Quantum scar affecting the motion of three interacting particles in a circular trap

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We theoretically propose a quantum scar affecting the motion of three interacting particles in a circular trap. We numerically calculate the quantum eigenstates of the system and show that some of them are scarred by a classically unstable periodic trajectory, in the vicinity of which the classical analog exhibits chaos. The few–body scar we consider is stabilized by quantum mechanics, and we analyze it along the lines of the original quantum scarring mechanism [Heller, Phys. Rev. Lett. 53, 1515 (1984)]. In particular, we identify towers of scarred quantum states which we fully explain in terms of the unstable classical trajectory underlying the scar. Our proposal is within experimental reach owing to very recent advances in Rydberg atom trapping.

I. INTRODUCTION

The thermalization of closed interacting quantum systems [1] may be impeded by various mechanisms [2, 3] whose investigation is strongly motivated by contemporary applications [4, 5]. Indeed, slowly–thermalizing systems retain memory of their initial state over longer times [6], making them useful for quantum simulation [4] and quantum information processing [5]. Atomic systems are an excellent test–bed for chaos [7–9], and techniques for the individual manipulation [10] of Rydberg atoms [11] have extended its exploration to interacting systems. A recent experiment on Rydberg atom arrays [12] has initiated the investigation of weak ergodicity breaking in many–body systems [13, 14]. Systems exhibiting this phenomenon thermalize rapidly for most initial conditions, but specific initial states yield non–ergodic dynamics. This behavior is analogous to the quantum scars initially predicted [15] and observed [16] in the absence of interactions, which also lead to weak ergodicity breaking [17] by impacting some [18, chap. 22] quantum eigenstates. Hence, it is also called ‘many–body scarring’ [19]. A similar phenomenon has been predicted in the context of the Dicke model [20], where the quantum scars are due to the collective light–matter interaction and impact many quantum eigenstates [21].

Despite the intense theoretical scrutiny [22], only two experiments [12, 23] and one explicit proposal [19] explore many–body scarring so far [12, 19, 23]. In all three cases, the observed non–ergodic behavior is linked to classical physics. The experiments of Refs. [12, 23] both probe the PXP model [24] in regimes where the classical analog system [23] explores the vicinity of classically stable periodic trajectories, so that the absence of thermalization may be traced back to the classical Kolmogorov–Arnold–Moser theorem [25, Sec. VI]. The proposal of Ref. [19] refers to spin helices in various geometries. Their classical limit is stable [27], and from the quantum point of view they generalize helices predicted [28] and observed [10] in the integrable XXZ chain. Hence, the proximity of integrable models is expected to play a key role.

In this article, we propose a three–body system hosting a quantum scar which relies on the interaction between particles. It may be realized experimentally owing to very recent advances in Rydberg atom trapping [29, 30]. It is simple enough to be fully analyzed by combining the numerical calculation of stationary states and well–established tools for the analysis of chaotic systems [31], in the spirit of Heller’s original proposal [15].

The system we consider exhibits “towers” of scarred states which are approximately evenly spaced in energy. These are a key feature of both quantum scars [32] and many–body scars [13, 14, 22, 23, 33]. In the present context, we explain them in terms of the classically unstable periodic trajectory causing the scar, in the spirit of Heller’s original argument [32, Fig. 22]. The phase space dimensionality of the few–body system we consider (4, see below) matches the maximum number of independent parameters introduced so far in the variational approaches applied to the many–body PXP model and its generalizations [26, §III.A]. In stark contrast to

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the many–body PXP model where approximate classical limits have to be cleverly constructed [25], our few–body system affords an exact reduction to four parameters and the identification of the classical analog is straightforward.

We formulate our proposal in terms of trapped Rydberg atoms [24, 30]. However, we expect other interacting systems with the same symmetries to exhibit similar quantum scars. We substantiate this claim in the Appendix (Sec. A.1) by identifying the quantum scar for systems with the same symmetries to exhibit similar classical ergodicity. We have characterized the periodic trajectories of the classically ergodic region.

II. THE CONSIDERED SYSTEM

We consider three identical bosonic particles of mass $m$ in a circular trap of radius $R$. The Hamiltonian reads:

$$H = \frac{l_1^2 + l_2^2 + l_3^2}{2mR^2} + v(d_{12}) + v(d_{23}) + v(d_{31}) ,$$

(1)

where $l_i$ is the component of the angular momentum of particle $i$ along the rotation axis, which is perpendicular to Fig. 1b. We assume that the interaction $v(d_{ij})$ between the particles $i$ and $j$ only depends on their distance $d_{ij} = 2R |\sin[(\theta_i - \theta_j)/2]|$. For circular Rydberg atoms whose electronic angular momenta are perpendicular to the plane, $v(d_{ij}) = C_6/d_{ij}^6$ with $C_6 > 0$ [30, App. A].

We introduce the Jacobi coordinates [38, §1.2.2] $x = [(\theta_1 + \theta_2)/2 - \theta_3 + \pi]/\sqrt{3}, y = (\theta_2 - \theta_1)/2 - \pi/3, z = (\theta_1 + \theta_2 + \theta_3)/3 - 2\pi/3,$ and their conjugate momenta $p_x, p_y, p_z$ (which carry the unit of action). In terms of these,

$$H = p_x^2/(3mR^2) + H_{2D},$$

(2)

$$H_{2D} = (p_x^2 + p_y^2)/(4mR^2) + V(x, y) .$$

III. CLASSICAL PHYSICS

We first analyze the classical dynamics described by the Hamiltonian $H$. Expressing momenta, energies, and times in units of $P_{ref} = (mC_6/R^4)^{1/2}, C_6/R^6,$ and $(mR^8/C_6)^{1/2}$, respectively, the classical results are independent of $m, C_6,$ and $R,$ leading to the scaled predictions in Figs. 1a–3. We choose the rotating reference frame such that $p_z = 0$ and $z = 0$. The divergence of $v(d_{ij})$ prevents the particles from crossing, so that we assume $\theta_1 < \theta_2 < \theta_3 < \theta_1 + 2\pi$ at all times. Hence, the classical problem is reduced to a point moving in the 2D plane $(x, y)$ within the equilateral triangle of Fig. 1a, in the presence of the potential $V(x, y)$.

We characterize the periodic trajectories of $V$ using our own C4+ implementation of the numerical approach of Ref. [41]. We find three families of periodic trajectories, existing for all energies $\epsilon > 0$: we label them...
A, B, C in analogy with the results for the HH potential \[12\]. We shall analyze them and their bifurcations in a forthcoming paper \[12\]. Here, we focus on family B, which yields the quantum scar. For a given \( \epsilon \), there are three trajectories of type B, due to the three–fold rotational symmetry of the potential \( V \). They are represented in the \((x, y)\) plane in Fig. 1b, and the one which is symmetric about the vertical axis is shown as a function of time in Fig. 2a. They are unstable for all energies, as shown by the Lyapunov exponent \( \lambda_3 > 0 \) in Fig. 2a. Figure 2b shows that trajectory B satisfies both conditions heralding a quantum scar: \( \lambda_3 T_B < 2\pi \) \[18, ch. 22\], and lower values of \( \lambda_3 T_B \) signal stronger scarring \[44\] \[§2.5\], so that the scar strengths associated with it for all \( E > 0 \) do not benefit from the classical enhancement due to the proximity of bifurcations \[44\]. This sets it apart from a previous proposal involving a scar hinging on this enhancement \[46\] so that, in stark contrast to ours, it is captured by Einstein–Brillouin–Keller quantization \[47\].

In order to visualize effects beyond the linear regime, Fig. 3 shows the surface of section \( \{ \psi(x, y) \} \) of \( H \) for \( \epsilon = 7C_6/R^6 \) and the conditions \( x = 0 \), \( p_x > 0 \) (allowing for a comparison with the HH potential \[48\]). It exhibits both non–ergodic regions comprising tori \[49\], App. 8] and an ergodic region, as is typical for a non–integrable system \[50, §1\]. The three fixed points corresponding to trajectories B are all located in the ergodic region. This precludes their stabilization by any classical mechanism.

### IV. QUANTUM PHYSICS

We seek the eigenfunctions of \( H \) in the form \( \Psi_n(\theta_1, \theta_2, \theta_3) = \psi_n(r) e^{i m z} \), where \( r = (x, y) \) and \( n = p_z/\hbar \). The wavefunction \( \psi_n \) is an eigenstate of \( H_{\text{HH}} \) with the energy \( \epsilon \). It is defined on the whole \((x, y)\) plane. Its symmetries are related to (i) angular periodicity, (ii) bosonic symmetry, and (iii) the point group \( C_{3v} \).

We first discuss (i) and (ii). (i) The \( 2\pi \)–periodicity of \( \Psi_n \) in terms of \( (\theta_i)_{1 \leq i \leq 3} \) yields \( \psi_n(r - BC) = \psi_n(r - CA) = \psi_n(r - AB) = \psi_n(r) e^{-i 2\pi n/3} \), so that \( n \) is an integer. (ii) The bosonic symmetry of \( \Psi_n \) leads to \( \psi_n(Sr) = \psi_n(r) \), where \( S \) is the symmetry about any of the lines (AB), (BC) or (CA) in the \((x, y)\) plane. Hence, we may restrict the configuration space to the inside of the triangle ABC of Fig. 1b. Along its edges, \( \psi(x, y) \) strongly diverges (e.g. \( \psi \approx (y + \pi/3)^{-6} \) near \( [AB] \)), so that \( \psi_n = 0 \) there. Combining (i) and (ii), and calling \( R \) the rotation of angle \( 2\pi/3 \) about \( O \), \( \psi_n(Rr) = \psi_n(r) e^{2i\pi n/3} \).

We analyze the role of the point group \( C_{3v} \). We classify the energy levels in terms of its three irreducible representations \( \rho = A_1, A_2, \) and \( E \) \[41, §95\]. Hence, Hilbert space is split into three unconnected blocks. These may be told apart through the behavior of \( \psi_n \) under two operations in the \((x, y)\) plane \[51\]: \( R \) and the reflection \( S_{\Delta} \) about the line \( \Delta = (CL) \) (see Fig. 1b). Wavefunctions pertaining to the 1D representations \( A_1 \) or \( A_2 \) satisfy \( \psi_n(Rr) = \psi_n(r) \), so that \( n = 0 \) modulo 3. Under reflection, \( \psi_n(S_{\Delta} r) = \pm \psi_n(r) \), where the + and − signs hold for \( A_1 \) and \( A_2 \), respectively. Wavefunctions pertaining to the 2D representation \( E \) satisfy \( \psi_n(Rr) = \exp(\pm 2i\pi/3) \psi_n(r) \), so that \( n = \pm 1 \) modulo 3 \[52\].

These symmetry considerations further reduce the configuration space to the green triangle \( OLB \) of Fig. 1b. We deal with representations \( A_1, A_2, \) and \( E \) separately by applying different boundary conditions on its edges (see Sec. A2 in the Appendix). We solve the resulting stationary Schrödinger equations using the finite–element software FreeFEM \[53\]. The classical scaling no longer holds.

Instead, the energy spectra and wavefunctions depend on the dimensionless ratio \( \eta = \hbar / P_{\text{ref}} = \hbar R^2 / (m C_6)^{1/2} \). Smaller values of \( \eta \) signal deeper quasiclassical behavior: we choose \( \eta = 0.01 \). We focus on energies \( \epsilon \sim 7C_6/R^6 \), which are large enough for the classical ergodic trajectory (brown dots on Fig. 3) to occupy a substantial part.
Table I. Parameters $T_B^{(\rho)}$, $S_B^{(\rho)}$, $\alpha_B^{(\rho)}$, and $k^{(\rho)}$ for Eq. 4 depending on the irreducible representation $\rho = A_1, A_2, or E$.

| $T_B^{(\rho)}$ | $S_B^{(\rho)}$ | $\alpha_B^{(\rho)}$ | $k^{(\rho)}$ |
|--------------|-------------|-----------------|---------|
| $A_1$ $T_B/2$ | $S_B/2$ | $\lambda_B T_B/2$ | $k$ |
| $A_2$ $T_B/2$ | $S_B/2$ | $\lambda_B T_B/2$ | $k - 1/2$ |
| $E$ $2T_B$ | $S_B$ | $\lambda_B T_B$ | $k + 1/2$ |

Figure 5 shows $\Delta \nu$ agree with the energies of the scarred states. Hence, they are directly related to the classical period $T_B(\epsilon)$ of unstable trajectories. We use Gutzwiller’s trace formula [31, 54, 55] to the tower of scarred many-body states with an approximating $n$ to the ergodic region of classical phase space (see Fig. 3).

V. SEMICLASSICAL ANALYSIS

For the majority of the calculated quantum states, the probability density $|\psi(x, y)|^2$ is unrelated to the periodic trajectories of type B. Nevertheless, for each representation, we find multiple scarred quantum states, represented by the vertical dashed lines in Fig. 3 whose energy spacings is approximately regular. This is analogous to the tower of scarred many-body states with an approximately constant energy separation found in a PXP chain [13], which is a recurrent feature in theoretical analyses of weak ergodicity breaking [14, 22, 23]. In the present context, we explain the series of scarred quantum states semiclassically. We use Gutzwiller’s trace formula [31] describing the impact of the classical periodic trajectories on the quantum density of states $n(\epsilon)$. We isolate the contribution $\Delta n_B^{(\rho)}$ to $n$ coming from the unstable trajectory $B$, which depends on the representation $\rho$ [54, 55]:

\[
(\Delta n_B^{(\rho)} 2\pi \hbar/T_B^{(\rho)} + 1)/\alpha_B^{(\rho)} = 
\sum_{k=0}^{\infty} \{[S_B^{(\rho)}/\hbar - 2\pi(k^{(\rho)} + 1/2)]^2 + \alpha_B^{(\rho)/2})^2\}^{-1}.
\]  

The parameters $T_B^{(\rho)}(\epsilon)$, $S_B^{(\rho)}(\epsilon)$, $\alpha_B^{(\rho)}(\epsilon)$, and $k^{(\rho)}$ in Eq. 4 are defined in Table I for each representation. They are directly related to the classical period $T_B(\epsilon)$ and action $S_B(\epsilon) = \oint p \cdot dx$ along one trajectory $B$, the product $\lambda_B(\epsilon) T_B(\epsilon)$, and the summation index $k$, respectively. Figure 4 shows $\Delta n_B^{(\rho)}(\epsilon)$ for each representation. Its maxima agree with the energies of the scarred states. Hence, the series of scarred states found in each representation reflects the multiple resonances in $n(\epsilon)$ due to the unstable trajectory $B$. The regularity in their energy spacing follows from the resonance maxima being evenly spaced in terms of the classical action, $S_B^{(\rho)}/\hbar = 2\pi(k^{(\rho)} + 1/2)$.

VI. EXPERIMENTAL PROSPECTS AND OUTLOOK

We consider e.g. $^{87}$Rb atoms in the 50C circular Rydberg state [36, 56], for which $C_6/\hbar = 3$ GHz $\mu$m$^6$. The value $\eta = 0.01$ corresponds to $R = 7 \mu$m. The ring-shaped trap may be realized optically using Laguerre–Gauss laser beams and light sheets [57, §II.C.2]. The energy $\epsilon = 7 C_6/\hbar = h \times 200$ kHz is within experimental reach [36]. For small angular momenta, the centrifugal energy, which is proportional to $(m \mu)^2/3$, is negligible compared to $\epsilon$. The position of the atoms may be detected at a given time by turning on a 2D optical lattice trapping individual Rydberg atoms [36, 58], which freezes the dynamics, followed by atomic deexcitation and site-resolved ground–state imaging [59].

Further investigation will be devoted to the stability of the quantum scar. Recent experiments [23, 60] have shown that it may be enhanced by periodically modulating the parameters. Depending on the stabilization mechanism (see e.g. Ref. [61] or Ref. [62, §27]), this may lead to a discrete time crystal [63] which is either quantum or classical.

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Appendix A

The goal of this Appendix is twofold. In Section A1 we identify novel quantum scars supported by the Hénon–Heiles Hamiltonian, and characterize them using the same semiclassical argument as in the main text. In Section A2 for each of the three irreducible representations of the group $C_{3v}$, we derive boundary conditions defining quantum stationary states within the reduced configuration space.

1. Quantum scars in the Hénon–Heiles model

In this Section, we briefly describe our results, analogous to those of the main text, for the Hénon–Heiles Hamiltonian $H_{HH} = (p_x^2 + p_y^2)/(2m) + V_{HH}$, where:

$$V_{HH} = mω_0^2(x^2 + y^2)/2 + α(x^2y - y^3/3).$$

(A1)

Equation [A1] is written in the dimensional form of Ref. [64 §5.6.4] which assumes that the coordinates $x$ and $y$ carry the unit of length. The quantities $p_x$, $p_y$ are their conjugate momenta, the parameters $m$ and $ω_0$ denote a mass and a frequency, and the coefficient $α$ sets the strength of the cubic term. If lengths, momenta, energies, and times are expressed in units of $L_{HH} = mω_0^2/α$, $P_{HH} = m^2ω_0^2/α$, $E_{HH} = m^3ω_0^2/α^2$, $T_{HH} = 1/ω_0$, the dimensionless form matches that of Ref. [34]. As in the main text, in terms of these units, the classical dynamics is independent of $m$, $ω_0$, $α$. As for quantum physics, the classical scaling no longer holds, and the energy spectra and wavefunctions depend on the dimensionless parameter $η_{HH} = ℏ/(L_{HH}P_{HH}) = ℏα^2/(m^3ω_0^5)$. Smaller values of $η_{HH}$ signal deeper quasiclassical behavior.

The Hénon–Heiles potential is related to our main discussion for two reasons. First, its symmetry group is $C_{3v}$ [52], which is the point group of the system analyzed in the main text. Second, expanding Eq. (3) there to third order in $x$ and $y$ near the equilibrium position $O$ shows that it reduces to Eq. (A1) in the low–energy limit.

The Hénon–Heiles Hamiltonian has been extensively studied (see e.g. Ref. [52 §1.4]). Our goal in revisiting it was twofold. First, we have calibrated our codes against published results for this potential. Second, we have identified quantum scars for the Hénon–Heiles Hamiltonian which, to the best of our knowledge, are novel. At the end of the section, we point out the relevance of the Hénon–Heiles potential in relation to a broad family of systems, which includes the case of dipolar particles.

a. Calibration

We have used our codes to reproduce the known classical periodic trajectories of $H_{HH}$, their periods and Lyapunov exponents [42], and its surfaces of section for various energies [48]. We have also recovered the quantum energy levels and wavefunctions, belonging to all three representations, in Refs. [51, 63] for $η_{HH} = 1/80$ and in Ref. [52] for $η_{HH} = 0.06^2$.

b. Quantum scars for the Hénon–Heiles potential

We now turn to the lower value $η = 0.04^2$, so as to consider the deep quasiclassical regime. We focus on energies $ε \sim 0.13 E_{HH}$: these are large enough for the ergodic region to occupy a substantial part of phase space [48], while remaining below the threshold energy $E_{HH}/6$ above which $H_{HH}$ supports trajectories that are not bound [42]. Figure 6 shows the probability density density for the scarred state with the energy $ε$ closest to 0.13$E_{HH}$ for each representation. It is maximal near the three trajectories $B$ for the energy $ε$, signaling the scar.

In each irreducible representation $ρ = A_1$, $A_2$, and $E$, we find multiple scarred quantum states for the Hénon–Heiles potential (vertical dashed lines in Fig. 4) whose energy spacing is approximately regular, in direct analogy with the results of the main text. They may be explained using the same semiclassical argument relying on Gutzwiller’s trace formula. We isolate the contribution $Δn^{(ρ)}$ to the density of states $n$ for each representation $ρ$ due to the unstable trajectory $B$. Both Eq. 4 and Table I in the main text are applicable to the Hénon–Heiles potential with no change. We have calculated the required period $T_B$, action $S_B$ and Lyapunov exponent $λ_B$ characterizing the periodic trajectory $B$ in the Hénon–Heiles potential as a function of the energy $ε$ using our codes. Figure 7 shows $Δn^{(ρ)}$ for each representation $ρ$. Just like in the main text, its maxima coincide with the energies of the scarred states. Hence, the same conclusion holds, and we may ascribe the regularity in their energy spacing to the resonance maxima being equally spaced in terms of the classical action $S_B$.

c. Generality of the Hénon–Heiles potential

The potential $V_{HH}$ combines a 2D isotropic harmonic trap with a two–variable cubic polynomial function. Hence, it may be seen as the simplest possible 2D potential exhibiting $C_{3v}$ symmetry. The three-body Hamiltonian given by Eq. 1 in the main text reduces to it near one of its (equivalent) minima for the repulsive pair–wise interaction $v(d_{ij}) = α d_{ij}^α$ regardless of the power law exponent $α > 0$. The presence of quantum scars in the Hénon–Heiles model leads us to expect similar scars in all of these systems. In particular, the dipole–dipole interaction [35] in the case where all three dipole moments are polarized perpendicular to the plane, corresponding to $α = 3$, is expected to yield the same phenomena.
2. Boundary conditions defining a basis of quantum stationary states

In this Section, we exploit the spatial symmetries of the point group $C_{3v}$ and time-reversal symmetry to state boundary conditions uniquely defining a basis of quantum stationary states. We state our reasoning in terms of the system considered in the main text, but it applies without change to the Hénon–Heiles Hamiltonian discussed in Sec. A1 above.

We expect the quantum states scarred by the classically unstable periodic trajectory $B$ to exhibit an enhanced probability density along all three trajectories $B$ at a given energy (red dotted lines in Figs. 1a and 4b–c in the main text for the system discussed there, and in Figs. 6b–c in the present Appendix for the Hénon–Heiles model). Hence, the probability density for the scarred states is expected to exhibit $C_{3v}$ symmetry. Therefore, we construct a basis of quantum stationary states whose corresponding density profiles all exhibit this symmetry. This property is not automatically satisfied and requires choosing appropriate basis functions. For example, Figs. 7a and 7b in Ref. 51 show probability densities corresponding to eigenstates of the Hénon–Heiles model which do not exhibit $C_{3v}$ symmetry despite the fact that the Hamiltonian does, see Sec. A1 above.

The group $C_{3v}$ admits 3 irreducible representations, $\rho = A_1$, $A_2$, and $E$ [40, §95]. Representations $A_1$ and $A_2$ are 1D, whereas representation $E$ is 2D. For each representation, we shall formulate a boundary condition defining basis functions belonging to it. All wavefunctions $\psi(r)$ are normalized according to $\int_{AB} d^2r |\psi(r)|^2 = 1$, the integral being taken over the triangle $ABC$.

\textbf{a. One–dimensional representations $A_1$ and $A_2$}

We first consider a 1D representation $\rho = A_1$ or $A_2$. Let $\psi$ be an eigenstate of $H_{\text{2D}}$ for the energy $\epsilon$ transforming according to $\rho$. We call $S_1$, $S_2$, $S_3 = S_\Delta$ the reflections about $(AH)$, $(BK)$, $(CL)$ in the $(x,y)$ plane (see Fig. 8). The wavefunction $\psi(S_\rho r)$ is also an eigenstate of $H_{\text{2D}}$ for the same energy $\epsilon$. Because $\rho$ is 1D, $\psi(S_\rho r) = \chi_\rho \psi(r)$ for some complex number $\chi_\rho$. The reflections $S_i$ satisfy $S_i^2 = 1$, so that $\chi_i = \pm 1$. They also satisfy $S_2 S_1 = S_3 S_2 = R$, with $R$ being the rotation of angle $2\pi/3$ about the point $O$. The transformation $R^3 = 1$, so that $(\chi_1 \chi_2)^3 = (\chi_2 \chi_3)^3 = 1$. Hence, $\chi_1 = \chi_2 = \chi_3 = \pm 1$.

The case $\chi_1 = \chi_2 = \chi_3 = 1$ leads to $\rho(R) = \rho(S_i) =...
The black triangle $ABC$ is the classical configuration space for the Hamiltonian $H_{2D}$ of the main text. The smaller green triangle $OBL$ is the reduced configuration space within which we solve for the quantum stationary states. The classically accessible region, limited by the dotted golden line, is shown for the energy $\epsilon = 7C_{6}/R^6$. We enforce the boundary condition $\psi = 0$ on the quantum wavefunctions along the horizontal dashed green line. The three classical periodic trajectories $B$ (dotted red lines) are also shown.

1, so that $\rho = A_1$ [40, §95, Table 7]. Then, $\psi(S_1r) = +\psi(r)$, leading to the boundary condition $\partial_n\psi = 0$ along the sides $[LO]$ and $[OB]$ of the green triangle $OBL$ in Fig. 8. Combined with the condition $\psi = 0$ along the side $[LB]$ derived in the main text, it defines a basis of wavefunctions $\psi$ for Representation $A_1$.

The case $\chi_1 = \chi_2 = \chi_3 = -1$ leads to $\rho(R) = 1$ and $\rho(S_1) = -1$, so that $\rho = A_2$. Then, $\psi(S_1r) = -\psi(r)$, leading to the condition $\psi = 0$ along the sides $[LO]$ and $[OB]$. Hence, imposing the Dirichlet boundary conditions on the three edges of the triangle $OBL$ defines a basis of wavefunctions $\psi$ for Representation $A_2$.

The energy levels transforming according to the 1D representations $A_1$ and $A_2$ are non-degenerate, hence, the time-reversal invariance of $H_{2D}$ allows us to choose all basis wavefunctions $\psi(r)$ to be real [40, §18]. Furthermore, $\psi(r) = \psi(R^{-1}r) = \chi_1\psi(S_1r)$ differ by a sign at most. Hence, the corresponding probability densities coincide, and $|\psi(r)|^2$ does exhibit $C_{3v}$ symmetry.

b. Two-dimensional representation $E$

We now turn to the 2D representation $\rho = E$. Let $\epsilon$ be a two–degenerate energy level of $H_{2D}$. The corresponding eigenspace is spanned by two complex wavefunctions, $\phi_+$ and $\phi_-$ which transform according to $\rho$:

$$\phi_\pm(R^{-1}r) = e^{\pm i2\pi/3}\phi_\pm(r), \quad (A2a)$$
$$\phi_\pm(S_\Delta r) = \phi_\mp(r), \quad (A2b)$$

where the transformations $R$ and $S_\Delta$ are defined as in Sec. A2a above and the main text.

The time–reversal invariance [40, §18] of $H_{2D}$ entails that the complex–conjugate wavefunctions $\phi_+(r)$ and $\phi_-^*(r)$ are also eigenstates of $H_{2D}$ with the same energy $\epsilon$. Complex–conjugating Eqs. (A2a), accounting for normalization, and writing $(\phi_+)^* = \phi_-$ lead to $\phi_\pm = e^{i\alpha}\phi_\mp$, where $e^{i\alpha}$ is a complex number of modulus 1. Introducing the new basis wavefunctions $\psi_+(r) = e^{i\alpha/2}\phi_+(r)$ and $\psi_-(r) = \psi_+^*(r)$, Eqs. (A2) reduce to 2 conditions on $\psi_+$:

$$\psi_+(R^{-1}r) = e^{i2\pi/3}\psi_+(r), \quad (A3a)$$
$$\psi_+(S_\Delta r) = \psi_+(r). \quad (A3b)$$

The probability densities $|\psi_+(r)|^2 = |\psi_+(R^{-1}r)|^2 = |\psi_+(S_\Delta r)|^2$ coincide. Hence, $|\psi_+(r)|^2$ exhibits $C_{3v}$ symmetry: this is the probability density plotted in Figs. 4a–c of the main text (three Rydberg atoms) and Figs. 6a–c (Hénon–Heiles model).

We seek $\psi_+(r)$ in the following form, which is more amenable to numerical computation:

$$\psi_+(r) = (x - iy)(u_1(r) + iu_2(r)), \quad (A4)$$

where $u_1$ and $u_2$ are two real functions satisfying coupled Schrödinger equations. In Eq. (A1), the factor $(x - iy)$ accounts for the fact that $\psi_+(0) = 0$, i.e. for the stationary states of the 2D isotropic harmonic oscillator carrying angular momentum [40, §112]. Equations (A3) yield the boundary conditions $u_1 = 0$, $\partial_nu_2 = 0$ on both $[LO]$ and $[OB]$ (see Fig. 8). Combined with the condition $\psi = 0$ along $[LB]$ derived in the main text, they define a basis of stationary states related to Representation $E$. For each of the twice-degenerate energy levels, $\psi_+(r)$ is given by Eq. (A4) and the second basis function is $\psi_-(r)$.

d. Spatial extent of the wavefunctions

For a given energy level $\epsilon$, the spatial extent of the stationary states defined in Secs. A2a and A2b barely exceeds the classically accessible region (limited by the dotted golden line in Fig. 8 for the Hamiltonian $H_{2D}$ of the main text and $\epsilon = 7C_{6}/R^6$). Therefore, we restrict the region within which we solve for the wavefunctions to a part of the triangle OLB which slightly exceeds this region. In other words, we enforce the condition $\psi = 0$ not on $[LB]$, but on the horizontal dashed line in Fig. 8.

d. Indices of the quantum states

We order the quantum states pertaining to a given irreducible representation $\rho$ by increasing energies. This gives rise to the state index $\nu^{(\rho)}$ appearing in Figs. 4 and 5 in the main text, and Figs. 6 and 7 in the present Appendix. The irreducible representations $A_1$ and $A_2$ have
dimension 1, so that, barring accidental degeneracies, the corresponding energy levels are non-degenerate. By contrast, the irreducible representation $E$ has dimension 2, meaning that each energy level is twice degenerate. For this representation, we consistently indicate one half of the state index, $\nu(E)/2$, and one half of the density of states contribution $\Delta n_B(E)/2$.

The relative level indices $\Delta \nu(p)$ given in Fig. 5 of the main text and in Fig. 7 of this Appendix are exact. The level indices of Fig. 5 concerning the Hénon–Heiles model are also exact. We obtain approximations to the level indices of Fig. 5 in the main text, concerning three Rydberg atoms moving along a circle, using the semiclassical approximation to the density of states, accounting for the role of discrete spatial symmetries [60].

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