

Split and overlapped binary solitons in optical latticesGolam Ali Sekh,^{1,2} Francesco V. Pepe,^{3,4,1} Paolo Facchi,^{4,1} Saverio Pascazio,^{4,1} and Mario Salerno⁵¹*Istituto Nazionale di Fisica Nucleare (INFN), Sezione di Bari, I-70126 Bari, Italy*²*Department of Physics, University of Kashmir, Hazratbal, Srinagar-190006, J & K, India*³*Museo Storico della Fisica e Centro Studi e Ricerche “Enrico Fermi,” I-00184 Roma, Italy*⁴*Dipartimento di Fisica and MECENAS, Università di Bari, I-70126 Bari, Italy*⁵*Dipartimento di Fisica “E. R. Caianiello,” INFN–Gruppo Collegato di Salerno, and CNISM, Università di Salerno, I-84084 Fisciano, Italy*

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We analyze the energetic and dynamical properties of bright-bright (BB) soliton pairs in a binary mixture of Bose-Einstein condensates subjected to the action of a combined optical lattice, acting as an external potential for the first species while modulating the intraspecies coupling constant of the second. In particular, we use a variational approach and direct numerical integrations to investigate the existence and stability of bright-bright (BB) solitons in which the two species are either spatially separated (split soliton) or located at the same optical lattice site (overlapped soliton). The dependence of these solitons on the interspecies interaction parameter is explicitly investigated. For repulsive interspecies interaction we show the existence of a series of critical values at which transitions from an initially overlapped soliton to split solitons occur. For attractive interspecies interaction only single direct transitions from split to overlapped BB solitons are found. The possibility to use split solitons for indirect measurements of scattering lengths is also suggested.

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I. INTRODUCTION

Bose-Einstein condensates (BECs) are fascinating tools for simulating different physical systems. Advanced laser technology and its successful applications to ultracold atoms have enabled us to engineer potentials of different geometries. A well-established technique consists in creating a linear optical lattice (LOL) by interfering pairs of counterpropagating laser beams [1]. On the other hand, laser beams can also be used to vary atomic interaction periodically in space with the help of optical Feshbach resonances [2]. Periodically modulated atomic interaction leads to a nonlinear optical lattice (NOL). BEC in LOL has been used to investigate different physical phenomena in condensed matter physics, including Bloch oscillations [3,4], generation of coherent atomic pulses (atom laser) [5], dynamical localization [6,7], Landau-Zener tunneling [8–10], and superfluid-Mott transition [11].

Interatomic interaction in BECs gives rise to a nonlinearity which permits localized bound states to remain stable for a long time, due to the balance between the effects of nonlinearity and dispersion. In the presence of a LOL, the interplay between lattice periodicity and interatomic interaction was shown to induce modulation instabilities of Bloch wave functions near the band edges [12], leading to the formation of localized excitations with chemical potentials inside band gaps, the so-called gap solitons (GSs). These excitations have been investigated both for continuous BECs, in one-dimensional [13–17] and multidimensional [18–20] settings, and for the discrete case (BEC arrays) [21,22] in the presence of attractive and repulsive interactions. NOL can also support special kinds of solitons both in 1D [23] and in multidimensional settings in combination with LOL [24,25]. NOLs have been used to avoid dynamical instabilities of gap solitons and to induce long-lived Bloch oscillations [26], Rabi oscillations [27], and dynamical localization [28] in the nonlinear regime. For comprehensive reviews on single-component BECs in linear and/or nonlinear optical lattices, see [1,23,29,30].

On the other hand, the analysis of the physical properties of binary mixtures of condensates still presents open issues and is an interesting research topic [31–37]. In past years some work has been done on the stability and dynamics of binary BEC mixtures with both components loaded in LOLs [38], NOLs [39], or combinations thereof [40–43].

However, BEC mixtures with one component loaded in a LOL and the other loaded in a NOL have not been investigated, to the best of our knowledge. This setting is particularly interesting because it may support new types of matter waves, due to the interplay between the different types of OL and the intrinsic nonlinearities. In particular, in the absence of any interaction (e.g., with all scattering lengths tuned to zero), the spectrum of the component in the LOL displays a band structure while that of the other component has free-particle features. It is known that for attractive intraspecies interactions, uncoupled mixtures will feature localized states. In this situation one can expect that a rich variety of bound states can be formed once the interspecies interaction is switched on.

The aim of the present paper is to study localized matter waves of binary BEC mixtures with one component loaded in a LOL and the other in a NOL. In particular, we concentrate on localized states which have chemical potentials of both components in the lower semi-infinite part of the spectrum. We call these states bright-bright (BB) solitons, or also “fundamental” solitons, because when intraspecies scattering lengths are both negatives (the case investigated in this paper) they coincide with the ground state of the system. We show that BB solitons can be classified according to the distance between the lattice sites where centers of their component densities are located. Denoting these distances by nL , $n = 0, 1, 2, \dots$, with L the spatial period of the lattices (assumed to be the same for both LOL and NOL), the $n = 0$ and $n \neq 0$ families are referred to as overlapped and split BB solitons, respectively. The existence and stability of these solitons are investigated both by a variational approach (VA) for the mean-field two-component Gross-Pitaevskii equation (GPE) and by direct numerical

integrations of the system. In particular, the dependence of the existence ranges of BB soliton pairs on the interspecies interaction parameter γ_{12} is investigated. As an interesting result, we find that one can pass from one soliton family to another by simply changing the strength of the interspecies interaction. In particular, by starting from an overlapped ($n = 0$) BB soliton, one finds a series of repulsive critical values of γ_{12} at which the transition from the n - to the $n + 1$ -split BB soliton occurs as γ_{12} is adiabatically increased away from the uncoupling limit ($\gamma_{12} = 0$). On the contrary, for attractive interspecies interaction only direct transitions from split to overlapped BB solitons are possible. Since critical values at which transitions occur depend on physical parameters of the mixture, these phenomena suggest that split BB solitons could be used for indirect measurements of scattering lengths in experiments.

The paper is organized as follows. In Sec. II, we introduce the mean-field equations for the coupled system and envisage a variational study for stationary localized states. We examine the linear stability of these states for attractive and repulsive intercomponent interaction. In Sec. III we introduce a time-dependent variational approach, with Gaussian trial solutions, to study different classes of BB soliton pairs. The stability of split and overlapped families of soliton pairs is checked by numerical integration of the mean-field equations. In Sec. IV, a numerical routine is employed to understand the role of interspecies interaction in the splitting mechanism, for both attraction and repulsion between different species. Finally, in Sec. V we make concluding remarks.

II. ANALYTICAL FORMULATION

Throughout this paper, we shall consider a quasi-one-dimensional binary mixture of BECs, in which the transverse motion is frozen into the ground state of a tight transverse trapping potential, with trapping frequency ω_{\perp} . The mean-field dynamics of a mixture in which the two species' particles have equal mass m is modeled by the coupled GPEs [44],

$$i\hbar \frac{\partial \Psi_j}{\partial \tau} = \left(-\frac{\hbar^2}{2m} \frac{\partial^2}{\partial s^2} + \mathcal{V}_j(s) + 2\sigma_j(s)\hbar\omega_{\perp}|\Psi_j|^2 + 2\sigma_{12}\hbar\omega_{\perp}|\Psi_{3-j}|^2 \right) \Psi_j, \quad (1)$$

where $j = 1, 2$ is the species index, \mathcal{V}_j 's are the external trapping potentials, σ_j 's the intraspecies scattering lengths (assumed to depend on position), and σ_{12} the interspecies scattering length. The wave functions are normalized to the numbers of particles,

$$\mathcal{N}_j = \int ds |\Psi_j|^2. \quad (2)$$

Since our system is subject to an external potential proportional to $\cos(2k_L s)$ generated by two counterpropagating laser beams, the inverse wave number k_L^{-1} and the recoil energy $E_r = (\hbar k_L)^2/2m$ provide natural units for length, energy, and time [29]. To simplify the notation, we introduce the

adiimensional quantities

$$\begin{aligned} x &:= k_L s, & t &:= \frac{2E_r}{\hbar} \tau, \\ V_j &:= \frac{\mathcal{V}_j}{2E_r}, & \psi_j &:= \sqrt{\frac{\hbar\omega_{\perp}}{2E_r k_L}} \Psi_j, \\ N_j &:= \frac{\hbar\omega_{\perp}}{2E_r} \mathcal{N}_j, & \tilde{\gamma}_j &:= 2\sigma_j k_L, & \gamma_{12} &:= 2\sigma_{12} k_L, \end{aligned} \quad (3)$$

yielding the GPEs,

$$i \frac{\partial \psi_j}{\partial t} = \left(-\frac{1}{2} \frac{\partial^2}{\partial x^2} + V_j(x) + \tilde{\gamma}_j(x)|\psi_j|^2 + \gamma_{12}|\psi_{3-j}|^2 \right) \psi_j, \quad (4)$$

and the constraint,

$$\int dx |\psi_j(x)|^2 = N_j. \quad (5)$$

(Note how the variables N_j , which will be used throughout this article, coincide with the actual numbers of particles \mathcal{N}_j only up to a factor.) In the physical case of interest, only the first species is subject to an external lattice potential:

$$V_1(x) = V_{01} \cos(2x), \quad V_2(x) = 0. \quad (6)$$

The loose longitudinal harmonic trapping will be neglected, since we will focus on states that are localized over a few lattice sites. As for the interspecies coupling constants, we shall assume that the interaction of the second-species particles depends on the lattice modulation:

$$\tilde{\gamma}_1(x) = \gamma_1, \quad \tilde{\gamma}_2(x) = \gamma_2 + V_{02} \cos(2x). \quad (7)$$

To allow the existence of bright solitons, which are stationary states with both species localized, the average coupling constants γ_j are assumed to be negative (so that $|\gamma_2| > |V_{02}|$). Matter-wave bright solitons have also been observed experimentally in trapped systems [45]. Since the presence of the linear and nonlinear lattice potentials in Eq. (4) is an obstruction to finding exact bright soliton solutions, our study will be based on a reasonable variational approach, with a subsequent numerical test.

Stationary solutions: Overlapped solitons

We are interested in the stationary solutions of the coupled GPEs (4). The form $\psi_j(x, t) = \phi_j(x) e^{-i\mu_j t}$ yields the stationary GPEs,

$$\left(-\frac{1}{2} \frac{\partial^2}{\partial x^2} + V_j(x) - \mu_j + \tilde{\gamma}_j(x)|\phi_j|^2 + \gamma_{12}|\phi_{3-j}|^2 \right) \phi_j = 0, \quad (8)$$

where the external potentials and coupling constants are given, respectively, by Eqs. (6) and (7). Each stationary state is characterized by the chemical potentials (μ_1, μ_2) , which are obtained by solving the nonlinear eigenvalue problem in Eq. (8) with fixed normalization conditions (2).

Let us assume that the intraspecies interactions are attractive ($\gamma_j < 0$). Moreover, we shall focus on the case $V_{01}, V_{02} < 0$: in this situation, due to the (linear and nonlinear) trapping mechanisms, density profiles peaked around the points where $\cos(2x) = 1$ are energetically favorable for both species. In

order to investigate the features of BB soliton pairs, we choose a Gaussian trial solution

$$\phi_j(x) = A_j \exp[-x^2/2a_j^2]. \quad (9)$$

Since the amplitudes A_j and the widths a_j are bound by the normalization conditions

$$N_j = \int dx |\phi_j(x)|^2 = \sqrt{\pi} a_j A_j^2, \quad (10)$$

the functions (9) have only one free parameter. Moreover, this class of trial solutions fits *overlapped* BB solitons, with the peak of their densities sitting at the same position, say, at $x = 0$. Since, due to attractive interspecies interactions, the superposition of densities lowers the energy of the system, we expect the most energetically favorable soliton pair to be overlapped.

At fixed numbers of particles, the Gross-Pitaevskii energy functional for ϕ_j in the class (9) can be viewed as a function of the soliton width:

$$\begin{aligned} E = \int dx \left[\frac{1}{2} \sum_{j=1,2} \left(\left| \frac{\partial \phi_j}{\partial x} \right|^2 + \tilde{\gamma}_j(x) |\phi_j|^4 \right) \right. \\ \left. + \gamma_{12} |\phi_1|^2 |\phi_2|^2 + V_1(x) |\phi_1|^2 \right] \\ = \frac{1}{\sqrt{8\pi}} \left(\frac{\sqrt{\pi} N_1}{\sqrt{2} a_1^2} + \frac{\sqrt{\pi} N_2}{\sqrt{2} a_2^2} + \sqrt{8\pi} V_{01} N_1 e^{-a_1^2} \right. \\ \left. + \frac{V_{02} N_2^2}{a_2} e^{-\frac{a_2^2}{2}} + \frac{\gamma_1 N_1^2}{a_1} + \frac{\gamma_2 N_2^2}{a_2} + \frac{\gamma_{12} N_1 N_2}{\sqrt{a_1^2 + a_2^2}} \right). \quad (11) \end{aligned}$$

The optimal width values $[\bar{a}_1(N_1, N_2), \bar{a}_2(N_1, N_2)]$ are determined by

$$\left. \frac{\partial E}{\partial a_j} \right|_{a_k = \bar{a}_k} = 0 \quad \text{for } j = 1, 2, \quad (12)$$

and fix the trial ground state. The corresponding energy will be denoted by

$$E_{\min}(N_1, N_2) = E|_{a_j = \bar{a}_j(N_1, N_2)}. \quad (13)$$

The chemical potentials $\mu_j = \partial E_{\min} / \partial N_j$ can be used to test the linear stability of the ground-state solution through the Vakhitov-Kolokolov criterion [46].

Relevant properties of the overlapped solitons can be inferred from the energy functional in Eq. (11) and the chemical potential. If one keeps constant the total number of atoms $N = N_1 + N_2$, the change in the energy $E_{\min}(N_1, N_2)$ of the trial ground state with N_1 (or equivalently N_2) can be analyzed. Let us fix for definiteness $\gamma_1 = \gamma_2 = -1$, $\gamma_{12} = -0.5$, $V_{01} = -0.5$, $V_{02} = -0.25$. Throughout the paper, numbers N_j of order 1 will be extensively used: recall that they are related to the actual numbers of atoms \mathcal{N}_j by the factor $2E_r/\hbar\omega_{\perp}$ through Eq. (3). In an experiment with ^7Li atoms ($m \simeq 1.2 \times 10^{-27}$ kg), with a transverse trapping potential $\omega_{\perp} \simeq 2\pi \times 700$ Hz [45,47] and a laser wave number $k_L = 5 \times 10^{-7} \text{ m}^{-1}$, the conversion factor is of the order of 10^4 .

In Fig. 1 we set the (scaled) number of particles to $N = 3$ and study the behavior of the minimal energy and

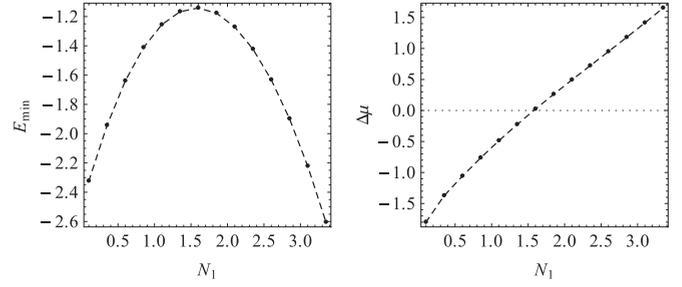


FIG. 1. (Left panel) Dependence of the minimum energy of the system, within ansatz (9), vs the number of atoms N_1 , with $N_1 + N_2 = 3$. (Right panel) Difference between chemical potentials $\Delta\mu = \mu_1 - \mu_2$ vs N_1 . The parameters are fixed to $\gamma_1 = \gamma_2 = -1$, $V_{01} = -0.5$, and $V_{02} = -0.25$. Energies and chemical potentials are in units of $2E_r$.

the difference in chemical potentials of the soliton pair. The evident asymmetry in the plot of the energy $E(N_1, N - N_1)$ (left panel) with respect to $N_1 = N/2$ is related to the inhomogeneity of the lattice potentials and the self-interactions for the two components. Moreover, the difference in chemical potentials $\Delta\mu = \mu_1 - \mu_2$ (right panel) can lead to changes in the numbers of particles if the system is in contact with a particle reservoir or if a transition mechanism between the two species is present.

Existence curves in the $\mu - N$ plane are plotted in Fig. 2 for the case of equal attractive intraspecies interactions and for both attractive (left panel) and repulsive (right panel) interspecies interaction. To reduce the number of parameters, we have considered the case $N_1 = N_2 \equiv N$. The dotted lines refer to the VA results obtained from the numerical minimization of the energy (11), while the solid lines represent the corresponding curves obtained from numerical relaxation method and self-consistent diagonalization of the stationary GPEs in Eq. (8) [15,39]. From these plots one can see that $d\mu_j/dN < 0$, both for the VA and for the GPE curves, which implies, according to the Vakhitov-Kolokolov criterion, that the BB solitons are linearly stable. This result is also confirmed by direct numerical time integrations of the two-component GPEs (4) (not shown here).

From Fig. 2 one can see that, while in the attractive case the agreement is quite good for a wide range of N (this is true also for relatively large values of γ_{12}), in the repulsive case deviations of the VA and numerical curves are larger. This discrepancy is due to the fact that the Gaussian ansatz becomes less accurate in the repulsive case, in which the interspecies interaction reduces the stability of the overlapped configuration, as one can see from Fig. 3 where VA and GPE soliton profiles are compared. However, the accuracy of the VA result increases as positive γ_{12} values are decreased towards the uncoupling limit $\gamma_{12} = 0$. Also note the existence of points where the chemical potential curves intersect, both for attractive and repulsive interspecies interactions. At these points, BB solitons, having the same number of atoms and the same chemical potentials (related to their width), will have equal VA profiles for the two components. Despite the discrepancy in the location of the intersection point, the equality of the profiles is well confirmed by numerical GPE results.

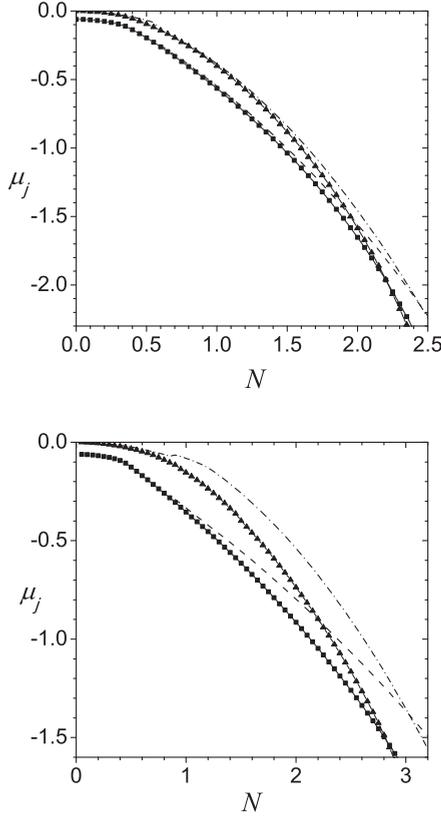


FIG. 2. Existence curves in the $\mu_j - N$ plane ($N \equiv N_1 = N_2$) for BB solitons of the GPEs with attractive ($\gamma_{12} = -0.5$, left) and repulsive ($\gamma_{12} = 0.1$, right) interspecies interaction. Other parameters are fixed as $\gamma_1 = \gamma_2 = -1, V_{01} = -0.5, V_{02} = -0.25$. Squares and triangles denote numerical results obtained by self-consistent diagonalization of Eq. (8) and refer to the first and the second component, respectively. Dashed and dot-dashed lines represent the corresponding results obtained from the minimization of the energy functional. Chemical potentials in units of $2E_r$.

III. SPLIT BB SOLITONS AND DYNAMICAL PROPERTIES

In the previous section, the choice of Gaussian trial wave functions (9) is aimed at studying overlapped BB soliton configurations, with the peaks of the two density profiles coinciding at $x = 0$. Due to attractive interspecies interaction, this configuration is expected to be the lowest-energy BB soliton pair. It is possible, however, to extend the analysis to *split BB solitons*, in which the centers of mass of the two species do not coincide.

Let us initially consider the uncoupled limit $\gamma_{12} = 0$. In the case $V_{01} < 0$ and $V_{02} < 0$, one expects an infinite set of degenerate energy-minimizing BB solitons, since the centers of mass of each species can be located at any point x_{0j} such that $\cos(2x_{0j}) = 0$, regardless of the other species' density profile. These minimizing configurations can be classified in families

$$BB_n(\gamma_{12} = 0), \quad \text{with } n \in \mathbb{N}, \quad (14)$$

according to the absolute distance between the centers of mass:

$$\Delta x := |x_{02} - x_{01}| \in \left[\left(n - \frac{1}{2}\right)\pi, \left(n + \frac{1}{2}\right)\pi \right]. \quad (15)$$

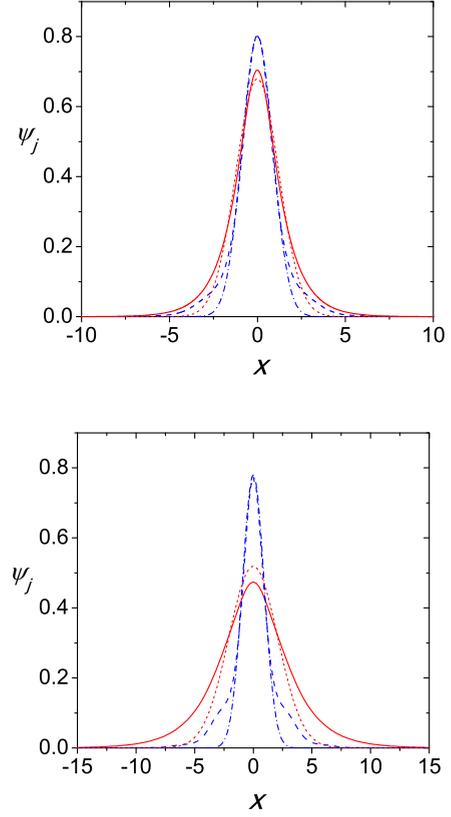


FIG. 3. (Color online) Profiles of BB solitons with $N_1 = N_2 = N$ corresponding to the $N = 1$ points on the curves of Fig. 2 for attractive ($\gamma_{12} = -0.5$, left panel) and repulsive ($\gamma_{12} = 0.1$, right panel) interspecies interaction. Dashed and dot-dashed (blue) lines and continuous and dotted (red) lines refer to first and second component, respectively. In both panels continuous and dashed lines refer to numerical solutions of Eq. (4), while dotted and dot-dashed lines to variational analysis. Position in units of k_L^{-1} .

Clearly, the energy of the BB soliton configurations is the same for all families, $E_n(\gamma_{12} = 0) = E_0$. When the effect of the interspecies coupling γ_{12} can be treated as a small perturbation, one expects the existence of stationary Gross-Pitaevskii solutions close to the ones at $\gamma_{12} = 0$, which can still be classified in families $BB_n(\gamma_{12})$ according to the criterion (15). However, if $\gamma_{12} < 0$, the interspecies attraction will break the energetic degeneracy in favor of the overlapped configuration BB_0 . Thus, the split solitons in $BB_{n>0}(\gamma_{12})$ become metastable. While configurations with a very large distance between the two species are almost unaffected by interspecies interactions, larger values of $|\gamma_{12}|$ weaken the (local) stability of solitons with small n , since attractive interactions can give a sufficient amount of energy to overcome the (linear or nonlinear) potential barrier and reduce the distance between the centers of mass. Thus one expects a critical value $\gamma_{cr}^{(n)}$, such that for $\gamma_{12} < \gamma_{cr}^{(n)}$ metastable solutions in BB_n would no longer exist.

In the following, we will numerically analyze the existence of split BB soliton pairs, as well as their energetic and dynamical behavior. To this end, we shall generalize our ansatz to include the positions of the component density centers as free parameters. We will also consider time-dependent parameters to investigate the dynamics of the system. The

Gross-Pitaevskii equations (4) can be restated as a variational problem [48]

$$\delta \int \mathcal{L} \left(\psi_j, \psi_j^*, \frac{\partial \psi_j}{\partial x}, \frac{\partial \psi_j^*}{\partial x}, \frac{\partial \psi_j}{\partial t}, \frac{\partial \psi_j^*}{\partial t} \right) dx dt = 0, \quad (16)$$

where the Lagrangian density reads

$$\begin{aligned} \mathcal{L} = & \frac{1}{2} \sum_{j=1}^2 \left[i \left(\psi_j^* \frac{\partial \psi_j}{\partial t} - \psi_j \frac{\partial \psi_j^*}{\partial t} \right) - \left| \frac{\partial \psi_j}{\partial x} \right|^2 - \gamma_j |\psi_j|^4 \right] \\ & - V_{01} \cos(2x) |\psi_1|^2 - \frac{V_{02}}{2} \cos(2x) |\psi_2|^4 \\ & - \gamma_{12} |\psi_1|^2 |\psi_2|^2. \end{aligned} \quad (17)$$

We generalize the Gaussian ansatz (9) to

$$\begin{aligned} \psi_j(x, t) = & A_j \exp \left[i \frac{dx_{0j}}{dt} (x - x_{0j}) + i \phi_j (x - x_{0j})^2 + i \theta_j \right] \\ & \times \exp \left[- \frac{(x - x_{0j})^2}{2a_j^2} \right], \end{aligned} \quad (18)$$

where the variational parameters $(A_j, a_j, x_{0j}, \phi_j, \theta_j)$ for $j = 1, 2$ are generally time-dependent and represent, respectively, the amplitude, width, center-of-mass position, frequency chirp, and overall phase of a soliton in the j th component. The Lagrangian L for the trial wave functions (18), obtained by integrating the Lagrangian density in (17), reads

$$\begin{aligned} L = & \int_{-\infty}^{+\infty} \mathcal{L} dx \\ = & \frac{\sqrt{\pi}}{4} \left(\sum_{j=1}^2 \left[\frac{A_j^2}{a_j} + \sqrt{2} \gamma_j a_j A_j^4 + 4a_j^3 A_j^2 \phi_j^2 \right. \right. \\ & \left. \left. - 2a_j A_j^2 \left(\frac{dx_{0j}}{dt} \right)^2 + 4a_j A_j^2 \frac{d\theta_j}{dt} + 2a_j^3 A_j^2 \frac{d\phi_j}{dt} \right] \right. \\ & \left. + 4V_{01} e^{-a_1^2} a_1 A_1^2 \cos(2x_{01}) + \sqrt{2} V_{02} e^{-\frac{1}{2}a_2^2} a_2 A_2^4 \cos(2x_{02}) \right. \\ & \left. + \frac{4\gamma_{12} e^{-\frac{(x_{01}-x_{02})^2}{a_1^2+a_2^2}} a_1 A_1^2 a_2 A_2^2}{\sqrt{a_1^2+a_2^2}} \right). \end{aligned} \quad (19)$$

Note the dependence of the interaction term on Δx , defined in Eq. (15).

The functional derivatives of L with respect to the variational parameters yield a set of Euler-Lagrange equations. After appropriate manipulation, we can obtain a picture for the dynamics of the soliton pairs. In particular, the equations

$$\frac{d}{dt} (\sqrt{\pi} a_j A_j^2) = 0, \quad (20)$$

$$-2a_j \phi_j + \frac{da_j}{dt} = 0, \quad (21)$$

can be interpreted as dynamical constraints on the amplitude (particle number conservation) and the frequency chirp. Taking into account Eqs. (20) and (21), the equations of motion for the centers of mass and the width can all be derived from the

TABLE I. Locally stable BB soliton pairs for $N_1 = 1$, $N_2 = 1$, $\gamma_1 = -1$, $\gamma_2 = -1$, $V_{01} = -0.5$, $V_{02} = -0.25$, belonging to the classes BB_n with $n = 0, 1, 2$. The optimal parameters Δx , a_1 , and a_2 (units of k_L^{-1}) are obtained by minimizing the global potential Π (units of $2E_r$).

γ_{12}	n	Δx	a_1	a_2	Π_{\min}
0	0,1,2	$0, \pi, 2\pi$	0.929	2.229	-0.177
-0.25	0	0	0.910	1.472	-0.249
-0.25	1		No local minimum		
-0.25	2	2.991π	0.930	2.260	-0.177

effective potential

$$\begin{aligned} \Pi(a_1, a_2, x_{01}, x_{02}) = & \Pi_1(a_1, x_{01}) + \Pi_2(a_2, x_{02}) \\ & + \Pi_{12}(a_1, x_{01}, a_2, x_{02}), \end{aligned} \quad (22)$$

with

$$\begin{aligned} \Pi_1(a_1, x_{01}) = & \frac{N_1}{4a_1^2} + \frac{\gamma_1 N_1^2}{2\sqrt{2\pi} a_1} \\ & + e^{-a_1^2} N_1 V_{01} \cos(2x_{01}), \end{aligned} \quad (23a)$$

$$\begin{aligned} \Pi_2(a_2, x_{02}) = & \frac{N_2}{4a_2^2} + \frac{\gamma_2 N_2^2}{2\sqrt{2\pi} a_2} \\ & + \frac{e^{-\frac{1}{2}a_2^2} N_2^2 V_{02} \cos(2x_{02})}{2\sqrt{2\pi} a_2}, \end{aligned} \quad (23b)$$

$$\Pi_{12}(a_1, x_{01}, a_2, x_{02}) = \frac{e^{-\frac{(x_{01}-x_{02})^2}{a_1^2+a_2^2}} \gamma_{12} N_1 N_2}{\sqrt{\pi} \sqrt{a_1^2+a_2^2}}, \quad (23c)$$

through the Newton-like equations

$$\frac{d^2}{dt^2} \begin{pmatrix} N_1 x_{01} \\ N_2 x_{02} \\ (N_1/2) a_1 \\ (N_2/2) a_2 \end{pmatrix} = - \begin{pmatrix} \partial_{x_{01}} \\ \partial_{x_{02}} \\ \partial_{a_1} \\ \partial_{a_2} \end{pmatrix} \Pi(a_1, a_2, x_{01}, x_{02}). \quad (24)$$

The problem of finding stationary solutions within the ansatz (18) thus reduces to searching the minima of the effective potential. Equation (24) can also be linearized around the stable equilibrium points for Π to obtain information on small center-of-mass and width oscillations. On the other hand, due to the (still restrictive) form of the trial wave functions, the far-from-equilibrium dynamics of Eq. (24) cannot be considered physically relevant. (See the discussion below.) In Table I, we show some illustrative results of the optimal parameters Δx , a_1 , and a_2 for solitons in the classes BB_n with $n = 0, 1, 2$. The numbers of particles are fixed to $N_1 = N_2 = 1$. At $\gamma_{12} = 0$, the three minima are degenerate. At $\gamma_{12} = -0.25$, the split solitons in the family BB_2 are locally stable [see Eq. (14)], while the global minimum is, as expected, the overlapped configuration. Instead, no local minimum can be found in the family BB_1 , indicating that $\gamma_{\text{cr}}^{(1)} > -0.25$.

A more systematic picture of the existence and behavior of solitons for small n with varying γ_{12} is given in Fig. 4: a numerical minimization procedure is used to find stable configurations with their centers of mass close to some initial points $(\tilde{x}_{01}, \tilde{x}_{02})$. A minimum with centers of mass around

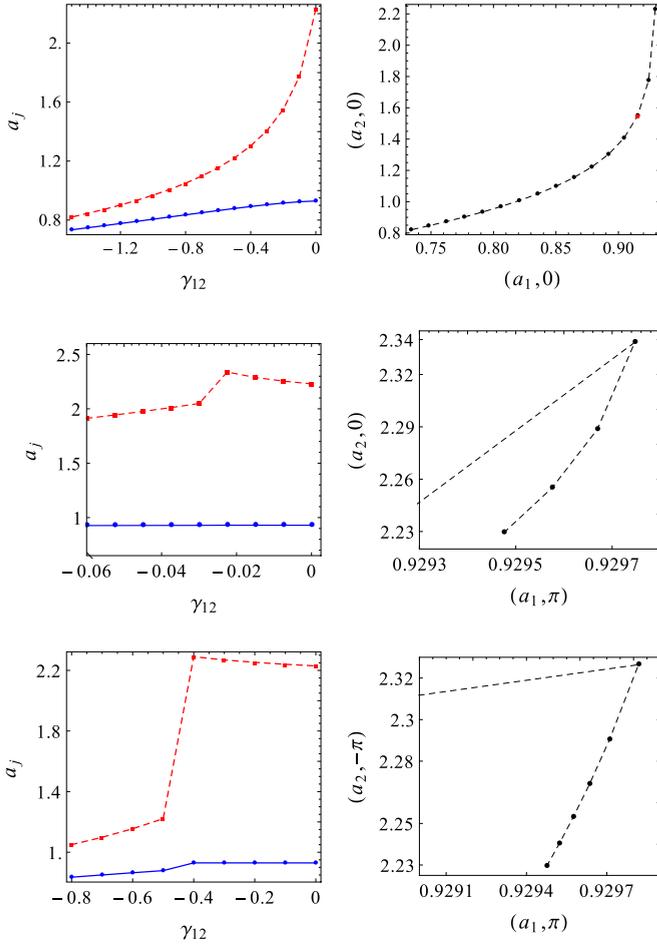


FIG. 4. (Color online) Numerical minimization of the pseudopotential Π [see (22)] with soliton peaks around $(x_{01}, x_{02}) = (0, 0)$ (top panels), $(x_{01}, x_{02}) = (\pi, 0)$ (central panels), and $(x_{01}, x_{02}) = (\pi, -\pi)$ (bottom panels). Plots in the left column represent the variation of the optimal soliton widths a_1 (solid blue lines) and a_2 (dashed red lines), in units of k_L^{-1} , with γ_{12} . Plots in the right column show the behavior of one optimal width with respect to the other. The jumps observed in the central and bottom lines (right column) are due to the disappearance of local minima with $\Delta x \simeq \pi$ and $\Delta x \simeq 2\pi$ as $|\gamma_{12}|$ decreases under a critical value. Minima found below the critical $|\gamma_{12}|$ coincide with the values in the top plots and are thus not shown in the right column.

$(\tilde{x}_{01}, \tilde{x}_{02}) = (0, 0)$ can be found for all negative γ_{12} (top panels). On the other hand, the search of minima whose centers of mass are close to $(\tilde{x}_{01}, \tilde{x}_{02}) = (\pi, 0)$ and $(\tilde{x}_{01}, \tilde{x}_{02}) = (-\pi, \pi)$ yields discontinuous behaviors in the optimal parameters (central and bottom panels). The discontinuities are present because solitons in BB_1 (respectively BB_2) exist only for $\gamma_{12} > \gamma_{cr}^{(1)}$ (respectively $\gamma_{12} > \gamma_{cr}^{(2)}$), while for more negative values the algorithm actually finds the global minimum belonging to BB_0 . It is also worth noticing that in Fig. 4 the optimal widths a_1 and a_2 of the overlapped solitons decrease with $|\gamma_{12}|$. In the case of split solitons, the amplitudes remain almost constant, with a slight increase (more evident in a_2) with $|\gamma_{12}|$, due to the attraction exerted between densities.

These findings are corroborated by the behavior of the center-of-mass positions displayed in Fig. 5, where the

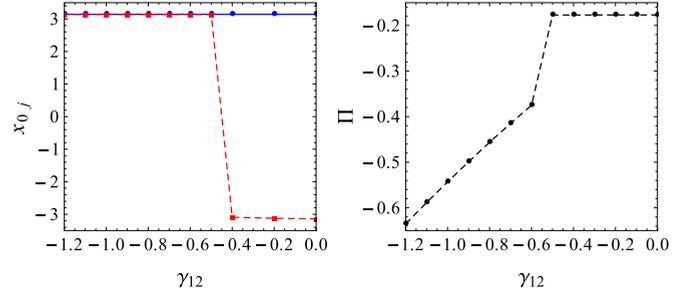


FIG. 5. (Color online) (Left panel) Peak positions x_{01} (solid blue line) and (x_{02}) (dotted red line), in units of k_L^{-1} , of optimal BB profiles, obtained by searching the local minimum around $(x_{01}, x_{02}) = (\pi, -\pi)$, for different values of γ_{12} . The sudden jump towards (π, π) is due to the disappearance of the local minimum (see also Fig. 4, bottom panels). (Right panel) Local minimum value of the effective potential energy in Eq. (22), in units of $2E_r$.

displacement of the center of mass of the second species with increasing interspecies interaction is observed. The situation is the same as that depicted in Fig. 4, bottom panels. In the left panel of Fig. 5, the jump of x_{02} as $|\gamma_{12}|$ is decreased signals the disappearance of the local energy minimum. The value of the local minimum of the effective potential energy (22) is shown in the right panel.

In order to check the existence and stability of BB soliton pairs as approximate solutions of the GPEs, we employ a numerical simulation of the dynamics generated by Eq. (4). First, we have checked the stationarity of overlapped soliton pairs, localized around $x = 0$. It is possible to verify, for different values of γ_{12} , that the soliton pair determined by the minimization procedure is stationary within very good approximation. In Fig. 6, the time evolution of the overlapped

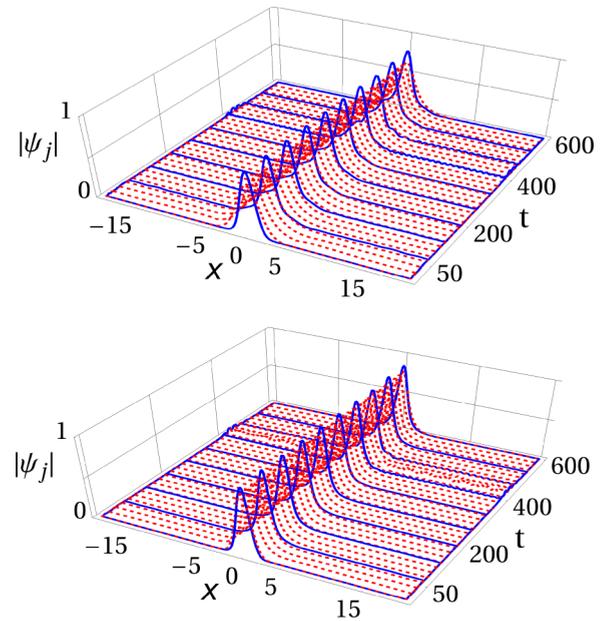


FIG. 6. (Color online) Time evolution of density profiles for overlapped solitons in the first (solid blue lines) and second (dashed red lines) species for $\gamma_{12} = -0.25$ (left panel) and $\gamma_{12} = -1$ (right panel). The results are obtained by direct numerical integration. Position in units of k_L^{-1} . Time in units of $\hbar/(2E_r)$.

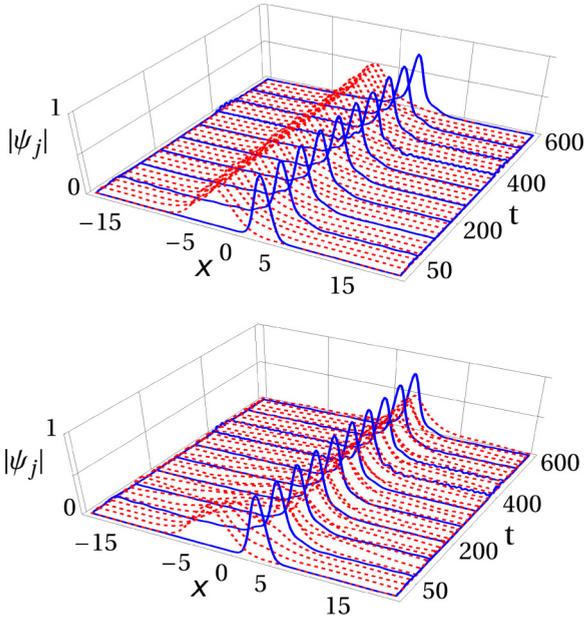


FIG. 7. (Color online) Time evolution of density profiles for split solitons in the first (solid blue lines) and second (dashed red lines) species for $\gamma_{12} = -0.25$ (left panel) and $\gamma_{12} = -0.38$ (right panel). In the latter case, numerical integration shows the instability of the soliton pair. Position in units of k_L^{-1} . Time in units of $\hbar/(2E_r)$

solitons is represented for $\gamma_{12} = -0.25$ (left) and $\gamma_{12} = -1$ (right). Then, we have tested the behavior of split soliton pairs in BB_2 in different regimes. In the case $\gamma_{12} = -0.25$, which is larger than the critical value $\gamma_{cr}^{(2)} \simeq -0.4$, the split configuration evolves in time with slight distortions, but it preserves the qualitative features of the initial state for all the time of the simulation (left panel of Fig. 7). When $\gamma_{12} \simeq \gamma_{cr}^{(2)}$, the energetic instability of the soliton pair in BB_2 is reflected by a *dynamical* instability: the second-species density distribution is gradually attracted by the first species (right panel of Fig. 7), ending with an overlapped configuration, which is eventually stabilized by radiating wave packets [49,50].

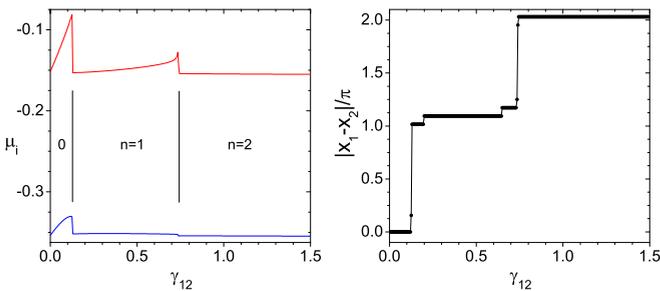


FIG. 8. (Color online) Dependence of chemical potentials (left panel) and distance between peaks (right panel) on repulsive interspecies interactions of a BB soliton with $N_1 = N_2 = 1$. Other parameters are fixed as $\gamma_1 = \gamma_2 = -1, V_{01} = -0.5, V_{02} = -0.25$. Top (blue) and bottom (red) curves in the left panel refer to chemical potentials of first and second component, respectively, while vertical lines separate γ_{12} regions for existence of BB solitons in BB_0 (overlapped), BB_1 , and BB_2 (split). Chemical potential in units of $2E_r$, distance in units of k_L^{-1} .

IV. SPLITTING DEPENDENCE ON INTERSPECIES INTERACTION

In the previous section we observed that split BB solitons can become unstable at some negative critical values of the interspecies scattering length. We shall now investigate these critical values in more detail by direct numerical integration of the GPEs, both for attractive and repulsive interatomic interactions.

Let us first discuss the repulsive case. We can consider as the initial state an overlapped BB soliton, centered at $x = 0$, with no interspecies coupling ($\gamma_{12} = 0$), and adiabatically switch on a repulsive interspecies interaction between components at $t > 0$. One expects that, due to the repulsive interspecies interaction, the initial BB_0 soliton will evolve into a split one belonging to the BB_1 family at some value $\gamma_{12} = \gamma_{rep}^{(1)}$ and then into the BB_2 family at $\gamma_{12} = \gamma_{rep}^{(2)}$. This picture coincides with the numerical results in Fig. 8, both in terms of the chemical

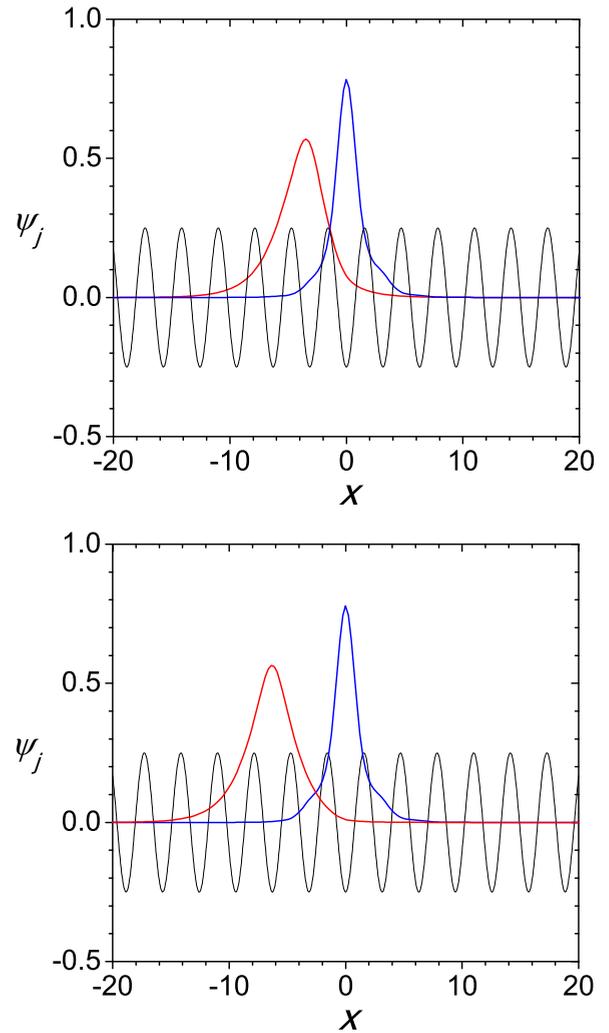


FIG. 9. (Color online) Split BB solitons inside the regions $n = 1$ and $n = 2$ of the left panel of Fig. 8, at $\gamma_{12} = 0.5$ (top panel) and $\gamma_{12} = 1.0$ (bottom panel). The (blue) profiles centered at $x = 0$ and the (red) ones centered at $x < 0$ refer to first and second component, respectively, while the sinusoidal (black) lines shows the periodicity of the optical lattice. Position in units of k_L^{-1} .

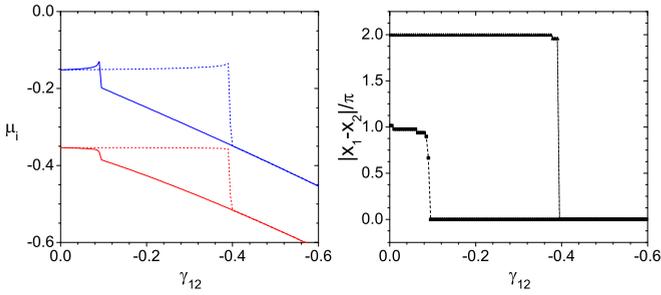


FIG. 10. (Color online) Dependence on attractive interspecies interaction of chemical potentials (left panel) and distance between density centers, x_1, x_2 , (right panel) of a BB soliton with $N_1 = N_2 = 1$. Bottom (blue) and top (red) curves in the left panel refer to the first and the second component, respectively, while dotted and continuous lines (left panel) and square and triangle symbols (right panel) refer to BB solitons with centers initially separated by π and by 2π , respectively. Other parameters are fixed as $\gamma_1 = \gamma_2 = -1$, $V_{01} = -0.5$, $V_{02} = -0.25$. Chemical potential in units of $2E_r$, distance in units of k_L^{-1} .

potentials and the distances between peaks Δx , normalized to π . The jumps in the distance are correlated with jumps in the chemical potentials at the critical values, which are uniquely fixed by the parameters of the system. In Fig. 9, the profiles of the split solitons with $n = 1$ and $n = 2$ are represented at two different γ_{12} values belonging to their existence curve. Despite the smaller value of γ_{12} , the two BB components in the $n = 1$ case appear to be more distorted in their overlapping region than in the $n = 2$ case. This is an evident consequence of the exponential decay of the soliton-soliton interaction with distance [see Eq. (19)]. Notice that, as in the attractive case, the normalized distance between soliton centers is not an integer number. This is a clear consequence of the existence of a repulsive force between components.

When attractive interactions $\gamma_{12} < 0$ are adiabatically turned on at $t > 0$, we expect that an initial split soliton in BB_{n_0} , with $n_0 > 0$, will undergo *only one jump* towards $n = 0$. Indeed, from the analysis in the previous section, we can deduce that the negative critical values are ordered as $\gamma_{cr}^{(n)} > \gamma_{cr}^{(n+1)}$. Thus, if $\gamma_{12} > \gamma_{cr}^{(n_0)}$, interactions give enough energy to overcome all the intermediate barriers from the n_0 -th down to $n = 0$. This intuitive result, based on energetic considerations, matches very well the results of the numerical simulation, as one can see from Fig. 10 where the cases $n_0 = 2$ and $n_0 = 1$ are represented.

Since the critical values of γ_{12} at which the transitions occur are uniquely fixed by the parameters of the mixture, including the number of atoms and intraspecies interactions, an experimental implementation of the above numerical simulations could be used for indirect measurements of the interspecies scattering length of BEC mixtures. The interspecies scattering length can also be measured from the oscillatory motion of coupled solitons as predicted in [51].

V. CONCLUSIONS

We have considered matter-wave bright-bright solitons in coupled Bose-Einstein condensates by assuming that the first component is loaded in a linear optical lattice and the second component in a nonlinear optical lattice. In particular, the existence and stability of split and overlapped BB solitons has been investigated by VA, by direct numerical integrations of the coupled GPEs, and by direct numerical integrations of the system. The dependence of the existence ranges of BB solitons on the interspecies interaction parameter has been also investigated. In particular, for repulsive interspecies interactions, we showed the existence of a series of critical values of γ_{12} at which transitions from the n - to the $n + 1$ -split BB soliton occur. For attractive interspecies interaction we showed that only direct transitions from a split BB soliton to the overlapped BB soliton are possible. Since critical values at which transitions occur depend on the physical parameters of the mixture, these phenomena suggest that split BB solitons could be used for indirect measurements of these parameters in experiments.

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