

Adiabatic splitting, transport, and self-trapping of a Bose-Einstein condensate in a double-well potential

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We show that the adiabatic dynamics of a Bose-Einstein condensate (BEC) in a double-well potential can be described in terms of a dark variable resulting from the combination of the population imbalance and the spatial atomic coherence between the two wells. By means of this dark variable, we extend, to the nonlinear matter-wave case, the recent proposal by Vitanov and Shore [Phys. Rev. A **73**, 053402 (2006)] on adiabatic passage techniques to coherently control the population of two internal levels of an atom or molecule. We investigate the conditions to adiabatically split or transport a BEC as well as to prepare an adiabatic self-trapping state by the optimal delayed temporal variation of the tunneling rate via either the energy bias between the two wells or the BEC nonlinearity. The emergence of nonlinear eigenstates and unstable stationary solutions of the system as well as their role in the breaking down of the adiabatic dynamics is investigated in detail.

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I. INTRODUCTION

Bose-Einstein condensates (BEC) in double-well potentials have drawn a lot of attention both theoretically and experimentally for the possibilities they offer to study fundamental quantum mechanical effects at the macroscopic level as well as for potential applications such as interferometry, high-precision measurements, or thermometry [1]. Most of the experimental realizations lead to the splitting of the condensate into two independent parts. Nevertheless, recently, weakly linked parts of a BEC in a double-well potential forming a single Josephson junction [2] have also been achieved [3]. In contrast to Josephson junctions realized in superconductors and superfluids [4], in BEC the nonlinear interatomic interactions play a crucial role. In the presence of the nonlinearity two dynamical regimes have been predicted: (i) anharmonic Josephson oscillations [5], if the initial population imbalance of the two wells is below a critical value, and (ii) macroscopic quantum self-trapping [6] (i.e., the inhibition of large-amplitude Josephson oscillations above the threshold for the population imbalance). This threshold corresponds to the population imbalance for which the difference between the two on-site interaction energies becomes larger than the tunneling energy splitting. Both dynamical regimes have been explored experimentally in a single Josephson junction [3] and in arrays [7].

Recently, a lot of attention has been devoted to explore techniques to coherently control the nonlinear dynamics of a BEC in double-well potentials by modulating in time either the potential [8] or the nonlinearity [9]. The two main studied regimes are the nonlinear Landau-Zener [10–16] and the Rosen-Zener [17] regimes. In the former the tunneling rate is fixed while the energy bias is linearly varied in time; in the latter the energy bias is varied in time while the tunneling rate is switching on and off following, for example, a temporal Gaussian profile. Both these regimes have been deeply investigated in double- [10–12, 15–17] and triple-well [13, 14] potentials, yielding a wide variety of dynamical scenarios ranging from robust population transfer to quantum blocking, nonlinear oscillations, and, even, breaking down of the adiabatic dynamics.

Following a different perspective, there have been several recent proposals to coherently manipulate single atoms [18–20] and BECs [13, 21, 22] in triple-well potentials by adiabatically following a particular energy eigenstate of the system, the so-called spatial dark state. This spatial dark state only involves the two ground states of the extreme traps in a close analogy to the well-known quantum optical stimulated Raman adiabatic passage (STIRAP) technique [23]. Accordingly, these tools were named three-level atom optics (TLAO) techniques [18]. In this paper, following the work by Vitanov and Shore [24] on the adiabatic passage on two-level atoms, we extend the TLAO techniques to the two-level matter-wave case showing that the adiabatic dynamics of a BEC, in a double-well potential, can be described in terms of a dark variable resulting from the combination of the population imbalance and the spatial atomic coherence. This dark variable can be tailored by varying in time the tunneling rate, the energy bias, and the nonlinearity with the goal of (i) adiabatically splitting of the BEC, (ii) achieving complete BEC transfer from one well to the other, and (iii) preparing an adiabatic self-trapping state.

The paper is organized as follows. In Sec. II we present the physical system consisting of a BEC in a double-well potential. We assume the two-level approximation and describe the BEC dynamics in terms of a dark variable. The conditions for the adiabatic control of tunneling by means of this dark variable are discussed in Sec. III from a nonlinear dynamics perspective. In Sec. IV we present detailed numerical simulations on the adiabatic splitting, transport, and trapping of a BEC. Section V summarizes the main conclusions of the paper and briefly discusses the validity of the two-mode approximation. It also presents some possible extensions of the present work such as its formulation in second quantization.

II. MODEL

We consider a BEC trapped in a double-well potential [Fig. 1(a)], whose dynamics at zero temperature is described by

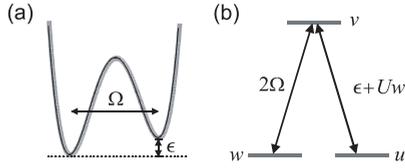


FIG. 1. (a) Two-level model for the BEC in a double-well potential with Ω the tunneling rate and ϵ the energy bias. (b) Three-level correspondence of (a) for the density matrix variables and coupling strengths. w is the population difference, $u/2$ ($v/2$) is the real (imaginary) part of the spatial coherence, and U is the nonlinear self-interaction energy ($\hbar = 1$).

the time-dependent Gross-Pitaevskii equation (GPE) ($\hbar = 1$)

$$i \frac{d\psi(\vec{r}, t)}{dt} = \left[-\frac{\Delta}{2m} + V(\vec{r}, t) + g|\psi(\vec{r}, t)|^2 \right] \psi(\vec{r}, t), \quad (1)$$

where $V(\vec{r}, t)$ is the external trapping potential and the nonlinearity is given by $g = 4N\pi a_s/m$, with N the total atom number, a_s the s -wave scattering length, and m the atomic mass. By assuming the two-level approximation [6,25], the wave function or classical order parameter of the BEC under study can be written as $\psi(\vec{r}, t) = c_L(t)\phi_L(\vec{r}) + c_R(t)\phi_R(\vec{r})$, with $c_{L(R)}(t) = \sqrt{N_{L(R)}}e^{i\theta_{L(R)}(t)}$ satisfying $|c_L|^2 + |c_R|^2 = 1$. $\phi_{L,R}(\vec{r})$ accounts for the ground state of the corresponding isolated trap. The amplitudes c_L and c_R obey the nonlinear two-mode dynamical equations given by

$$\mathcal{H} \begin{pmatrix} c_L \\ c_R \end{pmatrix} = i \frac{d}{dt} \begin{pmatrix} c_L \\ c_R \end{pmatrix}, \quad (2)$$

with the Hamiltonian

$$\mathcal{H} = \begin{pmatrix} \epsilon_L + U_L|c_L|^2 & \Omega \\ \Omega & \epsilon_R + U_R|c_R|^2 \end{pmatrix}, \quad (3)$$

where $U_{R,L}$ are the atomic self-interaction energies, Ω is the tunneling rate between the two wells, and $\epsilon_{R,L}$ are the on-site energies. In terms of the $\phi_{L,R}(\vec{r})$ overlaps, the former parameters read

$$\Omega = - \int d\vec{r} \left[\frac{1}{2m} \nabla \phi_L^* \nabla \phi_R + \phi_L^* V(\vec{r}, t) \phi_R \right], \quad (4)$$

$$U_{L,R} = g \int d\vec{r} |\phi_{L,R}|^4, \quad (5)$$

$$\epsilon_{R,L} = \int d\vec{r} \left[\frac{1}{2m} |\nabla \phi_{R,L}|^2 + \phi_{R,L}^* V(\vec{r}, t) \phi_{R,L} \right], \quad (6)$$

with $\int d\vec{r} [\phi_i^* \phi_j] = \delta_{ij}$ ($i, j = L, R$).

Notice that in the two-mode approximation one assumes that the parameters of the problem (tunneling rate, nonlinear interaction, and energy bias) can be varied independently. In general, however, the temporal modification of any of these parameters will result in the modification of the spatial mode functions affecting the rest of the parameters. From the experimental point of view, a double-well optical potential can be created with well separations in the range of few microns by means of two focused laser beams [26], or by the superposition of a 3D crossed beam dipole trap with a 1D optical lattice [3]. In the latter case, the intensity of the standing wave field that creates the optical lattice and its

displacement with respect to the center of the dipole trap could be used to manipulate both the tunneling and the energy bias. In addition, the scattering length could be controlled by adjusting a Feshbach resonance [27].

Within the density matrix formalism, assuming $U_R \sim U_L (\equiv U)$, and taking $\epsilon \equiv \epsilon_R - \epsilon_L$ as the energy bias between the two wells, one can write the coherent dynamics of the BEC in the two-well potential as follows [28]:

$$\frac{d}{dt} \begin{pmatrix} u \\ v \\ w \end{pmatrix} = \begin{pmatrix} 0 & -(\epsilon + Uw) & 0 \\ (\epsilon + Uw) & 0 & -2\Omega \\ 0 & 2\Omega & 0 \end{pmatrix} \begin{pmatrix} u \\ v \\ w \end{pmatrix}, \quad (7)$$

where we have introduced the real-valued variables $u = 2\text{Re}\{\sigma_{LR}\}$, $v = 2\text{Im}\{\sigma_{LR}\}$, and $w = \sigma_{RR} - \sigma_{LL}$, where $\sigma_{ii} = c_i c_i^*$ and $\sigma_{ij} = c_i c_j^*$ with $i, j = L, R$ the corresponding trap populations and spatial coherence [see Fig. 1(b)]. Note that the conservation of the norm implies $w^2 + u^2 + v^2 = 1$.

Inspired by the work of Vitanov and Shore [24] on STIRAP in a system of two internal atomic levels, we extend their proposal to the external degrees of freedom of matter waves. It is easy to verify that Eqs. (7) are equivalent to the discrete nonlinear Schrödinger equation (DNLS) for a BEC in a triple-well potential (see, for instance, Ref. [13]) under the following identifications: $u \Leftrightarrow C_R$, $v \Leftrightarrow -iC_M$, $w \Leftrightarrow C_L$, $\epsilon + Uw \Leftrightarrow \Omega_{MR}$, and $2\Omega \Leftrightarrow \Omega_{LM}$, where C_i with $i = L, M, R$ are the probability amplitudes of finding the BEC in the left, middle, or right trap, respectively. Ω_{LM} and Ω_{MR} denote, respectively, the tunneling interaction between the left-middle and middle-right traps. Note however that for the two-level system the variables u, v , and w are real quantities, while for a three-level system the variables C_i are, in general, complex numbers.

In the TLAO techniques within the three-level approximation with the left and right levels resonant, the transfer process is based on adiabatically following one of the three energy eigenstates of the system $|D(\Theta)\rangle = \cos \Theta |L\rangle - \sin \Theta |R\rangle$, with the mixing angle defined as $\Theta(t) = \tan^{-1}[\Omega_{LM}(t)/\Omega_{MR}(t)]$, and $|L\rangle$ and $|R\rangle$ being the ground states of the left and right wells. In the two-level system under investigation, the analog state will be given by a combination of the population difference and the coherence $d(\theta) = \cos \theta w - \sin \theta u$, with the mixing angle given by $\theta(t) = \tan^{-1}\{2\Omega(t)/[\epsilon(t) + U(t)w(t)]\}$. This analogy opens the possibility of extending the TLAO techniques to two-level systems, by appropriately engineering the time dependence of the tunneling rate $\Omega(t)$, the energy bias $\epsilon(t)$, and/or the nonlinear interaction $U(t)$. Note that the temporal variation of the scattering length a_s using either magnetic [29] or optical [30] Feshbach resonances or by varying the trap frequency, leading to a modification of the BEC spatial profile according to Eq. (5). For a review on the manipulation of Feshbach resonances see [31].

III. ADIABATIC CONTROL OF TUNNELING

In this section we study different scenarios for the coherent control of the external degrees of freedom of a BEC by adiabatically following the dark variable $d(\theta)$. In particular, we

show how to adiabatically split, transport, and inhibit tunneling of a BEC via the matter-wave analogs of both STIRAP and double-STIRAP techniques in two-level systems.

Let us assume that the BEC is initially prepared in the left trap with $\Omega(t = -\infty) = 0$. If so, note then that $\theta = 0$, meaning $w = -1$, and $u = v = 0$ will be the initial state. Then, to coherently split the condensate, θ should vary adiabatically from $\theta = 0$ to $\theta = \pi/2$, corresponding to $u = 1$ and $w = v = 0$ [see Fig. 1(b)] by appropriately changing Ω , ϵ , and U in time. To transfer the BEC from the left to the right trap, the mixing angle should be slowly increased to $\theta = \pi$. For the adiabatic self-trapping state case, the evolution will consist in varying θ from 0 to, for example, $\pi/2$ and then back to 0. In all cases, in order to guarantee that the BEC follows the dark variable $d(\theta)$ during the whole process two conditions must be fulfilled:

Condition 1: Adiabaticity criteria. $\Omega(t)$, $\epsilon(t)$, and $U(t)$ should be smoothly varied in time to adiabatically follow the dark variable $d(\theta)$, which, in turn, means that the mixing angle $\theta(t) = \tan^{-1} 2\{[\Omega(t)/[\epsilon(t) + U(t)w]]\}$ must be slowly changed from $\theta(t = -\infty) = 0$ to its expected final value. Gaussian profiles for the temporal variations of the control parameters are assumed:

$$\Omega = \Omega_g + \Omega_0 e^{-(t-t_\Omega)^2/\sigma_\Omega^2}, \quad (8)$$

$$\epsilon = \epsilon_g + \epsilon_0 [e^{-(t-t_\Omega + \Delta t_\epsilon)^2/\sigma_\epsilon^2} + n_\epsilon e^{-(t-t_\Omega - \Delta t_\epsilon)^2/\sigma_\epsilon^2}], \quad (9)$$

$$U = U_g + U_0 [e^{-(t-t_\Omega + \Delta t_U)^2/\sigma_U^2} + n_U e^{-(t-t_\Omega - \Delta t_U)^2/\sigma_U^2}], \quad (10)$$

with either the energy bias $\epsilon(t)$, or the BEC nonlinearity $U(t)$, preceding the tunneling interaction $\Omega(t)$ (i.e., the counter-intuitive sequences $\Delta t_\epsilon > 0$ or $\Delta t_U > 0$ will be assumed). Here $n_{\epsilon,U} = 0, \pm 1$ is a switch that takes the value 0 for the two-level matter-wave analog of the STIRAP sequence and $+1$ (-1) for the symmetric (antisymmetric) double-STIRAP sequences. Adiabaticity means that, at any time, $|\dot{\theta}(t)|$ should be much smaller than the energy separation between the selected eigenstate and the one energetically closest. For a weak nonlinear interaction, $|U(t)/\Omega(t)| \ll 1$, this adiabaticity condition reads $|\dot{\theta}(t)| \ll \sqrt{4\Omega(t)^2 + [\epsilon(t) + U(t)w]^2}$. It is worth remarking that, for $U \neq 0$, $\theta(t)$ is not a parameter of the system but a dynamical variable since it contains the population difference w in its definition.

Condition 2: Avoiding adverse bifurcation points. For large enough values of the nonlinearity, the interaction between the atoms of the BEC results in a nonlinear temporal coupling, producing additional nonlinear stationary states yielding loop structures and a rich variety of level-crossing scenarios [10, 13, 32, 33]. In the matter-wave STIRAP case for a triple-well potential [13], it has been shown that, even in the adiabatic limit given by $|\dot{\theta}(t)| \rightarrow 0$, the appearance of nonlinear stationary states breaks down, in some cases, the adiabatic evolution. To address this issue in the double-well potential, we start by first looking for those critical U values giving rise to nonlinear stationary states by solving the eigenvalue equation

$$\mathcal{H} \begin{pmatrix} c_L \\ c_R \end{pmatrix} = \mu \begin{pmatrix} c_L \\ c_R \end{pmatrix}, \quad (11)$$

with \mathcal{H} given in Eq. (3) and μ being the chemical potential. After some algebra one obtains the following fourth-order eigenvalue equation for μ :

$$[\epsilon^2 - (-2\mu + \epsilon + 2U)^2](-2\mu + \epsilon + U)^2 + 4\Omega^2(-2\mu + \epsilon + 2U)^2 = 0. \quad (12)$$

For $\epsilon = 0$ and $U^2/(2\Omega)^2 < 1$ (> 1) this quartic equation gives two (four) real roots [10].

Additional information on the BEC dynamics can be obtained by looking for the stationary solutions of the density matrix equations (7) and analyzing their stability (see the Appendix). The stationary solutions can be written as $\{u^{ss}, v^{ss} = 0, w^{ss}\}$, with $u^{ss}/w^{ss} = \tan \theta^{ss}$ and $(u^{ss})^2 + (w^{ss})^2 = 1$. The energy of these stationary solutions (for which there are up to four) is given by Eq. (12). Note that, as the dynamics is conservative, the eigenvalues of the linear stability matrix must satisfy $\sum_{i=1}^3 \lambda_i = 0$ for each stationary solution. In fact, a linear stability analysis (LSA) around any of these solutions yields one eigenvalue equal to zero together with either two pure imaginary eigenvalues, corresponding to a stable fixed point (or center), or two real eigenvalues, accounting for an unstable saddle point. In the limit of a large nonlinear interaction, that is, for $U^2/(2\Omega)^2 > 1$, the system presents four stationary solutions, one of them being an unstable saddle point while the other three are stable elliptical ones.

If the number of stationary solutions remains constant during the whole adiabatic dynamics, the system will successfully follow the selected energy eigenstate. However, depending on the interplay between the nonlinear interaction and the tunneling rate, during the dynamics the number of stationary solutions will change through a bifurcation point from four to two or vice versa. If the selected stationary solution is not involved in the bifurcation, the system will again adiabatically follow the corresponding eigenstate. For the opposite case, whether the adiabatic dynamics will be affected will depend on the particular bifurcation scenario:

Case 1. From four to two stationary solutions obtained through a backward pitchfork bifurcation point in which the stationary solution to be followed merges two more solutions (one unstable) and yields one stable solution. In this case, the system will evolve adiabatically. In what follows, this case will be termed the PB42 case, even if the pitchfork bifurcation is perturbed and becomes imperfect.

Case 2. From two to four solutions obtained through a forward pitchfork bifurcation point in which the stationary solution to be followed becomes unstable and yields two stable solutions corresponding, in the limit of vanishing tunneling, to $w = \pm 1$. In this case, after reaching the bifurcation point, the system will split into a combination of the two stationary solutions. In this case, named PB24 in what follows, the adiabatic dynamics will, in general, break down.

Case 3. From four to two solutions obtained through a saddle-node bifurcation in which the stationary solution to be followed is annihilated by an unstable solution. The adiabatic dynamics breaks down even in the adiabatic limit. After reaching the bifurcation the system will split into a combination of two nondegenerate energy eigenstates and,

therefore, will oscillate at the corresponding energy splitting. This is termed case SNB42 in what follows.

IV. NUMERICAL SIMULATIONS

This section is devoted to the numerical simulations of the BEC adiabatic splitting, transport, and self-trapping by means of the temporal variation of the energy bias, the nonlinearity, and the tunneling rate. We will show, for the most relevant cases, the eigenvalues and stationary states predicted by Eqs. (12) and (7) as well as their linear stability. The main goal of these numerical simulations will consist in illustrating the different dynamical scenarios to adiabatically control the BEC while looking for those parameter values that prevent the system from reaching the two unwanted bifurcation cases PB24 and SNB42 previously described.

A. Adiabatic splitting of a BEC

We start first by fixing the nonlinear interaction control parameter U to a constant value U_g while, counterintuitively, time-varying the energy bias and the tunneling rate [see the inset of Fig. 2(b)] with the goal of achieving equal population in the two wells, that is, $w(t_{\text{out}}) = 0$ starting with the BEC being in the left trap [i.e., $w(t_{\text{in}}) = -1$], or, equivalently, $\theta(t_{\text{out}}) = \pi/2$ from $\theta(t_{\text{in}}) = 0$. Figure 2(a) shows $w(t_{\text{out}})$, after the numerical integration of Eqs. (7) with $w(t_{\text{in}}) = -1$ and the rest of parameter values given in the figure caption. Clearly, there is a large set of parameters where the adiabatic splitting process takes place with a high fidelity [see the plateau in Fig. 2(a)]. Figure 2(b) shows the plateau $w(t_{\text{out}}) = 0$ for $U_g/\Omega_0 = (-0.12, 1.5)$ and $\epsilon_0 = 1.5\Omega_0$. To give more insight into the dynamics of the system, we have divided the nonlinear interaction range studied into five regions, from a to e , as shown at the top of Fig. 2(b). In the following we will discuss in detail the dynamics for each of these regions. Note that, in all cases, the dynamics starts and ends with $\Omega(t_{\text{in}}) = \Omega(t_{\text{out}}) = 0$, which, for $U \neq 0$, implies that for $t = t_{\text{in,out}}$ the system presents four fixed points or stationary solutions. For intermediate times and large enough tunneling rates, there are only two stationary solutions.

In Fig. 3 we show the temporal dynamics corresponding to the parameter values of case (i) in region c of Fig. 2(b) with $U_g = 0.5\Omega_0$ [see Fig. 3(a)]. The adiabatic energy eigenvalues and stationary solutions w^{ss} are shown in Figs. 3(b) and 3(c), respectively, with the dashed curve accounting for the unstable ones. Short arrows in Figs. 3(b) and 3(c) indicate the stationary solution to be adiabatically followed. The dynamical variable $w(t)$, after integration of Eqs. (7), is plotted in Fig. 3(d). For this set of parameter values the adiabaticity condition is fulfilled and the selected eigenstate is involved in a single bifurcation, at $\Omega_0 t = 777$, corresponding to the PB42 case previously described. After the bifurcation point, the system follows the eigenstate that, at the end of the process, yields $w = 0$ and $\mu = 0.25\Omega_0$. This nonlinear dynamical scenario holds for the whole plateau, region c in Fig. 2(b), where the adiabatic splitting succeeds.

For the parameter range $U_g/\Omega_0 = (1.5, 2.07)$, corresponding to region d in Fig. 2(b), the final state of the BEC strongly depends on the parameter values. For a slight modification of

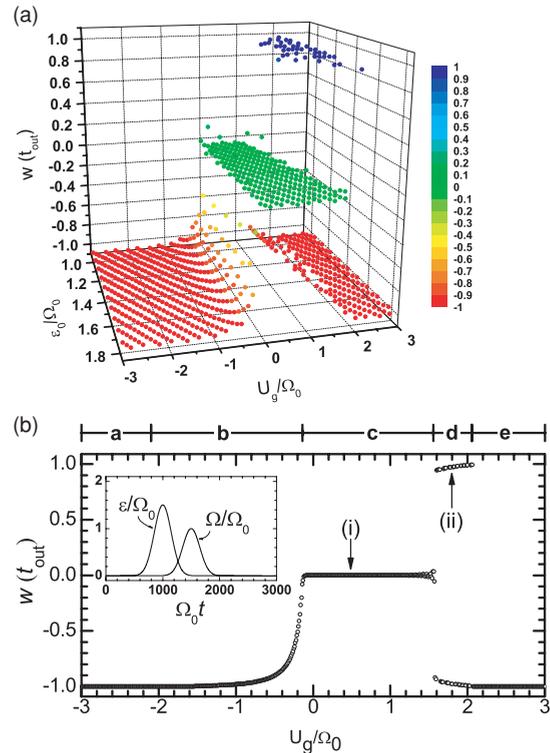


FIG. 2. (Color online) Adiabatic splitting of a BEC. (a) Population difference $w(t_{\text{out}})$ at the end of the splitting process in the parameter plane modulation amplitude of the energy bias ϵ_0 versus the nonlinear interaction parameter U_g . Initial conditions: $w(t_{\text{in}}) = -1$ and $u(t_{\text{in}}) = v(t_{\text{in}}) = 0$. Parameter setting: $\sigma_\epsilon = \sigma_\Omega = 212.8\Omega_0^{-1}$, $\Delta t_\epsilon = 500\Omega_0^{-1}$, $t_\Omega = 1500\Omega_0^{-1}$, and $\Omega_g = \epsilon_g = n_\epsilon = U_0 = n_U = 0$. (b) Same as in (a) for the fixed value $\epsilon_0 = 1.5\Omega_0$. The temporal variation of the energy bias and the tunneling rate are shown in the inset. From a nonlinear dynamics perspective, we have classified the results into five different regions, from a to e , depending on the number and type of bifurcations that the system suffers during its dynamics (see text). The detailed temporal dynamics corresponding to cases (i) and (ii) are plotted in Figs. 3 and 4, respectively. The adiabatic splitting succeeds in region c , corresponding to $U_g/\Omega_0 = (-0.12, 1.5)$.

U_g , the final state alternates between $w \sim +1$ and $w \sim -1$, accounting for the BEC being located at the right or left trap, respectively. To understand the origin of this behavior, we plot in Fig. 4 the detailed dynamics for $U_g = 1.7\Omega_0$, corresponding to case (ii) in Fig. 2(b). At $\Omega_0 t = 1374$, the solution that the system is adiabatically following is annihilated with an unstable one through a saddle-node bifurcation (i.e., case SNB42 previously described). At this point, the system splits into two nondegenerate components (being the largest and energetically closest stable solution) and starts to oscillate at a frequency corresponding to the energy separation of the two solutions. At $\Omega_0 t = 1588$, the system reaches a bifurcation point of the PB24 type. This bifurcation yields two energetically degenerate stable solutions, corresponding to the two final states $w = \pm 1$. Thus, at $\Omega_0 t = 1588$, the largest component of the system chooses one out of the two new stable solutions. We have numerically verified that the selected solution strongly depends on the previous oscillatory dynamics.

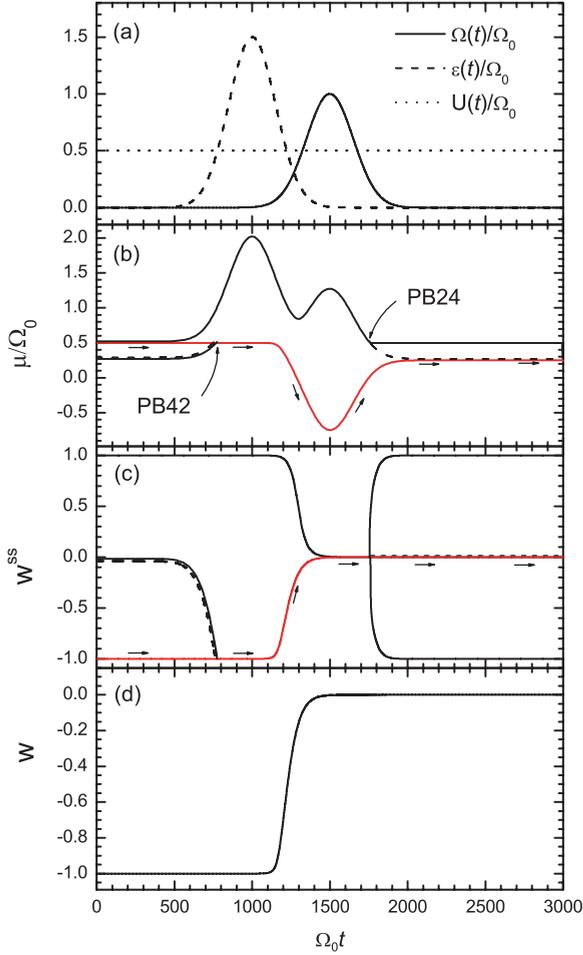


FIG. 3. (Color online) Adiabatic splitting of a BEC. $U_g = 0.5\Omega_0$ and the rest of the parameters are as in Fig. 2. Plotted as a function of time are (a) the nonlinear interaction, energy bias, and tunneling rate profiles; (b) energy eigenvalues; (c) stationary solutions for the population difference w^{ss} ; and (d) the population difference $w(t)$ after numerical integration of Eqs. (7). In (b) and (c), the dashed curve accounts for the unstable solution, while the short arrows indicate the adiabatic solution (in red) to be followed by the system. PB42 and PB24 denote the two pitchfork bifurcations that the system suffers at $\Omega_0 t = 777$ and $\Omega_0 t = 1750$, respectively. Note that the pitchfork bifurcation PB24 yields one unstable solution (dashed curve) plus two energetically degenerated stable solutions (solid curve), corresponding, in the limit of vanishing tunneling, to $w = \pm 1$.

For $|U_g| > 2.07\Omega_0$, corresponding to regions *a* and *e* of Fig. 2(b), there are four stationary solutions during the whole dynamics that do not present any bifurcation. In this case, the nonlinear coupling forces the adiabatic evolution from $\theta(t_{in}) = 0$ to $\theta(t_{out}) = 0$ instead of evolving to the desired $\theta(t_{out}) = \pi/2$. Then, the initial $(u, v, w) = (0, 0, -1)$ and the final state of the system coincide.

Finally, for $-2.07\Omega_0 < U_g < -0.12\Omega_0$, region *b* in Fig. 2, the energy eigenstate to be followed reaches, during the ramping down of the tunneling interaction $\Omega(t)$, a pitchfork bifurcation of the PB24 type. After this bifurcation point, the two stable stationary solutions are energetically degenerated and, therefore, the system splits into a nonoscillatory

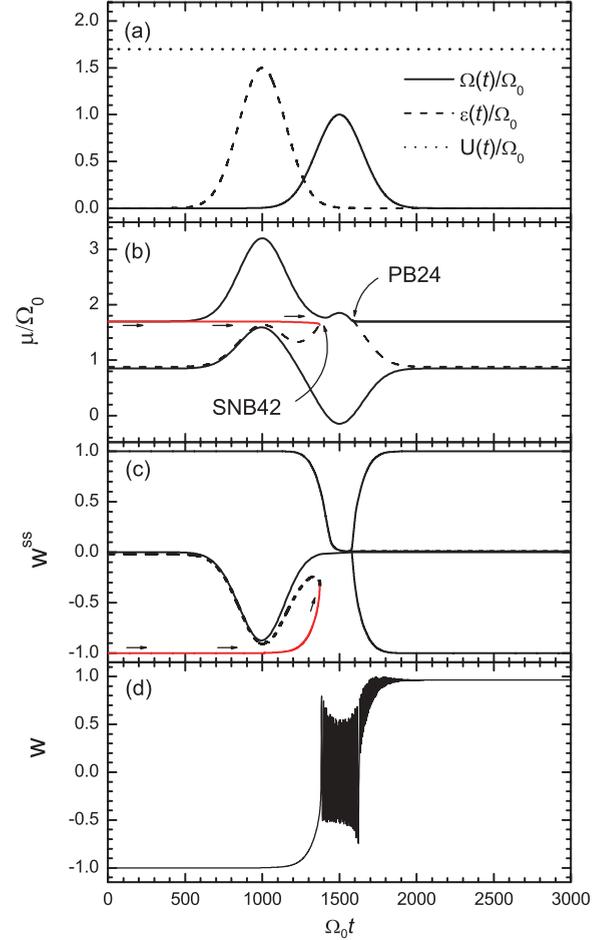


FIG. 4. (Color online) Same as in Fig. 2 for $U_g = 1.7\Omega_0$. SNB42 and PB24 account for a saddle-node and a pitchfork bifurcation at $\Omega_0 t = 1374$ and $\Omega_0 t = 1588$, respectively.

combination of them that for $U_g = -0.12\Omega_0$ ($U_g = -2.07\Omega_0$) yields $w(t_{out}) = 0$ [$w(t_{out}) = -1$].

Up to now, we have discussed the possibility of adiabatically splitting a BEC by fixing U and time-varying $\epsilon(t)$ and $\Omega(t)$. Note, however, that it is also possible to adiabatically split the BEC by appropriately time-varying the nonlinear interaction $U(t)$ and the tunneling rate $\Omega(t)$ with a fixed energy bias. Thus, for $U_g = -0.4\Omega_0$, $U_0 = 1.5\Omega_0$, $\epsilon = 0$, and the rest of the parameters given in the figure caption, Fig. 5 shows $w(t_{out})$ as a function of the time delay between the modulation of the tunneling rate and the nonlinear interaction. We again obtain a plateau of parameter values, $\Omega_0 \Delta t_U \simeq (250, 600)$, where the adiabatic splitting, corresponding to $w(t_{out}) \sim 0$, becomes successful. A detailed analysis of the temporal dynamics reveals that, as in Fig. 2, the plateau corresponds to the case where the selected solution reaches during its dynamics a single bifurcation point of the PB42 type.

B. Adiabatic transport

To adiabatically transfer the BEC from the left to the right trap, the mixing angle must be slowly varied from $\theta(t_{in}) = 0$ to $\theta(t_{out}) = \pi$. By taking $n_\epsilon = -1$ and the rest of

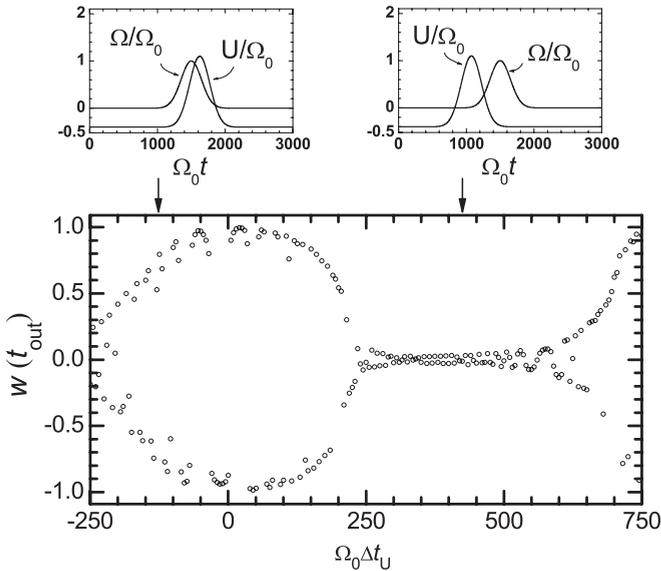


FIG. 5. Adiabatic splitting of a BEC. Plotted is the population difference $w(t_{out})$ at the end of the BEC splitting process as a function of the temporal delay Δt_U . The temporal variation of the nonlinear interaction and the tunneling rate is shown in the two upper figures for $\Delta t_U = -125\Omega_0^{-1}$ (left) and $\Delta t_U = 425\Omega_0^{-1}$ (right). Initial conditions: $w(t_{in}) = -1$ and $u(t_{in}) = v(t_{in}) = 0$. Parameter setting: $U_g = -0.4\Omega_0$, $U_0 = 1.5\Omega_0$, $\sigma_U = \sigma_\Omega = 212.8\Omega_0^{-1}$, $t_\Omega = 1500\Omega_0^{-1}$, and $\epsilon_g = \epsilon_0 = n_\epsilon = n_U = 0$. The adiabatic splitting succeeds in the parameter region $\Omega_0 \Delta t_U \simeq (250, 600)$.

parameters as in Fig. 2, Fig. 6 shows the population difference $w(t_{out})$ at the end of the antisymmetric double-STIRAP process (shown in the inset) as a function of the nonlinear interaction U_g . Clearly, there is a wide parameter setting, the plateau $U_g/\Omega_0 = (-2, 1.4)$, where the BEC transfer process takes place with a high fidelity. In the parameter region $U_g/\Omega_0 = (-2, 0)$, the system starts at $w(t_{in}) = -1$ following an energy eigenstate of the system that does not involve any bifurcation point. For $U_g/\Omega_0 = (0, 1.4)$ the system crosses first a PB42 bifurcation to later on reaches a PB24 bifurcation point that, for these parameters, yields eventually $w(t_{out}) = 1$. In regions

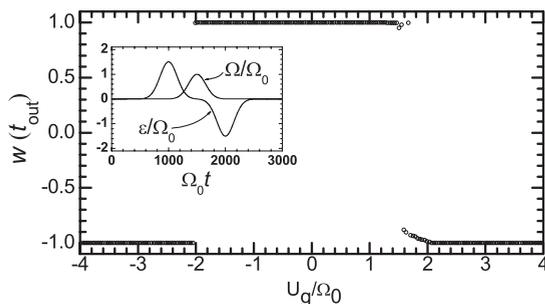


FIG. 6. Adiabatic transport of a BEC. Plotted is the population difference $w(t_{out})$ at the end of the BEC transport process as a function of the nonlinear interaction parameter U_g . The temporal variation of the energy bias and the tunneling rate are shown in the inset. Initial conditions: $w(t_{in}) = -1$ and $u(t_{in}) = v(t_{in}) = 0$. Parameter setting: $\epsilon_0 = 1.5\Omega_0$, $\sigma_\epsilon = \sigma_\Omega = 212.8\Omega_0^{-1}$, $\Delta t_\epsilon = 500\Omega_0^{-1}$, $t_\Omega = 1500\Omega_0^{-1}$, $n_\epsilon = -1$, and $U_0 = \epsilon_g = n_U = 0$. The adiabatic transport succeeds in the parameter region $U_g/\Omega_0 = (-2, 1.4)$.

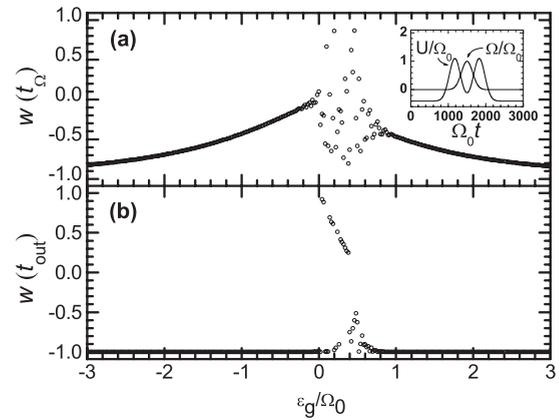


FIG. 7. Adiabatic self-trapping of a BEC. Plotted is the population difference at (a) the intermediate time $t = t_\Omega$ and (b) the end of the BEC transport process, as a function of the energy bias ϵ_g . The temporal variation of the nonlinear interaction and the tunneling rate are shown in the inset for $U_0 = 1.5\Omega_0$. Initial conditions: $w(t_{in}) = -1$ and $u(t_{in}) = v(t_{in}) = 0$. Parameter setting: $U_g = -0.4\Omega_0$, $\sigma_U = \sigma_\Omega = 212.8\Omega_0^{-1}$, $\Delta t_U = 325\Omega_0^{-1}$, $t_\Omega = 1500\Omega_0^{-1}$, $n_U = 1$, and $\epsilon_0 = U_0 = n_\epsilon = 0$. The adiabatic self-trapping state does not succeed in the parameter region $\epsilon_g/\Omega_0 = (0, 0.8)$.

$U_g/\Omega_0 = (-4, -2)$ and $U_g/\Omega_0 = (1.4, 4)$ the nonlinearity yields $\theta(t_{out}) = 0$, forcing the initial and final state to be the same.

C. Adiabatic self-trapping

It is also possible to completely inhibit the BEC transport between the two wells by means of the matter-wave analog to the double-STIRAP process (see inset of Fig. 7). Starting with the BEC located at the left trap, $w(t_{in}) = -1$, we show in Fig. 7 the population difference at the time when the tunneling rate is maximum, $w(t_\Omega)$, and at the end of the double-STIRAP sequence, $w(t_{out})$, as a function of the energy bias ϵ_g . The adiabatic self-trapping process succeeds for a wide domain of parameter values, except for the region $\epsilon_g/\Omega_0 = (0, 0.8)$. In this last region, the solution to be followed reaches during its dynamics a saddle-point bifurcation while, for the rest of the values of the energy bias, the selected solution does not involve any bifurcation point.

V. CONCLUSIONS

By means of a dark variable that results from the combination of the population imbalance and the spatial atomic coherence, we have investigated in detail the adiabatic dynamics of a BEC in a double-well potential in the framework of the two-level approximation. We have shown that it is possible to robustly split, transport, or trap a BEC by appropriate temporal variation of either the energy bias or the nonlinear interaction together with the tunneling rate. All these proposals have been studied from a nonlinear dynamics perspective by deriving the stationary solutions of the system, evaluating their linear stability, and discussing the bifurcation scenarios. For the eventual implementation of the techniques discussed here, we want to highlight the recent work by Bauer *et al.* [27], who

report an accurate control on magnetic Feshbach resonances by means of laser light.

In order to check the validity of the two-mode approximation, we have numerically integrated the one-dimensional Gross-Pitaevskii (GP) equation for the nonlinear adiabatic splitting of a BEC in a double-well potential, obtaining the characteristic “plateau” of the STIRAP protocol shown in Fig. 2(b). However, a detailed numerical investigation of the GP equation still remains to be performed to validate that the matter-wave nonlinear STIRAP techniques derived here in the two-mode approximation could be used for matter-wave interferometry or coherent transport of a BEC. In performing this analysis, the global landscape provided by the results of the two-mode approximation should be a useful roadmap.

Although it is beyond the scope of the present paper, it would be very interesting also to extend the present nonlinear matter-wave STIRAP techniques to the second-quantization formalism [34]. Within this formalism, we could elucidate whether the adiabatic splitting of the BEC results in a macroscopic superposition where the entire condensate localizes in one of the wells, that is, in a NOON state [35], or in a perfect 50% spatial splitting of all the atoms. Note that in both cases the final population difference reads $w(t_{\text{out}}) = 0$.

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APPENDIX: LINEAR STABILITY ANALYSIS

The stationary solutions of the density matrix equations (7) take the form $\{u^{\text{ss}}, v^{\text{ss}} = 0, w^{\text{ss}}\}$ with $u^{\text{ss}}/w^{\text{ss}} = 2\Omega/(\epsilon + Uw^{\text{ss}})$ and $(u^{\text{ss}})^2 + (w^{\text{ss}})^2 = 1$. Let us consider now a perturbation around any of these solutions in the form of

$$u(t) = u^{\text{ss}} + \delta u e^{\lambda t}, \quad (\text{A1})$$

$$v(t) = \delta v e^{\lambda t}, \quad (\text{A2})$$

$$w(t) = w^{\text{ss}} + \delta w e^{\lambda t}, \quad (\text{A3})$$

where the real part of λ determines the stability of the solution. The stability is governed by the equation $(\hat{M} - \lambda \hat{I})\delta\vec{s} = 0$, where $\delta\vec{s} = (\delta u, \delta v, \delta w)^T$ and \hat{M} is the linear stability matrix defined as

$$\hat{M} \equiv \begin{pmatrix} 0 & -(\epsilon + Uw^{\text{ss}}) & 0 \\ (\epsilon + Uw^{\text{ss}}) & 0 & Uu^{\text{ss}} - 2\Omega \\ 0 & 2\Omega & 0 \end{pmatrix}. \quad (\text{A4})$$

Solving the secular equation $\det(\hat{M} - \lambda \hat{I}) = 0$, one obtains the following three eigenvalues:

$$\lambda_0 = 0, \quad (\text{A5})$$

$$\lambda_{\pm} = \pm \sqrt{2\Omega(Uu^{\text{ss}} - 2\Omega) - (\epsilon + Uw^{\text{ss}})^2}. \quad (\text{A6})$$

As expected from a conservative system, $\sum_{i=1}^3 \lambda_i = 0$ for every stationary solution. One eigenvalue is equal to zero while the other two are either pure imaginary, corresponding to a stable fixed point (or center), or pure real, accounting for an unstable one (or saddle point).

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- [1] M. R. Andrews, C. G. Townsend, H.-J. Miesner, D. S. Durfee, D. M. Kurn, and W. Ketterle, *Science* **275**, 637 (1997); T. G. Tiecke, M. Kemmann, Ch. Buggle, I. Schvachuck, W. von Klitzing, and J. T. M. Walraven, *J. Opt. B: Quantum Semiclass. Opt.* **5**, S119 (2003); Y. Shin, M. Saba, A. Schirotzek, T. A. Pasquini, A. E. Leanhardt, D. E. Pritchard, and W. Ketterle, *Phys. Rev. Lett.* **92**, 150401 (2004); J. Estève, J.-B. Trebbia, T. Schumm, A. Aspect, C. I. Westbrook, and I. Bouchoule, *ibid.* **96**, 130403 (2006); B. V. Hall, S. Whitlock, R. Anderson, P. Hannaford, and A. I. Sidorov, *ibid.* **98**, 030402 (2007); U. Hohenester, P. K. Rekdal, A. Borzi, and J. Schmiedmayer, *Phys. Rev. A* **75**, 023602 (2007); J. Esteve, C. Gross, A. Weller, S. Giovanazzi, and M. K. Oberthaler, *Nature (London)* **455**, 1216 (2008).
- [2] B. D. Josephson, *Phys. Lett.* **1**, 251 (1962).
- [3] M. Albiez, R. Gati, J. Fölling, S. Hunsmann, M. Cristiani, and M. K. Oberthaler, *Phys. Rev. Lett.* **95**, 010402 (2005).
- [4] P. L. Anderson, and J. W. Rowell, *Phys. Rev. Lett.* **10**, 230 (1963); S. V. Pereverzev, A. Loshak, S. Backhaus, J. C. Davis, and R. E. Packard, *Nature (London)* **388**, 449 (1997); A. K. Sukhatme, Y. Mukharsky, T. Chui, and D. Pearson, *ibid.* **411**, 280 (2001).
- [5] J. Javanainen, *Phys. Rev. Lett.* **57**, 3164 (1986); M. W. Jack, M. J. Collett, and D. F. Walls, *Phys. Rev. A* **54**, R4625 (1996); I. Zapata, F. Sols, and A. J. Leggett, *ibid.* **57**, R28 (1998).
- [6] G. J. Milburn, J. Corney, E. M. Wright, and D. F. Walls, *Phys. Rev. A* **55**, 4318 (1997); A. Smerzi, S. Fantoni, S. Giovanazzi, and S. R. Shenoy, *Phys. Rev. Lett.* **79**, 4950 (1997); S. Raghavan, A. Smerzi, S. Fantoni, and S. R. Shenoy, *Phys. Rev. A* **59**, 620 (1999).
- [7] F. S. Cataliotti, S. Burger, C. Fort, P. Maddaloni, F. Minardi, A. Trombettoni, A. Smerzi, and M. Inguscio, *Science* **293**, 843 (2001); F. S. Cataliotti, L. Fallani, F. Ferlaino, C. Fort, P. Maddaloni, and M. Inguscio, *New J. Phys.* **5**, 71 (2003); T. Anker, M. Albiez, R. Gati, S. Hunsmann, B. Eiermann, A. Trombettoni, and M. K. Oberthaler, *Phys. Rev. Lett.* **94**, 020403 (2005).
- [8] F. Kh. Abdullaev and R. A. Kraenkel, *Phys. Rev. A* **62**, 023613 (2000); M. Holthaus, *ibid.* **64**, 011601(R) (2001); C. Lee, W. Hai, L. Shi, X. Zhu, and K. Gao, *ibid.* **64**, 053604 (2001); G. F. Wang, L. B. Fu, and J. Liu, *ibid.* **73**, 013619 (2006); Q. Xie and W. Hai, *ibid.* **75**, 015603 (2007).
- [9] F. K. Abdullaev and R. A. Kraenkel, *Phys. Lett. A* **272**, 395 (2000); F. K. Abdullaev, J. S. Shaari, and M. R. B. Wahiddin,

- ibid.* **345**, 237 (2005); C. Lee, *Phys. Rev. Lett.* **102**, 070401 (2009).
- [10] B. Wu and Q. Niu, *Phys. Rev. A* **61**, 023402 (2000).
- [11] J. Liu, B. Wu, and Q. Niu, *Phys. Rev. Lett.* **90**, 170404 (2003).
- [12] G. Theocharis, P. G. Kevrekidis, D. J. Frantzeskakis, and P. Schmelcher, *Phys. Rev. E* **74**, 056608 (2006).
- [13] E. M. Graefe, H. J. Korsch, and D. Witthaut, *Phys. Rev. A* **73**, 013617 (2006).
- [14] G.-F. Wang, D.-F. Ye, L.-B. Fu, X.-Z. Chen, and J. Liu, *Phys. Rev. A* **74**, 033414 (2006).
- [15] A. P. Itin and S. Watanabe, *Phys. Rev. E* **76**, 026218 (2007).
- [16] Q. Zhang, P. Hänggi, and J. Gong, *Phys. Rev. A* **77**, 053607 (2008).
- [17] D.-F. Ye, L.-B. Fu, and J. Liu, *Phys. Rev. A* **77**, 013402 (2008).
- [18] K. Eckert, M. Lewenstein, R. Corbalán, G. Birkel, W. Ertmer, and J. Mompart, *Phys. Rev. A* **70**, 023606 (2004); Th. Busch, K. Deasy, and S. Nic Chormaic, *J. Phys. Conf. Ser.* **84**, 012002 (2007).
- [19] K. Eckert, J. Mompart, R. Corbalán, M. Lewenstein, and G. Birkel, *Opt. Commun.* **264**, 264 (2006).
- [20] T. Opatrný and K. K. Das, *Phys. Rev. A* **79**, 012113 (2009).
- [21] M. Rab, J. H. Cole, N. G. Parker, A. D. Greentree, L. C. L. Hollenberg, and A. M. Martin, *Phys. Rev. A* **77**, 061602(R) (2008).
- [22] V. O. Nesterenko, A. N. Nikonov, F. F. de Souza Cruz, and E. L. Lapolli, *Laser Phys.* **19**, 616 (2009).
- [23] K. Bergmann, H. Theuer, and B. W. Shore, *Rev. Mod. Phys.* **70**, 1003 (1998).
- [24] N. V. Vitanov and B. W. Shore, *Phys. Rev. A* **73**, 053402 (2006).
- [25] I. Zapata, F. Sols, and A. J. Leggett, *Phys. Rev. A* **57**, R28 (1998); G. J. Milburn, J. Corney, E. M. Wright, and D. F. Walls, *ibid.* **55**, 4318 (1997); R. W. Spekkens and J. E. Sipe, *ibid.* **59**, 3868 (1999); S. Giovanazzi, A. Smerzi, and S. Fantoni, *Phys. Rev. Lett.* **84**, 4521 (2000); D. Ananikian and T. Bergeman, *Phys. Rev. A* **73**, 013604 (2006).
- [26] Y. Shin, M. Saba, T. A. Pasquini, W. Ketterle, D. E. Pritchard, and A. E. Leanhardt, *Phys. Rev. Lett.* **92**, 050405 (2004).
- [27] D. M. Bauer, M. Lettner, C. Vo, G. Rempe, and S. Dürr, *Nature Phys.* **5**, 339 (2009).
- [28] C. Lee, W. Hai, L. Shi, and K. Gao, *Phys. Rev. A* **69**, 033611 (2004).
- [29] S. Inouye, M. R. Andrews, J. Stenger, H.-J. Miesner, D. M. Stamper-Kurn, and W. Ketterle, *Nature (London)* **392**, 151 (1998); E. A. Donley, N. R. Claussen, S. L. Cornish, J. L. Roberts, E. A. Cornell, and C. E. Wieman, *ibid.* **412**, 295 (2001); H. Saito and M. Ueda, *Phys. Rev. A* **65**, 033624 (2002).
- [30] P. O. Fedichev, Yu. Kagan, G. V. Shlyapnikov, and J. T. M. Walraven, *Phys. Rev. Lett.* **77**, 2913 (1996); J. L. Bohn and P. S. Julienne, *Phys. Rev. A* **56**, 1486 (1997); F. K. Fatemi, K. M. Jones, and P. D. Lett, *Phys. Rev. Lett.* **85**, 4462 (2000); M. Theis, G. Thalhammer, K. Winkler, M. Hellwig, G. Ruff, R. Grimm, and J. H. Denschlag, *ibid.* **93**, 123001 (2004); R. Ciurylo, E. Tiesinga, and P. S. Julienne, *Phys. Rev. A* **71**, 030701(R) (2005).
- [31] C. Chin, R. Grimm, P. Julienne, and E. Tiesinga, e-print [arXiv:0812.1496](https://arxiv.org/abs/0812.1496) [quant-ph].
- [32] M. Holthaus, *Phys. Rev. A* **64**, 011601(R) (2001).
- [33] R. D'Agosta and C. Presilla, *Phys. Rev. A* **65**, 043609 (2002).
- [34] G.-B. Jo, Y. Shin, S. Will, T. A. Pasquini, M. Saba, W. Ketterle, D. E. Pritchard, M. Vengalattore, and M. Prentiss, *Phys. Rev. Lett.* **98**, 030407 (2007); G.-B. Jo, J.-H. Choi, C. A. Christensen, T. A. Pasquini, Y.-R. Lee, W. Ketterle, and D. E. Pritchard, *ibid.* **98**, 180401 (2007); J. Estève, C. Gross, A. Weller, S. Giovanazzi, and M. K. Oberthaler, *Nature (London)* **455**, 1216 (2008).
- [35] D. W. Hallwood, K. Burnett, and J. Dunningham, *New J. Phys.* **8**, 180 (2006); T. J. Haigh, A. J. Ferris, and M. K. Olsen, e-print [arXiv:0907.1333](https://arxiv.org/abs/0907.1333) [quant-ph]; D. W. Hallwood, A. Stokes, J. J. Cooper, and J. Dunningham, *New J. Phys.* **11**, 103040 (2009).