Anderson Localization of Solitons

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At low temperature, a quasi-one-dimensional ensemble of atoms with an attractive interaction forms a bright soliton. When exposed to a weak and smooth external potential, the shape of the soliton is hardly modified, but its center-of-mass motion is affected. We show that in a spatially correlated disordered potential, the quantum motion of a bright soliton displays Anderson localization. The localization length can be much larger than the soliton size and could be observed experimentally.

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At zero temperature, cold atoms interacting attractively in a one-dimensional (1D) system tend to cluster together, forming a bright soliton. Explicit solutions of the manybody problem can be found in some cases, for example, for contact interactions [1]. Using external potentials, it has been experimentally shown how to put solitons in motion [2]. What happens to a soliton exposed to a disordered potential? If the potential is strong, it will destroy the soliton. If it is sufficiently weak and smooth not to perturb the soliton shape, one expects the soliton to undergo multiple scattering, diffusive motion, and possibly Anderson localization [3]. Indeed, propagation of waves in a disordered potential is profoundly affected by Anderson localization. Multiple scattering on random defects yields exponentially localized density profiles and a suppression of the usual diffusive transport associated with incoherent wave scattering [4]. In 1D Anderson localization is a ubiquitous phenomenon [5], which has been recently observed for cold atomic matter waves [6]. It is important to understand how it is modified when interactions between particles are taken into account.

We consider a Bose-Einstein condensate in a quasi-1D geometry. Within mean-field theory, it is described by the Gross-Pitaevskii energy functional

$$E = \int dz \left[\frac{1}{2} |\partial_z \phi|^2 + \frac{g}{2} |\phi|^4 - \mu |\phi|^2 \right]$$
(1)

in units of $E_0 = 4m\omega_{\perp}^2 a^2$, $l_0 = \hbar/2|a|m\omega_{\perp}$, and $t_0 = \hbar/4a^2m\omega_{\perp}^2$ for energy, length, and time, respectively. Here, ω_{\perp} denotes the transverse harmonic confinement frequency, *a* the atomic *s*-wave scattering length, and μ the chemical potential. The cases of repulsive and attractive atomic interaction are covered by $g = \pm 1$.

The dynamics is extremely different in both cases, so that we first discuss the case of attractive interaction, g = -1. The ground state of (1) is the bright soliton [7]

$$\phi_0(z-q) = \sqrt{\frac{N}{2\xi}} \frac{e^{-i\theta}}{\cosh[(z-q)/\xi]},\tag{2}$$

normalized to the total number of particles N. The chemical potential is $\mu = -N^2/8$ and the soliton width is $\xi = 2/N$. This ground-state solution has an arbitrary center-ofmass (c.m.) position q and an arbitrary global phase θ that spontaneously break the translational and the U(1) gauge symmetry of the energy functional (1), respectively. These degrees of freedom appear as zero-energy modes of Bogoliubov theory, and their quantum dynamics requires special attention [8,9].

The energy functional (1) is no longer translation invariant when a potential term $\int dz V(z) |\phi|^2$ is added. If V(z) is sufficiently weak and smooth, the soliton shape remains unchanged to lowest order in *V*, and only its c.m. position *q* is affected. In this Letter, we show that the quantum dynamics of *q* in a disorder potential leads to Anderson localization of the soliton, over a localization length possibly longer than the soliton size.

An intuitive picture emerges using a simple ansatz in terms of collective coordinates (cf. [10] in the context of Bloch oscillations). Inserting $\phi(z; q, P_q) = e^{iP_q z/N} \phi_0(z - q)$ into (1) leads to an effective quantum Hamiltonian

$$\hat{H}_{q} = \frac{\hat{P}_{q}^{2}}{2N} + \int dz V(z) |\phi_{0}(z - \hat{q})|^{2}.$$
 (3)

It describes the entire soliton as an object of mass N evolving in an effective potential $\int dz V(z) |\phi_0(\hat{q} - z)|^2$ that is the convolution of the bare potential with the soliton density.

This simple ansatz yields no information on the remaining degrees of freedom. Therefore, we apply a more complete analysis, expanding the energy functional (1) to second order in deviations from the ground-state solution (2), as in, e.g., [8]. Diagonalization of the quadratic Hamiltonian results in the eigenvalue problem for the non-Hermitian operator

$$\mathcal{L} = \begin{pmatrix} -\frac{1}{2}\partial_z^2 - 2|\phi_0|^2 - \mu & -\phi_0^2 \\ \phi_0^{*2} & \frac{1}{2}\partial_z^2 + 2|\phi_0|^2 + \mu \end{pmatrix}.$$
 (4)

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Its right eigenvectors (u_n, v_n) and corresponding adjoint modes $(u_n^{\rm ad}, v_n^{\rm ad})$ build a basis that spans the functional space of (ϕ, ϕ^*) . For all nonzero eigenvalues E_n , the adjoint modes are left eigenvectors of \mathcal{L} . This is no longer true for the zero-energy modes. The first zero mode $(u_{\theta}, v_{\theta}) = i\partial_{\theta}(\phi_0, \phi_0^*)$, is related to the global U(1) gauge invariance $\phi e^{-i\theta} \rightarrow \phi e^{-i(\theta+\epsilon)}$ broken by the classical solution (2) and its adjoint mode is well known, $(u_{\theta}^{ad}, v_{\theta}^{ad}) =$ $\partial_N(\phi_0, \phi_0^*)$ [8,9]. The other zero mode, $(u_q, v_q) =$ $i\partial_a(\phi_0, \phi_0^*)$, originates from the translational invariance $q \rightarrow q + \epsilon$ also broken by the soliton solution. To find the adjoint mode, one solves $\mathcal{L}(u_q^{\mathrm{ad}}, v_q^{\mathrm{ad}}) = M^{-1}(u_q, v_q)$, where M is determined by the requirement $\langle u_q^{\rm ad} | u_q \rangle$ – $\langle v_q^{\rm ad} | v_q \rangle = 1$. This ensures that $(u_q^{\rm ad}, v_q^{\rm ad})$ is orthogonal to all eigenvectors of \mathcal{L} with $E_n \neq 0$ [8,9]. It is easy to verify that

$$\begin{pmatrix} u_q^{\text{ad}} \\ v_q^{\text{ad}} \end{pmatrix} = i \frac{z - q}{N} \begin{pmatrix} \phi_0 \\ -\phi_0^* \end{pmatrix},$$
(5)

and M = N, the mass of the system.

Following Dziarmaga [9] we may now perform an expansion around the ground state,

$$\phi = \phi_0 + P_\theta u_\theta^{\rm ad} + P_q u_q^{\rm ad} + \sum_{n, E_n > 0} (b_n u_n + b_n^* v_n^*), \quad (6)$$

where all modes are implicit functions of θ and q. Inserting (6) in the energy functional (1) and expanding to second order in P_{θ} , P_q , and b_n (requiring $P_q \xi \ll N$ for the c.m. momentum) results in the Bogoliubov Hamiltonian, whose quantum version reads

$$\hat{H}_{0} = -\frac{N}{8}\hat{P}_{\theta}^{2} + \frac{\hat{P}_{q}^{2}}{2N} + \sum_{n, E_{n} > 0} E_{n}\hat{b}_{n}^{\dagger}\hat{b}_{n}, \qquad (7)$$

where $\hat{P}_{\theta} = \hat{N} - N = -i\partial_{\theta}$ and $\hat{P}_{q} = -i\partial_{q}$.

Because θ can take nonperturbatively large values, we can work in a subspace of Hilbert space with definite $\hat{N} = N$. The minus sign in front of P_{θ}^2 in (7) arises because the bright soliton (2) is a saddle point of the energy functional (1). It has no consequences since the number of particles is fixed.

When the energy functional is supplemented with a potential term $\int dz V(z) |\phi|^2$, this perturbation commutes with \hat{N} , so that we can focus on its incidence on the second and third term in (7), i.e., the c.m. motion and Bogoliubov excitations. A smooth external potential can only slightly distort the soliton. If the potential energy variation across the soliton is of the order of the standard deviation $|V_0|$ of the disordered potential, then $|V_0| \ll |\mu| = N^2/8$ is a sufficient condition for the soliton shape to be only weakly modified by the external potential [11]. Moreover, the large energy gap $|\mu|$ above the ground state [12] makes Bogoliubov excitations by the potential negligible. The soliton shape thus follows adiabatically and reversibly, i.e., without heating, the variations of the external potential potential.

tial. Then, the only degree of freedom affected by the external potential is the soliton position q. Inserting the expansion (6) into the energy functional, expanding up to quadratic terms, and quantizing, we arrive at the effective Hamiltonian (3) to leading order in 1/N and V.

In the following, we study the case of $N = 100 \text{ Li}^7$ atoms with scattering length a = -3 nm in a transverse harmonic trap with $\omega_{\perp} = 2\pi \times 5$ kHz. Then, units for energy, length, and time are $E_0 = 1.28 \times 10^{-4}\hbar\omega_{\perp}$, $l_0 =$ 47.8 μ m and $t_0 = 0.25$ s. A soliton (2) of size $\xi = 2/N =$ 0.02 ($\approx 1 \mu$ m) is initially prepared in a large axial harmonic trap with $\omega_z = 100 (2\pi \times 64 \text{ Hz})$. The trapping potential is sufficiently small not to distort the soliton, whose c.m. occupies the ground state of a harmonic oscillator with frequency ω_z . When the trap is turned off, the soliton position starts its quantum dynamics with the corresponding momentum distribution

$$\pi_0(k) = \frac{1}{\sqrt{\pi}\Delta k} \exp[-k^2/\Delta k^2], \qquad \Delta k^2 = N\omega_z, \quad (8)$$

and begins to explore the disordered potential.

In 1D random potentials, Anderson localization is generic: the amplitude of every plane wave with wave vector k decreases asymptotically as $\exp\{-\gamma(k)|q|/2\}$. The inverse localization length $\gamma(k)$ can be calculated analytically in the weak-disorder limit (see below) [13]. Equivalently, the energy spectrum is discrete and dense (in the limit of infinitely large systems) with exponentially localized eigenstates.

Disorder potentials are completely characterized by correlation functions $V(z_1) \dots V(z_n)$ where the overbar denotes an ensemble average over disorder realizations. The average potential value shifts the origin of energy and can always be set to zero, V(z) = 0. The pair correlator can be written as $V(z')V(z' + z) = V_0^2C(z/\sigma_0)$, where V_0 measures the potential strength, and σ_0 the spatial correlation length. Higher-order correlations are required to fully describe non-Gaussian disorder such as the optical speckle potential considered in the following.

Optical speckle yields a light-shift potential $V(z) \propto \chi |E(z)|^2$ proportional to the intensity of the light field E(z) and to the atomic polarizability χ , whose sign depends on the detuning of the external light frequency from the atomic resonance. At fixed detuning, the potential features either random peaks (the "blue-detuned" case) or wells ("red-detuned"). The potential distribution is asymmetric, and the importance of odd moments can be probed by comparing the blue- and red-detuned cases for fixed $|V_0|$. We use $|V_0| = 8 \times 10^{-5} |\mu| = 0.1$ in the following. A 1D speckle potential has the pair correlation function $C(y) = [\sin(y)/y]^2$, with a correlation length that can be as short as $0.26 \ \mu m \ [6]$ or $\sigma_0 = 0.0056$ in our units. In *k*-space, the corresponding power spectrum reads $\mathcal{P}_V(k) = \pi \sigma_0 V_0^2 (1 - |k\sigma_0|/2) \Theta (1 - |k\sigma_0|/2)$.

Our Hamiltonian (3) shows that the bright soliton sees a convoluted disorder potential whose k-space components

are the product $(N\pi k\xi/2)/\sinh(\pi k\xi/2)V_k$ of Fourier components from soliton density and speckle potential. To second order (Born approximation) in the potential strength, the inverse localization length [13] then reads

$$\gamma(k) = \frac{N^4 \xi^2 \pi^3 \sigma_0 V_0^2 (1 - |k\sigma_0|)}{[\sinh(\pi k\xi)]^2} \Theta(1 - |k\sigma_0|).$$
(9)

For a short correlation range $\sigma_0 \ll \xi$, the *k* dependence due to the soliton convolution dominates and the bare speckle can be approximated by its white-noise limit $\mathcal{P}_V(0) = \pi \sigma_0 V_0^2$. The soliton width ξ takes over as the new effective correlation length scale. For $1/(\pi\xi) < k < 1/\sigma_0$ one can use the approximate expression

$$\gamma(k) \approx (2\pi N^2 \xi)^2 \mathcal{P}_V(0) e^{-2\pi \xi k}.$$
(10)

The lowest-order perturbation result (9) can be compared with exact data computed numerically, both by exact diagonalization and transfer-matrix methods [14] for the Hamiltonian (3). Figure 1 confirms that Anderson localization is observed for all k values, but also shows that the localization lengths for the blue- and red-detuned potential differ by up to an order of magnitude for $k\xi > 0.5$. Consequently, they also differ from the lowest-order perturbation results. Thus, perturbation theory cannot be expected to apply, even though the disorder potential is much smaller than the kinetic energy of the soliton. But the salient feature here is the rapid exponential decrease of $\gamma(k)$, clearly visible in the inset of Fig. 1. Thus, the inverse localization length still obeys

$$\gamma = \gamma_0 \exp(-\alpha k) \tag{11}$$

with $\alpha \propto \xi$, the natural length scale of the problem.

Let us return to the c.m. dynamics. The initial state is the superposition of various *k* components, with weights given



FIG. 1 (color online). Inverse localization length of a bright soliton versus its momentum k in units of its width $\xi = 0.02$ ($\approx 1 \ \mu$ m) in a speckle potential with correlation length $\sigma_0 = 0.28\xi$. Circles are numerical results obtained by diagonalizing the effective Hamiltonian (3), solid lines are numerical results from a transfer-matrix calculation. The bare potential amplitude is $V_0 = \pm 0.1$ [blue (dark gray) and red (gray) curve] in our units, corresponding to $\pm 8 \times 10^{-5} |\mu|$. Inset: same data in logarithmic scale. The lowest-order perturbation result, Eq. (9) is shown as a black dashed line.

by Eq. (8). Each k component localizes at long time into the asymptotic form $\exp\{-\gamma(k)|q|/2\}$, implying that the superposition will also localize. If the extension of the initial state is small compared to the typical localization length (which is the case here), one can neglect the phase correlations between the various k components at long time. The final density distribution, ensemble averaged over disorder realizations, is then given by

$$W(q) = \overline{|\psi(q)|^2} = \int dk \frac{\gamma(k)}{2} \exp\{-\gamma(k)|q|\}\pi_0(k).$$
(12)

Using (11) this integral can be evaluated by a saddle-point argument from which it becomes apparent that at distance q, one finds the soliton with an initial momentum k_q such that $|q| = 1/\gamma(k_q)$ provided $\alpha \Delta k \gg 1$. The probability distribution for the soliton position then takes the form

$$W(q) = \frac{\sqrt{2}}{e\alpha\Delta k} \frac{\exp[-k_q^2/\Delta k^2]}{|q|},$$
(13)

where $k_q = \ln(\gamma_0 |q|)/(\alpha)$ from (11). In the regime of interest, this is an almost perfect algebraic decrease as $|q|^{-1}$ with a small logarithmic correction [15].

These predictions have been tested by numerical integration of the Schrödinger equation with the Hamiltonian (3) starting from the initial Gaussian wave function corresponding to (8), and averaging over 250 disorder realizations. As shown in Fig. 2, the final probability distribution for the soliton position follows quite well the predicted algebraic decay. The time scale required to observe a stationary localized state around position q is given by $\tau(q) = |q|/v(k_q) = N|q|/k_q = \alpha |q|/\ln(\gamma_0|q|)$.

Finally, we turn to the case of repulsive interactions [16] by taking g = +1 in (1). The Bose-Einstein condensate ground state in a harmonic trap extends over the Thomas-Fermi radius. How such a condensate wave packet expands in a disordered potential has been studied for different dimensionalities [17–19]. Because of repulsive interaction, the entire condensate is an extended object that requires a fieldlike description, in sharp contrast with a bright soliton that features particlelike properties.

There exists, however, a dark soliton solution $\phi_0 \propto \tanh[(x-q)/\xi]$ where $\xi = 1/\sqrt{\rho}$ and ρ is the background density. Then, one can use the nonperturbative description presented in [9] or a collective-coordinate method [20] to obtain an effective Hamiltonian for the dynamics of the dark soliton in the presence of a weak disorder potential:

$$H'_{q} = -\left[\frac{P_{q}^{2}}{2|M|} + \frac{|M|}{4\xi} \int dz \frac{V(z)}{\left[\cosh((z-q)/\xi)\right]^{2}}\right].$$
 (14)

The mass of the dark soliton is negative, $M = -4\rho\xi$, and its modulus equals twice the number of particles missing in the soliton notch. This effective Hamiltonian is valid only when the velocity of the c.m. of the soliton is much smaller than the sound velocity $c = \sqrt{\rho}$ in the condensate. At



FIG. 2 (color online). Probability distribution of soliton position q in units of its width $\xi = 0.02 \ (\approx 1 \ \mu m)$. Blue (dark gray) and red (gray) lines: Data from numerical integration of the Schrödinger equation with the Hamiltonian (3) starting in the ground state of an axial harmonic trap with $\omega_z = 100 \ (2\pi \times 64 \ Hz)$ and lasting $\Delta t = 20 \ (5 \ s)$ for a blue (dark gray) and red (gray)-detuned speckle potential of amplitude $V_0 = \pm 8 \times 10^{-5} \mu \ (2\pi \times 0.064 \ Hz)$. Data averaged over 250 different realizations of the disorder potential. Black dashed line: saddle-point result (13) for parameters in the Born approximation. The full integral (12) yields the same curve for $|q| \gg \xi$. The double-logarithmic inset plot shows the dominantly algebraic decrease as $|q|^{-1}$.

velocity comparable to c, the shape of the soliton changes, making the analysis more difficult.

Equation (14) shows that we obtain the same form of the effective Hamiltonian as for the bright soliton. At first sight, it predicts similar Anderson localization effects. However, there is a fundamental difference between dark and bright solitons. The bright soliton (2), being the ground state of the N particle system, is protected by a large gap from quasiparticle excitations. In that respect, entire solitons can undergo quantum dynamics and localize like rubidium atoms [6]. For the dark soliton a gap can only be imposed by the boundary conditions in a finite-size system, and is inversely proportional to the system size, making radiation of Bogoliubov excitations [21] and decoherence much easier. Furthermore, the dark soliton is an excited state, as signaled by the global minus sign of the effective Hamiltonian (14). Interactions with the thermal cloud may accelerate the soliton and make it disappear [22]. Estimating the relevant time scale and evaluating its effect on the localization dynamics is left for future work.

In summary we have shown that the center of mass of a bright soliton may undergo Anderson localization in a smooth disorder potential. In realistic situations where the soliton wave packet is prepared in a small region of space, this leads to an essentially algebraic localization of the ensemble-averaged atomic density at long times. We emphasize that the effects discussed here are beyond the standard mean-field description: while the soliton's shape is described by a mean-field theory, its center of mass is treated quantum mechanically. This has important consequences: although the one-body density matrix will display Anderson localization as shown in Fig. 2, a single realization of the experiment is expected to find a single soliton at a given random position, with a probability density given by W(q), Eq. (13).

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