Bose-Einstein condensates with attractive interactions on a ring

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Considering an effectively attractive quasi-one-dimensional Bose-Einstein condensate of atoms confined in a toroidal trap, we find that the system undergoes a phase transition from a uniform to a localized state, as the magnitude of the coupling constant increases. Both the mean-field approximation, as well as a diagonalization scheme are used to attack the problem.

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Effectively attractive Bose-Einstein condensates have concentrated a lot of attention recently, in connection with the experimental formation of bright solitons in them [1,2]. More specifically, in the experiments of Refs. [1,2], 7Li atoms were confined in quasi-one-dimensional traps. With use of the Feschbach resonances [3] the effective coupling constant that describes the atom interactions was then tuned and as it became negative – corresponding to an effective attraction between the atoms – localized states, “bright soliton trains” were observed to form in Ref. [1], while a single bright soliton was observed in Ref. [2]. References [4,5] have examined theoretically these systems. In the limit where transversely to the long axis of the trap the gas is in the lowest harmonic-oscillator level, the transverse degrees of freedom are frozen out and the system is essentially one-dimensional [6].

Motivated by these developments, we study here an effectively attractive one-dimensional Bose-Einstein condensate of atoms confined in a toroidal trap. In a recent theoretical paper Kanamoto, Saito, and Ueda have investigated the ground state and the low-lying excited states of such a system [5] (see also Ref. [7] for a detailed discussion of this problem). As shown, the Gross-Pitaevskii mean-field theory predicts a quantum phase transition between a uniform state and a localized state as the absolute value of the strength of the interaction increases. Furthermore, numerical diagonalization of the Hamiltonian for a finite number of bosons shows that the transition in this case is smeared out, as expected in finite systems.

In our study we examine the same problem using different techniques. Initially we use the mean-field approximation with a properly chosen variational wavefunction to study the phase transition and the order parameter in the two phases. Then, working in the same truncated space we use a Bogoliubov transformation to diagonalize the Hamiltonian. Having diagonalized the problem, we examine the lowest state of the system, the low-lying excited states, as well as the depletion of the condensate at the region of the transition. Our results are consistent with those of Ref. [5].

Let us therefore consider a Bose-Einstein condensate in a toroidal trap. Following Ref. [5], we assume that the system contains N bosons, that the radius of the torus is R and its cross section is $S = \pi r^2$, with $r \ll R$. If $\hat{\psi}(\theta)$ is the field operator, the Hamiltonian is

$$\hat{H} = \int_0^{2\pi} d\theta \left[ -\hat{\psi}^\dagger(\theta) \frac{\partial^2}{\partial \theta^2} \hat{\psi}(\theta) + \frac{U_0}{2} \hat{\psi}^\dagger(\theta) \hat{\psi}(\theta) \hat{\psi}(\theta) \hat{\psi}(\theta) \right],$$

(1)

where $U_0 = 8\pi aR/S$, with $a$ being the scattering length for elastic atom-atom collisions, and $\theta$ is the azimuthal angle. Here the length is measured in units of $R$ and the energy in units of $\hbar^2/2mR^2$, with $m$ being the atom mass.

Let us start with the mean-field approach. Within this approximation the system is described by a single wavefunction, the order parameter $\psi(r)$, and the many-body state is the product $\Pi_i \psi(r_i)$, with $i = 1, \ldots, N$, where $N$ is the number of atoms. Therefore this approximation ignores correlations between the atoms and in general it has a higher energy than the exact solution that one can get by diagonalizing the Hamiltonian.

In this problem it is natural to work in the basis of plane-wave states $\phi_\ell(\theta) = e^{i\ell\theta}/\sqrt{2\pi}$ and according to the analysis of Ref. [5], the order parameter close to the transition consists of the state $\phi_0$ (the dominant component), and the states $\phi_{\pm 1}$. To get a simple physical picture, we thus develop a variational approach, expanding $\psi(\theta)$ in the basis of the $\phi_\ell$ states and keeping only these three components. This is a reasonable assumption, since states with higher values of $l$ have higher kinetic energy, and indeed as shown in Ref. [5] this approximation gives qualitatively (but not quantitatively) good results. We thus write

$$\psi(\theta) = c_{-1}\phi_{-1} + c_0\phi_0 + c_1\phi_1.$$ 

(2)

Since the dominant component of $\psi$ is $\phi_0$, therefore $|c_0| \gg |c_{-1}|$, and $|c_0| \gg |c_1|$. Because of the symmetry of the problem, $|c_{-1}| = |c_1|$, which also guarantees that the total angular momentum of $\psi$ is zero, as it should. The normalization condition imposes the further constraint $|c_{-1}|^2 + |c_0|^2 + |c_1|^2 = 1$.

To proceed, we express the energy per particle $\epsilon_0$ in terms of the coefficients $c_i$, which are then determined by minimizing $\epsilon_0$ with respect to them [8],

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\[ \epsilon_0 = 2|c_1|^2 + \frac{\gamma}{2}(|c_0|^4 + |c_{-1}|^4 + |c_1|^2) + 4|c_0|^2|c_{-1}|^2 + 4|c_1|^2|c_{-1}|^2 + 2\gamma c^*_{-1}c_1 + 2(c_0^2) c_{-1}c_1, \]

where \( \gamma = NU_0/(2\pi) = 4N\alpha R/S \) is essentially the ratio between the interaction energy and the kinetic energy. If \( c_j = |c_j|e^{i\theta_j} \), then the last two terms in Eq. (3) are equal to \( 4|c_0|^2|c_1|^2 \cos(\theta_{-1} + \theta_1 - 2\theta_0) \). To minimize \( \epsilon_0 \) one has to choose \( \theta_{-1} + \theta_1 - 2\theta_0 = 0 \). Since \(|c_0|^2 = 1 - 2|c_1|^2\),

\[ \epsilon_0 - \frac{\gamma}{2} = 2|c_1|^2(1 + 2\gamma) - 7\gamma|c_1|^4. \]

Here \( \gamma/2 \) is the interaction energy per particle of the uniform state. The above equation implies that for \( \gamma > \gamma_{cr} \), where \( \gamma_{cr} = -1/2 \), the minimum of the energy occurs for \( c_0 = 1 \), and \( c_{-1} = c_1 = 0 \). In this regime the density of the system is uniform and the energy per particle is \( \gamma/2 \). On the other hand, if \( \gamma < \gamma_{cr} \), \( \epsilon_0(|c_1|^2) \) is like a mexican hat, and its minimum occurs for all three \( c_i \neq 0 \), the energy per particle is lower than \( \gamma/2 \) and the cloud develops a non-uniform density. Our approach implies that the transition takes place for the same critical value of \( \gamma \) as in the exact solution of the mean-field approach [5,7]. It is interesting that within our scheme including more basis states in the order parameter does not affect \( \gamma_{cr} \). More precisely, including the states with \( |l| \leq m \), the quadratic terms in the energy have the form

\[ \epsilon_0 - \frac{\gamma}{2} = 2|c_1|^2(1 + 2\gamma) + 2|c_1|^4(4 + 2\gamma) + \ldots + 2|c_m|^2(m^2 + 2\gamma). \]

From the above expression it is obvious that \( \gamma_{cr} \) is always \(-1/2\), essentially because of the high kinetic energy of atoms in states with high values of \( |l| \), that scales as \( l^2 \).

Let us now examine the non-uniform state when \( \gamma \geq \gamma_{cr} \). Demanding that \( \partial\epsilon_0/\partial|c_1|^2 = 0 \), we get

\[ |c_{\pm 1}|^2 = \frac{2}{\gamma_{cr}}(\gamma - \gamma_{cr}); |c_0|^2 = 1 - \frac{4}{\gamma_{cr}}(\gamma - \gamma_{cr}). \]

Figure 1 shows the corresponding order parameter. In addition

\[ \epsilon_0 - \frac{\gamma}{2} = \frac{4}{\gamma_{cr}}(\gamma - \gamma_{cr})^2. \]

If \( \gamma = -1/2 - \delta \), where \( \delta \to 0 \) and \( \delta > 0 \), the above equation implies that \( \epsilon_0 = -1/4 - \delta/2 - 8\delta^2/7 \), which should be compared with the exact mean-field solution [5], \( \epsilon_0^* \approx -1/4 - \delta/2 - 2\delta^2 \). Therefore our trial wavefunction reproduces the linear correction, but it fails in the quadratic term and it has higher energy. This is not a surprise, however, since our approach is variational and it does not necessarily capture all these features. Including more basis states, one can improve the wavefunction and the energy to the desired order in \( \delta \).

Turning to our alternative approach of the problem, where we will go beyond the mean-field approximation and we will calculate terms in the energy of order \( 1/N \), let us expand the field operator \( \hat{\psi} \),

\[ \hat{\psi}(\theta) = \hat{\phi}_{-1} \hat{c}_{-1} + \hat{\phi}_0 \hat{c}_0 + \hat{\phi}_1 \hat{c}_1, \]

where \( \hat{c}_l \) is now the annihilation operator of an atom with angular momentum \( l \). Again we have restricted ourselves to the states with \( l = 0, \pm 1 \). In this basis the Hamiltonian (1) can be written as

\[ \hat{H} = \sum_l l^2 \hat{c}_l^\dagger \hat{c}_l + \frac{U_0}{4\pi \sum_{klnm}} \hat{c}_k^\dagger \hat{c}_l^\dagger \hat{c}_m \hat{c}_n \delta_{m+n-k-l}. \]

A similar Hamiltonian has been studied in Ref. [8] in the context of weakly-interacting Bose-Einstein condensates under rotation. Denoting the basis vectors as

\[ |m\rangle \equiv |N_{-1}, N_0, N_1\rangle \equiv |(-1)^m, 0^{N-2m}, (+1)^m\rangle, \]

where \( N_l \) is the occupancy of the state \( \phi_l \), with \( \sum_l N_l = N \), and \( \sum_l l N_l = 0 \), it is straightforward to calculate the diagonal matrix elements,

\[ \langle m|\hat{H}|m\rangle = 2m + \frac{U_0}{4\pi}[(N - 2m)(N - 2m - 1) + 2m(m - 1) + 8m(N - 2m) + 4m^2]. \]

Assuming that the system is close to the transition point, but on the side of the “uniform state” (in the infinite-\( N \) limit of the mean-field approximation), then \( m \) is of order unity and thus \( m \ll N \). In this limit,
$$\langle m|\hat{H}|m+1\rangle = \frac{U_0}{4\pi^2}\sqrt{(N-2m)(N-2m-1)(m+1)^2},$$  \hspace{1cm} (13)

and in the limit \(m \ll N\),

$$\langle m|\hat{H}|m+1\rangle = \gamma(m+1).$$  \hspace{1cm} (14)

From Eqs. (12) and (14) \(\hat{H}\) can be written as

$$\hat{H} - \frac{\gamma}{2}(N-1) = (1+\gamma)(\hat{c}_{-1}^{\dagger}\hat{c}_{-1} + \hat{c}_{1}^{\dagger}\hat{c}_{1}) + \gamma(\hat{c}_{-1}^{\dagger}\hat{c}_{1} + \hat{c}_{1}^{\dagger}\hat{c}_{-1}).$$  \hspace{1cm} (15)

The above Hamiltonian can be diagonalized with use of a Bogoliubov transformation [9],

$$\hat{b} = \lambda_{1}\hat{c}_{-1} + \lambda_{2}\hat{c}_{1},$$

$$\hat{d} = \lambda_{2}\hat{c}_{-1} + \lambda_{1}\hat{c}_{1}^{\dagger}.\hspace{1cm} (16)$$

For \(\hat{b}\) and \(\hat{d}\) to satisfy bosonic commutation relations, \(\lambda_{2}^{2} - \lambda_{1}^{2} = 1\). Following the usual tricks, \(\hat{H}\) can be written in the diagonal form

$$\hat{H} - \frac{\gamma}{2}(N-1) = E_0 + E(\hat{b}^{\dagger}\hat{b} + \hat{d}^{\dagger}\hat{d}),$$  \hspace{1cm} (17)

where

$$E_0(\gamma) = 2[(1+\gamma)\lambda_{1}^{2} - \gamma\lambda_{1}\lambda_{2}],$$  \hspace{1cm} (18)

and

$$E(\gamma) = (1+\gamma)(\lambda_{1}^{2} + \lambda_{2}^{2}) - 2\gamma\lambda_{1}\lambda_{2}.\hspace{1cm} (19)$$

In addition, to eliminate the off-diagonal terms one has

$$\frac{2\lambda_{1}\lambda_{2}}{\lambda_{1}^{2} + \lambda_{2}^{2}} = \frac{\gamma}{\gamma + 1}.\hspace{1cm} (20)$$

Parametrizing, i.e., writing \(\lambda_{1} = \sinh(\theta)\) and \(\lambda_{2} = \cosh(\theta)\), Eq. (20) can be written as

$$\tanh(2\theta) = \gamma/\gamma + 1,$$  \hspace{1cm} (21)

and in order for a solution to exist, \(\gamma > -1/2\). Equation (21) implies that \(\theta = \ln(2\gamma + 1)\), which can be used to express all the quantities in terms of \(\gamma\). Thus we get for the eigenvalues of \(\hat{H}\),

$$E_{n_b,n_d} = -\frac{\gamma}{2}(N-1) = -(\gamma + 1) + \sqrt{1 + 2\gamma(1 + n_b + n_d)},$$  \hspace{1cm} (22)

where \(n_b, n_d = 0, 1, 2, \ldots\) are the eigenvalues of the number operators \(\hat{b}^{\dagger}\hat{b}\) and \(\hat{d}^{\dagger}\hat{d}\) respectively. For \(N \to \infty\), \(E_{0,0} = \gamma/2 = c_0\), in agreement with mean field.

Having diagonalized \(\hat{H}\), we can easily get the occupancy of the \(\phi_{\pm 1}\) states, which is

$$\langle \hat{c}_{\pm 1}^{\dagger}\hat{c}_{\pm 1} \rangle = \lambda_{1}^{2}/N.\hspace{1cm} (23)$$

Expanding we get that close to the transition,

$$\langle \hat{c}_{\pm 1}^{\dagger}\hat{c}_{\pm 1} \rangle \approx \frac{\sqrt{2}}{8}(\gamma + 1/2)^{-1/2}N.\hspace{1cm} (24)$$

Since the occupancy of the \(\phi_{\pm 1}\) states is a smooth and continuous function for any value of \(\gamma\) when \(N\) is finite, and since in the limit \(N \to \infty\), \(\langle \hat{c}_{\pm 1}^{\dagger}\hat{c}_{\pm 1} \rangle / \hat{N}\) has to equal \(|c_{\pm 1}|^{2}\) [10], a simple geometric construction shows that the value of the function in Eq. (24) at \(\gamma = -1/2 + \delta\), minus the value of \(|c_{\pm 1}|^{2}\) at \(\gamma = -1/2 - \delta\) has to be proportional to \(\delta\), and thus

$$\delta = \eta N^{-2/3},$$  \hspace{1cm} (25)

in agreement with Ref. [5], where it was found numerically that \(\eta \approx 1.077\). Therefore to leading order

$$|c_{\pm 1}|^{2} = \frac{4}{7}(\gamma + 1/2 + \eta N^{-2/3}),\hspace{1cm} (26)$$

and

$$\gamma_{cr} = -1/2 - \eta N^{-2/3}.\hspace{1cm} (27)$$

Turning to the energy spectrum, in the lowest state \(n_b = n_d = 0\), and thus the ground state energy is

$$E_{0}(\gamma) = -(\gamma + 1) + (2\gamma + 1)^{1/2},$$  \hspace{1cm} (28)

measured from the energy of the state with uniform density, \(\gamma(N-1)/2\). Making use of Eq. (27) we plot the function \(E_{0}(\gamma + \eta N^{-2/3})\) in Fig. 2 for \(N = 100\) (top curve), 500 (middle curve), and \(N \to \infty\) (bottom curve), with \(\eta = 1.077\). From Eq. (27) we get that to leading order

$$E_{0}(\gamma = -1/2) = -(1/2 + 2\eta)^{1/2}N^{-1/3}.$$  \hspace{1cm} (29)

The low-lying excited states of \(\hat{H}\) lie above the lowest state by \(E(\gamma)(n_b + n_d)\), where

$$E(\gamma) = (2\gamma + 1)^{1/2}.$$

The energy of the lowest excited state is thus \(E(\gamma)\), in agreement with Ref. [5]. Figure 3 shows \(E(\gamma + \eta N^{-2/3})\) for \(N = 100\) (top curve), 500 (middle curve), and \(N \to \infty\) (bottom curve). From Eq. (27), \(E(\gamma = -1/2) = (2\eta)^{1/2}N^{-1/3}\).

Finally the depletion of the condensate \(\Delta N\) is,

$$\Delta N = \langle \hat{c}_{-1}^{\dagger}\hat{c}_{-1} \rangle + \langle \hat{c}_{1}^{\dagger}\hat{c}_{1} \rangle = 2\lambda_{1}^{2} = \left(\frac{\gamma + 1}{\sqrt{2\gamma + 1}} - 1\right).\hspace{1cm} (30)$$

Figure 4 shows \(\Delta N(\gamma + \eta N^{-2/3})\) for \(N = 100\) (bottom curve), 500 (middle curve), and \(N \to \infty\) (top curve).
According to Eq. (27), to leading order $\Delta N(\gamma = -1/2) = (\sqrt{2}/4)\eta^{-1/2}N^{1/3}$.

To summarize, we examined a Bose-Einstein condensate in a toroidal trap, with an effective attractive interaction between the atoms. Using a variational approach and working within an appropriately chosen set of basis states, we demonstrated that within the mean-field approximation for a strong enough coupling constant a condensate of uniform density becomes unstable against the formation of a localized state. Going beyond the mean-field approximation, we diagonalized the Hamiltonian in the same truncated space just above the transition and we calculated the energy of the low-lying states, as well as the depletion of the condensate. While the energy in this case is lower than that of the mean-field, the two methods give the same results in the $N \to \infty$ limit, assumed implicitly in mean-field.

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