Three-wave mixing of Bogoliubov quasi-particles in a Bose condensate

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A dressed basis is used to calculate the dynamics of three-wave mixing between Bogoliubov quasi-particles in a Bose condensate. Due to the observed oscillations between different momenta modes, an energy splitting, analogous to the optical Mollow triplet, appears in the Beliaev damping spectrum of the excitations from the oscillating modes.

Since the experimental realization of Bose-Einstein condensation in trapped atomic gases, which serves as a mono-energetic and dense atomic source, experiments in non-linear atom optics have become feasible. Atomic four-wave mixing (4WM) [1], Superradiance [2], and matter-wave amplification [3], [4] are all examples of non-linear atom optics, which involve the mixing of several atomic fields or the mixing of atomic and electromagnetic fields.

In the case of atomic 4WM, bosonic amplification directs one of the products of a collision between an excitation atom and a condensate atom into an, initially, largely populated mode [1], [5], [6]. Thus, a new mode is macroscopically populated through the second collision product. When the energy, $\varepsilon_k$, of an excitation mode $k$ is low compared to the condensate chemical potential, the excitation can no longer be described as a free atom moving with momentum $k$, but rather as a collective phonon excitation, which involves a large number of atoms [7]. Atomic 4WM is therefore inadequate for the description of phonon decay. Phonon excitations are described in the framework of Bogoliubov theory [8]. As Bogoliubov excitations are bosons, bosonic amplification will also direct one of the damping products of a phonon into an, initially, largely populated mode, leading to three-wave-mixing (3WM) of Bogoliubov quasi-particles.

In this letter we introduce a basis of states which are dressed by the interaction between three, largely populated, modes of Bogoliubov quasi-particles in a Bose-Einstein condensate at zero temperature. Using this basis of quantum states we calculate the time evolution of the system, which exhibits non-linear oscillations between the different momentum modes. The use of the wave-mixing eigen-basis enables the calculation of the system dynamics efficiently and with no approximations. In contrast to other theoretical approaches to wave-mixing, which lead to non-linear differential equations [5], [9] the dressed state approach turns the problem into a linear one, for which propagation in time is trivial. We show that, due to relative number squeezing, the variance in the number difference between the two low momentum modes remains constant.

In analogy to the treatment of spontaneous photon scattering in the atom-laser dressed system [10], we treat damping into the quasi-continuum of empty modes, during the 3WM dynamics, as transfer between dressed state manifolds. Thus, the effects of 3WM on the damping process, are calculated. The damping energy spectrum is presented. A transition from elastic to inelastic damping is observed, which, in analogy to the optical Mollow triplet, leads to a splitting of the spectrum into a doublet of resonance energies.

Previously, nonlinear mixing of quasi-particles in a Bose condensate was studied, between the two lowest energy excitation modes in the discrete regime, where the energy of the excited mode is of the order of the energy separation between modes [9], [11], [12].

Our model system is a homogenous condensate of finite volume $V$, with $N_0$ atoms in the ground state. The Hamiltonian for the system in the Bogoliubov basis, taken to the $\sqrt{N_0}$ order, is given by [13]

$$H = H_0 + H_{\text{int}},$$

where $H_0 = \frac{1}{2} g n N_0 - \sum_k \varepsilon_k b_k^\dagger b_k + \frac{1}{2} \sum_{k} \varepsilon_k b_k^\dagger b_k^\dagger b_k b_k$ is the part of $H$ which is diagonalized by the Bogoliubov transformation. $N_0$ is the number of atoms in the $k = 0$ mode. $\varepsilon_k = \sqrt{\varepsilon_k^2 + 2gn}$ is the recently measured Bogoliubov energy [14] and $\varepsilon_k^0 = \frac{\hbar^2 k^2}{2m}$ is the free particle energy. $g = 4\pi a_s/m$ is the coupling constant, $a$ is the s-wave scattering length, $m$ is the mass of the atoms. $b_k^\dagger$ and $b_k$ are the creation and annihilation operators respectively, of quasi-particle excitations with momentum $k$. $u_k$ and $v_k$ are the Bogoliubov quasi-particle amplitudes. This part of the Hamiltonian describes excitations in the condensate with momentum $k$ and energy $\varepsilon_k$.

$H_{\text{int}}$ is the part of the Hamiltonian which is responsible for the interaction between excitations

$$H_{\text{int}} = \frac{g}{2V} \sqrt{N_0} \sum_{k,q} A_{k,q} \left( b_k^\dagger b_q b_{-q}^\dagger + b_k^\dagger b_q^\dagger b_{-q} b_k \right).$$

The first term in parentheses is referred to as Landau damping and is analogous to photon up conversion, such as second harmonic generation in optics [15], whereas the second term is referred to as Beliaev damping, and is analogous to photon down conversion [16]. $A_{k,q}$ is the many-body suppression factor.
\[ A_{kq} = 2u_k (u_q u_{k-q} - v_q u_{k-q} - u_q v_{k-q}) - 2v_k (v_q v_{k-q} - u_q v_{k-q} - v_q u_{k-q}). \] (3)

The main damping mechanism from a single, largely populated mode \( k \), will be elastic Beliaev damping of the excitations into empty modes, which are on a monoenergetic surface in momentum space [17].

We now consider the case where the condensate is excited with \( N \) excitations of momentum \( k \) and \( M \) excitations of momentum \( q \), \(|N_k, M_q, 0_{k-q}\rangle\), such that \( k \) and \( q \) fulfill the Bragg condition, i.e. \( \varepsilon_k = \varepsilon_q + \varepsilon_{k-q} \). Two-photon Bragg transitions can be used to excite the condensate, and populate different momentum modes with a variable number of excitations [14]. The amplitude given by \( H_{int} \) for Beliaev damping of the \( k \) momentum excitation into two excitations with momenta \( q \) and \( k - q \), \(|(N-1)_k, (M+1)_q, 1_{k-q}\rangle\), will be \( \sqrt{N} \) or \( \sqrt{M} \) fold larger than any of the other damping channels. This will result in 3WM into a newly populated \( k - q \) momentum mode. We assume that \( N_0 >> N, M >> \Gamma t_0 \), where \( \Gamma \) is the total Beliaev damping rate of the excitations into the quasi-continuum of empty modes and \( t_0 \) is the time of the experiment. Thus, during the experiment time, 3WM dynamics dominates over damping into empty modes. In general the set of \( N + 1 \) excitation Fock states \(|(N-i)_k, (M+i)_q, i_{k-q}\rangle\), where \( i \) varies between 0 and \( N \), spans a degenerate subspace of the eigenstates of \( H_0 \). \( H_{int} \) couples between pairs of states in this subspace which have a difference of 1 in \( i \), and can be represented by the \( (N+1) \times (M+1) \) tri-diagonal matrix

\[ \langle N-s, M+s, s \mid H_{int} \mid N-i, M+i, i \rangle = \frac{g}{2V} \sqrt{N_0} A_{kq} \sqrt{N-i} \sqrt{M+i} + 1 \delta_{i+1,s} + \sqrt{N-i} + 1 \sqrt{M+i} + 1 \delta_{i-1,s}. \] (4)

When \( H_{int} \) is diagonalized, we get a new set of \( N + 1 \) eigenstates, \(|j\rangle\), where \( j \) varies between 1 and \( N + 1 \), that are dressed by the interaction. Note that since the dressed states are superpositions of degenerate eigenstates of \( H_0 \), they are eigenstates of the complete Hamiltonian in Eq. (1).

The filled circles in Fig. 1 show the calculated energy spectrum of the dressed basis for \( N = M = 100 \) in units of \( g \sqrt{N_0} / 2V \) \( A_{kq} \). The degeneracy between the excitation Fock states is removed by the interaction. The spectrum appears to be roughly linear with an average energy spacing of \( dE \approx 2.48 \sqrt{N} \) in the above units [18]. The hollow circles in Fig. 1 show the value of the energy difference between each two dressed states. Due to the non-linearity of the problem, energy differences between each pair of dressed states are slightly different.

![FIG. 1. Filled circles show the spectrum of the \( M = N = 100 \) manifold in units of \( \sqrt{N_0} A_{kq} \). Hollow circles show the energy difference between each two dressed states in the same units. The spectrum is not linear, and the energy spacing varies parabolically around an average of \( 2.48 \sqrt{N} \).](image)

![FIG. 2. The square of the transfer matrix between the excitation Fock state basis and the dressed state basis for \( N = M = 100 \). Darker areas correspond to larger probability. The white dashed-dotted line draws the solution to \( \varepsilon_{j} = 2(\sqrt{(N^2 - i^2)}, \) \( i \) The Fock states are numbered by \( i \). The dressed states are numbered by \( j \) from lowest to highest energy.](image)

Figure 2 shows the absolute value squared of the transfer matrix between the bare excitation Fock state basis and the dressed basis for \( N = M = 100 \). Since \( H_{int} \), after being applied a sufficient number of times, can couple each two bare states, most dressed states are spanned by a large number of bare states [19].

In the limit of \( M = N \rightarrow \infty \) we can use the tri-diagonality of \( H_{int} \) to approximate the dressed state as the solution to the differential equation

\[ \frac{\partial a_j^2(x)}{\partial x^2} = \left( \frac{\lambda_j}{\sqrt{1 - x^2}} - f_0^2 \right) a_j(x), \] (5)

with the boundary
conditions \( da_j(x)/dx = \left( \lambda_j/\sqrt{N} - N \right) a_j(x) \) at \( x = 0 \),
and \( da_j(x)/dx = \left( N - \lambda_j/\sqrt{2N} \right) a_j(x) \) at \( x = 1 \). Here
\( dx = 1/N \), \( x = i dx \), and \( a_j(x) = \langle N - i, M + i, i|j \rangle \)
is the projection of \( |j \rangle \) on the Fock state with \( i \) excitations removed from the \( k \) momentum mode. Also,
\( \lambda_j = E_j/\sqrt{N} \), \( E_j \) is the energy eigenvalue of \( |j \rangle \), and
\( f_0^2 = 2N^2 \). The main contribution to the dressed state superposition of Fock states comes from the two states
which solve \( \frac{|\lambda_i|}{\sqrt{1-x^2}} - f_0^2 = 0 \). The dashed-dotted line in Fig. 2 draws the solution of this equation. In or-
der for this equation to have real solutions the resulting spectrum must satisfy \( |E_j| < 1.24N^{3/2} \). Since there are
\( N + 1 \) dressed states we get an average energy difference between dressed states of \( dE \simeq 2.48\sqrt{N} \), consistent with
our numerical observation.

|FIG. 3. The expectation value \( N_k \). \( k = 0.7 \ h/\xi \) and \( q = k/\sqrt{2} \), as a function of time. \( N = M = 100 \), for a conden-
sate of \( 3 \times 10^5 \) Rb atoms in the \( F = 2 \), \( m_f = 2 \) ground state. The(condensate is homogeneous, with a density of \( 3 \times 10^{14} \)
\( \text{atoms/cm}^3 \). The oscillation frequency is roughly the average energy difference in the dressed state spectrum. Insets (a) and
(b) show \( N_k \) and the standard deviation of \( N_k \) respectively, during a longer time.

We choose as a model system a condensate of \( 3 \times 10^{5} \) \( \text{Rb} \) atoms in the \( F = 2 \), \( m_f = 2 \) ground state. The condensate, which is similar to the experimental param-
eters of [14] is homogeneous and has a density of \( 3 \times 10^{14} \)
\( \text{atoms/cm}^3 \). The \( k = 0.7 \ h/\xi \) mode and \( q = k/\sqrt{2} \) mode, where \( \xi \) is the healing length of the condensate given by
\( \xi = \sqrt{8\pi a} \), are populated with a 100 excitations each.

We start from the above excitation Fock state, written as a linear superposition of dressed states, which are ob-
tained by diagonalizing (4). The state of the system is then readily propagated in time by evolving each of the
dressed states phases according to its energy. Figure 3 shows \( N_k \), the expectation value of the number of excita-
tions with momentum \( k \), as a function of time. Excita-
tions oscillate between the \( k \) momentum mode and the \( q \)
and \( k - q \) momenta modes. Inset (a) of Fig. 3 shows \( N_k \)
during a much longer time. Since the energy spectrum is not precisely linear, at longer times beating between
the different oscillation frequencies gives rise to a slow amplitude modulation of the oscillations.

Even though we start from an excitation Fock state, the system immediately evolves into a superposition of Fock states. Inset (b) of Fig. 3 shows the standard devi-
ation of \( N_k \) vs. time. After a time scale which is set by the
non-linearity of the spectrum, the average value of \( N_k \)
and its standard deviation are of similar size. However,
the expectation value of the number difference between
mode \( q \) and mode \( k - q \) equals \( M \) and is constant in time.
The standard deviation of that difference always remains
zero, which implies relative number squeezing between
the two modes.

Thus far we have discussed the time evolution of the
system within the dressed state manifold, which is de-
\( \text{defined by the initial population of the \( k \) and \( q \) modes, and}
\( \text{and have neglected damping into empty modes. In analogy to the}
\text{treatment of spontaneous photon scattering as trans-
\( \text{fer between dressed states manifolds of the dressed atom-
\text{laser system [10], we now consider scattering into empty}
\text{modes as transfer between dressed state manifolds. Thus,
\text{the damping of excitations from mode \( k \) is no longer elas-
\text{tic, but rather carries the energy difference between the}
\text{dressed states among which it occurred.}

\text{Evaluating} \ H_B = \frac{2\pi}{\hbar} \sqrt{N_0} \sum_{\mathbf{q}} A_{\mathbf{k}\mathbf{q}} \left( b_{\mathbf{k}+\mathbf{q}}^+ b_{\mathbf{k}}^+ b_{\mathbf{k}-\mathbf{q}} b_{\mathbf{k}} \right),
\text{between every pair of states in the two manifolds reveals the}
\text{spectral structure of the damping process. We find that,}
\text{the spectrum of the \( N' = N - 1 \), \( M' = N \) manifold, which}
\text{has one less energy eigenvalue, is shifted by roughly half}
\text{the energy difference with respect to the \( N = M \) energy}
\text{spectrum.} \ H_B \text{significantly couples only dressed states
\text{with neighboring energies. This results in a structure of a}
\text{doublet in the Beliaev damping spectrum [20].}

\text{The finite lifetime of the dressed state results from the}
\text{fact that there is a quasi-continuum of \( N' = N - 1 \), \( M' = N \) manifolds, all with an identical energy spectrum, to
\text{which a dressed state in the \( N = M \) manifold can couple}
\text{to via} \ H_B \text{. To determine the width of each transition}
\text{between two dressed states, we use the Fermi golden rule.}
\text{The damping rate between a state} |j \rangle_{N,M}
\text{in the \( N, M \) manifold and a state} |i \rangle_{N-1,M}
\text{in the \( N - 1, M \) manifold is then given by}
\Gamma = \frac{2\pi}{\hbar} \sum_{\mathbf{q}} g^2 N_0 |A_{\mathbf{k}\mathbf{q}}|^2 \delta \left( \varepsilon_0 - \varepsilon_{\mathbf{k}+\mathbf{q}} - \varepsilon_{\mathbf{k}-\mathbf{q}} \right),
\text{where} \varepsilon_0 \text{is the energy difference between} |j \rangle_{N,M}
\text{and} |i \rangle_{N-1,M} \text{. Using momentum conservation the energy con-
\text{}servation} \delta \text{-function becomes a geometrical condition on}
\text{the angle} \theta \text{between} \mathbf{k} \text{and} \mathbf{q} \)
\begin{align}
\cos(\theta) &= \frac{1}{2kq'} \left[ k^2 + q'^2 + 1 - \sqrt{(\varepsilon_k + \varepsilon_0 - \varepsilon_q')^2 + 1} \right]. \tag{7}
\end{align}

where momentum is in units of \( h/\xi \), and energy is in units of \( gn \). Equation (6) is then simplified to

\begin{align}
\Gamma &= \frac{g}{2\xi^3h} \left| A_{kQ} \right|^2 \left| n_{-1,M} \langle i| b_k | j \rangle_{N,M} \right|^2 \times \frac{q'}{2k} \left( \frac{\varepsilon_k + \varepsilon_0 - \varepsilon_q'}{\sqrt{(\varepsilon_k + \varepsilon_0 - \varepsilon_q')^2 + 1}} \right) dq'. \tag{8}
\end{align}

The spectrum of each transition is taken as a Lorenzian with a width of \( \Gamma/2\pi N_k \) and a normalization of \( \Gamma/N_k \) around \( \varepsilon_0 \). \( N_k \) is the average occupation of mode \( k \) for the two dressed states involved. Averaging over all of the possible transitions between the manifolds, Fig. 4 shows the damping spectrum between the \( N = M = 5 \times 10^3 \) and the \( N = 5 \times 10^3 - 1 \), \( M = 5 \times 10^3 \) manifolds, for the same model system as in Fig. 3 for \( k = 3, 2, 1, 6 \), and \( 0.7 \ h/\xi \), and \( q = k/\sqrt{2} \) [21]. A clear doublet structure is evident. The inset of Fig. 3 shows the energy-conserving surfaces for the two center energies of the \( k = 0.7 \ h/\xi \) curve (solid line), and the energy conserving surface for elastic damping from the same mode (dashed line).

![FIG. 4. Damping spectrum between the \( N = M = 5 \times 10^3 \) manifold and the \( N = 5 \times 10^3 - 1 \), \( M = 5 \times 10^3 \) manifold, for the same condensate as in Fig. 2. The three curves are for \( k = 3.2 \) (dashed), \( k = 1.6 \) (dotted) and \( k = 0.7 \ h/\xi \) (solid line), \( q = k/\sqrt{2} \). The inset shows the energy-conserving surfaces for the two center frequencies of the \( k = 0.7 \ h/\xi \) curve (solid line) and the energy conserving surface for elastic damping from mode \( k \) (dashed line).](image)

The separation between the peaks in Fig. 4, which is equal to the oscillation frequency, decreases with \( k \) as \( A_{kQ} \). The decrease in the width of each peak as a function of \( k \) comes from two contributions. Firstly, since damping into empty modes is incoherent, it decreases as \( |A_{kQ}|^2 \). Secondly, as \( k \) decreases there are less allowed empty modes on the energy-conserving surface. Therefore, the doublet structure is more resolved for the lower \( k \) values. The difference in width between the positive and negative energy peaks is mainly due to the difference in the number of allowed empty modes on the energy conserving surfaces. For small enough \( k \) values, with respect to the oscillation frequency, the energy conserving surface for the negative energy peak completely disappears and only the positive energy peak remains. Another contribution to the width of the resonance in Fig. 4 is due to the non-linearity of the dressed state spectrum, which gives a slightly different energy for transitions between different dressed state pairs.

Several other mechanisms contribute to the broadening of the two peaks which are not included in Fig. 4. The fact that only the first scattering event occurs between the \( N = M = 5 \times 10^3 \) and the \( N = 5 \times 10^3 - 1 \), \( M = 5 \times 10^3 \) manifolds will further broaden the resonances. Since the energy splitting scales as \( \sqrt{N} \), in an experiment where one scatters \( dN \) atoms from mode \( k \), this will result in a relative broadening of \( \frac{dN}{\sqrt{N}} \). According to the same scaling, an initial coherent, rather than Fock, state will cause a relative broadening of \( \frac{dN}{N} \). The condensate finite size or inhomogeneous density profile, will further contribute to the width of the resonance. We estimate that for the experimental parameters of [14] a doublet structure in the Beliaev damping spectrum can be resolved. Experimentally, the energy doublet can be observed by computerized tomography analysis of time of flight absorption images of the 3WM system [22].

Another process which will transfer the system between manifolds is that of a two-photon Bragg transition from the condensate to either the \( k \) or \( q \) modes [14]. The line-shape of the Bragg transition into the \( k \) mode also splits into a doublet structure [23].

The dressed state formalism can be applied to other processes which involve bosonic amplification. We have applied this method to a strongly Bragg-driven condensate, with large momentum transfer, by diagonalizing \( H_{int} = \frac{\hbar 2k}{2} \left( a_k^\dagger a_0 + a_0^\dagger a_k \right) \). \( \Omega_R \) is the two-photon Rabi frequency, and to atomic 4WM with large relative momentum by diagonalizing \( H_{int} = \frac{\hbar}{2} \left( a_k^\dagger a_0^\dagger a_k - a_0 a_k^\dagger a_0 \right) \). Both systems show a doublet structure in the spectrum of collisions from a certain momentum mode, due to rapid oscillations in the population of that mode [23].

In conclusion, we calculate the wave-mixing dynamics between three, low \( k \), Bogoliubov quasi-particles in a Bose condensate. The Hamiltonian of this system is diagonalized to the next order in \( \sqrt{N} \). The resulting basis of dressed states allows for the efficient, linear, propaga-
tion of the system in time. Non-linear oscillations between the different momentum modes are observed. Relative number squeezing between the \( q \) and the \( k - q \) momentum modes is shown. Beliaev damping of excitations from these modes is treated as a transfer between dressed states manifolds. The damping process is shown to become inelastic and, similarly to the optical Mollow triplet, exhibits a doublet-like energy spectrum, which is more resolved for the low \( k \) values.

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[18] This scaling law was verified numerically over several \( N = M \) values. A similar scaling law can be written for the \( N \neq M \) case.
[19] Since excitations will usually bear an uncertainty of \( \sqrt{N} \) in their population, the choice of an excitation Fock state as the system initial condition is somewhat unrealistic. This will smear the system over an ensemble of manifolds. However, for large \( N, M \) the relative spread in the excitation population becomes less and less significant.
[20] Damping from the \( q \) momentum mode can be treated in a similar manner. Since the population in the \( q \) mode does not reach 0, the damping spectrum has the structure of a triplet.
[21] Since we cannot diagonalize matrices of size \( 5 \times 10^3 \) by 5 \times 10^3 \), we diagonalize smaller \( N, M \) matrices and scale the resulting energies by a factor of \( \sqrt{N} \).