zhang, et al., 2022; Zhang et al., 2022). Therefore,
 374 for the mapping technique, we would need to evaluate $10^2 - 10^3$ realizations of 2D $F(\Delta E, E_0)$
 375 distributions. Additional system dimensions can be introduced due to the dependence
 376 of $\mathcal{P}_w(B_w, \omega)$ on the geomagnetic activity and geophysical coordinates, (MLT , L -shell).
 377 Therefore, although the mapping technique essentially reduces computations relative to
 378 the full test particle simulations, this approach requires significant resources for simu-
 379 lation of electron dynamics during long-term global events (where wave and background
 380 characteristics can vary significantly), e.g., geomagnetic storms. More suitable applica-
 381 tion of the mapping technique is simulations of localized (spatially and temporally) events,
 382 like plasma sheet injections (Artemyev, Neishtadt, & Angelopoulos, 2022) or strong pre-
 383 cipitation bursts (Artemyev, Zhang, et al., 2022; Zhang et al., 2022). For global simu-
 384 lations with SDE, which substitute 2D $F(\Delta E, E_0)$ distributions by 1D diffusion coeffi-
 385 cients $D_{EE}(E_0)$, should be much more realistic, although this approach does not account
 386 for nonlinear resonant effects.

387 Both the SDE approach and mapping technique assume the evaluation of $F(\Delta E, E_0)$
 388 distributions before simulating the electron dynamics, and thus such distributions are
 389 usually evaluated for a prescribed background magnetic field. This simplification may
 390 work for the inner magnetosphere, where the background dipole field does not vary too
 391 much (see Orlova & Shprits, 2010; Ni et al., 2011, for discussions on when this assump-
 392 tion does not work well), but cannot be well justified for plasma injections characterized
 393 by rapid, significant variations of the magnetic field configuration (see discussion in Ashour-
 394 Abdalla et al., 2013; Sorathia et al., 2018; Birn et al., 2022). The background magnetic
 395 field configuration determines the resonant condition and thus should control the effi-
 396 ciency of electron scattering by waves. In principle, the effect of the magnetic field re-
 397 configuration can be included into SDE and mapping approaches, but this would require
 398 the evaluation of $F(\Delta E, E_0)$ distributions for multiple magnetic field configurations, which
 399 is usually unlikely with available computational resources. A possible solution will be
 400 to analytically evaluate $F(\Delta E, E_0)$ distributions (e.g., Vainchtein et al., 2018; Artemyev
 401 et al., 2021), but this solution requires more developed theoretical models of wave-particle
 402 resonant interactions including effects of wave-field deviations from a simple plane wave
 403 model (see discussions in Mourenas et al., 2018; Artemyev et al., 2023). So far such ef-
 404 fects have been evaluated (e.g., Tao et al., 2011; Zhang et al., 2020; Allanson et al., 2021;
 405 Gan et al., 2022; An et al., 2022) and incorporated into $F(\Delta E, E_0)$ distributions (e.g.,
 406 Kubota & Omura, 2018; Hsieh et al., 2020, 2022) only via numerical integration of a large
 407 test particle ensemble.

408 Both the SDE approach and mapping technique are based on numerical evaluations
 409 of the probability distribution function of energy jumps, $F(\Delta E, E_0)$. Although we have
 410 adopted the plane wave approximation in this study, the procedure of $F(\Delta E, E_0)$ eval-
 411 uations is independent of such approximations and F can be calculated for more real-
 412 istic modes of wave-packets (see observations and simulations in Zhang et al., 2018, 2021).
 413 This generalization for the finite wave-packet size is quite important for constraining the
 414 efficiency of nonlinear resonant effects (see, e.g., Kubota & Omura, 2018; Zhang et al.,
 415 2020; An et al., 2022). Therefore, in the next step, one would need to incorporate real-
 416 istic distributions of whistler-mode wave-packet sizes.

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Open Research

No data sets were used in this work.

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