Quantized Fields in a Nonlinear Dielectric Medium: A Microscopic Approach

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Abstract

Theories which have been used to describe the quantized electromagnetic field interacting with a nonlinear dielectric medium are either phenomenological or derived by quantizing the macroscopic Maxwell equations. Here we take a different approach and derive a Hamiltonian describing interacting fields from one which contains both field and matter degrees of freedom. The medium is modelled as a collection of two-level atoms, and these interact with the electromagnetic field. The atoms are grouped into effective spins and the Holstein-Primakoff representation of the spin operators is used to expand them in inverse powers of the total spin. When the lowest order term of the interaction is combined with the free atomic and field Hamiltonians, a Hamiltonian describing a theory of noninteracting polaritons results. When higher order terms are expressed in terms of polariton operators standard nonlinear optical interactions emerge. These are then compared to the results of phenomenological and macroscopic theories. The theory is also used to derive an effective Hamiltonian describing counterpropagating modes in a nonlinear
medium.
1 Introduction

The study of the propagation of quantized electromagnetic fields in nonlinear
dielectric media is an area in which there is presently considerable theoretical
and experimental activity [1,2]. This interest is being driven by both funda-
mental and practical considerations. On a fundamental level nonlinear optics
provides an example of an experimentally accessible nonlinear quantum field
theory, and the light which is produced can exhibit distinctly quantum effects
such as squeezing and quantum phase diffusion [3-7]. The practical interest
comes from the proposed use of optical solitons in fibers for long-distance
communication [8]. This has made necessary a better understanding of the
role of quantum noise in such systems.

The first step which must be taken in these problems is the quantization
of the electromagnetic field in the presence of a nonlinear dielectric. This can
be done in two ways which we shall call the macroscopic and microscopic ap-
proaches. In the first, the macroscopic approach, the medium is completely
described by its linear and nonlinear susceptibilities. No matter degrees of
freedom appear explicitly in this treatment. One then finds a Lagrangian
which produces the macroscopic Maxwell equations for the field in a nonlinear
medium. From it one finds the canonical momenta and the Hamiltonian.
Quantization is accomplished by imposing the standard equal-time commutation relations. We explored this approach was in a previous paper [9]. In the second approach, the microscopic, a model for the medium is constructed, and both the field and matter degrees of freedom appear in the theory. Both are quantized. This is what we shall explore here. As we shall see, the result is a theory of mixed matter-field modes which are coupled by a nonlinear interaction.

A number of problems arise in the quantized macroscopic theory which lead us to examine a microscopic one. The first is the difficulty in including dispersion. Dispersion is a result of the fact that the medium is nonlocal in time, i.e. the polarization at a given time depends not only on the electric field at that time but also on the field at earlier times. It is not clear whether nonlocal theories can be encompassed within the canonical formalism. (Drummond has shown how it can be done if one is considering a narrow frequency interval [10]. He then showed that additional intervals could be considered at the expense of introducing additional fields.) Another issue is operator ordering. Because the vacuum state of the theory now includes the effects of the medium, it is no longer clear that normal ordering is the appropriate form to use. Finally, it is expected that the inclusion of
losses in the theory will be most easily accomplished at the microscopic level [11].

The macroscopic and microscopic approaches are complementary; each has its own range of application and illuminates the other. The microscopic theory we present here, being more fundamental, is more appropriate for addressing basic problems such as dispersion and operator ordering. It is, however, at the moment restricted to rather simple models of the medium. Real media are more easily described by the macroscopic theory in that the medium is characterized by only a few parameters. It has been used, for example, by Drummond and Carter to study the propagation of quantized fields in silica fibers [4].

While Drummond and Carter have developed the macroscopic theory into a useful tool, much of the work done so far on the nonlinear optics of quantized fields has used a phenomenological approach. One assumes that the field in a $\chi^{(3)}$ medium is described by the nonlinear Schroedinger equation and the theory is quantized by imposing equal-space, rather than equal-time, commutation relations [5]. A study of the use of equal-space commutations has been made by Deutsch [12]. He concluded that for linear theories they lead to the same results as the more standard equal-time relations, but for
nonlinear theories it is expected that for some quantities this will no longer be the case. These issues are in need of further clarification.

We shall consider a medium consisting of two-level atoms. This is a model which we treated before, but in that case the field was restricted to a single mode [13]. This did not allow the inclusion of propagation. We now extend that work to include an arbitrary number of modes. This extension is not straightforward, because in the single-mode case all of the two-level atoms can be treated as one large spin. This is no longer possible when more than one mode is present. This necessitates modification of the methods used in our first paper. To accomplish this, we divide the medium into little boxes whose dimensions are smaller than the shortest wavelength which appears in the theory. The atoms in each of these boxes can be treated as a large spin and the Holstein-Primakoff representation of the resulting spin operators can then be used. This allows us to make a $1/N$ expansion of the Hamiltonian where $N$ is the number of atoms in a box [14]. A continuum version of the theory is obtained by letting the box size go to zero.

What results is a Hamiltonian containing electromagnetic and matter fields. The linear part of the Hamiltonian is essentially the same as that obtained by Hopfield in his treatment of linear dielectric media [15]. It can
be diagonalized by transforming from the matter and field modes to mixed matter-field modes, known as polaritons. When the nonlinear part of the Hamiltonian is expressed in terms of polariton operators the result consists of interaction terms which are similar to those which are familiar from non-linear optics, i.e. terms describing an intensity-dependent index of refraction and third harmonic generation. The effects of dispersion are automatically included because the polaritons have a dispersion relation which differs from that of photons. The operator ordering in the Hamiltonian is also dictated by the theory. The end result is a microscopic theory to which the macroscopic and phenonenological theories can be compared. The microscopic theory is derived in Sections II-IV and the comparisons with other theories are made in Section V.

2 Hamiltonian for a Medium of Two-Level Atoms

Let us consider a large number, $N$, of atoms distributed uniformly throughout a volume $V$. These atoms interact with the radiation field whose quantization volume is also taken to be $V$. For simplicity we shall assume that each atom
has a single optically active electron. The Hamiltonian describing this system is

\[ H = \sum_{j=1}^{N} \left[ \frac{1}{2m} (\vec{p}_j + e\vec{A}(\vec{x}_j))^2 + V(\vec{x}_j - \vec{R}_j) \right] + \frac{1}{2} \int d^3r [\dot{\vec{A}}^2 + (\nabla \times \vec{A})^2]. \]  

(1)

Here \( \vec{x}_j \) and \( \vec{p}_j \) are the position and momentum of the electron on the jth atom which is located at \( \vec{R}_j \). The potential energy of the electron in the jth atom is \( V(\vec{x}_j - \vec{R}_j) \). The charge on the electron is \(-e\) where \( e > 0 \). Finally, we are assuming that we are in the Coulomb gauge so that \( \nabla \cdot \vec{A} = 0 \).

We now make two approximations. First we wish to specialize to the case where only two levels for each atom are included; an upper level \(|a⟩\) with energy \( E_0/2 \) and a lower level \(|b⟩\) with energy \(-E_0/2 \). The projection operator onto the two-state space is

\[ P = \prod_{j=1}^{N} (|a_j⟩⟨a_j| + |b_j⟩⟨b_j|), \]  

(2)

and the effective Hamiltonian for the two-level atoms is obtained by applying it to both sides of Eq.(1). Our Hamiltonian now becomes

\[ H = P \left\{ \sum_{j=1}^{N} \left[ \frac{1}{2} E_0 \sigma_j^{(3)} + \frac{e^2}{2m} \vec{A}(\vec{x}_j)^2 \right] + \frac{1}{2} \int d^3r [\dot{\vec{A}}^2 + (\nabla \times \vec{A})^2] \right\} P \]

\[ + \frac{e}{m} \left[ \sum_{j=1}^{N} \vec{p}_j \cdot \vec{A}(\vec{x}_j) \right] P. \]  

(3)

Because we are dealing with optical wavelengths we are also in a position to make the dipole approximation and replace \( \vec{x}_j \), the electron coordinate, by
\( \vec{R}_j \), the position vector of the atom, in the argument of \( \vec{A} \). This yields

\[
H = P \left\{ \sum_{j=1}^{N} \left[ \frac{1}{2} E_0 \sigma_j^{(3)} + \frac{e^2}{2m} \vec{A}(\vec{R}_j)^2 \right] + \frac{1}{2} \int d^3r \left[ \dot{\vec{A}}^2 + (\nabla \times \vec{A})^2 \right] \right\} P
\]

\[
+ \frac{e}{m} P \left[ \sum_{j=1}^{N} \vec{p}_j \cdot \vec{A}(\vec{R}_j) \right] P.
\]

(4)

Let us now examine the last term of the above equation which represents the dipole interaction between the atoms and the field. As a first step we expand the vector potential in plane wave modes

\[
\vec{A}(\vec{R}) = \sum_{\vec{k},\lambda} \sqrt{\frac{1}{2\omega_k V}} (a_{\vec{k},\lambda} e^{i\vec{k} \cdot \vec{R}} + a_{\vec{k},\lambda}^\dagger e^{-i\vec{k} \cdot \vec{R}}) \hat{\epsilon}_\lambda(\vec{k}),
\]

(5)

where \( \omega_k = |\vec{k}| \) and the polarization vectors, \( \hat{\epsilon}_\lambda(\vec{k}) \), where \( \lambda = 1, 2 \), satisfy \( \vec{k} \cdot \hat{\epsilon}_\lambda(\vec{k}) = 0 \) and are real. We also adopt the convention

\[
\hat{\epsilon}_\lambda(-\vec{k}) = (-1)^\lambda \hat{\epsilon}_\lambda(\vec{k}).
\]

(6)

If we now define

\[
i \mu_\lambda(\vec{k}) = \frac{e}{m} (a|\vec{p}|b) \cdot \hat{\epsilon}_\lambda(\vec{k}) = ieE_0 \langle a|\vec{x}|b \rangle \cdot \hat{\epsilon}_\lambda(\vec{k}),
\]

(7)

then the interaction term becomes

\[
\frac{e}{m} \sum_{j=1}^{N} P[\vec{A}(\vec{R}_j) \cdot \vec{p}_j] P = P \sum_{j=1}^{N} \sum_{\vec{k},\lambda} \sqrt{\frac{1}{2\omega_k V}} (a_{\vec{k},\lambda} e^{i\vec{k} \cdot \vec{R}_j} + a_{\vec{k},\lambda}^\dagger e^{-i\vec{k} \cdot \vec{R}_j})
\]

\[
i (\mu_\lambda(\vec{k}) \sigma_j^{(+)}) - \mu_\lambda^*(\vec{k}) \sigma_j^{(-)} \right\} P
\]

(8)
At this point we shall make a number of simplifications. First, we shall not write the projection operators explicitly and shall assume that all matter operators act only on the two levels, $|a\rangle$ and $|b\rangle$, of each atom. We shall also assume that the phases of the atomic wave functions have been chosen so that $\mu_\lambda(\vec{k})$ is real. Finally, we shall look at the case when only one polarization of the field is significantly populated so that the polarization index, $\lambda$, will be dropped. The remaining polarization will be assumed to obey Eq. (6) with $\lambda = 2$.

We now wish to rephrase the theory in terms of effective spins. To this end we divide the total volume, $V$, into blocks of volume $\Delta V$. The center of the $l$th block is located at the position $\vec{r}_l$. We suppose that there are $n_0$ atoms in each block, where $n_0 >> 1$, and that the dimensions of each block are much smaller than an optical wavelength. This means that if we are looking at optical phenomena all of the atoms in the $l$th block can be treated as if they are at $\vec{r}_l$. The atoms can then be treated as a spin $s = n_0/2$ object located at $\vec{r}_l$. Expressing the Hamiltonian in terms of the block variables we have

$$H = \sum_{l=1}^{N_b} E_0 S_i^{(3)} + \sum_\vec{k} \omega_\vec{k} a_\vec{k}^\dagger a_\vec{k} + \frac{e^2 n_0}{2m} \sum_{l=1}^{N_b} \vec{A}(\vec{r}_l)^2 + H_{int}, \quad (9)$$
where

\[ H_{\text{int}} = \sum_{l=1}^{N_b} \sum_{\vec{k}} \sqrt{\frac{1}{2\omega_k V}} (a_{\vec{k}} e^{i\vec{k} \cdot \vec{r}_l} + a_{\vec{k}}^\dagger e^{-i\vec{k} \cdot \vec{r}_l}) i\mu(\vec{k}) (S_l^{(\uparrow)} - S_l^{(\downarrow)}). \]  

(10)

We have designated the \( z \) component of the spin at \( \vec{r}_l \) by \( S_l^{(z)} \) and the raising and lowering operators by \( S_l^{(\uparrow)} \) and \( S_l^{(\downarrow)} \), respectively. \( N_b \), which is equal to \( V/\Delta V \), is the total number of blocks.

We can further simplify the \( \vec{A}^2 \) term in the Hamiltonian. If we multiply and divide the sum by \( \Delta V \) we have

\[ \frac{e^2}{2m \Delta V} \sum_{l=1}^{N_b} \vec{A}(\vec{r}_l)^2 \approx \frac{e^2 \rho}{2m} \int_V d^3r \vec{A}(\vec{r})^2 \]

\[ \approx \frac{e^2 \rho}{2m} \sum_{\vec{k}} \frac{1}{2\omega_k} [a_{\vec{k}} a_{-\vec{k}}^\dagger + a_{\vec{k}}^\dagger a_{-\vec{k}}^\dagger] + a_{\vec{k}} a_{\vec{k}}^\dagger + a_{\vec{k}}^\dagger a_{-\vec{k}}]. \]  

(11)

We have denoted by \( \rho \) the density of atoms which is equal to either \( N/V \) or \( n_0/\Delta V \).

We now have our Hamiltonian in the desired form. The medium is described by spin variables and the field by creation and annihilation operators. The next step is to expand it so that we can extract the linear interaction and the different orders of nonlinear interaction. We accomplish this with a semiclassical, or \( 1/s \), expansion.
3 Expansion and Continuum Limit

Before proceeding to the actual expansion we must restrict the $\vec{k}$ sum- 

tations by imposing a high-frequency cutoff. This is necessary so that the 

wavelengths in the theory do not become so small that they violate the con- 

ditions under which the macroscopic theory is valid. In particular, we have 

assumed that the mean spacing between atoms is much less than any wave- 

length in the theory. Therefore, we shall restrict $|\vec{k}|$ to be less than $k_u$ where 

$k_u$ corresponds to a wavelength shorter than those in the optical regime but 

considerably larger than the interatomic separation.

We would expect that the ground state of the field-atom system would 

be the state in which all of the spins are down, corresponding to all of the 

atoms in their ground states, and no photons present. This is actually the 

case in the semiclassical approximation. A derivation of this fact is presented 

in Appendix A. Here we shall assume this to be the case and shall expand 

our Hamiltonian about the no-photon, all-spin-down state.

In the case of the spin operators this expansion is implemented by means 

of the Holstein-Primakoff representation [16] in which the spin $s$ operators 

$S^{(3)}, S^{(+)}$, and $S^{(-)}$ are represented in terms of boson creation and annihila-
tion operators, $\zeta^\dagger$ and $\zeta$ as

$$S^{(-)} = (2s - \zeta^\dagger\zeta)^{1/2}\zeta \quad S^{(+)} = \zeta^\dagger(2s - \zeta^\dagger\zeta)^{1/2}$$

$$S^{(3)} = -s + \zeta^\dagger\zeta,$$  \hspace{1cm} (12)

where $[\zeta, \zeta^\dagger] = 1$. Our convention for the square roots is that when the argument is positive so is the square root, and when the argument is negative, the square root is $i$ times a positive number. Eqs. (12) then give us the proper commutation relations for the spin operators, i.e.

$$[S^{(3)}, S^{(\pm)}] = \pm S^{(\pm)} \quad [S^{(+)}, S^{(-)}] = 2S^{(3)}.$$  \hspace{1cm} (13)

The excitation number for the boson operators, i.e. the eigenvalue of $\zeta^\dagger\zeta$, corresponds to $s_3 + s$ where $s_3$ is the eigenvalue of $S^{(3)}$. Therefore, the boson vacuum state is the spin state with the spin pointing down. If we are only considering states whose excitation number is small then we can expand the square roots

$$S^{(-)} \cong \sqrt{2s} \left(1 - \frac{1}{4s}\zeta^\dagger\zeta\right)\zeta \quad S^{(+)} = \sqrt{2s}\zeta^\dagger \left(1 - \frac{1}{4s}\zeta^\dagger\zeta\right).$$  \hspace{1cm} (14)

In our model of a nonlinear medium the fraction of atoms in a block which is excited will be small because we are off resonance. This corresponds to small excitation numbers. Therefore, the use of the expansions in Eq. (14)
is justified. With these approximations our Hamiltonian becomes

$$H = H_0 + H_{int}^{(1)} + H_{int}^{(2)},$$

(15)

where

$$H_0 = \sum_{l=1}^{N_b} E_0 (-s + \zeta_l^{\dagger} \zeta_l) + \sum_{|k|<k_u} \omega_k a_k^{\dagger} a_k + \frac{e^2 \rho}{2m} \int_V d^3r \bar{A}(\vec{r})^2$$

(16)

$$H_{int}^{(1)} = i \sum_{l=1}^{N_b} \sum_{|k|<k_u} \sqrt{\frac{s}{\omega_k V}} \mu(\vec{k}) (a_k e^{i\vec{k}\cdot\vec{r}_l} + a_k^{\dagger} e^{-i\vec{k}\cdot\vec{r}_l}) (\zeta_l^{\dagger} - \zeta_l)$$

(17)

$$H_{int}^{(2)} = -\frac{i}{4s} \sum_{l=1}^{N_b} \sum_{|k|<k_u} \sqrt{\frac{s}{\omega_k V}} \mu(\vec{k}) (a_k e^{i\vec{k}\cdot\vec{r}_l} + a_k^{\dagger} e^{-i\vec{k}\cdot\vec{r}_l})$$

$$((\zeta_l^{\dagger})^2 \zeta_l - \zeta_l^{\dagger} \zeta_l^2),$$

(18)

and $[\zeta_l, \zeta_l^{\dagger}] = \delta_{ll'}$. $H_{int}^{(1)}$ represents the linear part of the interaction between the atoms and the field while $H_{int}^{(2)}$ represents the nonlinear interaction.

We now want to go to a continuum representation of the matter operators. Instead of the operators $\zeta_l$ and $\zeta_l^{\dagger}$ we wish to employ operators, $\zeta(\vec{r})$ and $\zeta^{\dagger}(\vec{r})$ which are functions of a continuous position variable and whose commutation relations are

$$[\zeta(\vec{r}), \zeta^{\dagger}(\vec{r}')] = \delta^3(\vec{r} - \vec{r}').$$

(19)

Note that the operators $\zeta_l/\sqrt{\Delta V}$ and $\zeta_l^{\dagger}/\sqrt{\Delta V}$ have the commutation relations

$$\left[ \frac{1}{\sqrt{\Delta V}} \zeta_l, \frac{1}{\sqrt{\Delta V}} \zeta_l^{\dagger} \right] = \frac{1}{\Delta V} \delta_{ll'},$$

(20)
which in the limit $\Delta V \to 0$ limit goes to Eq. (19). This gives us the identification

$$\frac{1}{\sqrt{\Delta V}} \zeta_l \to \zeta(\vec{r}).$$  \hspace{1cm} (21)

This relation can be used to find the continuum representation of the different terms in the Hamiltonian. For example, the $\zeta_l^\dagger \zeta_l$ term becomes

$$\sum_{l=1}^{N_b} \zeta_l^\dagger \zeta_l = \sum_{l=1}^{N_b} \frac{1}{\sqrt{\Delta V}} \zeta_l^\dagger \frac{1}{\sqrt{\Delta V}} \zeta_l \Delta V \to \int_V d^3r \zeta^\dagger(\vec{r}) \zeta(\vec{r}),$$  \hspace{1cm} (22)

which, after noting that $N = 2sN_b$, gives us

$$H_0 = -\frac{1}{2}NE_0 + E_0 \int_V d^3r \zeta^\dagger(\vec{r}) \zeta(\vec{r}) + \sum_{|k|<k_u} \omega_k a_k^\dagger a_k^\vphantom{\dagger}
+ \frac{e^2\rho}{2m} \sum_{|k|<k_u} \frac{1}{2\omega_k} [a_k^\vphantom{\dagger} a_{-k}^\dagger + a_k^\dagger a_{-k}^\vphantom{\dagger} + a_k^\dagger a_k^\dagger + a_{-k}^\vphantom{\dagger} a_{-k}^\dagger].$$  \hspace{1cm} (23)

Similarly, for $H^{(1)}_{\text{int}}$ and $H^{(2)}_{\text{int}}$ we find

$$H^{(1)}_{\text{int}} = \frac{i}{\sqrt{V}} \sum_{|k|<k_u} \int_V d^3r (a_k e^{i\vec{k} \cdot \vec{r}} + a_k^\dagger e^{-i\vec{k} \cdot \vec{r}}) g_k \zeta^\dagger(\vec{r}) - \zeta(\vec{r}))$$  \hspace{1cm} (24)

$$H^{(2)}_{\text{int}} = -\frac{i}{2\rho \sqrt{V}} \sum_{|k|<k_u} \int_V d^3r (a_k e^{i\vec{k} \cdot \vec{r}} + a_k^\dagger e^{-i\vec{k} \cdot \vec{r}}) g_k \zeta^\dagger(\vec{r})^2 \zeta(\vec{r})$$

$$- \zeta^\dagger(\vec{r})\zeta(\vec{r})^2),$$  \hspace{1cm} (25)

where $g_k = \mu(\vec{k}) \sqrt{\rho/2\omega_k}$.  

16
4 Diagonalization of the Quadratic Part

So far we have derived a Hamiltonian which contains two interacting fields, the electromagnetic field and the matter field. These fields have both linear and nonlinear interactions with each other. We have yet to see an interaction of the type found in nonlinear optics, that of a field interacting nonlinearly with itself. For example, it is not immediately clear how something that looks like the usual description of self-phase modulation would be described by our Hamiltonian.

Something which does look more like the usual theory emerges if we diagonalize the quadratic part of the Hamiltonian, i. e. $H_0 + H_{int}^{(1)}$. What results is a description in terms of mixed matter-field modes, or polaritons [15]. If $H_{int}^{(2)}$ is expressed in terms of polariton operators, it describes a nonlinear interaction between polaritons. Therefore, instead of a Hamiltonian describing the interaction of a photon field with itself, which phenomenological effective Hamiltonians in nonlinear optics do, we have a Hamiltonian which describes the interaction of a polariton field with itself.

In order to diagonalize $H_0 + H_{int}^{(1)}$ we introduce the operators

$$\zeta_\mathbf{\vec{k}} = \frac{1}{\sqrt{V}} \int_V d^3r e^{-i\mathbf{\vec{k}} \cdot \mathbf{\vec{r}}} \zeta(\mathbf{\vec{r}}),$$

(26)
which implies that
\[ \zeta(\vec{r}) = \frac{1}{\sqrt{V}} \sum_{\vec{k}} e^{i\vec{k} \cdot \vec{r}} \zeta_{\vec{k}}. \] (27)

These operators have the commutation relations
\[ [\zeta_{\vec{k}}, \zeta_{\vec{k}'}^\dagger] = \delta_{\vec{k},\vec{k}'}. \] (28)

In terms of these operators \(H_0 + H_{\text{int}}^{(1)}\) become
\[
H_0 = -\frac{1}{2} N E_0 + E_0 \sum_{|\vec{k}| < k_u} \zeta_{\vec{k}}^\dagger \zeta_{\vec{k}} + \sum_{|\vec{k}| < k_u} \omega_k a_{\vec{k}}^\dagger a_{\vec{k}} \\
+ \frac{e^2 \rho}{2m} \sum_{|\vec{k}| < k_u} \frac{1}{2\omega_k} [a_{\vec{k}} a_{-\vec{k}} + a_{\vec{k}}^\dagger a_{-\vec{k}} + a_{\vec{k}}^\dagger a_{\vec{k}} + a_{-\vec{k}}^\dagger a_{-\vec{k}}] \] (29)

and
\[
H_{\text{int}}^{(1)} = i \sum_{|\vec{k}| < k_u} g_{\vec{k}} [a_{\vec{k}} (\zeta_{\vec{k}}^\dagger - \zeta_{-\vec{k}}) + a_{\vec{k}}^\dagger (\zeta_{\vec{k}} - \zeta_{-\vec{k}}^\dagger)]. \] (30)

We first note that only the modes \(\vec{k}\) and \(-\vec{k}\) are coupled to each other. Therefore, we can diagonalize \(H_0 + H_{\text{int}}^{(1)}\) by considering modes a pair at a time and adding up the results. Let \(H_{\vec{k}}\) be the part of \(H_0 + H_{\text{int}}^{(1)}\) containing operators for the modes \(\vec{k}\) and \(-\vec{k}\). It can be expressed as
\[
H_{\vec{k}} = \nu^\dagger B \nu, \] (31)
where

\[
\nu = \begin{pmatrix}
a_{\mathbf{k}} \\
\zeta_{\mathbf{k}} \\
a^\dagger_{-\mathbf{k}} \\
\zeta^\dagger_{-\mathbf{k}}
\end{pmatrix}
\]  \hspace{1cm} (32)

and

\[
B = \begin{pmatrix}
\omega_k + \frac{e^2 \rho}{2m \omega_k} & -ig_k & \frac{e^2 \rho}{2m \omega_k} & ig_k \\
ig_k & E_0 & ig_k & 0 \\
\frac{e^2 \rho}{2m \omega_k} & -ig_k & \omega_k + \frac{e^2 \rho}{2m \omega_k} & ig_k \\
-ig_k & E_0 & -ig_k & E_0
\end{pmatrix}
\]  \hspace{1cm} (33)

In arriving at Eq.(31) we have dropped all constants, such as those arising from commutation relations, from the Hamiltonian, and used the fact that for our choice of polarization, \( g_{-\mathbf{k}} = g_{\mathbf{k}} \).

Following Hopfield we want to introduce operators which are linear combinations of the components of \( \nu \). We shall let \( \alpha_{\mathbf{k}} \) and \( \beta_{\mathbf{k}} \) be these new operators where we require

\[
[\alpha_{\mathbf{k}}, \alpha^\dagger_{\mathbf{k}}] = [\beta_{\mathbf{k}}, \beta^\dagger_{\mathbf{k}}] = 1 \]  \hspace{1cm} (34)

and

\[
[\alpha_{\mathbf{k}}, \beta_{\mathbf{k}}] = [\alpha_{\mathbf{k}}, \beta^\dagger_{\mathbf{k}}] = 0. \]  \hspace{1cm} (35)
Defining the vector
\[
v = \begin{pmatrix}
\alpha_k \\
\beta_k \\
\alpha^\dagger_{-k} \\
\beta^\dagger_{-k}
\end{pmatrix},
\]  
(36)
we have that
\[
v = M\nu,
\]  
(37)
where \( M \) is a 4x4 matrix yet to be determined. The commutation relations obeyed by the elements of \( v \) and \( \nu \) imply that \( M \) has the property that
\[
M^{-1} = GM^\dagger G,
\]  
(38)
where the matrix \( G \) has the elements \( G_{ij} = \delta_{ij} \) if \( i, j = 1, 2 \), and \( G_{ij} = -\delta_{ij} \) if \( i, j = 3, 4 \), with all other matrix elements being zero. From Eq. (38) we see immediately that \( M \) is not unitary.

Returning to \( H_k \) we have that
\[
H_k = v^\dagger(M^{-1})^\dagger BM^{-1}v,
\]  
(39)
which we want to be diagonal. This implies that
\[
(M^{-1})^\dagger BM^{-1} = D
\]  
(40)
where $D$ is a diagonal matrix, or, making use of Eq. (38) and the fact that $G^2 = 1$,
\[ M(GB)M^{-1} = GD. \] (41)

Noting that if $D$ is diagonal so is $GD$ we see that $M$ must diagonalize the matrix $GB$.

We can reduce finding $M$ to finding the eigenvectors of $GB$. Let us denote the eigenvectors of $GB$ by $\eta_i$, where $i = 1, \ldots, 4$, and the unit vector whose only nonzero component is the $i$th by $c(i)$, i.e. $c(i)_j = \delta_{ij}$. If $M$ satisfies the equation
\[ M\eta_i = c(i), \] (42)
then we find that
\[ \langle c(i)|M(GB)M^{-1}|c(j)\rangle = \lambda_i \delta_{ij}, \] (43)
where $\lambda_i$ is the eigenvalue of $GB$ corresponding to $\eta_i$. Therefore, if $M$ satisfies Eq. (42) it will diagonalize $GB$. This equation also gives us immediately an explicit representation for $M^{-1}$
\[ M^{-1} = \sum_{i=1}^{4} |\eta_i\rangle \langle c(i)|. \] (44)

It is then possible to use Eq. (38) to find $M$
\[ M = G(M^{-1})^\dagger G = G \sum_{i=1}^{4} |c(i)\rangle \langle \eta_i| G \] (45)
Finally, let us note that Eq. (38) imposes a condition on the eigenvectors \( \eta_i \).

We have from this equation that

\[
(M^{-1})^\dagger GM^{-1} = G. \quad (46)
\]

Substitution of Eq.(42) into this result gives us that

\[
\langle \eta_j | G | \eta_i \rangle = G_{ij}. \quad (47)
\]

For \( i \neq j \) this condition is automatically satisfied because \( \eta_i \) and \( \eta_j \) are eigenvalues of \( GB \) corresponding to different eigenvalues. For \( i = j \) it can be imposed as a normalization condition.

The only task remaining is to find the eigenvectors and eigenvalues of \( GB \). Explicit expressions are given in Appendix B along with those for the matrix elements of \( M \). Those results along with Eqs. (31) and (37) imply that

\[
H_\vec{k} = E_1(k)\alpha_\vec{k}^\dagger\alpha_\vec{k} + E_2(k)\beta_\vec{k}^\dagger\beta_\vec{k} + E_1(k)\alpha_{-\vec{k}}^\dagger\alpha_{-\vec{k}} + E_2(k)\beta_{-\vec{k}}^\dagger\beta_{-\vec{k}},
\]

where

\[
E_1(k) = \frac{1}{\sqrt{2}} \left( [E_0^2 + \omega_k(\omega_k + 2C_0)]
+ \left[ [E_0^2 - \omega_k(\omega_k + 2C_0)]^2 + 16E_0\omega_kg_{\vec{k}}^2 \right]^{1/2} \right)^{1/2} \quad (49)
\]
\begin{align*}
E_2(k) &= \frac{1}{\sqrt{2}} ([E_0^2 + \omega_k(\omega_k + 2C_0)] \\
&\quad - [[E_0^2 - \omega_k(\omega_k + 2C_0)]^2 + 16E_0\omega_kg_k^2]^{1/2})^{1/2},
\end{align*}
(50)

with \( C_0 = (e^2\rho)/(2m\omega) \). All of the square roots in the above equations are taken to be positive. The energies \( E_1(k) \) and \( E_2(k) \) are just the two branches of the polariton energy curve. For large values of \( \omega_k \) we find that \( E_1(k) \to \omega_k \) and \( E_2(k) \to E_0 \). As \( \omega_k \to 0 \) we find that \( E_1(k) \) goes to a value slightly larger than \( E_0 \)
\begin{equation}
E_1(k) \to E_0 + 2\frac{\omega_kg_k^2}{E_0^2},
\end{equation}
(51)

and \( E_2(k) \) goes to
\begin{equation}
E_2(k) = \left(2\omega_kC_0 - \frac{4\omega_kg_k^2}{E_0^2} + \omega_k^2 \right)^{1/2}.
\end{equation}
(52)

It should be noted that \( \omega_kC_0 \) and \( \omega_kg_k^2 \) are independent of \( \omega_k \), and that \( \omega_kg_k^2 << E_0^3 \) and \( \omega_kC_0 << E_0^2 \). In Hopfield’s analysis the small constant term inside the parentheses vanished due to an atomic sum rule. This tells us that keeping only one excited atomic level is not a good approximation near \( \omega_k = 0 \). Because we are interested in optical phenomena this does not present a problem.

We can finally express \( H_0 + H_{int}^{(1)} \) in terms of the polariton operators.
Dropping constant terms we have

\[ H_0 + H_{\text{int}}^{(1)} = \sum_{|\mathbf{k}| < k_u} (E_1(k)\alpha_\mathbf{k}^\dagger \alpha_\mathbf{k} + E_2(k)\beta_\mathbf{k}^\dagger \beta_\mathbf{k}). \]  

(53)

The effect of the medium appears in two ways in this Hamiltonian. First, the operators are mixed matter-field operators, i.e. polariton, not photon, operators. Second, the effects of dispersion appear through the energies \(E_1(k)\) and \(E_2(k)\) which are not of photon form. Therefore, by including the matter degrees of freedom in the theory dispersion emerges naturally, and we avoid the problems of trying to quantize a theory which is nonlocal in time.

5 Nonlinear Interaction

We now turn our attention to the nonlinear part of the Hamiltonian, \(H_{\text{int}}^{(2)}\), which can be interpreted as describing an interaction between polaritons. With the interaction expressed as in Eq. (25), however, this interpretation is not obvious. In order to bring it out we begin by expressing it in terms of the operators \(\zeta_\mathbf{k}\)

\[ H_{\text{int}}^{(2)} = -\frac{i}{2\rho V} \sum_{|\mathbf{k}| < k_u} \cdots \sum_{|\mathbf{k}_3| < k_u} g_\mathbf{k} (\delta_{\mathbf{k}+\mathbf{k}_3,\mathbf{k}_1+\mathbf{k}_2} a_\mathbf{k} \zeta_\mathbf{k}_2^\dagger \zeta_\mathbf{k}_3^\dagger \zeta_\mathbf{k}_1^\dagger \\
- \delta_{\mathbf{k}+\mathbf{k}_2+\mathbf{k}_3,\mathbf{k}_1} a_\mathbf{k} \zeta_\mathbf{k}_2^\dagger \zeta_\mathbf{k}_3^\dagger + \delta_{\mathbf{k}+\mathbf{k}_2+\mathbf{k}_3,\mathbf{k}_1} a_\mathbf{k}^\dagger \zeta_\mathbf{k}_2^\dagger \zeta_\mathbf{k}_1^\dagger \\
- \delta_{\mathbf{k}+\mathbf{k}_3,\mathbf{k}_1+\mathbf{k}_2} a_\mathbf{k}^\dagger \zeta_\mathbf{k}_3^\dagger \zeta_\mathbf{k}_1^\dagger \zeta_\mathbf{k}_2^\dagger). \]  

(54)
The next step is to express the photon and atomic operators in terms of polariton operators. This leads to a great many terms so it is perhaps best to consider a specific physical process and then to select the terms which are relevant to it.

Let us first consider the situation when a single polariton mode, for example, the one corresponding to the operator \( \alpha_{\vec{k}_0} \) in the regime where \( \omega_{k_0} > E_0 \), is highly excited. The dominant terms in \( H_{int}^{(2)} \), at least for times which are not too long, will be those in which each of the four operators refers to the excited mode. The other terms will have a smaller effect on the time evolution because they contain at least one operator for a mode which is initially in the vacuum state. Examining the operator transformations in Eq. (37) we see that if we only keep the polariton operators \( \alpha_{\vec{k}_0} \) and \( \alpha_{\vec{k}_0}^{\dagger} \), then

\[

t_{\vec{k}_0} \rightarrow A_{11} (k_0) \alpha_{\vec{k}_0} \\
\zeta_{-k_0} \rightarrow A_{31} (k_0) \alpha_{\vec{k}_0}^{\dagger}
\]

where have set \( A = M^{-1} \). In the regime we have chosen, \( \omega_{k_0} > E_0 \) on the \( \alpha \) branch, we find that \( A_{11} (k_0) \) is of order one, \( A_{21} (k_0) \) and \( A_{41} (k_0) \) are of order \( g_{k_0}/E_0 \), and \( A_{31} (k_0) \) is of order \( C_0/E_0 \) which is considerably smaller than \( g_{k_0}/E_0 \). Therefore, we shall drop terms containing \( A_{31} (k_0) \). Keeping this in mind and making the substitutions indicated in Eq. (55) in Eq. (54)
we find

\[ H_{\text{int}}^{(\text{dom})} = -\frac{i g_{k_0}}{2\rho V} A_{11} [2 |A_{21}|^2 A_{21}^* (\alpha_{k_0}^\dagger)^2 (\alpha_{k_0})^2 + 2 |A_{41}|^2 A_{41}^* (\alpha_{k_0}^\dagger)^2 (\alpha_{k_0})^2 \]

\[ + |A_{21}|^2 A_{21}^* \alpha_{k_0}^\dagger \alpha_{k_0} (\alpha_{k_0})^2 + (|A_{41}|^2 A_{41}^* - |A_{21}|^2 A_{21}^*) \alpha_{k_0} (\alpha_{k_0}^\dagger)^2 \alpha_{k_0} \]

\[ + (|A_{21}|^2 A_{41} - |A_{41}|^2 A_{21}) \alpha_{k_0}^\dagger (\alpha_{k_0}^\dagger)^2 + 2 |A_{41}|^2 A_{41} \alpha_{k_0} \alpha_{k_0}^\dagger \alpha_{k_0} \alpha_{k_0}^\dagger \]

\[ + 2 |A_{41}|^2 A_{41}^* (\alpha_{k_0})^2 (\alpha_{k_0}^\dagger)^2, \quad (56) \]

where all of the matrix elements are evaluated at \( k_0 \), and we have made use of the fact that \( A_{11}(k_0) \) is real and both \( A_{21}(k_0) \) and \( A_{41}(k_0) \) are imaginary.

Let us note several things about this expression. First, it looks similar to what we would expect an interaction which describes self-phase modulation to look like. Therefore, by going to a description in terms of polariton operators we have recovered a familiar nonlinear optical interaction. Second, it is not normally ordered. If we use commutators to bring it into normal order we will pick up terms proportional to \( \alpha_{k_0}^\dagger \alpha_{k_0}^\dagger \). These represent small shifts to the polariton frequency and can be neglected in most applications. Neglecting them gives us

\[ H_{\text{int}}^{(\text{dom})} \approx \chi (\alpha_{k_0}^\dagger)^2 (\alpha_{k_0})^2 \]

\[ \quad (57) \]

where

\[ \chi = -\frac{i g_{k_0}}{2\rho V} A_{11} [2(|A_{21}|^2 + 2|A_{41}|^2) A_{21}^* + 2(|A_{41}|^2 + 2|A_{21}|^2) A_{41}]. \]

\[ \quad (58) \]
The effect of self-phase modulation on a single mode is often described by the Hamiltonian

\[ H_{\text{int}} = \lambda (a^\dagger)^2 a^2. \]  

which is superficially similar to Eq. (57). However, in most treatments the operators appearing in Eq. (59) are assumed to be photon creation and annihilation operators. As is shown by Eq. (57), they should be interpreted as polariton operators instead.

Now let us consider a more complicated situation. Suppose we initially have a pulse which is made up of modes on the \( \alpha \) branch with wave vectors near \( \vec{k}_0 \). In particular, let us assume that all of the wave vectors of the modes present in the pulse lie in a small region \( S \) about \( \vec{k}_0 \). Keeping only terms in \( H_{\text{int}}^{(2)} \) which contain four excited modes (after transforming to polariton operators) we obtain a rather complicated interaction. Each term in it contains two creation and two annihilation operators, and four elements of the matrix \( A(\vec{k}) \), each evaluated at a different wave vector. The interaction simplifies considerably if we make two approximations. First, we ignore operator ordering, which, as we saw, is tantamount to neglecting small frequency shifts. Second, we approximate each matrix element of \( A \) by its value at \( \vec{k}_0 \). Because the spread in wave vectors is small this is a good approximation. With these...
approximations we find

\[ H^{(\text{dom})}_{\text{int}} \cong \chi \sum_{\vec{k}_1 \in S} \sum_{\vec{k}_2 \in S} \sum_{\vec{k}_3 \in S} \sum_{\vec{k}_1, \vec{k}_2, \vec{k}_3} \delta_{\vec{k} + \vec{k}_1 + \vec{k}_2 + \vec{k}_3} \alpha_{\vec{k}} \alpha_{\vec{k}_1} \alpha_{\vec{k}_2} \alpha_{\vec{k}_3}, \]  

(60)

This situation is often treated by using a phenomenological Hamiltonian which describes pulse propagation in a \(\chi^{(3)}\) medium [5]. The Hamiltonian is

\[ H = \int dx \left[ \frac{\partial \phi^\dagger}{\partial x} \frac{\partial \phi}{\partial x} + c(\phi^\dagger)^2 \phi^2 \right], \]

(61)

where the field \(\phi(x, t)\) is the field envelope of the pulse in a frame moving at the group velocity of the pulse, and it obeys the commutation relations

\[ [\phi(x, t), \phi^\dagger(x', t)] = \delta(x - x'). \]

(62)

The equation of motion for the field operator \(\phi(x, t)\) resulting from this theory is the nonlinear Schroedinger equation. It can also be expressed in terms of creation and annihilation operators. Defining the annihilation operator

\[ a(\beta, t) = \frac{1}{\sqrt{2\pi}} \int dx e^{i\beta x} \phi(x, t), \]

(63)

we find for the equation of motion [5]

\[ i \frac{\partial a(\beta, t)}{\partial t} = \beta^2 a(\beta, t) + 2c \int d\beta_1 \int d\beta_2 a^\dagger(\beta_1, t) a(\beta_2, t) a(\beta + \beta_1 - \beta_2, t). \]  

(64)

It should be pointed out that the \(t\) in these equations is not really time and the \(x\) is not really space. The above formalism is derived by an analogy to
the classical theory which describes the propagation of a pulse in a nonlinear dispersive medium. The pulse is assumed to be propagating in the $z$ direction. The variable $x$ is $v_g t - z$, where $v_g$ is the group velocity, and $t$ is proportional to $z$. This means that what look like equal-time commutation relations in Eq. (62) are actually equal-space commutation relations. The question of when a theory using equal-space commutation relations gives the same results as a canonically quantized theory has not been fully answered. Deutsch has done a preliminary investigation and found that for a linear theory they are equivalent [12]. However, for a nonlinear theory he found indications that field correlation functions in which the fields are evaluated at different spatial points will not be the same in the two theories. Because of the different commutation relations, it is rather difficult to directly compare Eq. (64) to an equation of motion derived from our interacting polariton theory.

Instead we shall compare it to one due to Carter and Drummond who derived it by quantizing the macroscopic theory [2]. Their theory describes a pulse, consisting of modes with wave numbers centered about $k_c$ and frequencies centered about $\omega_c$, propagating in a medium of length $L$ in the positive $z$ direction. The basic field in this theory is

$$\Psi(z, t) = e^{-i k_c z + i \omega_c t} \frac{1}{\sqrt{L}} \sum_k e^{i k z} a_k(t), \quad (65)$$
which, because of the initial exponential factor, is slowly varying in space and time. The annihilation operators are defined in terms, not of the vector potential and the electric field, but in terms of the dual potential and the displacement field [2,13]. They, therefore, implicitly contain matter degrees of freedom and are, in a sense, polariton operators. The field $\Psi(z,t)$ obeys the equal-time commutation relations

$$[\Psi(z,t), \Psi^\dagger(z',t)] = \delta(z - z').$$

Like ours, the Carter-Drummond theory describes the pulse in the lab frame, i.e. the frame in which the medium is at rest. This removes one of the difficulties we had in trying to compare our theory with the phenomenological nonlinear-Schrödinger equation theory which is formulated in a moving frame. Carter and Drummond also reformulate their theory in a moving frame in order to make a comparison to the phenomenological theory, but we shall use their original lab frame results. In this frame their Hamiltonian is

$$H = \frac{1}{2} \int_0^L dz \left[ i v \left( \frac{\partial \Psi^\dagger}{\partial z} \Psi - \Psi^\dagger \frac{\partial \Psi}{\partial z} \right) \right.\
+ \omega'' \frac{\partial \Psi^\dagger}{\partial z} \frac{\partial \Psi}{\partial z} - v^2 \chi^E (\Psi^\dagger)^2 \Psi^2 \left. \right],$$

where $v$ is the group velocity of the pulse, $\chi^E$ is proportional to the third
order nonlinear susceptibility, and $\omega''$ is the second derivative of the frequency with respect to the wave number evaluated at $k_c$. The Hamiltonian and the commutation relations give, for the equation of motion for the operator $a_k(t)$,

$$
\frac{ia_k}{dt} = [\omega_c + v(k - k_c) + \frac{1}{2}\omega''(k - k_c)^2]a_k
- v^2 \chi \frac{1}{L} \sum_{k_1} \sum_{k_2} a_{k_1}^\dagger a_{k_2} a_{k+k_1-k_2}.
$$

(68)

It is this equation which we wish to compare to the corresponding equation derived from our nonlinear polariton theory.

Let us find the equation of motion for $\alpha_{\vec{k}}(t)$. We shall assume that we have a pulse made up of wave vectors near $\vec{k}_0$ so that the interaction is Eq. (60) is appropriate. Because all wave vectors in the pulse are close to $\vec{k}_0$, we shall expand the polariton energy $E_1(\vec{k})$ about $\vec{k}_0$

$$
E_1(\vec{k}) = E_1(\vec{k}_0) + (\delta\vec{k} \cdot \hat{k}_0) \frac{dE_1}{d\vec{k}} + \frac{1}{2k_0}(\delta\vec{k}^2 - (\delta\vec{k} \cdot \hat{k}_0)^2) \frac{d^2E_1}{d\vec{k}^2}
+ \frac{1}{2}(\delta\vec{k} \cdot \hat{k}_0)^2 \frac{d^2E_1}{d\vec{k}^2},
$$

(69)

where $\delta\vec{k} = \vec{k} - \vec{k}_0$, $\hat{k}_0 = \vec{k}_0/k_0$, and all derivatives are evaluated at $k_0 = |\vec{k}_0|$. Setting $v = dE_1/dk$, we find

$$
\frac{id\alpha_{\vec{k}}}{dt} \cong [E_1(k_0) + v\hat{k}_0 \cdot \delta\vec{k} + \frac{1}{2k_0}(\delta\vec{k}^2 - (\delta\vec{k} \cdot \hat{k}_0)^2)v + \frac{1}{2}(\delta\vec{k} \cdot \hat{k}_0)^2 \frac{d^2E_1}{d\vec{k}^2}]\alpha_{\vec{k}}
+ 2\chi \sum_{\vec{k}_1 \in S} \sum_{\vec{k}_2 \in S} \alpha_{\vec{k}_1}^\dagger \alpha_{\vec{k}_2} \alpha_{\vec{k}+\vec{k}_1-\vec{k}_2}.
$$

(70)
If the pulse is one dimensional, i.e. if $\delta \vec{k}$ is always parallel to $\hat{k}_0$, then Eqs. (68) and (70) have the same form. Thus we recover from the microscopic theory an equation of motion of the same form as that which arises from the quantized macroscopic theory. Note that extending the treatment here to broadband pulses is straightforward for microscopic theory; one simply does not expand the polariton energy as a function of $\vec{k}$. This extension is more complicated for the quantized macroscopic theory.

We can also examine the more complicated situation in which two modes are initially highly excited. Let us look at two cases, one when the modes are counterpropagating and one when they are not. The counterpropagating case is the more complicated of the two so we shall consider it second.

Suppose that the two polariton modes on the $\alpha$ branch with wave vectors $\vec{k}_a$ and $\vec{k}_b$ are initially highly excited. Let us assume that $\vec{k}_a$ and $\vec{k}_b$ are both in the $x$-$y$ plane, and that both modes are polarized in the $z$ direction. We shall also assume that $\vec{k}_a \neq -\vec{k}_b$. Keeping only the two excited modes, and making the substitutions in Eq. (55) for each mode, the Hamiltonian becomes

$$H_{\text{int}}^{(\text{dom})} \cong \chi_a (\alpha^{\dagger}_{k_a})^2 \alpha_{k_a}^2 + \chi_b (\alpha^{\dagger}_{k_b})^2 \alpha_{k_b}^2 + \chi_{ab} \alpha^{\dagger}_{k_a} \alpha^{\dagger}_{k_b} \alpha_{k_a} \alpha_{k_b}.$$  (71)

Here $\chi_a$ is simply $\chi$ with all of the matrix elements evaluated at $\vec{k}_a$, with a
similar definition for $\chi_b$, and $\chi_{ab}$ is given by

$$\chi_{ab} = -\frac{i}{\rho v}(F - F^*), \quad (72)$$

where

$$F = A_{11}^{(a)} g_{\vec{k}_a} [(|A_{21}^{(b)}|^2 + |A_{41}^{(b)}|^2)(A_{21}^{(a)*} + A_{41}^{(a)}) + A_{21}^{(b)*} A_{41}^{(b)} (A_{41}^{(a)*} + A_{21}^{(a)})]$$

$$A_{11}^{(b)} g_{\vec{k}_b} [(|A_{21}^{(a)}|^2 + |A_{41}^{(a)}|^2)(A_{21}^{(b)*} + A_{41}^{(b)})$$

$$+ A_{21}^{(a)*} A_{41}^{(a)} (A_{41}^{(b)*} + A_{21}^{(b)})] \quad (73)$$

In the above equation the superscript on the matrix element of $A$ indicates whether it is evaluated at $\vec{k}_a$ or $\vec{k}_b$. In Eq. (71) one has the usual terms which describe cross- and self-phase modulation of the two modes. It should be noted that if $|\vec{k}_a| = |\vec{k}_b|$, then with the above stated conditions on $\vec{k}_a$ and $\vec{k}_b$, and the polarizations, we have that $A(\vec{k}_a) = A(\vec{k}_b)$. This in turn implies that $\chi_{ab} = 4\chi_a = 4\chi_b$, and that the interaction in Eq. (73) can be derived from the kind of interaction appearing in the Hamiltonian in Eq. (61) or Eq. (67). If $|\vec{k}_a| \neq |\vec{k}_b|$, then this is no longer the case, but will be approximately true if $|\vec{k}_a| - |\vec{k}_b|$ is small and the dependence of $A(\vec{k})$ on $|\vec{k}|$ is weak.

Now let us look at the case of two counterpropagating modes. We shall assume that two modes on the $\alpha$ branch with wave vectors $\vec{k}_0$ and $-\vec{k}_0$, and with the same polarization, are initially highly excited. We again want to
keep only these two modes in our Hamiltonian, but the situation is now considerably more complicated than in our previous cases. This is because each of the matter or field operators will now be replaced by a sum of two polariton operators, e.g.,

\[ a_{\vec{k}_0} \rightarrow A_{11}(\vec{k}_0)\alpha_{\vec{k}_0} + A_{13}(\vec{k}_0)\alpha_{-\vec{k}_0}^\dagger, \]

\[ \zeta_{\vec{k}_0} \rightarrow A_{21}(\vec{k}_0)\alpha_{\vec{k}_0} + A_{23}(\vec{k}_0)\alpha_{-\vec{k}_0}^\dagger, \] (74)

where contributions from initially unpopulated modes have been dropped. We find that

\[
H_{\text{int}}^{(2)} \simeq \frac{ig_{\vec{k}_0}}{2\rho V}A_{11}(A_{21}-A_{23})\{6A_{21}A_{23}[2(\alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0})^2 + (\alpha_{\vec{k}_0} \alpha_{-\vec{k}_0})^2] \\
+ 3(A_{21}-A_{23})^2[(\alpha_{\vec{k}_0}^\dagger)^2 \alpha_{-\vec{k}_0}^\dagger \alpha_{\vec{k}_0} + (\alpha_{-\vec{k}_0}^\dagger)^2 \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0} + \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0} + \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \\
+ \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \} - 2(A_{21}^2 - A_{21}A_{23} + A_{23}^2)[(\alpha_{\vec{k}_0}^\dagger)^2 \alpha_{\vec{k}_0}^2 \\
+ (\alpha_{-\vec{k}_0}^\dagger)^2 \alpha_{-\vec{k}_0}^2 + 4\alpha_{\vec{k}_0}^\dagger \alpha_{-\vec{k}_0}^\dagger \alpha_{\vec{k}_0} \alpha_{-\vec{k}_0}], \] (75)

where all of the matrix elements are evaluated at \( \vec{k}_0 \). In deriving Eq. (75) we have made use of the fact that \( A_{11} \) and \( A_{13} \) are real, that \( A_{21} \) and \( A_{23} \) are imaginary (see Appendix B), and that \( A_{13} \) can be neglected in comparison to the other three. We have also dropped terms due to commutators. These are similar in form to the terms in Eq. (75) except that they contain only two \( \alpha_\vec{k} \) operators instead of four. They are, therefore, smaller than the terms we
have kept by a factor of the order of the number of photons in the $\alpha_{\vec{k}_0}$ and $\alpha_{-\vec{k}_0}$ modes. If we were to examine the population of initially unpopulated modes by the two counterpropagating beams, these terms should be kept. In particular, these terms would play a role in the four-wave mixing process which is responsible for phase conjugation and squeezing.

The interaction in Eq. (75) cannot be derived from an effective Hamiltonian like that in Eq. (61). Eq. (75) contains terms with unequal numbers of creation and annihilation operators while Eq. (61) does not. This means that Eq. (61) cannot be used to treat counterpropagating pulses in a nonlinear medium, but that a more complicated Hamiltonian, such as that in Eq. (75), must be used.

6 Conclusion

The justification of the Hamiltonians which are used in the quantum theory of nonlinear optics is an important part of placing this theory on a firmer basis. We have presented a derivation of an effective Hamiltonian for the interaction of light and a medium consisting of two-level atoms which contains a number of the standard nonlinear optical interactions. This is accomplished
by expanding the atomic operators in terms of boson creation and annihila-
tion operators and diagonalizing the part of the Hamiltonian which describes
linear interactions. The result is a theory of interacting polaritons which is
a nonlinear extension of Hopfield’s theory. Because the dispersion relation
for polaritons is different from that of photons, the effects of dispersion are
automatically included in this theory.

The theory presented here provides further justification for the Hamil-
tonians which emerge from the quantized macroscopic theory. In addition,
it can be used directly to provide effective Hamiltonians for more complica-
ted situations than those which have so far been considered. We saw an
example of this in the case of counterpropagating beams. The microscopic
theory serves as a useful counterpart to the quantized macroscopic theory in
providing a description of quantized fields in nonlinear dielectric media.

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Appendix A

Here we examine the ground state of the Hamiltonian in the lowest order of the semiclassical expansion. The spin operators are replaced by the c-number quantities

\[
S^{(3)}_l \rightarrow s \cos \theta_l \\
S^{(+)}_l \rightarrow se^{i\phi_l} \sin \theta_l \\
S^{(-)}_l \rightarrow se^{-i\phi_l} \sin \theta_l,
\]

giving us the Hamiltonian

\[
H_{sc} = \sum_{l=1}^{N_b} E_0 s \cos \theta_l + \sum_{|\vec{k}|<k_u,\lambda} \omega_k a_{\vec{k}\lambda}^\dagger a_{\vec{k}\lambda} \\
- \sum_{|\vec{k}|<k_u,\lambda} \sum_{l=1}^{N_b} (a_{\vec{k}\lambda}^\dagger e^{i\vec{k}\cdot\vec{r}_l} + a_{\vec{k}\lambda} e^{-i\vec{k}\cdot\vec{r}_l}) D_{l\vec{k}\lambda} \\
+ \frac{e^2 \rho}{2m} \sum_{|\vec{k}|<k_u,\lambda} \frac{1}{2\omega_k} \left[ (-1)^\lambda a_{\vec{k}\lambda}^\dagger a_{-\vec{k}\lambda} + a_{\vec{k}\lambda} a_{-\vec{k}\lambda}^\dagger \right] \\
+ a_{\vec{k}\lambda}^\dagger \left( a_{\vec{k}\lambda} + (-1)^\lambda a_{\vec{k}\lambda}^\dagger \right) ,
\]

where

\[
D_{l\vec{k}\lambda} = \frac{2s}{\sqrt{2\omega_k}} \sin \theta_l \sin \phi_l \mu_\lambda(\vec{k}).
\]
We have assumed that $\mu_\lambda(\vec{k})$ is real which implies that $D_{\vec{k}\lambda}$ is also real. In principle, we should diagonalize this Hamiltonian which would allow us to determine the values of $\theta_\lambda$ and $\phi_\lambda$ as well as the field in the ground state to lowest order. In practice, we shall be able to obtain what we need after completing part of the procedure, and the rest of the diagonalization can be completed when additional terms from $H_0$ and $H^{(1)}_{int}$ are included.

We begin by eliminating the terms linear in the field operators. This can be accomplished by shifting the creation and annihilation operators, i.e. by setting

$$a_{\vec{k}\lambda} = b_{\vec{k}\lambda} + z_{\vec{k}\lambda},$$  

(79)

where $z_{\vec{k}\lambda}$ is a c number, and

$$b_{\vec{k}\lambda} = D_{\vec{k}\lambda}(-z_{\vec{k}\lambda})a_{\vec{k}\lambda}D_{\vec{k}\lambda}(-z_{\vec{k}\lambda})^{-1}. \quad (80)$$

The displacement operator $D_{\vec{k}\lambda}(z)$ is equal to $\exp(z a_{\vec{k}\lambda}^\dagger - z^* a_{\vec{k}\lambda})$. We now substitute Eq. (79) into $H_{sc}$ and collect the terms linear in $b_{\vec{k}\lambda}$ and $b_{\vec{k}\lambda}^\dagger$. These are

$$\sum_{|\vec{k}|<k_a,\lambda} \left\{ \omega_k (z^*_{\vec{k}\lambda} b_{\vec{k}\lambda} + z_{\vec{k}\lambda} b_{\vec{k}\lambda}^\dagger) - L^*_{\vec{k}\lambda} b_{\vec{k}\lambda} - L_{\vec{k}\lambda} b_{\vec{k}\lambda}^\dagger \right\}$$

$$+ \frac{e^2 \rho}{4m\omega_k} \left(-1\right)^\lambda \left(z_{\vec{k}\lambda} b_{-\vec{k}\lambda} + z_{-\vec{k}\lambda} b_{\vec{k}\lambda}^\dagger\right) + 2 \left(z_{\vec{k}\lambda}^* b_{\vec{k}\lambda} + z_{-\vec{k}\lambda}^* b_{-\vec{k}\lambda}^\dagger\right)$$

$$+ \left(-1\right)^\lambda \left(z_{\vec{k}\lambda} b_{-\vec{k}\lambda}^\dagger + z_{-\vec{k}\lambda} b_{\vec{k}\lambda}^\dagger\right); \quad (81)$$
where
\[ L_{\vec{r} \lambda} = \sum_{l=1}^{N_b} e^{i \vec{k} \cdot \vec{r}_l} D_{l \vec{r} \lambda}. \] (82)

By grouping the terms for \( \vec{k} \) and \(-\vec{k}\) together we find that the expression in Eq. (81) will vanish if
\[ \omega_k z_{\vec{k} \lambda} \left( 1 + \frac{e^2 \rho}{2m \omega_k^2 + e^2 \rho} \right) = L_{\vec{k} \lambda} - (-1)^\lambda \frac{e^2 \rho L_{\vec{k} \lambda}^*}{2m \omega_k^2 + e^2 \rho} \] (83)

Making use of the fact that \( L_{-\vec{k} \lambda}^* = (-1)^\lambda L_{\vec{k} \lambda} \) this simplifies to
\[ z_{\vec{k} \lambda} = \frac{m L_{\vec{k} \lambda}^* \omega_k}{m \omega_k^2 + e^2 \rho}. \] (84)

Substitution of this result back into \( H_{sc} \) gives us
\[
H_{sc} = \sum_{l=1}^{N_b} E_0 s \cos \theta_l - \varepsilon_0 + \sum_{\vec{k}<\vec{k}_u, \lambda} \omega_k b_{\vec{k} \lambda}^\dagger b_{\vec{k} \lambda}
+ \frac{e^2 \rho}{2m} \sum_{\vec{k}<\vec{k}_u, \lambda} \frac{1}{2 \omega_k} \left[ (-1)^\lambda b_{\vec{k} \lambda}^\dagger b_{-\vec{k} \lambda} + b_{-\vec{k} \lambda}^\dagger b_{\vec{k} \lambda} \right]
+ \sum_{\vec{k}<\vec{k}_u, \lambda} \varepsilon_0 = \sum_{\vec{k}<\vec{k}_u, \lambda} \frac{m|L_{\vec{k} \lambda}|^2 \omega_k}{m \omega_k^2 + e^2 \rho}. \] (85)

where
\[
\varepsilon_0 = \sum_{\vec{k}<\vec{k}_u, \lambda} \frac{m|L_{\vec{k} \lambda}|^2 \omega_k}{m \omega_k^2 + e^2 \rho}. \] (86)

We can find the semiclassical ground state spin configuration by minimizing the sum of the first two terms in Eq. (85) with respect to \( \phi_l \) and \( \theta_l \).

That is, we minimize the expression
\[
\sum_{l=1}^{N_b} E_0 s \cos \theta_l - \varepsilon_0. \] (87)
We can obtain an estimate of what the minimum configuration will be by examining the size of $\varepsilon_0$. If $\varepsilon_0 << N_B s E_0$, then the first term will be the dominant one, and the minimum will occur approximately at $\theta_l = \pi$, for all $l$. This we shall, in fact, find to be the case.

In order to estimate $\varepsilon_0$ we first express it in the form

$$\varepsilon_0 = \frac{2(s e E_0)^2}{V} \sum_{\vec{k} < k_u, \lambda} \frac{m|\langle a|\vec{x}|b\rangle \cdot \hat{e}_\lambda(\vec{k})|^2 |c_{\vec{k}}|^2}{m \omega_{\vec{k}}^2 + e^2 \rho}, \quad (88)$$

where

$$c_{\vec{k}} = \sum_{l=1}^{N_b} e^{i\vec{k} \cdot \vec{r}_l} \sin \theta_l \cos \phi_l. \quad (89)$$

Taking $k_u = 2\pi/(\Delta V)^{1/3}$ as the momentum cut off we find that

$$\sum_{\vec{k} < k_u} |c_{\vec{k}}|^2 = N_b \sum_{l=1}^{N_b} \sin^2 \theta_l \cos^2 \phi_l \leq N_b^2. \quad (90)$$

This implies that

$$\varepsilon_0 \leq N E_0^2 |\langle a|\vec{x}|b\rangle|^2 m. \quad (91)$$

Dividing this by $E_0 N_b s$, setting the matrix element equal to the Bohr radius, and taking $E_0$ to be the energy of a 500 nm wavelength photon we have that

$$\frac{\varepsilon_0}{E_0 N_b s} \sim 5 \times 10^{-3}. \quad (92)$$

Therefore, the minimum should be achieved when $\theta_l \simeq \pi$. 

40
We can, in fact, do better. Let us set \( \theta_l = \pi - \delta \theta_l \) and expand the expression in Eq. (87) in \( \delta \theta_l \). We then have that

\[
\sum_{l=1}^{N_b} E_0 s \cos \theta_l \approx -N_b E_0 s + \frac{E_0 s}{2} \sum_{l=1}^{N_b} (\delta \theta_l)^2,
\]

(93)

and

\[
\varepsilon_0 \leq E_0^2 s m |\langle a | \vec{r} | b \rangle|^2 \sum_{l=1}^{N_b} (\delta \theta_l)^2 \\
\leq E_0 s (5 \times 10^{-3}) \sum_{l=1}^{N_b} (\delta \theta_l)^2.
\]

(94)

From this equation and inequality it is clear that the minimum occurs at \( \delta \theta_l = 0 \). Any deviation of \( \delta \theta_l \) from zero causes a larger increase in the first term of Eq. (87) than can be compensated for by the decrease in the second. This also implies that the shifts \( z_{\vec{k} \lambda} \) are zero and that the operators \( a_{\vec{k} \lambda} \) and \( b_{\vec{k} \lambda} \) are the same. The lowest order ground state consists of all the spins pointing down and the field in the vacuum state. Given this information the diagonalization of the rest of the Hamiltonian can proceed as in Section 4.

**Appendix B**

The four eigenvalues of the matrix \( GB \) are given by \( \lambda_1 = E_1(\vec{k}) \), \( \lambda_2 = E_2(\vec{k}) \), \( \lambda_3 = -E_1(\vec{k}) \), and \( \lambda_4 = -E_2(\vec{k}) \). The eigenvector of \( GB \) corresponding to
the eigenvalue $\lambda_j$ and satisfying Eq. (47) as a normalization condition is

$$
\eta_j = \begin{pmatrix}
  x_{1j} \\
  x_{2j} \\
  x_{3j} \\
  x_{4j}
\end{pmatrix},
$$

(95)

where

$$
x_{1j} = \frac{|\omega_k + \lambda_j|}{2(\omega_k|\lambda_j|)^{1/2}} \left[ \frac{(E_0^2 - \lambda_j^2)^2}{(E_0^2 - \lambda_j^2)^2 + 4\omega_k E_0 g_k^2} \right]^{1/2}
$$

$$
x_{2j} = -\frac{2i\omega_k g_k^2}{(E_0 - \lambda_j)(\omega_k + \lambda_j)} x_{1j}
$$

$$
x_{3j} = \frac{\omega_k - \lambda_j}{\omega_k + \lambda_j} x_{1j}
$$

$$
x_{4j} = -\frac{E_0 - \lambda_j}{E_0 + \lambda_j} x_{2j}.
$$

(96)

Finally, we can use Eq. (45) to give us that

$$
M_{ij} = G_{ii} G_{jj} \langle \eta_i | c(j) \rangle = G_{ii} G_{jj} x_{ji}^*,
$$

(97)

which, with Eq. (44), implies that

$$
A_{ij} = x_{ij}.
$$

(98)

Let us examine, in particular, the matrix elements which enter into the calculation describing two counterpropagating beams (Section 5). We assume
that the beams both have wave vectors of magnitude \( k_0 \) and are on the \( \alpha \) branch. Defining

\[
d = \left[ \frac{(E_0^2 - E_1^2)^2}{(E_0^2 - E_1^2)^2 + 4\omega_{k_0} E_0 g_k^2} \right]^{1/2},
\]

we have that

\[
A_{11} = \frac{|\omega_{k_0} + E_0|d}{2\sqrt{\omega_{k_0} E_1}}
\]

\[
A_{13} = \frac{|\omega_{k_0} - E_1|d}{2\sqrt{\omega_{k_0} E_1}}
\]

\[
A_{21} = -\frac{ig_{k_0} d}{E_0 - E_1} \sqrt{\frac{\omega_{k_0}}{E_1}}
\]

\[
A_{23} = \frac{ig_{k_0} d}{E_0 + E_1} \sqrt{\frac{\omega_{k_0}}{E_1}},
\]

and \( A_{31} = -A_{13}, A_{33} = -A_{11}, A_{41} = A_{23}, \) and \( A_{43} = A_{21}. \)
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