Diffusion transitions in a 2D periodic lattice

Matheus J. Lazarotto* and Iberê L. Caldas
Instituto de Física, Universidade de São Paulo, Rua do Matão 1371, São Paulo 05508-090, Brazil

Yves Elskens
Aix-Marseille Université, CNRS, UMR 7345 PIIM, F-13397 Marseille, France
(Dated: May 16, 2022)

Spatial diffusion of particles in periodic potential models has provided a good framework for studying the role of chaos in global properties of classical systems. Here a bidimensional “soft” billiard, classically modeled from an optical lattice Hamiltonian system, is used to study diffusion transitions under variation of the control parameters. Sudden transitions between normal and ballistic regimes are found and characterized by inspection of topological changes in phase-space. Transitions correlated with increases in global stability area are shown to occur for energy levels where local maxima points become accessible, deviating trajectories approaching them. These instabilities promote a slowing down of the dynamics and an island myriad bifurcation phenomenon, along with the suppression of long flights within the lattice. Other diffusion regime variations occurring within small intervals of control parameters are shown to be related to the emergence of a set of orbits with long flights, thus altering the total average displacement for long integration times but without global changes in phase-space.

Keywords: Diffusion; Periodic Lattices; Hamiltonian System; Chaos

I. INTRODUCTION

From a wide range of experimental contexts, periodic potentials have been used to model diverse physical systems, from optical lattices, where ultra-cold gases or Bose-Einstein condensates are confined by a set of laser beams [1–4], to ionic particles submitted to guided waves in \( \mathbf{E} \times \mathbf{B} \) fields and plasma physics [5, 6], to wave propagation in photonic crystals [7], and to xenon atoms diffusion over platinum surfaces [8].

These systems have also been used as models for the quantum-classical correspondence of chaos for many-body systems. As shown by Thommen [9], particles in a quantum tilted lattice can present a behavior similar to classical chaos. Prants [10, 11] proposes scenarios where the effect of chaos in particles displacement could be observed experimentally in a periodic potential with semiclassical effects when the field interacts with two-level atoms.

Particularly within the theoretical perspective of classical dynamical systems, periodic potentials commonly appear in Sinai billiards and hard wall scatterers models for Lorentz gases [12–14]. Analogous Hamiltonian models with smooth periodic potentials, the so-called “soft” billiards, were also considered, providing a good framework to study chaotic dynamics and control theory [15–17]. Such a feature is particularly relevant for the transport of particles through lattices, since anomalous diffusion regimes can occur without the application of random forces, purely due to inner instabilities in the system motion [5, 18–24].

Recent works on periodic potentials as those on soft billiards based on optical lattices display their mixed Hamiltonian phase-space structure and some scenarios for different parameters. We particularly mention Horsley et al., for a study on the same Hamiltonian model used here [25] (introduced in section II), although no considerations on diffusion and its dependency on the system parameters were included. Regarding classical aspects on quantum models, Porter et al. studied bidimensional (2D) hamiltonian lattices for different topologies, such as a honeycomb lattice [26], including a quantum treatment for the energy bands and the effect of symmetry breaking [27]. In [20], Prants studies a Hamiltonian model analogous to the one considered here but modelling a dipole-field interaction by including into the coherent dynamics the interaction of internal atomic degrees of freedom with translational ones.

Regarding diffusion itself, Zaslavsky performed some of the prominent works for continuous flows and Hamiltonian models considering the variation of control parameters [12, 21, 23]. Particularly in [23], for a periodic Q-model, were shown the existence of long flights within the chaotic motion, the consequent anomalous diffusion regime, the tail thickening effect in the power-law distribution, and the occurrence of sudden transitions in diffusion rate as a function of a control parameter. Similarly, Argonov and Prants [18, 19, 28] showed the fractal structure of escape time basins and the effect of chaos in different diffusion regimes due to Lévy flights in a fully coherent semiclassical model of a 1D lattice. Kleva [5] considered in detail the structure of chaos close to movable separatrix curves of a periodic potential displaying these same Lévy flights. The connection between these flights and anomalous regimes has been shown for a wide variety of models [12, 29].

From this diffusion background, the present work aims to a detailed description of the dynamics behind the
transport of particles in a lattice Hamiltonian system. A bidimensional periodic classical potential, based on an optical lattice system, is used to numerically study transitions in the spatial diffusion as the main control parameters change, namely the total particle energy and the coupling between the waves that form the lattice. Generally, for either varying energy or coupling, the disruption of stability islands and further bifurcation of periodic orbits (PO), both stable (SPO) and unstable (UPO), results in a non-trivial dynamics with mixed phase-space. Due to such bifurcations, the system’s spatial diffusion regime, measured by a long time exponent rate, undergoes sudden variations between normal and ballistic regimes or even peaks of intermediate rate with long transient times. The lack of pattern or predictability of such transitions thus motivates an inspection of the topological changes within phase-space. In order to conduct this inspection as the control parameters vary, a series of numerical techniques are applied, namely, the measurement of the ratio of chaotic/regular area within phase-space, the automated search for POs and identification of their stability and period, the manifolds from UPOs, and the escape time (or range) basins.

From the general finding of non-trivial changes in diffusion, a particular transition was studied in more detail. An increase (decrease) of stability (chaotic) area was found to correlate with a sudden diffusion change from ballistic to normal regime as energy increases. We verified that such behavior occurs when particles energy is enough to reach local instability points (potential local maxima), promoting a slowing down of dynamics and giving rise to new regular structures amidst a web-like manifold pattern in phase-space. This myriad of islands soon vanishes as energy increases and is replaced with a global chaotic sea, determining a more restrained chaotic dynamics with suppressed long flights through the lattice. For diffusion variations uncorrelated with changes in chaotic/stable areas, increments in the particles average displacement are found to be caused by the onset of a set of long flight orbits. These transitions exist for a short energy interval and represent a small part of the total stable area but still alter the average displacement calculation as a transient behavior, perceptible for long times ($t \propto 10^3$).

In this paper, sections II and III respectively present the lattice model used in this study and how the transport of particles is evaluated. The main findings are discussed along section IV, starting by the aforementioned diffusion transitions (section IV A) and followed by an initial analysis of the onset of chaos and transport in the system (section IV B). A particular transition is detailed in sections IV C and IV D by showing the phase-space changes and POs bifurcations. For completeness, in section IV E we also present results for diffusion variations not related to global modifications in the dynamics. Final remarks are given in section V. An appendix section is provided to enhance the discussion on integration error, statistical convergence, further PO analysis and extra results. Each topic is mentioned through the text whenever it may be of interest.

II. MODEL

The bidimensional periodic lattice model studied in this work stems from a classical Hamiltonian motivated by the trapping of a single neutral particle in the superposition of standing laser waves [1]. Experimentally, this setup is used as an optical lattice to trap and control cold atom gases or Bose-Einstein condensates, although its treatment is made quantically. In such an arrangement, an electrical field $E$ from a standing wave, obtained by the superposition of multiple counter-propagating monochromatic waves, induces a parallel dipole $d$ into a neutral particle. The re-interaction between field and dipole yields the potential

$$V_{\text{dip}}(r) = - d \cdot E(r) = - \sum_i (d \cdot \hat{e}_i) E_i^0 \cos(\mathbf{k}_i \cdot r + \theta_i),$$

with $\mathbf{k}_i$ as the wave vector and $\theta_i$ the phase of each individual standing wave; $\hat{e}_i$ and $E^0_i$ are the respective polarization direction and amplitude of the waves, and $r$ the particle spatial position. The induced dipole given by $d = \gamma(\omega)E$ holds for the case where the incident radiation frequency $\omega$ does not resonate with the particle inner energy levels, avoiding absorption and further re-emission of photons and thus ensuring a linear polarization by a polarizability factor $\gamma(\omega)$. This aspect allows for an approximation of spatial trajectories as classical. Alternatively, Prants [10, 11, 20] considered a semiclassical effect of particle-field interaction by allowing photons absorption and emission for particles with two inner energy levels.

The generic form of potential (1) allows the construction of different lattices by the combination of multiple waves and frequencies. For a 2D lattice, at least two linearly independent wave vectors are selected, constraining the particle movement along two cartesian axes, although an extra wave propagating in the same ($\mathbf{k}_1, \mathbf{k}_2$) plane can be placed in order to achieve different topologies, as done by Porter et al. for a honeycomb lattice [27]. Indeed, even non-harmonic fields can be used for the lattice construction, particularly if one aims to consider symmetry breaking effects [26].

Here, two orthogonal waves within the $x - y$ plane are taken with equal amplitude and wavelength ($\mathbf{k}_x = k_x \hat{x}; \mathbf{k}_y = k_y \hat{y}$), yielding the periodic potential for the lattice

$$V(x', y') = U' \left( \cos^2(kx') + \cos^2(ky') \right) + 2\alpha \cos(kx') \cos(ky'),$$

with

$$U' = \gamma(\omega)E^2_0 > 0 \quad \text{and} \quad \alpha = (\ddot{\varepsilon}_x \cdot \ddot{\varepsilon}_y) \cos(\theta_x - \theta_y).$$
Therewith, for a single particle the lattice Hamiltonian can be written as

\[ H = p_x^2 + p_y^2 + U(\cos^2(x) + \cos^2(y) + 2\alpha \cos(x) \cos(y)), \]

with space units scaled to \((x = kx'; y = ky')\) and Hamiltonian to \(H = 2mH'\), so that the energy scale is \(U = 2mU'\). The classical dynamics of a particle is thus described by the surface potential shown in figure 1, where a particle can be trapped in the field wells for energies lower than those of the saddle point between pits or otherwise diffuse through the lattice cells above this threshold.

![3D representation of particles (in green) over the lattice potential surface \(V(x, y)\) displaying its periodic structure.](image)

The energy scale \(U\) is of no relevance in the classical regime, in the sense that it does not alter the topology of solutions whatsoever, and can be set to 1 by rescaling time. In the quantum regime on the other hand, this energy scale relates to the accessible eigenstates and thus has further relevance. For this study, we fix \(U = 20\) following Horsley et al. [25] since it corresponds to a feasible value obtainable in experiments. In [20], Prants considered a similar Hamiltonian to (4) but taking into account the coupling of internal and translational degrees of freedom of the atom via a coherent semiclassical model for laser frequencies close to optical resonance.

As can be seen from potential (2), the coupling parameter \(\alpha\) stands for the product of the fields polarizations and phase difference, acting as a perturbation to the integrable Hamiltonian of two pendula potentials along \(x\) and \(y\) (with period \(\pi\)), coupling them for any \(\alpha \neq 0\). Although \(\alpha\) values may vary within \([-1, 1]\), one can notice that it is only required to consider solutions for \([0, 1]\), since the negative counterpart is equivalent to a spatial translation by \(\pi\) in one of the cartesian directions, thus not altering solutions properties.

Figure 2 shows how the periodic potential structure changes as \(\alpha\) increases from the separable case (\(\alpha = 0\)) to the maximum superposition amplitude (\(\alpha = 1\)). As the saddle points move towards the local maxima, they finally merge when \(\alpha = 1\), forming two trenches with degenerate minima on the lines \(y + x \equiv \pi \mod (2\pi)\) and \(y - x \equiv \pi \mod (2\pi)\) (see table 1). For increasing \(\alpha\), the path between lattice pits for particles to diffuse through lattice cells gets wider and wider, since the diffusion’s energy threshold at the saddle points is given by \(V^{\text{saddle}}(\alpha) = U(1 - \alpha^2)\).

![Color plot of the top view of potential \(V(x, y)\) minimal unit cell for different values of the coupling \(\alpha\). The PSS (Poincaré Surface Section) used for phase-space display is shown as the horizontal dashed green line at \(y = \frac{\pi}{2}\).](image)

| Equilibrium points \((x^*, y^*)\) | \(V(x^*, y^*)\) |
|-------------------------------|----------------|
| **Minima**                     |                |
| \((\frac{\pi}{2}, \frac{\pi}{2})\) | 0              |
| \((-\frac{\pi}{2}, -\frac{\pi}{2})\) | \(2U(1 + \alpha)\) |
| \((\frac{\pi}{2}, -\frac{\pi}{2})\) | \(2U(1 - \alpha)\) |
| \((-\frac{\pi}{2}, \frac{\pi}{2})\) |                |
| **Maxima (global)**            |                |
| \((0, 0)\)                    | \(2U(1 + \alpha)\) |
| \((\pi, \pi)\)                | \(2U(1 - \alpha)\) |
| **Maxima (local)**             |                |
| \((\pi, 0)\)                  |                |
| \((0, \pi)\)                  |                |
| **Saddle**                     |                |
| \((0, \pm \cos^{-1}(-\alpha))\) |                |
| \((\pm \cos^{-1}(-\alpha), 0)\) | \(U(1 - \alpha^2)\) |
| \((\pi, \pm \cos^{-1}(\alpha))\) |                |
| \((\pm \cos^{-1}(\alpha), \pi)\) |                |

TABLE I. Equilibrium points position \((x^*, y^*)\) and energy value \(V(x^*, y^*)\) within a unit cell of the periodic lattice; positions are taken with modulo \(2\pi\). At \(\alpha = 1\), saddle points merge with local maxima and form minimum trench lines given by \(\cos(y_0(x)) = -\cos(x)\).

For the results shown along this work, all phase-space portraits will be made over the same Poincaré surface section (PSS) – highlighted by green lines in figure 2,
namely the one defined as the oriented surface along two of the lattice pits, i.e.,
\[
\Sigma = \left\{ (x, y, p_x, p_y) \in \mathbb{R}^4 : \left( y = \frac{\pi}{2} ; \ p_y > 0 \right) \right\}. \tag{5}
\]

Since Hamiltonian (4) is autonomous, energy \((E = H)\) is an immediate constant of motion, constraining trajectories in a three-dimensional surface, which can thus be pictured by a 2D section. The oriented surface \(\Sigma\) is particularly convenient for this potential since it intersects all possible solutions within a single unit cell pit, except for the UPOs along the stable direction of the saddle equilibrium points located at \(y = 0\) and \(y = \pi\). Indeed, bounded solutions around the minima points with \(y < 0\) will occur, but nonetheless the \(\frac{\pi}{2}\) rotation invariance implies that their symmetrical counterpart solutions will intersect \(\Sigma\) at \(y = \frac{\pi}{2}\). The potential symmetry allows us to consider the motion in a restricted phase-space if one modulates trajectories with periodic boundary conditions \((x, y \in \left[ -\pi, \pi \right))\), but it also allows for the scattering or diffusion of particles if one lets spatial variables run freely.

The PSS \(\Sigma\) will also be used as reference for representing periodic orbits. Along any PSS, POs appear as fixed points with discrete period, i.e., an \(n\)-periodic orbit will cross the PSS \(n\) times until it return to its initial point. However, the number of these crossings depends on the orientation of the surface and how it intercepts the orbit path. Therefore, along the text we refer to discrete periods (labeled as \(T\)) only defined relative to \(\Sigma\) (5). Moreover, we differentiate the discrete period from the real-valued dynamical period (labeled as \(\tau\)), which is simply the total time elapsed for the orbit to returns to its initial point.

III. DIFFUSION CALCULATION

The quantitative characterization of the diffusion regime of an ensemble of particles through the lattice was straightly achieved from the asymptotic power law [23]
\[
\langle R^2 \rangle(t) \propto t^\mu, \tag{6}
\]
where \(R = \sqrt{(x-x_0)^2 + (y-y_0)^2}\) is the spatial displacement of a particle from its initial position, with the average taken over all particles. The exponent \(\mu\) thus indicates the diffusion rate, being normal \((\mu = 1)\), ballistic \((\mu = 2)\) or anomalous for different results within \(0 < \mu < 2\). One can immediately assert from energy conservation and the boundedness of the potential (4) that super-diffusivity \((\mu > 2)\) cannot occur in the system.

To numerically compute \(\mu\) for a given pair of parameters (\(\alpha, E\)), an ensemble of \(N\) random initial points covering the 3D energy shell in phase-space is evolved for long times \((t = 2000.0)\). From the time series data for \(\langle R^2 \rangle(t)\), one can retrieve the rate exponent \(\mu\) by fitting it with the power-law (6). Since the power-law is expected to hold true only asymptotically to long times, the data fitting is performed over the last 30% of time interval data. The random initial points are sampled as a Monte-Carlo procedure by randomly generating a position \((x_0, y_0)\) within the constraint \(V(x_0, y_0, \alpha) < E\) and a momentum vector \(p = (p_x, p_y)\) with random direction given by an angle uniformly distributed within \([0, 2\pi]\) and modulus defined by the conservation of energy (4).

Besides errors of statistical order, the long-time integration required raises the issue of numerical error in the solutions. A Runge-Kutta-Cash-Karp (RKCK) [30] method is used, and therefore it does not conserve the symplectic 2-form (even for regular solutions). However, energy deviations did not exceed the order of \(10^{-9}\), implying that the solutions obtained are very well bounded within the same energy hyper-surface, even though individual trajectories present small divergence from the real solution. This divergence is surmounted in this case since only the average over a uniformly filled phase-space is required, given that no stickiness was found to be relevant for the system’s dynamics. Nevertheless, a direct comparison to a symplectic method, developed by M. Tao [31], was made and is inserted in appendix B, showing that the lack of symplecticity of the RKCK method does not impact the results obtained here.

IV. RESULTS

A. Diffusion exponent

The study of the transport of particles through the lattice is made by the procedure described in section III while varying the main control parameters, i.e., the total energy of the particle \((E)\), and the coupling between radiation waves \((\alpha)\). Figure 3 shows the profile found for different \(\alpha\) values and varying energy. When the system is integrable \((\alpha = 0)\), it is spatially separable and, for energies allowing for diffusion, its behavior is completely ballistic, since this case is similar to a pendulum with rotation energy, moving unimpeded through the lattice. When integrability is broken \((\alpha \neq 0)\), generally the diffusion exponent \(\mu(E)\) undergoes a series of short and sudden transitions, with changes from normal \((\mu = 1)\) to free \((\mu = 2)\) regimes occurring abruptly and intermediate variations to intermediary regimes \((1 < \mu < 2)\), mostly seen as sharp peaks or valleys.

As exposed in section II, the coupling \(\alpha\) relates to the broadening of spatial channels through which a particle travels to neighbour lattice pits (stable minima points). Both the saddle points between the pits and local maxima decrease in energy, thus allowing a wider space for flights. Thus, one could hope to primarily find a continuous transition for the transport exponent as a function of both \(\alpha\) and \(E\). However, inner instabilities from chaotic trajectories and a mixed phase-space, with the emergence or destruction of Kolmogorov-Arnold-Moser (KAM) islands (PO bifurcations), imply a more intricate profile for
the particles diffusion and its exponent $\mu$. Indeed, when conducting the equivalent diffusion calculation for fixed energies and $\alpha$ as the free parameter, similar behavior is found (results shown in appendix C), given that changes in dynamics occur in a similar way, namely the bifurcation of stability islands or UPOs, for the variation of both control parameters.

For energies close to global maximum points ($E \approx V_{g\text{max}} = 2U(1+\alpha)$), a common plateau at ballistic regime occurs for all values of $\alpha$. At this energy level, the particle dynamics covers a wider space region with large momenta, where instabilities from the coupling can be seen as small perturbations. Therefore, the particle movement is that of a point moving rapidly through the lattice with smaller deviations, thus yielding a ballistic diffusion regime. Furthermore, this is also verified by an emergent dominance of large chains of stability islands in phase-space.

Given the statistical reliance on the calculation for $\mu(E)$, an immediate concern with its convergence and error is raised, particularly regarding the total integration time $t$ and ensemble size $N$. Details on this error analysis are given in appendix D, while here we summarize that the transitions undergone by $\mu(E)$ are found for either time and ensemble size of different magnitudes, whereas absolute deviations in their values do not exceed 15%. However, transitions composed of single thin peaks or intermediary rate values ($1 < \mu < 2$) may correspond to long transient behavior due to small stability islands composed of long flights, slowly converging towards ballistic rate $\mu \to 2$, although only manifested for longer times; a further discussion is made in section IV E.

FIG. 3. Diffusion exponent $\mu(E)$ for different values of the coupling $\alpha$. Any non-integrable scenario ($\alpha \neq 0$) presents sudden transitions between diffusion regimes. Energy scales range ($E_{\text{min}} \to E_{\text{max}} = (0 \to 2U(1+\alpha))$).

For energies close to global maximum points ($E \approx V_{g\text{max}} = 2U(1+\alpha)$), a common plateau at ballistic regime occurs for all values of $\alpha$. At this energy level, the particle dynamics covers a wider space region with large momenta, where instabilities from the coupling can be seen as small perturbations. Therefore, the particle movement is that of a point moving rapidly through the lattice with smaller deviations, thus yielding a ballistic diffusion regime. Furthermore, this is also verified by an emergent dominance of large chains of stability islands in phase-space.

FIG. 4. Normalized chaotic/regular area $A$ over the PSS $\Sigma$ for varying energy and different fixed coupling values calculated via SALI algorithm.

One may immediately expect that the transitions shown in figure 3 depend entirely on the mixed dynamics of the system’s solutions. Therefore, we firstly inquired whether these transitions correlate to the appearance or destruction of stability islands. For this purpose, the
chaotic area was measured over the PSS Σ (5) via Smaller Alignment Index (SALI) method [32, 33]. Briefly putting it, the algorithm considers two deviation vectors \((\hat{\omega}_1, \hat{\omega}_2)\) which are integrated by the first order variation of motion equations along with a reference orbit (their norms being kept constant). These deviation vectors thus behave differently for each kind of orbit. In case the reference orbit is chaotic, they align or anti-align to each other since they orient themselves towards the unstable manifold direction, whereas for a regular orbit they are kept at finite angle while only orienting themselves towards the tangent plane of the stability torus in which the orbit is contained. It is possible that for regular orbits the tangent vectors still align/anti-align due to shear between close torus layers, however this may occur for times much longer than the one for alignment in chaotic orbits.

From this principle, the defined index

\[
\text{SALI}(t) = \min (\|\hat{\omega}_1 + \hat{\omega}_2\|, \|\hat{\omega}_1 - \hat{\omega}_2\|)
\]

provides a quantitative way to discriminate the orbit’s nature, given that aligned vectors imply SALI\((t) \to 0\) while regular ones keep SALI at constant non-zero values (assuming non-parallel initial vectors). The algorithm does not rely on any particularity of the present model, thus being applicable in general to Hamiltonian systems or symplectic discrete maps.

Using this index to differentiate between regular and chaotic orbits, one can mesh the surface \(\Sigma\) and sum over the sub-areas from each initial condition assigned to each grid cell. Equivalently, the same procedure could be made for a 4D grid over the whole phase-space. However, the selected PSS intersects all possible orbits within a unit cell, except the UPOs along the saddle equilibrium points at \(y = 0\) and \(y = \pi\), thus ensuring that the section provides a good portrait of the chaos/stability ratio of the whole phase-space. Figure 4 shows the area portions of orbits along \(\Sigma\) normalized by the total area \(A_{\text{tot}} = A_{\text{chaos}} + A_{\text{regular}}\) for the same coupling values displayed in figure 3.

Figure 4 shows a correlation between a main transition for \((\alpha = 0.1; E = 36)\), where a sudden increase of stable area is found to occur, and the transition in diffusion exponent \(\mu(E)\) shown in figure 3, where a free transport plateau drops to normal regime. On the other hand, a series of peaks in diffusion rate, either from increasing or decreasing \(\mu(E)\), do not correlate with any pronounced changes in chaotic/stable area. This is particularly seen along the intervals at \((\alpha = 0.5; E \in [30, 50])\) and \((\alpha = 1.0; E \in [30, 60])\) where the system is dominantly chaotic along the whole energy interval although peaks of diffusion transition are seen for \(\mu(E)\). Even though no direct implication requires the diffusion exponent to correlate directly with the chaotic area, it is compelling to check whether the transitions found are related to the emergence of chaos or inner changes in its domain, such as the appearance of Lévy flights, which may not alter the area but change chaos properties. Given these two scenarios for transport variation, we aim to inspect the phase-space for energy values around these transitions and characterize the dynamical changes occurring along with it.

B. Diffusion onset

As previously mentioned, the diffusion transition occurring at \((\alpha = 0.1; E = 36.0)\) comprises a sudden change from ballistic \((\mu = 2)\) to normal \((\mu = 1)\) regime as seen from figure 3, correlated with a simultaneous drop/growth of chaotic/regular area. At the particular energy \(E = 36.0\), the potential surface does not undergo any sudden change but its local maxima points, located in the unit cell at \((x = \pm \pi, y = 0)\) and \((x = 0, y = \pm \pi)\), become accessible to trajectories. Counterintuitively, although spatially the path for transport widens, these new equilibrium points act as an instability source, changing the chaotic dynamics properties by disrupting long flights through the lattice. It is therefore interesting to use such

![FIG. 5. PO bifurcation process for increasing energy seen from the PSS Σ (left column), with fixed points drawn as colored circles, along with its correspondent spatial trajectories in the right column. Each PO is named after its stability – stable (S) or unstable (U) – and an index, related to its color. In the portraits, top row has \(E = 5\), center row \(E = 9\) and bottom row \(E = 17.6\); in all of them \(\alpha = 0.1\).]
an orbit as reference for changes in the system as energy increases.

The PO along the local maxima direction exists for all energy values \( E > 0 \), becoming unstable for \( E \approx 6.9 \) \((S_g \rightarrow U_g)\) whilst branching into two other SPOs \((S_r\) and \(S_y)\), as shown in figure 5. This bifurcation process is the first considerable emergence of chaos in phase-space, as the separatrix around the new islands (related to \(S_r\) and \(S_y\)) is disrupted and replaced by a chaotic layer with hyperbolic point at \(U_g\) fixed point. In figure 6, portrait A shows that for energy values slightly above the diffusion onset \((E = 22)\), the chaotic layer increases and forms a connected piece, surrounding the three main stability islands. One may notice that the bottommost island, for \(\frac{p_x}{\sqrt{E}} \approx 1\), with \(S_r\) at its center, is highly compressed along the energy shell border, but still exists.

Using the UPO \(U_g\) and the ones along the saddle equilibrium points \((U_x\) and \(U_y\), shown in portrait B of figure 6), their manifold structure indicates that for transient times the chaotic region is separately occupied by an outer layer, seen in portrait C as the unstable manifold of \(U_g\) (in red) and an inner layer, from the stable manifold branch of \(U_g\) (in green). These initially separated regions occur due to small turnstiles between the the manifold branches, although they appear infinitely many times for long periods, filling the whole chaotic sea as a single connected region. The unstable manifold is mostly located along regions of high \(p_x\) momentum, related to direct flights travelling horizontally along the lattice. Due to the \(x \leftrightarrow y\) symmetry, one can find the same division for vertical flights, from the unstable manifold of \(U_x\) (figure 6, portrait D), where it is placed inside the innermost lobe of \(U_g\) manifold, corresponding to high values of \(p_y\).

### C. Local maxima transition

In order to detail the transition at \(E = 36\) and \(\alpha = 0.1\), phase-space portraits over the PSS \(\Sigma\) (5) are given

![FIG. 6. A) PSS \(\Sigma\) for \(E = 22\). B) UPOs used for the calculation of the manifolds displayed at frames C and D. C) Stable (green) and unstable (red) manifolds from \(U_g\) and \(U_x\) respectively. D) Stable (green) and unstable (red) manifolds from UPO \(U_y\) and \(U_x\) respectively. In frame B, the blue dashed line depicts the spatial location of the PSS \(\Sigma\) and the blue dot in the remaining frames the fixed point from UPOs within it.](image)

![FIG. 7. Phase-space portraits for \((\alpha, E) = (0.1, 32)\) with ballistic diffusion regime, before transition. (Top) PSS with orbits crossing points in black. (Center) Escape time basins color map. (Bottom) Stable manifolds from \(U_g\).](image)
for energy values before \((E = 32.0 - \text{Fig. 7})\), around \((E = 36.1 - \text{Fig. 8})\) and after \((E = 38.0 - \text{Fig. 9})\) the transition. These figures show the PSS with its orbit crossings (black points), the stable manifold from the \(U_g\) UPO (Fig. 6), and a color map of escape time basins. In this context, escape time basins are simply defined as a map of the time required for initial conditions on the PSS \(\Sigma\) to reach outside the square box with \(n\) unit cells of size, \(i.e., x, y \in [-n\pi, n\pi]\) (here \(n = 10\)).

Before the transition, for \(E = 32\) (Fig. 7), phase-space stability regions are still given by the three main islands with center given by \(S_y, S_r\) and \(S_b\) SPOs (Fig. 5). As indicated by the escape time basins, the bottom islands \((S_y\) and \(S_r\)) are related to direct flights through the lattice, given their small escape time. Besides, channels of low escape time infiltrate the chaotic sea by stretching themselves from the instability region. These channels are also visible in the manifold portrait, while penetrating between the finger-like structure of the folded lobes that delimit the island. The upper island, related to orbits bounded to inside the unit cell, is the only set with “trapped” trajectories, whereas the chaotic sea alone presents an average escape time around \(\langle t \rangle \approx 20\). As expected, the escape time basins reflect the manifolds structure, since these are the invariant lines that approach/depart the PSS single fixed point from the UPO \(U_g\). Although not shown in figure 7, the unstable branch is perfectly symmetric, under the reflection \(y’ = y; x’ = \pi - x\), to the stable branch. The men-

**FIG. 8.** Same as figure 7 but for \((\alpha, E) = (0.1, 36.1)\) energetically slightly above the transition.

**FIG. 9.** Same as figures 7 and 8 but for \((\alpha, E) = (0.1, 38)\) after the transition, at normal diffusion regime.
tioned channels may fade for higher box sizes or integration time, however they imply a transient behavior lasting for at least hundreds of time units, thus considerably long.

While the system exceeds the transition energy ($E = 36$), the manifold finger-like lobes fold into the bottom stability islands (from $S_r$ and $S_p$), merging exactly at the fixed point position of the UPO $U_g$ with the stable regions vanishing in the process (figure 8). At this point, a myriad of islands emerge from the chaotic sea, forming centered chains around the unstable fixed point related to $U_g$, located at $(x, \frac{p}{\sqrt{E}}) \approx \left( \frac{\pi}{2}, -0.71 \right)$ and also around the upper stability island related to $S_b$. These structures last for a narrow energy interval, approximately $E \in [36.0, 36.3]$. For the group around $S_b$ island, a higher escape time (constrained orbits) can be seen, indicating a stickiness behavior for this region. The bottom chain structure, around $U_g$, presents an alternated layered structure (indicated by the yellow and purple islands in figure 8). Every layer has even discrete period, with each one increasing its period by 2 as they grow englobing the inner layers, forming an onion-like structure with the web-like manifold folding through them. More details on the alternating escape time range of this structure are given in appendix A, where the SPOs related to the chains are shown. Briefly, this alternation occurs due to the spatial “closure” of SPOs at the center of yellow chains, i.e., they return to their initial point, considering an unbounded dynamics, without periodic boundary conditions (PBC), thus having limited range of transport, whereas the chains with fast escape time (in purple) are related to SPOs that only close themselves when one considers spatial PBC, meaning that when unbounded, they travel as direct flights through the lattice. Moreover, these islands are discrete isochronous, being formed by multiple SPOs, with either odd or even discrete period.

After the transition, with the vanishing of the island myriad, a uniform chaotic sea mostly invades all phase-space, with only the upper island still remaining but in reduced size (figure 9). Amidst this chaotic region, the previous manifold web structure loses its “horizontal” tangling lines, leaving only a swirling structure, with a “knot” at $(x, p_x) = (0, 0)$. The vanishing of the two previous bottom islands and their channels within the chaotic region contributes to the suppression of long flights and therefore transport in the system ceases from being ballistic.

### D. Periodic orbits bifurcations

In order to analyse POs bifurcations and their modification in phase-space, a period-energy diagram was calculated with a monodromy algorithm. As developed by Baranger et al. [34] and further detailed by Simonović [35], one can obtain a periodic orbit, either stable or unstable, from a given initial condition attempt and iteratively applying a Newton-Raphson algorithm to approach a periodic solution. This technique is generically applicable to any conservative Hamiltonian system of $N$ degrees of freedom or symplectic map. Running this procedure extensively along the PSS $\Sigma$, a series of POs can be found in an automated way for different energies, allowing the construction of a diagram displaying the dynamical period $\tau$ of the orbits found as a function of the energy value.

The result for such a diagram calculated for $\alpha = 0.1$ is shown in figure 10, where POs with discrete period $T = 1, 2$ and 3 on the section $\Sigma$ are considered. The horizontal dashed lines mark two key energy values, namely the diffusion onset ($E = 19.8$) and the potential local maxima ($E = 36.0$). As the energy approaches either one of these values, a slowing down of the dynamics occurs, with the dynamical period of orbits asymptotically diverging as they reach the horizontal lines. This is easily understandable as the case of orbits reaching an unstable equilibrium point for the exact energy value of access, taking an infinite time span to reach it, analogous to the dynamics of a simple pendulum at the exact libration-rotation threshold. In this case, the saddle point between energy unit cells marks the energy for diffusion onset and the local maxima to the transition we studied here. Since the discrete periods considered in the diagram are low ($T = 1, 2, 3$), the diagram does not display the island myriad appearance, given its higher period chains. However, it can be seen that the presence of an unstable point, even though it spreads trajectories and introduces chaos, is also related to a slowing down of the dynamics and the vanishing of a series of POs, along with the creation of completely new ones. Therewith, the scenario for the transition studied here is that the effect of a new instability source suppresses long flights within

![FIG. 10. Period energy diagram for orbits with PSS discrete period $T = 1, 2$ and 3.](image-url)
the lattice while it permeates the phase-space with chaos of more constrained trajectories.

E. Local flights

To fully explore the aspects of the diffusion profiles obtained here (figure 3), we briefly describe in this section the short peaks and valleys comprising sudden changes in them. These are variations in diffusion exponent uncorrelated with considerable changes in chaotic area, as exemplified by the energy intervals around $E \approx 24$ and $E \approx 24.5$. 

FIG. 11. Displacement range color map for different energies around the sudden peak for $E \approx 24$. The total integration time is $t = 1000$ for each point in a $850 \times 830$ grid.

FIG. 12. PSS portraits of the stable manifold branch from UPO $U_\alpha$ for different energies along the diffusion variation. Integration time is $t = 6.9$. 

As we did previously, we look at phase-space portraits for energies in these intervals searching for alterations in transport characteristics. Therewith, to clearly display regions in phase-space with long flights, we use a displacement color map, similar to the escape-time ones but instead coloring each initial point with its displacement for a given integration time \( t \) (here \( t = 1000 \)). Figure 11 exemplifies this result for the transition at \( E \approx 24 \) and \( \alpha = 0.5 \), where a peak from normal diffusion to an intermediary value (\( \mu \approx 1.3 \)) is found. It can be seen that the only modification is the emergence of three small islands (shown in dark blue) around each of the twin islands (shown in light red) in the bottom region of the PSS, along with the bifurcation of these twin islands. The emergent island triads that appear present a very high displacement when compared to the chaotic sea surrounding them or the other stability islands composed of trajectories with limited range. It is clear then that the ballistic transport due to these small islands increases the global average diffusion rate, making it increase quadratically but with smaller amplitude, thus taking longer times to increase \( \mu \) but still converging towards \( \mu \approx 2 \). Although not shown here, the same behavior is seen for the other peaks and variations in transport rate mentioned that are not correlated with changes in chaotic/regular areas.

To emphasize that the variations for the energy values mentioned above are not global transitions, in the sense that they are not related to major changes in the chaotic domain, figure 12 shows the manifold structure for the case discussed above (\( E \approx 24, \alpha = 0.5 \)). Even though the island triads are not visible in the figure, one may notice that they appear between the manifold lobes without disturbing them. After the island triad disappears, for the highest energy value \( E = 24.5 \), the new bifurcated islands from the bottommost twin islands present a small stickiness around them. Here this is indicated by the region unfilled by manifold lobes where these islands previously existed. However, this behavior is not very pronounced and is not related to the increase in displacement seen for the transport rate, given that the bifurcated island pairs are composed of orbits with limited range.

The effect of ballistic islands over the global diffusion rate is in agreement with the diffusion profiles simulated for different integration times. Whenever these islands are present, as the evolution time increases, the diffusion regime slowly converges towards ballistic, as shown in appendix D. Therewith, we point out that, for the diffusion evaluation method used here, the presence of even small portions of phase-space with quadratic rate will imply a dominance over the total regime rate. Therefore the method use requires care for analysis of long times; however, it still serves as an indicator of the existence of long flights, given that they are indeed effects of the system dynamics. It is also interesting that islands with long displacement range appear amidst a chaotic sea with transport in a different regime, without major changes in the sea.

V. CONCLUSIONS AND PERSPECTIVES

This work presented a study of classical transport of particles for a 2D lattice model based on the periodic potential resulting from a dipole-field interaction. It was shown that as the system control parameters change, namely the coupling \( \alpha \) and total energy \( E \), the diffusion exponent \( \mu \), evaluated from an asymptotic law, presents sudden variations between ballistic and normal regimes due to the mixed nature of the system dynamics and its series of POs bifurcations.

A sudden drop of diffusion rate from ballistic (\( \mu = 2 \)) to normal (\( \mu = 1 \)) regime is found to occur in correlation with a decrease in chaotic area for \( (\alpha = 0.1, E = 36.0) \). A detailed analysis indicated that this transition occurs at an energy level for which orbits reach local maxima points. Even though the transport of particles seems facilitated by the wider spatial channel, the new instability source promotes a large bifurcation process with the emergence of multiple stable structures. Moreover, a global slow-down of the dynamics occurs due to orbits reaching the unstable equilibrium position, as seen in a period-energy diagram for periodic orbits.

Before the transition, long flights occur due to two main stability islands that vanish as \( E \rightarrow 36 \). At transition, phase-space becomes populated by an island myriad amidst a web-like manifold structure, with multiple isochronous chains of even discrete period. These chains form layers with increasing period and alternate displacement range in the lattice due to SPOs with closed or open spatial periodic topology. After a narrow energy interval (\( E \approx 36.3 \)), the island myriad vanishes and the remaining phase-space is dominated by a single chaotic region with long flights suppressed, presenting a normal diffusion regime. This characteristic transition may occur on the system for any coupling \( \alpha \neq 0 \), although it shall be more pronounced for small values, since local maxima energy levels decrease as a function of \( \alpha \), thus being less influential in the dynamics.

Diffusion variations that do not present correlation with chaotic area are shown to be caused by the emergence of small stability islands composed of long flights rather than global changes in dynamics. These long flights increase the average displacement towards ballistic rate but appearing only for long integration times, given their small weight within the orbits ensemble.

Further investigations can be made on the island myriad structure; it is possible that this phenomenon may occur due to a superposition of a fast dynamics, related to bounded orbits oscillating inside a single lattice pit, and a slow dynamics, related to the period divergence of UPOs reaching the unstable point \[36\]. Moreover, different lattice topologies could be experimented and the local maxima transition studied for a different symme-
try, such as in a honeycomb lattice, in order to analyse changes on the spatial closeness of the arising POs of such transitions. Further investigations could include an analysis dedicated solely to the chaotic dynamics and its main UPOs, better detailing the influence of the unstable point deviation in bifurcations of POs that suppress long flights. Alternatively, Hamiltonian perturbations feasible within experiments, such as potential amplitude variation, extra monochromatic waves or noise, could be applied and compared with the conservative case shown here, in order to enhance the control of particles in the lattice.

ACKNOWLEDGMENTS

M. Lazarotto would like to acknowledge Vitor M. de Oliveira for fruitful discussions and comments, particularly on manifold calculation and periodic orbits bifurcations. We acknowledge the financial support from the scientific agencies: São Paulo Research Foundation (FAPESP) under Grant No 2018/03211-6; Coordenação de Aperfeiçoamento de Pessoal de Nível Superior (CAPES) and Comité Français d’Evaluation de la Coopération Universitaire et Scientifique avec le Brésil (COFECUB) under Grant CAPES/COFECUB 8881.143103/2017-1; COFECUB under Grant No. 40273QA-Pb908/18; Conselho Nacional de Desenvolvimento Científico e Tecnológico (CNPq) under Grant Nos 407299/2018-1, 302665/2017-0 and 141750/2019-7.

APPENDIX

This appendix presents supplementary material to the main article. It is meant to provide extra results to the ones shown in the main text and details on the numerical methods used for diffusion calculations. In section A we exemplify the difference between the island chains layers that form the myriad structure by showing the different periodic orbits related to each one. Section B shows the comparison between the diffusion calculation with orbits integrated with a Runge-Kutta method and a symplectic method, ensuring that the average displacement used...
for the calculation of \( \mu \) is accurately made. Section C presents extra results of diffusion profiles calculated for fixed energy values and \( \alpha \) as variable parameter, as opposed to the cases shown in the main text. Section D discusses the statistical convergence of the diffusion rate calculation and the dependence of the values of exponent \( \mu \) on the simulation parameters.

Appendix A: Periodic orbit analysis

As shown by the escape time basins of figure 8, amidst the emergent myriad structure, island chains with even discrete period present an alternated layer structure of high (in yellow) and low (in purple) escape times. The new emergent SPOs are of two different kinds: one related to orbits with spatial closure, in the sense that they return to their initial position even when PBC is disabled, thus forming closed loop paths; the other group comprises orbits without spatial closure, meaning that when PBC is not considered, they evolve with repeating patterns without returning, thus forming long flights through the lattice. Therewith, it becomes clear that closed SPOs will have limited range and therefore higher escape time (or no escape at all). The displacement range of these orbits may increase with their discrete period but will still remain bounded. Open SPOs on the other hand will travel unboundedly through space, in an approximately ballistic way.

Figures 13 and 14 respectively display closed and open SPOs belonging to the myriad chains. It is worth mentioning that all these orbits are perfectly periodic when considering PBC. The periodicity of open SPOs is allowed in this case due to the periodic nature of the potential itself, given that a particle can return to a symmetric point in a neighbour cell, thus repeating the same dynamical evolution. Besides, one may notice that every orbit will have a symmetric counterpart, obtained by a rotations of \( \pi \), with the same properties regarding stability (Lyapunov exponent), period and “closedness” but with different discrete period, given the different possibilities of intersection with the PSS \( \Sigma \).

Moreover, the island chains are isochronous, implying that for an \( N \)-periodic chain, their “links” are not sequentially populated by a single SPO as usual, but instead they are composed of multiple SPOs with discrete period of divisors of \( N \), with fixed points alternating between the links [37]. This may be due to the rotation symmetry of the lattice, making periodic solutions to occur in pairs and therefore have the same winding number. Therefore they will occur superposed in phase-space and form the isochronous chain structure. As shown here, since the PSS is restricted to \( x \in [0, \pi] \), some fixed points of these SPOs may be located in \( x \in [-\pi, 0] \), where the PSS map is identical to the one seen for \( x > 0 \) but inverted, given the reflection translational symmetry of \( \pi/2 \).

Appendix B: Symplectic integration

In order to assess the quality of integration over long times required in diffusion exponent calculations, figure 15 shows the comparison between two methods: the Runge-Kutta-Cash-Karp (RKCK) and a symplectic one developed by M. Tao [31]. The particles ensemble is randomly generated in each case, but all the remaining simulation parameters are equivalent, i.e., the total integration time \( t = 1000 \) and ensemble size \( N = 10000 \). The final values for \( \mu(E) \) present a divergence of less than 0.15 when compared between the two methods. This indicates that even though RKCK method is not symplectic, it still yields a good average result, mostly due to its great energy conservation, which in this case was conserved up to order \( 10^{-9} \), regardless of deviations in particular trajectories. Besides, RKCK method presents faster CPU performance, which allowed the extensive calculations required in other analyses shown in this work.

As an adaptive time-step method, the RKCK absolute and relative precisions were selected as: \( \epsilon_{\text{abs}} = \epsilon_{\text{rel}} = 10^{-13} \). For Tao’s method, time-step is fixed and was selected as \( dt = 10^{-5} \). Also, Tao’s method requires a binding factor parameter, which was set to \( \omega = 100 \), since it provided the best performance regarding energy conservation (up to \( 10^{-5} \)) and symplectic 2-form (up to \( 10^{-7} \)).

![FIG. 15. Comparison between diffusion exponent \( \mu(E) \) calculation with RKCK method and Tao’s symplectic method.](image)

Appendix C: Diffusion transitions

Figure 16 shows the calculation of diffusion exponent \( \mu \) as a function of the coupling parameter \( \alpha \) and fixed values of energy \( E \), analogous to the results shown in the
main text. The obtained profiles show the generality of the sudden transitions phenomena, showing that the variation of either $\alpha$ or $E$ generate the same non-trivial variations. One may also notice that for considerably large energies (here exemplified for $E = 60$), diffusion regime is mostly ballistic, with a sudden change to normal regime for $\alpha \approx 1.0$. As pointed out in the article, for high energies or small perturbations, the system behaves as two uncoupled pendula with rotation energy, with little impediment to diffusion. As the coupling reaches its maximum value, its perturbation effect becomes sufficient to influence the dynamics, since global maxima become energetically higher (since $V_{\text{max}} = 2U(1 + \alpha)$) than the energy level of the particle. In this condition, the particle is more deflected as it moves towards maxima unstable points, instead of passing through them as it would if its energy was higher. These strong deflections thus may induce normal diffusivity through the lattice.

Appendix D: Statistical convergence

In order to ensure the statistical reliability of the calculated diffusion profiles, this section discusses the dependence of the values of $\mu$ with the total integration time $t$ and the ensemble size $N$. This is done by comparing the simulation in the best case scenario (highest $t$ or $N$) with multiple other ones with lower values of $t$ and $N$. As mentioned, the best case scenario here corresponds to the parameters used in the results presented in the main article, namely $t = 3 \times 10^3$ and $N = 3 \times 10^4$. It can be asserted then that the fluctuations and transitions found in the diffusion rate as a function of either $\alpha$ or $E$ are not due to statistical imprecision, but indeed caused by changes in the dynamics. It can also indicate that the values selected for $t$ and $N$ ensure good convergence of the final result.

Figure 17 shows the plot for the difference $\Delta \mu = \mu_N - \mu_n$ between different values of ensemble size $n$ and the best case for $N = 3 \times 10^4$. In this case, $t = 3000$ and $\alpha = 0.1$. 
\[ \Delta \mu = \mu_T - \mu_t \] 

between different values of integration time \( t \) and the best case for \( T = 3000 \). In all cases: \( \alpha = 0.1 \) and \( t = 3 \times 10^3 \). One can notice that the amplitudes of \( \Delta \mu \) for different \( n \) did not show to be large, with the average difference being of order \( \Delta \mu \approx 0.05 \). A particularly higher deviation occurs around the energy value for diffusion onset, corresponding to \( E = 19.8 \) for \( \alpha = 0.1 \), although this is expected given the absolute amplitude variation of \( \mu \) itself. Even though the test presented here comprises only a single value of \( \alpha \), it can be expected that for other values this result still holds. One may also notice that simulations could be optimized by selecting \( N \) at least one order of magnitude smaller than the value used in the main article, with the same qualitative final result.

Similarly, figure 18 shows the same procedure but for different integration times. Here the comparison is made as \( \Delta \mu = \mu_T - \mu_t \), with \( \mu_T \) obtained for the longest integration and \( \mu_t \) for any smaller value. The results indicate a higher amplitude variation for the different \( t \), but still below 0.15 in average. As for the case of different \( N \), the same sensitivity is shown for energy values around transitions or high fluctuations of \( \mu \). Nonetheless, the data for different integration times present a trend of increasing values of \( \mu(E) \) towards ballistic regime as the integration time increases. This is due to the presence of particles moving at ballistic rate that dominate the average displacement of the ensemble. However, the same result points out that convergence is close and the results found here are valid, given the small amplitude for \( \Delta \mu \).

In general, regardless of the sensitivity of \( \Delta \mu \) around regions with high variation of \( \mu \), one can see that they occur at the same energy values. This allows us to assert that the transitions found in the exponent \( \mu \) as we change the system parameters are not mere statistical variation artifacts. Also, some fluctuations become apparent just after a long transient; in fact they are not transitions properly said, in the sense that topological changes occur in phase-space, but just a long time transient behavior due the appearance of small islands composed of long flights that end up dominating the ensemble average displacement.

---

[1] I. Bloch. Ultracold quantum gases in optical lattices. Nature Physics, 1:23–30, 2005.
[2] I. Bloch; J. Dalibard and W. Zwerger. Many-body physics with ultracold gases. Reviews of modern physics, 80:885–964, 2008.
[3] A. Hemmerich; D. Schropp Jr. and T. W. Hänisch. Light forces in two crossed standing waves with controlled time-phase difference. Physical Review A, 44(3):1911–1921,
