Relativistic Brownian motion on a graphene chip

A. Pototsky\textsuperscript{1}, F. Marchesoni\textsuperscript{2}, F.V. Kusmartsev\textsuperscript{3}, P. Hänggi\textsuperscript{4}, and S.E. Savel’ev\textsuperscript{3,\textsuperscript{a}}

\textsuperscript{1} Department of Mathematics, University of Cape Town, 7701 Rondebosch, South Africa
\textsuperscript{2} Dipartimento di Fisica, Università di Camerino, 62032 Camerino, Italy
\textsuperscript{3} Department of Physics, Loughborough University, Loughborough LE11 3TU, UK
\textsuperscript{4} Institut für Physik, Universität Augsburg, 86135 Augsburg, Germany

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Abstract. Relativistic Brownian motion can be inexpensively demonstrated on a graphene chip. The interplay of stochastic and relativistic dynamics, governing the transport of charge carrier in graphene, induces noise-controlled effects such as (i) a stochastic effective mass, detectable as a suppression of the particle mobility with increasing the temperature; (ii) transverse harmonic mixing, whereby electron transport can be controlled by two orthogonal, commensurate ac drives; (iii) a transverse ratchet effect, measurable as a net current orthogonal to an ac drive on an asymmetric substrate, and (iv) chaotic stochastic resonance. Such properties can be of practical applications in the emerging graphene technology.

1 Introduction

As astroparticle data often require a relativistic analysis, the challenge of a consistent formulation of relativistic statistical thermodynamics is so much intriguing as timely \cite{1}. Such a challenge encompasses the notion of relativistic Brownian motion \cite{2}, as well. In particular, issues like the relativistic generalization of inherently non-local thermodynamic quantities, such as heat and work, or the very concept of an equilibrium heat bath, have not been unambiguously settled, yet \cite{1,2}. The state of the art in this field thus calls for experiments capable to assess and validate the existing theoretical approaches. This presently constitutes a nearly impossible task, due to the scarcity of the available relativistic data sets, which often also lack the necessary accurateness. Typically, such data either are being obtained from cosmic rays \cite{3} or could be, in the future, from expensive high-energy experiments. To overcome this difficulty, we propose here an alternative and affordable route to demonstrate the physics of relativistic Brownian motion under controllable laboratory conditions, namely on a suitable graphene chip.

Graphene is essentially a monolayer of carbon atoms packed in a honeycomb lattice and isolated from the bulk \cite{4}. The fact that graphene has a 2D structure makes its electronic properties unique. Charge carriers propagating through such a lattice are known to behave as relativistic massless Dirac fermions \cite{4–6}. On the quantum mechanical level, they are described by the 2D analog of the Dirac equation with the Fermi velocity, $v_F = 10^6 \text{ cm/s}$, replacing the speed of light in the Dirac description. Close to the so-called Dirac points, the energy-momentum relation of the elementary excitations, often associated with quasi-particles, is linear: the velocity of the charge carriers is thus always collinear with the momentum, its modulus being constant and equal to $v_F$.

Our proposal relies on the observation that for strong applied electric fields or at high enough temperatures, quantum transport effects, otherwise dominant in graphene, are effectively suppressed and the motion of charge carriers can be satisfactorily described by classical (i.e., non-quantum) relativistic equations. Such a semiclassical description has been invoked \cite{7}, for instance, to successfully interpret graphene’s unusually broad cyclotron resonance \cite{8,9} and is consistent with the relativistic hydrodynamical approach proposed in reference \cite{10,11}.

As detailed with this paper, the combination of stochastic and relativistic dynamical effects governing the carrier motion in graphene results in some unexpected transport phenomena. The source of stochasticity is provided by the finite temperature, whereby the charge carriers are being scattered by phonons, lattice defects and buckling, and the sample boundaries, which leads to an equilibrium redistribution of their energies. We emphasize, however, that a noisy environment in the chip can also be generated by external noise sources, controllable or not, independent of the temperature, e.g., in the form of current or voltage fluctuations. Moreover, in graphene the Fermi velocity, which plays the role of the light speed in the relativistic equations for its charge carriers, is relatively low; one can then analyze driven relativistic Brownian
motion in the presence of electromagnetic fields by ignoring otherwise hardly tractable relativistic retardation effects. Under these conditions, driven relativistic Brownian particles become accessible to both theory and experiment [12,13].

This paper is organized as follows. In Section 2 the dynamics of charge carriers in a graphene chip is formulated, in quasi-classical approximation, as a 2D relativistic Langevin equation. A viscous damping force and a fluctuating source have been added so as to establish a high-temperature relativistic equilibrium Jüttner distribution. The phenomenological motivations for such a model and certain restrictions to its validity are discussed in Section 3. The combination of the stochastic and relativistic properties of the model is numerically shown to induce unexpected noise-controlled effects such as: (i) a temperature dependent effective mass, responsible for the suppression of the particle mobility at high temperatures (Sect. 4.1), (ii) transverse harmonic mixing, with electron rectification being sustained by two orthogonal, commensurate ac drives (Sect. 4.2), (iii) transverse ratcheting, signaled by a net current orthogonal to an ac drive applied on an asymmetric substrate (Sect. 4.3), and (iv) a chaos controlled manifestation of stochastic resonance (Sect. 4.4). All relativistic stochastic effects reported in this paper are then summarized in Section 5, where, for reader’s convenience, the different behaviors of relativistic (massless) and nonrelativistic (massive) Brownian particles confined to 2D geometries, are compared case by case.

2 The model

The scheme of a simple graphene chip is sketched in Figure 1a. An undoped graphene sheet sits on a periodic sequence of parallel electrodes with alternate constant potentials, respectively, $U^{+}$ and $U^{-}$, with $U^{+} > U^{-}$. By tuning the distance between electrode pairs within a fixed spatial period, $L$, the ensuing staggered electric potential, $U(x)$, can be modulated at will, symmetrically or asymmetrically, with amplitude $U$ in the $x$ direction, while being identically zero in the $y$ direction. An example of the directed potential $U(x)$ is drawn in Figure 1a. Additionally, the charge carriers are driven by a spatially homogeneous electric field, $E(t) = (E_{x}(t), E_{y}(t))$, with sinusoidal components $E_{x}(t)$ and $E_{y}(t)$ acting along the $x$ and the $y$ axis, respectively.

The random dynamics of the charge carriers in a graphene sheet can be described (see Sect. 3 for details) by a set of coupled Langevin equations for the components of the 2D momentum, $p = (p_{x}, p_{y})$, subject to the relativistic dispersion relation $\varepsilon = v_{F}p_{0}$, where $p_{0} = (p_{x}^{2} + p_{y}^{2})^{1/2}$, $\varepsilon$ is the particle energy and $v_{F}$ the Fermi velocity. The form of such phenomenological Langevin equations is determined by the condition that the chosen particle-reservoir coupling must lead to the same equilibrium momentum distribution as predicted by the fully microscopic (Hamiltonian) theory. At low temperatures, the equilibrium distribution function of the quasiparticle (electron) energy obeys the Fermi-Dirac statistics, $\rho(\varepsilon) = 1/(\exp[\varepsilon/kT] + 1)$, where $T$ is the temperature. However, the operation of the graphene chips considered here typically requires carrier energies $\varepsilon \sim U$, so that for sufficiently high energies, $kT \ll \varepsilon \ll U$, $\rho(\varepsilon)$ is conveniently approximated by the relativistic Jüttner distribution, $\rho(p) \sim \exp(-\varepsilon/p/kT)$ [2,14]. A viable set of Langevin equations proven to be consistent with a 2D Jüttner distribution reads [2]

$$
\begin{align*}
\dot{p}_{x} &= -\gamma v_{F}p_{x}/p_{0} - dU(x)/dx + E_{x}(t) + \sqrt{2\gamma kT}\xi_{x}(t),
\dot{p}_{y} &= -\gamma v_{F}p_{y}/p_{0} + E_{y}(t) + \sqrt{2\gamma kT}\xi_{y}(t),
\end{align*}
$$

where $\gamma$ denotes the phenomenological constant damping introduced in Section 3. The random forces $\xi_{x}(t)$ and $\xi_{y}(t)$ are two white Gaussian noises with $\langle \xi_{i}(t) \rangle = 0$ and $\langle \xi_{i}(t)\xi_{j}(0) \rangle = \delta_{ij}\delta(t)$, for $i, j = x, y$, which ensures proper thermalization at temperature $T$. We recall here that the modulus of $V$ is constant and equal to $v_{F}$ [7]. These equations describe a relativistic Brownian dynamics, thus suggesting new settings for the experimental and theoretical investigation of relativistic thermodynamics and relativistic Brownian motion. In Section 4 we report results from extensive numerical simulations we performed by integrating equations (1).

We stress here that the validity of equations (1) for electrons in graphene is restricted to the quasi-classical limit, where quantum mechanical effects can be safely neglected. In particular, as detailed in the forthcoming section, the testing ground for relativistic phenomena on a graphene chip is subject to the following physical restrictions: (i) the de Broglie wavelength across the barrier, $\lambda_{\varepsilon} = h/\varepsilon$, must be much smaller than the period of substrate potential, $\lambda_{s} \ll L$, so as to neglect either miniband or discrete energy levels; (ii) the de Broglie wavelength parallel to the barrier, $\lambda_{v} = h/p_{0}$, must be much smaller than $L$, so as to suppress the probability of chiral tunneling (Klein’s paradox [12,15]). The condition $\lambda_{\varepsilon} < L$ allows neglecting, in particular, the electron-hole tunneling; (iii) temperature (or noise strength) must be high enough to further ensure that the chiral tunneling at the top of the $U(x)$ barriers is negligible with respect to the competing noise-activated particle hopping, i.e.,

$$\min \{h\nu_{F}/L; \hbar \omega_{0}^{2}U^{2}/v_{F}A^{2}L\} \ll kT \ll U,$$

where $A$ is the amplitude of the field component parallel to the barrier.

Condition (iii) on $T$ is introduced because noiseless carriers moving along the $x$-axis with $\lambda_{\nu} \ll L$ would still be subject to Klein’s mechanism (barrier transparency) [16]. Note, however, that in 2D geometries with smooth potentials $U(x)$, the fraction of electrons undergoing Klein’s tunneling is negligibly small, not only at high temperatures [as guaranteed by condition (iii)], but also in the presence of sufficiently strong transverse drives. Indeed, drives oriented along the $y$ axis produce an optical
gap in the energy spectrum proportional to the ac electric field amplitude. This tends to suppress chiral tunneling even for particles with $p_y = 0$ [17,18], thus corroborating our quasiclassical model.

Conditions (i)–(iii) are formulated in Section 3; the reader not interested in those technical details can jump directly to Section 4.

3 Quasi-classical dynamics of charge carriers in graphene

In the low-energy limit and in the absence of randomness (fluctuations and disorder, alike), the graphene charge carriers, electrons and holes, obey the Dirac equations,

$$(P_x - iP_y) \psi_B + \frac{U(x,y,t)}{v_F} \psi_A = \frac{i\hbar}{v_F} \partial_x \psi_A,$$

$$(P_x + iP_y) \psi_A + \frac{U(x,y,t)}{v_F} \psi_B = \frac{i\hbar}{v_F} \partial_y \psi_B, \quad (3)$$

where $U$ is a chip-dependent scalar potential, the two spinor components of the carrier wave function, $(\psi_A, \psi_B)$, refer to the two triangular graphene sublattices and the momentum operators are defined as

$$(P_x, P_y) = (-i\hbar \partial / \partial x, -i\hbar \partial / \partial y) \equiv -i\hbar \nabla.$$

Applying the operators $(P_x - iP_y)$ and $(P_x + iP_y)$, respectively, to the first and the second equation leads to the coupled second-order partial differential equations,

$$
\begin{align*}
-\hbar^2 \nabla^2 + \frac{\hbar^2}{v_F^2} \partial_y^2 + 2 \frac{i\hbar}{v_F} \partial_t \frac{U^2}{v_F^2} \psi_B = \\
-\frac{i\hbar}{v_F} \partial_t \psi_B + \frac{i\hbar}{v_F} \psi_A \left( \frac{\partial U}{\partial x} + i \frac{\partial U}{\partial y} \right),
\end{align*}
$$

$$
\begin{align*}
-\hbar^2 \nabla^2 + \frac{\hbar^2}{v_F^2} \partial_y^2 + 2 \frac{i\hbar}{v_F} \partial_t \frac{U^2}{v_F^2} \psi_A = \\
-\frac{i\hbar}{v_F} \partial_t \psi_A + \frac{i\hbar}{v_F} \psi_B \left( \frac{\partial U}{\partial x} - i \frac{\partial U}{\partial y} \right), \quad (4)
\end{align*}
$$

where $\nabla^2 \equiv \partial^2 / \partial x^2 + \partial^2 / \partial y^2$. Equivalently, the effective actions $(S_A, S_B)$, defined as $(\psi_A, \psi_B) = (\exp(iS_A/v_F), \exp(iS_B/v_F))$, satisfy the coupled equations,

$$
\begin{align*}
(\nabla S_A)^2 &= -\frac{1}{v_F^2} \left( \frac{\partial S_A}{\partial t} - U \right)^2 - i\hbar \nabla^2 S_A + \frac{i\hbar}{v_F} \partial_y S_A = \\
&= -\frac{i\hbar}{v_F} \frac{\partial U}{\partial t} + \frac{\hbar}{v_F} \frac{\partial}{\partial x} (S_A - S_B) \left( \frac{\partial U}{\partial x} - \frac{\partial U}{\partial y} \right),
\end{align*}
$$

$$
\begin{align*}
(\nabla S_B)^2 &= -\frac{1}{v_F^2} \left( \frac{\partial S_B}{\partial t} - U \right)^2 - i\hbar \nabla^2 S_B + \frac{i\hbar}{v_F} \partial_y S_B = \\
&= -\frac{i\hbar}{v_F} \frac{\partial U}{\partial t} + \frac{\hbar}{v_F} \frac{\partial}{\partial x} (S_B - S_A) \left( \frac{\partial U}{\partial x} - \frac{\partial U}{\partial y} \right). \quad (5)
\end{align*}
$$

In the limit $\hbar \to 0$, $S_A$ and $S_B$ tend to the unique classical action, $S$, solution of the classical Hamilton-Jacobi equation,

$$
\left( \frac{\partial S}{\partial x} \right)^2 + \left( \frac{\partial S}{\partial y} \right)^2 - \frac{1}{v_F} \frac{\partial S}{\partial t} - U^2 = 0. \quad (6)
$$
On making use of the well-known relation, $H = \partial S/\partial t$, one eventually derives from equation (6) the classical Hamiltonian for a relativistic massless particle,

$$H = \pm v_F \sqrt{p_x^2 + p_y^2 + U},$$

(7)

where, in notation of Section 2, $U(x,y,t) = U(x) - E_y(t)y - E_z(t)x$ and $\pm$ referring respectively to electrons and holes, whose densities are related by electroneutrality and are controlled by the chemical potential.

Note that the lowest order corrections to the classical limit, equation (6), are $O(\hbar)$, at variance with the standard WKB formalism for the Schrödinger equation. Such corrections, for $|S_A - S_B| \sim \hbar$, are due to the electron-hole transitions, so that a systematic quasiclassical treatment would require a two particle species description [19,20].

Thus far, the interactions of a single charge carrier with, e.g., phonons and impurities have been neglected. If the relevant collision times are short compared with the operation time constants of the graphene chip (periods of the ac-drivers, carriers drift times across the substrate potential cells, etc.), then such interactions can be accounted for by adding to the potential function, $U$, an appropriate fluctuating term [21], $U_\xi$, namely, $U(x,y,t) = U(x) - E_y(t)y - E_z(t)x + U_\xi$, which corresponds to an isotropic white noise force in the plane $(x,y)$.

The same interactions are also responsible for the damping term in the Langevin equations (1). The appropriate form of the damping and the fluctuating terms are determined by the fluctuation-dissipation theorem. In Section 2 we assumed a damping term of the viscous type, $-\gamma V$ [22], and then determined the random forces accordingly [2]. The damping coefficient $\gamma$ contains contributions from all possible collision mechanisms a charge carrier may undergo when moving through a graphene chip. To determine the relative importance of such mechanisms would require a detailed quantum kinetic theory (see, e.g., Ref. [23]), which rests beyond the purposes of our semi-classical approach, and, more importantly, direct access to experimental parameters, such as spatial disorder (impurities), internal lattice stress in the presence of a substrate (phonon distribution), local temperature and magnetic field fluctuations, which are hardly controllable.

The semiclassical approach outlined above provides a meaningful phenomenological description of charge carrier transport in a graphene chip as long as its characteristic length scales (e.g., the chip size and the electrode spacings) are much larger than the de Broglie wavelengths, (i) $\lambda_z = \hbar/v_F$ and (ii) $\lambda_y = \hbar/p_y$, with $(p_x, p_y) = (\partial S/\partial x, \partial S/\partial y)$. Condition (i) states that for $\lambda_z$ much shorter than the characteristic length scale of the substrate potential along the $x$-axis, $L \sim |U/(dU/dx)|$ for smooth functions $U(x)$, the mini-zone structure or discreteness of the carrier energy spectrum becomes negligible.

Condition (ii), $\lambda_y = \hbar/p_y \ll L$, follows from requiring that the probability of chiral tunneling is distinctively smaller than unity, i.e., that the particle must be reflected many times by the potential barrier before it may tunnel across. Chiral tunneling is exponentially suppressed for non-zero transverse momenta, $p_y \neq 0$, and smooth time-independent potentials, $U(x)$. In this case, the classical action has the form $S = p_y u - Et + S_0(x)$, where $E$ is the conserved energy and $S_0(x)$ can be readily derived from the the Hamilton-Jacobi equation (6), that is, $S_0 = \pm \int (E - U)^{1/2}/v_F \sqrt{p_x^2 + p_y^2} \, dx$. Chiral tunneling occurs through the classically forbidden region, $|E - U| < v_F|p_y|$; therefore, the smoother the potential, the larger the size of the forbidden region, $2v_F|p_y|/|dU/dx|$. As a consequence, chiral tunneling is characterized by an exponentially small probability, $w = \exp(-k_D p_y^2/h|dU/dx|)$ [16], with $k_D$ denoting a constant of the order of unity, which is appreciable only for a small fraction of electrons with $p_y \gtrsim \sqrt{h|dU/dx|}/v_F$. By introducing the estimate $p_y \sim U/v_F$ for the transverse momentum, we derived condition (ii). It is important to stress that a time dependent electric field acting along the barrier, $E_y(t)$, tends to exponentially suppress also the chiral tunneling of electrons moving perpendicularly to the barrier, $p_x = 0$, that is, $w = \exp(-\pi e^2 v_F A^2/(4h|dU/dx|\omega_0^2))$ [17,18]. This remark further corroborates our quasiclassical approach in the presence of strong enough drives. As an additional consistency requirement for our semiclassical description of graphene charge carrier transport over the substrate potential, we must further ensure that chiral barrier tunneling is negligible with respect to the thermal barrier hopping, that is, $w \ll \exp(-U/kT)$; hence, the inequalities (2) of condition (iii). Correspondingly, the lowest temperature, $T_c$, where the semiclassical approach applies, can be set to $T_c \sim h e_F/\hbar L$. For the standard estimates of the Fermi velocity, $v_F \sim 10^6$ m/s, and of the characteristic length of graphene superlattices, $L > 1$ $\mu$m (Ref. [24]), $T_c \lesssim 50$ K.

Accordingly, the frequencies $E(t)$ in equations (1) are required to satisfy the condition $\hbar \omega < U \sim k_B T_c$, namely, to fall below the THz range. All these estimates suggest that the effects presented in the next section rest well within the reach of present nano-technology [24].

4 Relativistic stochastic effects

The interplay of stochastic and relativistic dynamics, governing the transport of charge carrier in graphene, induces unexpected noise-controlled effects. In this section we focus on the most remarkable differences between relativistic and nonrelativistic Brownian motion in 2D.

4.1 Stochastic effective mass

The momentum components $p_x$ and $p_y$ in equations (1) are coupled via the dispersion relation, $\varepsilon = v_F p_0$, even in the absence of a substrate potential, $U(x) \equiv 0$. This gives rise to a peculiar phenomenon, which characterizes the relativistic Brownian motion with respect to its nonrelativistic counterpart (Fig. 1b). By numerically solving equations (1) for electrons driven, say, along the $y$ axis by $E = (0, A \cos(\omega_0 t))$, we determined the strength, $S$, of the delta-like peak of the $y$ power spectral density
at the driving frequency, $\omega_y$, that is, $S = 4M_1^2$, where $\lim_{t \to -\infty} y(t) \simeq M_1 \cos(\omega_y t - \phi_0)$ for an appropriate phase-lag, $\phi_0$ (not shown) [25]. Contrary to the non-relativistic limit, where it is $T$ independent, for relativistic particles at high temperatures $S$ decreases according to the approximate power law $T^{-2}$. This result can be quantitatively explained by noticing that according to the high-$T$ Jüttner distribution, the particle energy distribution is $\rho(\epsilon) \sim \epsilon \exp(-\epsilon/kT)$; hence, $\langle \epsilon \rangle = v_F p_0 = 2kT$. Upon replacing $p_0$ by $2kT/v_F$, equation (1) is reduced to a set of decoupled Langevin equations describing a non-relativistic particle of effective mass $m_{\text{eff}} = 2kT/v_F^2$. Accordingly, $S/S_0 = 1/[1 + (\omega_y/\omega_0)^2]$ with $S_0 = (A/\gamma \omega_0)^2$ and $\omega_0 = \gamma/m_{\text{eff}} = v_F^2/2kT$ (dashed line in Fig. 1b). Stochastic mass renormalization provides a simple validation check for our phenomenological model in equation (1).

### 4.2 Relativistic harmonic mixing

Consider next the case of a relativistic particle driven by two orthogonal harmonic electric fields in a symmetric periodic potential $U(x)$. In the non-relativistic limit, two harmonic drives with commensurate frequencies can induce rectification (i.e., a net electron drift or directed current) only if both drives are aligned along the $x$-axis [26–28]. For the relativistic Langevin equation (1), the coupling between the $x$ and $y$ degrees of freedom produces an unusual manifestation of harmonic mixing, whereby rectification occurs even if the two harmonic drives are orthogonal. To demonstrate this genuine relativistic effect we integrated equations (1) for the ac electric field $E = (B \cos(\omega_z t + \phi), A \cos(\omega_y t))$ and the symmetric sublattice displacement potential $U(x) = -\sin(2\pi x)$. The $x$ and $y$ displacements for a non-relativistic particle subjected to such force fields are statistically uncorrelated, thus yielding zero average currents in both direction. In distinct contrast, for a relativistic particle the coupling between $p_x$ and $p_y$ causes a mixing of the orthogonal components of the ac electric field $E$, and, as a result, a net current in the $x$-direction. By inspecting Figure 2, we concluded that the net current $J_x = \langle \dot{x} \rangle/L$ is non-zero only for certain ratios of the two drive frequencies, namely, for $\omega_y/\omega_z = (2p-1)/2q$ with $p, q$ integer numbers. As expected [27,28], for fixed frequencies $J_x$ depends sinusoidally on the phase $\phi$ (see inset of Fig. 2). This intriguing effect yields a measure of the degree of coupling between particle diffusion in the $x$ and $y$ directions, and, therefore, a simple criterion to assess the validity of the relativistic Langevin equations (1) to model charge carriers in graphene.

### 4.3 Relativistic ratcheting

We consider next charge carrier transport due to the rectification of non-equilibrium perturbations on a substrate with asymmetric potential $U(x)$ (ratchet or Brownian motor effect [29]). The working principle of a Brownian motor is that, under certain conditions, asymmetric devices are capable of rectifying random (i.e. noisy) and/or deterministic (periodic) modulations. In the non-relativistic regime, ratcheting occurs only in the $x$ direction, as a mere effect of the $x$ component of the periodic drive, $E_x(t)$. The rectification current weakens if $E(t)$ is rotated at an angle with the $x$-axis, until it drops to zero for ac drives parallel to the substrate valleys. Indeed, the component $E_y(t)$ of the ac-field does keep the system out of equilibrium, but cannot be rectified, since the substrate potential is uniform in the $y$ direction. Stated otherwise, the $x$ and $y$ dynamics keep being decoupled. In the relativistic Langevin equations (1), instead, the orthogonal ac-drive components, $E_x(t)$ and $E_y(t)$, are nonlinearly coupled through the dispersion relation, so that both can be rectified by the asymmetric potential $U(x)$. Most remarkably, apart from a special parameter range dominated by relativistic chaotic effects (see Sect. 4.4), transverse rectification induced by $E_y(t)$ in the $x$ direction only occurs at finite temperatures, thus implying a *bona fide* noise-sustained Brownian ratchet [29]. Noise is required to force particle fluctuations in the $x$-direction, around the asymmetric minima of $U(x)$; as $p_y$ is driven by $E_y(t)$ toward values of the order of $\sqrt{T}$, or smaller, correspondingly $v_x \approx \pm v_F p_x$ jumps to a maximum, $v_x \approx v_F p_x/|p_y| = \pm v_F$, so that the particle is kicked in the $x$ direction, either to the right or to the left; this is how spatial asymmetry comes into play. Naturally, such a rectification mechanism becomes ineffective when the particle sits at a potential minimum with $p_y = 0$, which only occurs in the absence of noise, i.e., for $T \equiv 0$.

To demonstrate the relativistic transverse ratchet effect, it suffices to arrange the electrodes on the graphene chip so as to generate an asymmetric periodic potential, $U(x)$, directed along $x$ and ac drive forcing the charge carriers along $y$ with $E(t) = (0, A \cos(\omega_y t))$. By numerically integrating the corresponding equations (1), we observed that a relativistic particle tends, indeed, to drift in

![Fig. 2.](image-url)
the x direction with net current $J_x = \langle \dot{x} \rangle / L$ and diffusion coefficient $D_x = \lim_{t \to \infty} \langle (x(t)^2) - \langle x(t) \rangle^2 \rangle / 2t$. The dependence of $J_x$ and $D_x$ on the temperature and the driving frequency are displayed in Figure 3. In panel (a) it is apparent that at constant $\omega_y$ the absolute value of the current, $|J_x|$, hits a maximum for an optimal $T$ and tends to zero in the limits $T \to 0$ and $T \to \infty$, thus underscoring the key role played by noise. This result supports our conclusion that we deal with a new category of noise-sustained (no current at zero noise) and relativistic (no ratchet phenomenon in the non-relativistic limit) Brownian motors. The direction of the net rectification current depends on both the driving frequency and the temperature; its sign can be (multiply) reversed by changing either $T$, panel (a), or $\omega_y$ (not shown). Most remarkably, in panel (b) one can locate a finite temperature interval, where, for a given $\omega_y$, the scaled diffusion coefficient $D_x / D_0$ decreases with increasing $T$; indeed, $D_x$ approaches the expected asymptotic value $D_0 = kT / \gamma$ only after going through a minimum. This $T$-interval approximately coincides with the $T$-interval in panel (a), where the absolute value of the current and, therefore, the mobility of the charge carriers is the largest. Analogously, we checked that the absolute maximum of $D_x$ as a function of $\omega_y$ corresponds to a local minimum of $J_x$ versus $\omega_y$. This observation suggests that for appropriate system parameters, elementary charge excitations in graphene can be regarded as Brownian particles with large mobility and low diffusivity.

4.4 Chaos and stochastic resonance

To investigate the role of anomalous signal amplification we consider relativistic Brownian motion driven along $y$ by the time-periodic electric field $E(t) = (0, A \cos(\omega_y t))$, and confined along $x$ by the double-well potential $U(x) = x^4 / 4 - x^2 / 2$, sketched in Figure 4a. In the deterministic, or noiseless regime, $T \equiv 0$, the motion of a classical particle in the $x$-direction would be frozen, i.e. the particle would sit in one of the two $U(x)$ minima, regardless of the drive applied in the $y$ direction. In clear contrast, the response of the relativistic particle in the $x$-direction is strongly affected by the nonlinear coupling of the $x$- and $y$-dynamics, controlled, respectively, by $U(x)$ and $E_y(t)$. In fact, by appropriately tuning the drive parameters $A$ and $\omega_y$, the relativistic particle can execute either small amplitude intra-well oscillations within one confining well, or large-amplitude inter-well oscillations between the two confining wells, or even exhibit chaotic-like switching between the two wells. The phase diagram of these three different dynamical regimes is depicted in Figure 4b with samples of the corresponding $x$-trajectories reported in the insets.

Our numerical simulations showed that in the chaotic regime a relativistic particle performs small amplitude oscillations around one $U(x)$ minimum for an unpredictable time interval (residence time) before switching over into the other minimum, where it resumes its small amplitude oscillations until the next switching event. These switchings are due to the intrinsic instability of the particle dynamics and not to thermal fluctuations, as in our simulations the temperature was initially set to zero.

On raising the temperature, the inter-well dynamics induced by the transverse field, $E_y(t)$, becomes increasingly noise dominated. In Figure 4c we display the normalized distribution density of the residence times in either $U(x)$ well, $\tau$, at different $T$. The peak split structure, characteristic of the noiseless chaotic dynamics, merges into a single broad peak centered at around $\tau = 2\pi / \omega_y$. The corresponding peak strength, $S$, defined here as the area encircled by the peak, attains a maximum for $kT$ smaller than, but close to the barrier height of $U(x)$, $\Delta U = 0.25$. Such a behavior can be regarded as an instance of stochastic resonance, a phenomenon well established for bistable non-relativistic Brownian motion [25], with two additional peculiarities: here the periodic drive acts perpendicularly to the hopping direction and the resonance is related with chaotic rather than periodic switches between potential wells.

5 Conclusions

The relativistic stochastic effects reported in this paper are summarized in Table 1, where for reader’s convenience we compare case by case the different behaviors of relativistic (massless) and non-relativistic (massive, or overdamped) Brownian particles in 2D. One has to bear in mind that the orthogonal coordinates of the relativistic particle in equations (1) are dynamically coupled even in the absence of a nonlinear substrate or external gradients. Such a coupling vanishes in the non-relativistic limit. As a consequence, first, the response of graphene charge carriers to a time-periodic drive is temperature dependent, see
Fig. 4. (Color online) Relativistic stochastic resonance. (a) The confining potential, \( U(x) = x^4/4 - x^2/2 \), bistable in the \( x \)-direction and uniform in the transverse direction \( y \); (b) corresponding phase diagram for \( E = (0, A \cos(\omega_y t)) \). Insets: trajectory samples for points in the plane \((A, \omega_y)\) (denoted by arrows) belonging to different dynamical phases. (c) Residence time distributions, \( \rho(\tau) \), for different \( T \) (as indicated) and constant drive, \( A = 15 \) and \( \omega_y = 0.5 \). Relevant simulation parameters: integration step, \( 1.2 \times 10^{-3} \); run time length, \( 5 \times 10^5 \); distribution time bin, \( 2.5 \times 10^{-2} \). Inset: integrated distribution peak, \( S = \int_{\tau=14}^{\tau=11} \rho(\tau) d\tau \), vs. \( T \) (SR signature). Other simulation parameters are: \( v_F = \gamma = k = 1 \).

### Table 1. Driven stochastic dynamics of relativistic and non-relativistic particles: a comparison.

|                          | Relativistic massless particles | Classical massive or overdamped particles |
|--------------------------|--------------------------------|------------------------------------------|
| **Experimental data sources** | Cosmic rays, graphene chip                  | Nano-particles in liquids and nano-channels, Vortices in superconductors, bio-molecules in bio-systems etc. |
| **Possible applications** | THz graphene-based electronics              | Nanotechnology, Noise-driven nano-robots, Drug delivery etc. |
| **Current state of the art** | Conflicting formulations limited applicability | Well established working tool |

| Coupling between orthogonal coordinates | Yes | No |
|-----------------------------------------|-----|----|
| **Rectification in a spatially symmetric periodic potential \( U(x) \), induced by mixing of two harmonic signals** | Occurs if the signals are applied in the same as well as in the orthogonal directions | Occurs only if the two signals are applied in the same direction |
| **Rectification in a spatially asymmetric periodic potential \( U(x) \) induced by a single center-symmetric unbiased signal** | Occurs if the signal is applied along the \( x \), as well as along the \( y \) direction | Occurs only if the signal is applied along the \( x \) direction |
| **Motion in a double-well potential with orthogonal harmonic drive; chaos and resonance** | Zero temperature lateral response can be constant, periodic or chaotic in time. Finite temperature response in the chaotic regime: non-exponential double-peak distribution of the residence times; stochastic resonance | No zero temperature lateral response Finite temperature response: random switchings with exponential distribution of the residence times |
Figure 1b, whereas the response of a classical free particle only depends on the drive. Second, on applying a periodic symmetric spatial modulation in a given direction, a relativistic massless particle can be rectified in that direction, when driven by two center-symmetric unbiased time periodic fields acting in orthogonal directions (transverse harmonic mixing, Fig. 2). Rectification occurs only for a certain combination of the relevant driving frequencies. In the case of a classical particle, harmonic mixing would require that both harmonic signals act in the direction of the spatial modulation. Third, in our approach may apply not only to graphene but also to other materials, like cerium and its dispersion relation. In this regard our approach may apply to certain types of semiconductors. Moreover, the results discussed in this work, besides suggesting an inexpensive technique to assess the current stochastic formalisms for relativistic particles, might well find application in ultrafast electronics. Indeed, the implementation of the effects reported above is expected to become a viable technology, should one eventually succeed in replacing GaAs- with graphene-based electronics [30,31].

We conclude by stressing once more that, irrespective of the phenomenological details of equations (1), the spatial coordinates of a relativistic particle in 2D are nonlinearly coupled via a non-separable energy-momentum dispersion relation. In this regard our approach may apply not only to graphene but also to other materials, like certain types of semiconductors. Moreover, the results discussed in this work, besides suggesting an inexpensive technique to assess the current stochastic formalisms for relativistic particles, might well find application in ultrafast electronics. Indeed, the implementation of the effects reported above is expected to become a viable technology, should one eventually succeed in replacing GaAs- with graphene-based electronics [30,31].

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