

Input-output formalism for few-photon transport in one-dimensional nanophotonic waveguides coupled to a qubit

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We extend the input-output formalism of quantum optics to analyze few-photon transport in waveguides with an embedded qubit. We provide explicit analytical derivations for one- and two-photon scattering matrix elements based on operator equations in the Heisenberg picture.

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I. INTRODUCTION

In the context of quantum information technology, including quantum computing devices, understanding the interaction between a few-photon state and a two-level atom plays an important role [1–3]. The photons are a possible candidate for the “flying qubit” that carries the information, and the two-level atom constitutes the “stationary qubit,” where the flying qubits are generated on demand and correlated with each other.

Recently, there has been an increased activity in analyzing the properties of photons propagating in a waveguide coupled to a qubit—a two-level quantum mechanical system. Experimental demonstration of the control of single photons was made in a waveguide coupled to an optical cavity with an atom in its near field [4]. Similar effects were observed in the microwave domain, when low-frequency photons in a transmission line were coupled to a superconducting qubit [5,6], which later was shown to act as a photon amplifier [7].

To theoretically model such systems one needs to consider a continuous set of waveguide modes that are free to propagate in one dimension, either directly coupled to a multilevel system (referred to as an “atom” in the present article), or indirectly coupled through an optical cavity with a discrete set of modes. Photon transport properties are nontrivial in these structures [8–11], which can be tailored to perform logic operations [12] or form a diode [13]. Exact solutions of one- and two-photon scattering have been reported in [9,11].

The most widely used theoretical approach is to treat the set of equations in the Schrödinger picture and apply the Lippmann-Schwinger formalism to calculate the reflection and transmission properties of the single and multiphoton states [10,14–17]. An alternative technique is to use the reduction formulas from field theory to calculate the scattering matrix of the system [18,19]. Time-domain simulations that take the waveguide dispersion into account are also possible, and an interesting radiation trapping mechanism was recently predicted [20].

In this article, we extend the input-output formalism [21,22] of quantum optics—a Heisenberg picture approach originally introduced to analyze the interaction between an atom in a cavity and a continuous set of electromagnetic states outside

of the atom-cavity system—to analyze the transport of few-photon states in a waveguide with an embedded qubit. In the input-output formalism one obtains a nonlinear set of operator equations based on the Hamiltonian of the system. For a coherent- or a squeezed-state input, this formalism has been extensively used to calculate various coherence properties of the output state of light. Here, we show that one can adopt this formalism to obtain exact results regarding one- or two-photon properties. To do so, we establish a relationship between the input-output formalism and the scattering matrix elements of the system. Our approach complements the existing theoretical literature and bridges different analytical techniques.

This article is organized as follows: In Sec. II we introduce the Hamiltonian of the system. In Sec. III we build the link between the scattering theory and the input-output formalism and continue in Sec. IV with the derivation of the one-photon transport properties. In Sec. V we show how to extend the calculations to the two-photon case. In Sec. VI we make observations on correlation function calculations based on coherent-state inputs and end with our conclusions in Sec. VII.

II. HAMILTONIAN

We start by discussing the model Hamiltonian that we will use in this article. As an illustration of the formalism, we consider a two-level atom coupled to a single polarization, single-mode waveguide [9] and treat the transport properties of few-photon states in such a system (Fig. 1). The Hamiltonian \tilde{H} is defined as ($\hbar = 1$)

$$\tilde{H} = \tilde{H}_0 + \tilde{H}_1.$$

Here, \tilde{H}_0 describes a chiral (i.e., one-way) waveguide where photons propagate in only one direction:

$$\tilde{H}_0 = \int_0^\infty d\beta \tilde{\omega}(\beta) \tilde{a}_\beta^\dagger \tilde{a}_\beta,$$

where \tilde{a}_β and \tilde{a}_β^\dagger are the annihilation and creation operators for the photons with wave vector β , respectively. In Appendix A we calculate the reflection and transmission probabilities for photons in a waveguide where the fields propagate in both directions and show that the results are straightforward extensions of the chiral case. The operators obey the commutation

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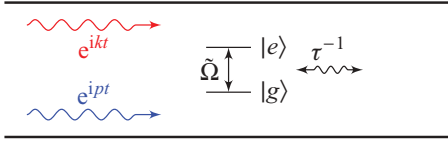


FIG. 1. (Color online) Schematic representation of two photons in a waveguide at frequencies k and p moving to the right toward a two-level atom with energy levels $|g\rangle$ and $|e\rangle$, respectively. $\tilde{\Omega}$ is the separation between the energy levels. Coupling of the two-level atom to the modes in the waveguide is proportional to τ^{-1} . Long horizontal lines denote the waveguide geometry.

relation $[\tilde{a}_\beta, \tilde{a}_{\beta'}^\dagger] = \delta(\beta - \beta')$. \tilde{H}_1 describes the atom as well as the atom-waveguide interaction:

$$\tilde{H}_1 = \frac{1}{2}\tilde{\Omega}\sigma_z + V \int_0^\infty d\beta(\sigma_+ \tilde{a}_\beta + \tilde{a}_\beta^\dagger \sigma_-).$$

Here, $\tilde{\Omega}$ is the atomic transition frequency, σ_\pm are the raising and lowering operators, respectively, for the two-level atom, and $\sigma_z = 2\sigma_+ \sigma_- - 1$. V denotes the coupling strength between the atomic states and the waveguide modes. The derivation of the Hamiltonian is based on the dipole and rotating wave approximations [23] and on taking the continuum limit for field operators. The details of taking the continuum limit are discussed in Appendix B.

It will be useful to have \tilde{H} in terms of the frequency of the photons instead of their wave vector; therefore, we linearize the waveguide dispersion around (β_0, ω_0) as $\tilde{\omega}(\beta) = \omega_0 + v_g(\beta - \beta_0)$ (see Fig. 2). Notice that the total excitation operator

$$N_E = \int_0^\infty d\beta \tilde{a}_\beta^\dagger \tilde{a}_\beta + \frac{1}{2}\sigma_z$$

commutes with \tilde{H} (i.e. $[\tilde{H}, N_E] = 0$). We could thus equivalently solve a system described by

$$H = \tilde{H} - \omega_0 N_E = H_0 + H_1, \quad (1)$$

where

$$H_0 = \int_{-\infty}^\infty d\beta v_g(\beta - \beta_0) \tilde{a}_\beta^\dagger \tilde{a}_\beta,$$

$$H_1 = \frac{1}{2}\Omega\sigma_z + V \int_{-\infty}^\infty d\beta(\sigma_+ \tilde{a}_\beta + \tilde{a}_\beta^\dagger \sigma_-).$$

Here, $\Omega = \tilde{\Omega} - \omega_0$, and we also extended the lower limit of integration to $-\infty$ so that we can define the Fourier transform of operators in the next section. Since we will be dealing with states with wave vectors around β_0 , the extension of the integration limit is well justified [24,25]. Finally, we complete our transition from wave vectors to frequencies by defining $\omega \equiv v_g \beta$ and the operator $a_\omega \equiv \tilde{a}_{\beta+\beta_0}/\sqrt{v_g}$, which satisfies the commutation relation $[a_\omega, a_{\omega'}^\dagger] = \delta(\omega - \omega')$. As a result of all these changes, we have

$$H_0 = \int_{-\infty}^\infty d\omega \omega a_\omega^\dagger a_\omega, \quad (2)$$

$$H_1 = \frac{1}{2}\Omega\sigma_z + \frac{V}{\sqrt{v_g}} \int_{-\infty}^\infty d\omega(\sigma_+ a_\omega + a_\omega^\dagger \sigma_-). \quad (3)$$

Throughout this article, the labels for photon degrees of freedom, for example k and p , refer to photon frequency.

III. CONNECTION BETWEEN SCATTERING THEORY AND INPUT-OUTPUT FORMALISM

In a typical scattering experiment, various input states are prepared and sent toward a scattering region. After the scattering takes place, the outgoing states of the experiment are observed, and information about the interaction is deduced. The mathematical formulation of such a process is commonly made using the scattering matrix with elements of the form

$$S_{p_1 p_2, k_1 k_2} = \langle p_1 p_2 | S | k_1 k_2 \rangle,$$

where $|k_1 k_2\rangle$ denotes the input states—here given as a two-particle state with energies (frequencies) k_1 and k_2 —and $|p_1 p_2\rangle$ denotes the outgoing states. These input and output states are assumed to be free states in the interaction picture and exist long before ($t \rightarrow -\infty$) and long after ($t \rightarrow \infty$) the interaction takes place. The S operator, then, is equal to the evolution operator U_I in the interaction picture from time $-\infty$ to $+\infty$:

$$\begin{aligned} S &= \lim_{\substack{t_0 \rightarrow -\infty \\ t_1 \rightarrow \infty}} U_I(t_1, t_0) \\ &= \lim_{\substack{t_0 \rightarrow -\infty \\ t_1 \rightarrow \infty}} e^{iH_0 t_1} e^{-iH(t_1 - t_0)} e^{-iH_0 t_0}, \end{aligned}$$

where H_0 is the noninteracting part of the Hamiltonian, and $H = H_0 + H_1$ is the total Hamiltonian.¹ In order to have a more compact notation, we will drop the limits and imply $t_0 \rightarrow -\infty$ and $t_1 \rightarrow \infty$.

An equivalent way to describe the scattering is in terms of the scattering eigenstates $|k_1 k_2^\pm\rangle$ that evolve in the interaction picture from a free state either in the distant past or the distant future:

$$\begin{aligned} |k_1 k_2^+\rangle &\equiv U_I(0, t_0) |k_1 k_2\rangle = e^{iH t_0} e^{-iH_0 t_0} |k_1 k_2\rangle \equiv \Omega_+ |k_1 k_2\rangle, \\ |k_1 k_2^-\rangle &\equiv U_I(0, t_1) |k_1 k_2\rangle = e^{iH t_1} e^{-iH_0 t_1} |k_1 k_2\rangle \equiv \Omega_- |k_1 k_2\rangle. \end{aligned}$$

The interaction picture time evolution operators that relate scattering and free states are called the Møller wave operators Ω_\pm . The scattering operator can equivalently be written as $S = \Omega_-^\dagger \Omega_+$.² It is also possible to write the scattering matrix elements as

$$\langle p_1 p_2 | S | k_1 k_2 \rangle = \langle p_1 p_2^- | k_1 k_2^+ \rangle.$$

We should note that scattering eigenstates and the free states with the same quantum numbers have the same energies; that is, $H_0 |k_1 k_2\rangle = E_{k_1 k_2} |k_1 k_2\rangle$ and $H |k_1 k_2^\pm\rangle = E_{k_1 k_2} |k_1 k_2^\pm\rangle$ [26].

¹See [26] for more information about stationary scattering theory. Reference [27] provides a historical account of the developments related to the S matrix.

²There is also an alternative definition of the scattering operator $S' = \Omega_+ \Omega_-^\dagger$, which relates the incoming and outgoing scattering eigenstates, $|k^+\rangle = S' |k^-\rangle$, such that $\langle p | S | k \rangle = \langle p^- | k^+ \rangle = \langle p^+ | S' | k^- \rangle = \langle p^+ | S' | k^+ \rangle$. See [28,29] for details.

It is possible to denote the scattering matrix elements by an appropriate definition of input and output operators such that

$$\langle p_1 p_2^- | k_1 k_2^+ \rangle = \langle 0 | a_{\text{out}}(p_1) a_{\text{out}}(p_2) a_{\text{in}}^\dagger(k_1) a_{\text{in}}^\dagger(k_2) | 0 \rangle, \quad (4)$$

where

$$a_{\text{in}}(k) \equiv \Omega_+ a_k \Omega_+^\dagger = e^{iHt_0} e^{-iH_0 t_0} a_k e^{iH_0 t_0} e^{-iHt_0}, \quad (5)$$

$$a_{\text{out}}(k) \equiv \Omega_- a_k \Omega_-^\dagger = e^{iHt_1} e^{-iH_0 t_1} a_k e^{iH_0 t_1} e^{-iHt_1}, \quad (6)$$

have the property of creating input and output scattering eigenstates:

$$\begin{aligned} a_{\text{in}}^\dagger(k) | 0 \rangle &= | k^+ \rangle, \\ a_{\text{out}}^\dagger(p) | 0 \rangle &= | p^- \rangle, \end{aligned}$$

and the commutation relations

$$[a_{\text{in}}(k), a_{\text{in}}^\dagger(p)] = [a_{\text{out}}(k), a_{\text{out}}^\dagger(p)] = \delta(k - p).$$

We now relate the scattering theory, as briefly sketched above, to the input-output formalism [21,22] of quantum optics. To do so, we start by recalling the definition of the input field operator [21]:

$$a_{\text{in}}(t) = \frac{1}{\sqrt{2\pi}} \int dk a_k(t_0) e^{-ik(t-t_0)},$$

where $a_k(t_0) \equiv e^{iHt_0} a_k e^{-iHt_0}$ is an operator in the Heisenberg picture. The relationship between $a_{\text{in}}(t)$ —which is defined in the input-output formalism—and $a_{\text{in}}(k)$ —which is defined above in (5) as a result of the scattering theory—can be determined by noting that

$$\begin{aligned} a_{\text{in}}(t) &= \frac{1}{\sqrt{2\pi}} \int dk e^{iHt_0} a_k e^{-iHt_0} e^{-ik(t-t_0)} \\ &= \frac{1}{\sqrt{2\pi}} \int dk e^{iHt_0} e^{-iH_0 t_0} a_k e^{iH_0 t_0} e^{-iHt_0} e^{-ikt} \\ &= \frac{1}{\sqrt{2\pi}} \int dk a_{\text{in}}(k) e^{-ikt}, \end{aligned} \quad (7)$$

where in the second line we used the fact that $[H_0, a_k] = -ka_k$ to convert the $a_k e^{ik t_0}$ term into $e^{-iH_0 t_0} a_k e^{iH_0 t_0}$. As a result, $a_{\text{in}}(k)$ provides the spectral representation of $a_{\text{in}}(t)$ in the limit $t_0 \rightarrow -\infty$. Similarly, the output field operator in the input-output formalism

$$a_{\text{out}}(t) = \frac{1}{\sqrt{2\pi}} \int dk a_k(t_1) e^{-ik(t-t_1)},$$

is related to $a_{\text{out}}(k)$ in the scattering theory through

$$a_{\text{out}}(t) = \frac{1}{\sqrt{2\pi}} \int dk a_{\text{out}}(k) e^{-ikt}, \quad (8)$$

in the limit $t_1 \rightarrow \infty$. We have thus established a direct connection between the input-output formalism and the scattering theory. We should note that a different set of input and output operators were defined in [30] with an aim to make a connection to correlation functions. In [31], a similar set of input-output operators were defined in order to relate two different quantization schemes in dielectric media. However, the connection we point out here differs from the previous literature.

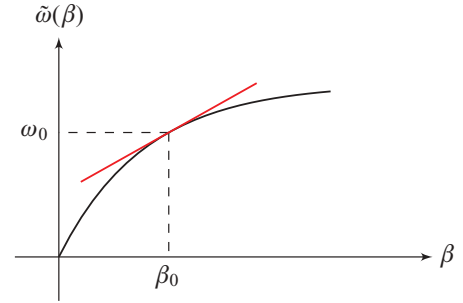


FIG. 2. (Color online) Linearization of a surface-plasmon-like waveguide dispersion relation $\tilde{\omega}(\beta)$ around a wave vector β_0 . The slope of the line is equal to the group velocity v_g . The photon states in the text are assumed to have frequencies in the vicinity of ω_0 so that the linearization is justified.

IV. SINGLE-PHOTON TRANSPORT

Now that we know the relationship between the input-output formalism and the scattering theory, let us now calculate the S -matrix elements $\langle p | S | k \rangle$ between two single-photon states $|k\rangle$ and $|p\rangle$. Following the standard procedure (see Appendix C), the input-output equations appropriate for the Hamiltonian in (1) are

$$\frac{dN}{dt} = -i\sqrt{\frac{2}{\tau}}(\sigma_+ a_{\text{in}} - a_{\text{in}}^\dagger \sigma_-) - \frac{2}{\tau} N, \quad (9)$$

$$\frac{d\sigma_-}{dt} = i\sqrt{\frac{2}{\tau}}\sigma_z a_{\text{in}} - \frac{1}{\tau}\sigma_- - i\Omega\sigma_-, \quad (10)$$

$$a_{\text{out}} = a_{\text{in}} - i\sqrt{\frac{2}{\tau}}\sigma_-, \quad (11)$$

where all operators are in the Heisenberg picture and hence they are all time dependent. The quantity $\tau^{-1} = \pi V^2/v_g$ is proportional to the spontaneous emission rate, and $N = \sigma_+ \sigma_- = (\sigma_z + 1)/2$ describes the probability of having the atom in the excited state.

The single-photon transport properties are described by the single-photon S matrix, which is related to the input and output operator by

$$\begin{aligned} \langle p | S | k \rangle &= \langle 0 | a_{\text{out}}(p) a_{\text{in}}^\dagger(k) | 0 \rangle \\ &= \frac{1}{\sqrt{2\pi}} \int dt \langle 0 | a_{\text{out}}(t) | k^+ \rangle e^{ipt}, \end{aligned}$$

where we used (8) to write $a_{\text{out}}(p)$ in terms of $a_{\text{out}}(t)$. It is therefore sufficient to first calculate $\langle 0 | a_{\text{out}}(t) | k^+ \rangle$ and then perform an inverse Fourier transformation to determine the single-photon S matrix. In the calculations to follow in this and the next section, we will go back and forth between Fourier transforms of the operators, and we will explicitly use t and t' to imply time-dependent operators and $k_{1,2}$ and $p_{1,2}$ to denote the time-independent Fourier-transformed pairs.

The quantity $\langle 0 | a_{\text{out}}(t) | k^+ \rangle$ can be obtained by sandwiching (10) and (12) between the two states $\langle 0 |$ and $| k^+ \rangle$. We have

$$\begin{aligned} \frac{d}{dt} \langle 0 | \sigma_- | k^+ \rangle &= i\sqrt{\frac{2}{\tau}} \langle 0 | \sigma_z a_{\text{in}} | k^+ \rangle - \frac{1}{\tau} \langle 0 | \sigma_- | k^+ \rangle \\ &\quad - i\Omega \langle 0 | \sigma_- | k^+ \rangle, \end{aligned} \quad (12)$$

$$\langle 0|a_{\text{out}}|k^+\rangle = \langle 0|a_{\text{in}}|k^+\rangle - i\sqrt{\frac{2}{\tau}}\langle 0|\sigma_-|k^+\rangle. \quad (13)$$

Note that

$$\langle 0|a_{\text{in}}(t)|k^+\rangle = \langle 0|a_{\text{in}}(t)a_{\text{in}}^\dagger(k)|0\rangle = \frac{1}{\sqrt{2\pi}}e^{-ikt}, \quad (14)$$

by the use of (7), and

$$\langle 0|\sigma_z a_{\text{in}}(t)|k^+\rangle = -\langle 0|a_{\text{in}}(t)|k^+\rangle, \quad (15)$$

since $|0\rangle$ has an atomic part that is in the ground state. Using (14) and (15) in (12) and (13) results in a first-order ordinary differential equation. By solving it, we get

$$\langle 0|\sigma_-|k^+\rangle = \frac{1}{\sqrt{2\pi}}e^{-ikt}\frac{\sqrt{2/\tau}}{(k-\Omega)+i/\tau}, \quad (16)$$

$$\langle 0|a_{\text{out}}|k^+\rangle = \frac{1}{\sqrt{2\pi}}e^{-ikt}\frac{(k-\Omega)-i/\tau}{(k-\Omega)+i/\tau}. \quad (17)$$

After Fourier transforming (17), we obtain the single-photon S matrix:

$$\langle p|S|k\rangle = t_k\delta(k-p), \quad (18)$$

where

$$t_k \equiv \frac{(k-\Omega)-i/\tau}{(k-\Omega)+i/\tau}$$

is the single-photon transmission coefficient. For subsequent calculations, we also define

$$s_k \equiv \frac{\sqrt{2/\tau}}{(k-\Omega)+i/\tau},$$

which measures the excitation of the atom by the single-photon wave when normalized against the incident wave amplitude. The quantities t_k and s_k are related by

$$t_k = 1 - i\sqrt{\frac{2}{\tau}}s_k.$$

These results for single-photon transport agree with [9,11], where the scattering wave function was directly calculated through a real-space formalism.

The crucial step in the derivation above is (15), which takes advantage of the single-excitation nature of the input state. Formally, the same result can also be obtained by approximately setting $\sigma_z = -1$ in (10), and thus linearizing the operator equation. Such a procedure has been commonly adopted in many quantum optics calculations [32–34]. Typically, such an approximation is justified by assuming a so-called weak excitation limit, where the atom is assumed to be mostly in the ground state. Physically, in the case of single-photon transport, the weak-excitation limit is valid when a single-photon pulse has a duration that is much longer than the spontaneous lifetime of the atom. However, we emphasize that the weak-excitation limit is not always valid in general, even for a single-photon pulse. It has been shown that, for the Hamiltonian in (1), a single-photon pulse with a duration comparable to the spontaneous emission lifetime can in fact completely invert an atom [35].

The formalism here removes the need for the assumption of the weak-excitation limit when calculating single-photon properties. In fact, we can directly calculate the excitation probability $\langle k^+|N|k^+\rangle$ for the scattering eigenstate $|k^+\rangle$. With $N = \sigma_+\sigma_-$ and using (16) we have

$$\begin{aligned} \langle k^+|N|k^+\rangle &= \langle k^+|\sigma_+\sigma_-|k^+\rangle = \langle k^+|\sigma_+|0\rangle\langle 0|\sigma_-|k^+\rangle \\ &= \frac{1}{2\pi}|s_k|^2 = \frac{1}{2\pi}\frac{2/\tau}{(k-\Omega)^2+(1/\tau)^2}. \end{aligned}$$

Here, we again have taken advantage of the fact that $|k^+\rangle$ is a single-excitation state, whereas σ_+ acting on any state except $|0\rangle$ would result in a multi-excitation state leading to a zero overlap with $\langle k^+|$. More generally, we have

$$\begin{aligned} \langle k^+|\sigma_z(t)|p^+\rangle &= \langle k^+|(2\sigma_+\sigma_- - 1)|p^+\rangle \\ &= 2\langle k^+|\sigma_+|0\rangle\langle 0|\sigma_-|p^+\rangle - \delta(k-p) \\ &= \frac{1}{\pi}e^{-i(p-k)t}s_k^*s_p - \delta(k-p), \end{aligned} \quad (19)$$

which will be useful when deriving the two-photon S matrix.

V. TWO-PHOTON TRANSPORT

Our aim in this section is to calculate the two-photon S matrix based on the results we obtained for the single-photon case. We first introduced the two-photon S -matrix element in (4). We will begin by inserting an identity operator in between $a_{\text{out}}(p_1)$ and $a_{\text{out}}(p_2)$:

$$\begin{aligned} \langle 0|a_{\text{out}}(p_1)a_{\text{out}}(p_2)a_{\text{in}}^\dagger(k_1)a_{\text{in}}^\dagger(k_2)|0\rangle \\ = \langle 0|a_{\text{out}}(p_1)\left(\int dk|k^+\rangle\langle k^+|\right)a_{\text{out}}(p_2)a_{\text{in}}^\dagger(k_1)a_{\text{in}}^\dagger(k_2)|0\rangle, \end{aligned}$$

and use the Fourier transform of (17) to simplify the expression as

$$= t_{p_1}\langle p_1^+|a_{\text{out}}(p_2)a_{\text{in}}^\dagger(k_1)a_{\text{in}}^\dagger(k_2)|0\rangle.$$

Using the Fourier transform of (11) we get

$$\begin{aligned} &= t_{p_1}\langle p_1^+|\left[a_{\text{in}}(p_2) - i\sqrt{\frac{2}{\tau}}\sigma_-(p_2)\right]a_{\text{in}}^\dagger(k_1)a_{\text{in}}^\dagger(k_2)|0\rangle \\ &= t_{p_1}\delta(p_1 - k_1)\delta(p_2 - k_2) + t_{p_1}\delta(p_1 - k_2)\delta(p_2 - k_1) \\ &\quad - i\sqrt{\frac{2}{\tau}}t_{p_1}\langle p_1^+|\sigma_-(p_2)|k_1k_2^+\rangle, \end{aligned}$$

where we used the orthogonality of the scattering eigenstates. Thus, to determine the two-photon S matrix, we will need to calculate $\langle p_1^+|\sigma_-(t)|k_1k_2^+\rangle$ and take its Fourier transform.

Using (10), we obtain the differential equation that describes $\langle p_1^+|\sigma_-(t)|k_1k_2^+\rangle$:

$$\begin{aligned} \frac{d}{dt}\langle p_1^+|\sigma_-(t)|k_1k_2^+\rangle &= i\sqrt{\frac{2}{\tau}}\langle p_1^+|\sigma_z(t)a_{\text{in}}(t)|k_1k_2^+\rangle \\ &\quad - \left(\frac{1}{\tau} + i\Omega\right)\langle p_1^+|\sigma_-(t)|k_1k_2^+\rangle. \end{aligned} \quad (20)$$

If we can simplify the part that depends on $\sigma_z a_{\text{in}}$, we can then solve the differential equation. Since a_{in} is an annihilation operator for scattering states, by using (7) we can write

$$\begin{aligned} & \langle p_1^+ | \sigma_z(t) a_{\text{in}}(t) | k_1 k_2^+ \rangle \\ &= \frac{1}{\sqrt{2\pi}} [\langle p_1^+ | \sigma_z(t) | k_2^+ \rangle e^{-ik_1 t} + \langle p_1^+ | \sigma_z(t) | k_1^+ \rangle e^{-ik_2 t}], \end{aligned}$$

which, using (19), results in

$$\begin{aligned} &= \frac{1}{\sqrt{2\pi}} \frac{1}{\pi} e^{-i(k_1+k_2-p_1)t} s_{p_1}^* (s_{k_1} + s_{k_2}) \\ &\quad - \frac{1}{\sqrt{2\pi}} \delta(k_2 - p_1) e^{-ik_1 t} - \frac{1}{\sqrt{2\pi}} \delta(k_1 - p_1) e^{-ik_2 t}, \end{aligned}$$

which is what we were after. We can now solve the first-order ordinary differential equation (20) in a way very similar to the derivation that led to (18). After some algebra and rearrangement we get

$$\begin{aligned} & \langle p_1^+ | \sigma_-(t) | k_1 k_2^+ \rangle \\ &= -\frac{1}{\sqrt{2\pi}} \frac{1}{\pi} s_{k_1+k_2-p_1} s_{p_1}^* (s_{k_1} + s_{k_2}) e^{-i(k_1+k_2-p_1)t} \\ &\quad + \frac{1}{\sqrt{2\pi}} \delta(k_2 - p_1) s_{k_1} e^{-ik_1 t} + \frac{1}{\sqrt{2\pi}} \delta(k_1 - p_1) s_{k_2} e^{-ik_2 t}. \end{aligned}$$

Taking the Fourier transform of the expression above gives us

$$\begin{aligned} & \langle p_1^+ | \sigma_-(p_2) | k_1 k_2^+ \rangle \\ &= -\frac{1}{\pi} \delta(k_1 + k_2 - p_1 - p_2) s_{p_2} s_{p_1}^* (s_{k_1} + s_{k_2}) \\ &\quad + s_{k_1} \delta(k_2 - p_1) \delta(k_1 - p_2) + s_{k_2} \delta(k_1 - p_1) \delta(k_2 - p_2). \end{aligned}$$

Lastly, using the relationship $t_{p_1} s_{p_1}^* = s_{p_1}$, we obtain

$$\begin{aligned} & \langle 0 | a_{\text{out}}(p_1) a_{\text{out}}(p_2) a_{\text{in}}^\dagger(k_1) a_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &= t_{k_1} t_{k_2} [\delta(k_2 - p_1) \delta(k_1 - p_2) + \delta(k_1 - p_1) \delta(k_2 - p_2)] \\ &\quad + i \frac{1}{\pi} \sqrt{\frac{2}{\tau}} \delta(k_1 + k_2 - p_1 - p_2) s_{p_1} s_{p_2} (s_{k_1} + s_{k_2}). \quad (21) \end{aligned}$$

This final result agrees with previous calculations using advanced techniques such as the Bethe ansatz³ in real space [10], the algebraic Bethe ansatz [14], and the Lehmann-Symanzik-Zimmermann (LSZ) formalism in quantum field theory [18,19]. The derivation here, however, is perhaps more elementary, and thus may serve to make such results more accessible. In addition, the results relate the presence of the background fluorescence to the excitation of the atoms.

VI. COHERENT-STATE COMPUTATION

A traditional use of the input-output formalism is to calculate the correlation function when the input is in a coherent state. Here we briefly outline such a calculation for our system in order to contrast it with the single-

two-photon calculations of the previous two sections. For this purpose, we consider a coherent input state $|\alpha_k\rangle$, such that

$$a_{\text{in}}(t) |\alpha_k^+\rangle = \alpha e^{-ikt} |\alpha_k^+\rangle,$$

and calculate, as an example, the $G^{(1)}$ correlation function

$$G^{(1)}(t', t) = \frac{\langle \alpha_k^+ | a_{\text{out}}^\dagger(t') a_{\text{out}}(t) | \alpha_k^+ \rangle}{\langle \alpha_k^+ | \alpha_k^+ \rangle}.$$

Using (11), we have

$$\begin{aligned} G^{(1)}(t, t') &= |\alpha|^2 e^{-ik(t-t')} + i\alpha e^{-ikt} \sqrt{\frac{2}{\tau}} \langle \sigma_+(t') \rangle \\ &\quad - i\alpha^* e^{ikt'} \sqrt{\frac{2}{\tau}} \langle \sigma_-(t) \rangle + \frac{2}{\tau} \langle \sigma_+(t') \sigma_-(t) \rangle, \quad (22) \end{aligned}$$

where for any operator O , $\langle O \rangle \equiv \langle \alpha_k^+ | O | \alpha_k^+ \rangle$.

Each of the expectation values in (22) can be calculated using the input-output formalism. Taking the expectation values in (9) and (10) results in

$$\begin{aligned} \frac{d}{dt} \langle \sigma_z(t) \rangle &= -i2\sqrt{\frac{2}{\tau}} [\alpha e^{-ikt} \langle \sigma_+(t) \rangle - \alpha^* e^{ikt} \langle \sigma_-(t) \rangle] \\ &\quad - \frac{2}{\tau} \langle \sigma_z(t) + 1 \rangle, \end{aligned}$$

$$\frac{d}{dt} \langle \sigma_-(t) \rangle = \left(-i\Omega - \frac{1}{\tau} \right) \langle \sigma_-(t) \rangle + i\alpha e^{-ikt} \sqrt{\frac{2}{\tau}} \langle \sigma_z(t) \rangle,$$

$$\frac{d}{dt} \langle \sigma_+(t) \rangle = \left(i\Omega - \frac{1}{\tau} \right) \langle \sigma_+(t) \rangle - i\alpha^* e^{ikt} \sqrt{\frac{2}{\tau}} \langle \sigma_z(t) \rangle.$$

Directly solving the equations above provides the values of $\langle \sigma_+(t') \rangle$ and $\langle \sigma_-(t) \rangle$ in (22), while the $\langle \sigma_+(t') \sigma_-(t) \rangle$ term can be computed using the quantum regression theorem. These calculations can be found in standard textbooks [22,23] in sections related to the properties of resonance fluorescence, and we will not repeat them here. Instead, based on the outline above, we make a few observations about the coherent-state computations, as commonly done, and the one- and two-photon computations as carried out in this article.

(1) The input-output formalism provides a set of nonlinear operator equations. Therefore, all computations, by necessity, involve the conversion of such operator equations into ordinary differential equations for various operator matrix elements. While the coherent-state computations typically involve taking expectation values in terms of the input states, the one- and two-photon computations involve matrix elements that have different photon numbers.

(2) It is certainly reasonable to expect that the one- or two-photon S matrices can be obtained by analyzing various correlation functions for a weak coherent-state input. Indeed, the connection between the two-photon out wave function and the $g^{(2)}$ correlation function has been pointed out in [10], and it is likely that stronger connections exist. This will be carried out in future work. However, if the aim is to determine the S matrix in the few-photon-Fock-state Hilbert space, the computation as discussed here should be far more direct.

(3) We emphasize that the few-photon computations yield the S matrix in the few-photon Hilbert space, and thus provide a complete description of all physical processes in the few-photon-Fock-state Hilbert space. In contrast, computing

³Equations (118)-(119) in [10] and equation (21) in this article are the same with the following notational correspondence: $\Gamma = 2/\tau$, $\Delta_1 = (k_1 - k_2)/2$, $\Delta_2 = (p_1 - p_2)/2$, $E_1 = k_1 + k_2$, $E_2 = p_1 + p_2$.

$G^{(1)}$ or $G^{(2)}$ correlation functions alone do not in general completely specify the out state for a given incident coherent state. Certainly, in the majority of quantum optics experiments at present, one probes a quantum system with a coherent input state and obtains information about the system by measuring different correlation functions. The coherent-state computations, as briefly reproduced above, are adequate to describe these experiments. However, these quantum systems are beginning to be considered as prospective devices which will eventually process quantum states [36,37]. In such an engineering context, one ultimately has to be able to completely specify the output quantum states. It is in this respect that we hope the few-photon transport computations will prove to be valuable for future engineering applications.

VII. CONCLUSION

In this article, we extend the input-output formalism of quantum optics to analyze one- and two-photon scattering in waveguides with a two-level atom inside. We develop the relationship between the input-output operators and the scattering theory, which in turn enables us to analytically calculate the photon scattering matrix elements with a minimum amount of algebra. We also contrast our calculations for few-photon Fock-state transport with the conventional application of input-output formalism for coherent-state transport. This work helps us go beyond the correlation function analysis in input-output formalism and leads to exact solutions for the scattering matrix elements.

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APPENDIX A: TWO-MODE MODEL

In this section we write the Hamiltonian for the case when photons are allowed to propagate in both directions within the waveguide. We refer to this case as the two-mode model. After introducing the Hamiltonian, we use the results of Secs. IV and V to calculate one- and two-photon reflection and transmission coefficients for right-to-left moving fields.

When photons propagate to the right (r) and to the left (ℓ), we will need to add extra terms to the Hamiltonian. We begin as we did in Sec. II and write

$$\tilde{H}_0 = \int_0^\infty d\beta \omega_r(\beta) r_\beta^\dagger r_\beta + \int_{-\infty}^0 d\beta \omega_\ell(\beta) \ell_\beta^\dagger \ell_\beta$$

for the waveguide part of the Hamiltonian. The dispersion relation for the left-moving modes $\omega_\ell(\beta)$ is the mirror image of the one for the right-moving modes. We linearize the left and right branches of the dispersion relationship at $\beta = \pm\beta_0$ to get $\omega_r \approx \omega_0 + v_g(\beta - \beta_0)$ and $\omega_\ell \approx \omega_0 - v_g(\beta + \beta_0)$. Following linearization, we extend the limits of integration to $\pm\infty$, make a change of variables $\beta \mapsto \beta \mp \beta_0$ for the right and left waveguides, respectively, and define $\omega = v_g\beta$, $r_\omega \equiv r_{\beta+\beta_0}/\sqrt{v_g}$, and $\ell_\omega \equiv \ell_{\beta-\beta_0}/\sqrt{v_g}$ to get

$$H_0 = \int_{-\infty}^\infty d\omega \omega (r_\omega^\dagger r_\omega - \ell_\omega^\dagger \ell_\omega). \quad (\text{A1})$$

The interaction part of the Hamiltonian is given by

$$H_1 = \frac{1}{2} \Omega \sigma_z + \frac{V}{\sqrt{v_g}} \int_{-\infty}^\infty d\omega [\sigma_+ (r_\omega + \ell_\omega) + (r_\omega^\dagger + \ell_\omega^\dagger) \sigma_-]. \quad (\text{A2})$$

Since the total excitation operator

$$N_E = \int_0^\infty d\beta r_\beta^\dagger r_\beta + \int_{-\infty}^0 d\beta \ell_\beta^\dagger \ell_\beta + \frac{1}{2} \sigma_z$$

commutes with the Hamiltonian, we subtracted the term $\omega_0 N_E$ from the Hamiltonian and set $\Omega = \tilde{\Omega} - \omega_0$ in the derivation, mimicking the steps in Sec. II.

Now that we have the Hamiltonian, we can write down the Heisenberg equations of motion and define the input-output operators for the fields as illustrated in detail for a chiral model in Appendix C. The equations for the annihilation operators are

$$\begin{aligned} \frac{dr_\omega(t)}{dt} &= -i[r_\omega, H] = -i\omega r_\omega - i\tilde{V}\sigma_-, \\ \frac{d\ell_\omega(t)}{dt} &= -i[\ell_\omega, H] = +i\omega \ell_\omega - i\tilde{V}\sigma_-, \end{aligned}$$

where $\tilde{V} = V/\sqrt{v_g}$. The definitions for the input and output operators for right-going fields are the same as in Appendix C, and we get

$$r_{\text{out}}(t) = r_{\text{in}}(t) - i\sqrt{\frac{2}{\tau}}\sigma_-(t),$$

where τ is defined in (C6). Left-going modes have a group velocity which is the negative of the right-going modes and that leads to a negative sign in (A1). As a result, starting from the definition of the input and output operators in (5)–(6), the input and output operators for left-going modes have the form

$$\begin{aligned} \ell_{\text{out}}(t) &= \frac{1}{\sqrt{2\pi}} \int d\omega \ell_\omega(t_1) e^{i\omega(t-t_1)}, \\ \ell_{\text{in}}(t) &= \frac{1}{\sqrt{2\pi}} \int d\omega \ell_\omega(t_0) e^{i\omega(t-t_0)}, \end{aligned}$$

$$\ell_{\text{out}}(t) = \ell_{\text{in}}(t) - i\sqrt{\frac{2}{\tau}}\sigma_-(t),$$

where we note the change of sign in the frequency variable. Using these results we can show that

$$\frac{d\sigma_-}{dt} = i\sqrt{\frac{2}{\tau}}\sigma_z r_{\text{in}} + i\sqrt{\frac{2}{\tau}}\sigma_z \ell_{\text{in}} - \frac{2}{\tau}\sigma_- - i\Omega\sigma_-,$$

which is in a form similar to those that we get in temporal coupled-mode theory [38].

We now have all the tools to solve for the scattering that takes place in the two-mode model. Let us define even and odd combinations of the operators for the right- and left-propagating modes as

$$a_\omega = \frac{r_\omega + \ell_{-\omega}}{\sqrt{2}} \text{ (even)}, \quad \hat{a}_\omega = \frac{r_\omega - \ell_{-\omega}}{\sqrt{2}} \text{ (odd)}. \quad (\text{A3})$$

Using these definitions in (A1)–(A2), we can show

$$H_0 = \int d\omega a_\omega^\dagger a_\omega + \hat{a}_\omega^\dagger \hat{a}_\omega \equiv H_{e,0} + H_{o,0},$$

$$H_1 = \frac{1}{2} \Omega \sigma_z + \frac{\sqrt{2}V}{\sqrt{v_g}} \int d\omega (\sigma_+ a_\omega + a_\omega^\dagger \sigma_-) \equiv H_{e,1},$$

where we see that the interaction part of the Hamiltonian depends only on the even combination of modes. In Secs. IV and V we solved for $H = H_{e,0} + H_{e,1}$ for a rescaled value of V . The odd part $H_{o,0}$ is interaction free and hence is also solved. From (A3) we get

$$r_{\text{in/out}}(\omega) = \frac{a_{\text{in/out}}(\omega) + \hat{a}_{\text{in/out}}(\omega)}{\sqrt{2}},$$

$$\ell_{\text{out}}(\omega) = \frac{a_{\text{out}}(-\omega) - \hat{a}_{\text{out}}(-\omega)}{\sqrt{2}},$$
(A4)

where we wrote the Fourier transforms of two-mode input-output operators in terms of the combinations of even and odd fields.

To get the one-photon reflection probability, we look at the scattering matrix element which corresponds to a right-propagating input photon and a left-propagating output photon

$$\begin{aligned} \langle 0 | \ell_{\text{out}}(p) r_{\text{in}}^\dagger(k) | 0 \rangle &= \frac{1}{2} \langle 0 | [a_{\text{out}}(-p) - \hat{a}_{\text{out}}(-p)] [a_{\text{in}}^\dagger(k) + \hat{a}_{\text{in}}^\dagger(k)] | 0 \rangle \\ &= \frac{1}{2} \langle 0 | a_{\text{out}}(-p) a_{\text{in}}^\dagger(k) | 0 \rangle - \frac{1}{2} \langle 0 | \hat{a}_{\text{out}}(-p) \hat{a}_{\text{in}}^\dagger(k) | 0 \rangle \\ &= \frac{1}{2} (t_k - 1) \delta(p + k) \equiv \bar{r}_k \delta(p + k). \end{aligned}$$

Here we used (A4) and (18) to get the one-photon reflection coefficient \bar{r}_k . Similarly, the one-photon transmission coefficient \bar{t}_k is given by

$$\langle 0 | r_{\text{out}}(p) r_{\text{in}}^\dagger(k) | 0 \rangle = \frac{1}{2} (t_k + 1) \delta(p - k) \equiv \bar{t}_k \delta(p - k).$$

Two-photon calculations require adding another input-output pair. For instance, the scattering matrix element associated with one photon scattering to the right, another to the left when two photons initially propagate to the right is given by

$$\begin{aligned} \langle 0 | r_{\text{out}}(p_1) \ell_{\text{out}}(p_2) r_{\text{in}}^\dagger(k_1) r_{\text{in}}^\dagger(k_2) | 0 \rangle &= \frac{1}{4} [\langle 0 | a_{\text{out}}(p_1) a_{\text{out}}(-p_2) a_{\text{in}}^\dagger(k_1) a_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &\quad - \langle 0 | a_{\text{out}}(p_1) \hat{a}_{\text{out}}(-p_2) a_{\text{in}}^\dagger(k_1) \hat{a}_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &\quad - \langle 0 | a_{\text{out}}(p_1) \hat{a}_{\text{out}}(-p_2) \hat{a}_{\text{in}}^\dagger(k_1) a_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &\quad + \langle 0 | \hat{a}_{\text{out}}(p_1) a_{\text{out}}(-p_2) a_{\text{in}}^\dagger(k_1) \hat{a}_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &\quad + \langle 0 | \hat{a}_{\text{out}}(p_1) a_{\text{out}}(-p_2) \hat{a}_{\text{in}}^\dagger(k_1) a_{\text{in}}^\dagger(k_2) | 0 \rangle \\ &\quad - \langle 0 | \hat{a}_{\text{out}}(p_1) \hat{a}_{\text{out}}(-p_2) \hat{a}_{\text{in}}^\dagger(k_1) \hat{a}_{\text{in}}^\dagger(k_2) | 0 \rangle] \\ &= \bar{t}_{k_1} \bar{r}_{k_2} \delta(k_1 - p_1) \delta(k_2 + p_2) + \bar{r}_{k_1} \bar{t}_{k_2} \delta(k_1 + p_2) \\ &\quad \times \delta(k_2 - p_1) + \frac{1}{4} B \delta(k_1 + k_2 - p_1 + p_2), \end{aligned}$$

where from (21)

$$B = i \frac{1}{\pi} \sqrt{\frac{2}{\tau'}} s_{p_1} s_{-p_2} (s_{k_1} + s_{k_2}).$$

We note that $\tau' = \tau/2$ due to an extra factor of $\sqrt{2}$ before V in the definition of H_1 . These results agree with equations (52) and (130) in [10].

APPENDIX B: HAMILTONIAN IN THE CONTINUUM LIMIT

This section will summarize the steps taken to obtain the continuum form of the Hamiltonian from its discrete version. We will follow the approach in [24,25].

The discrete variables are assumed to be for those in a one-dimensional (1D) cavity of length L . The mode spacing in the cavity is given by $\Delta\beta = 2\pi/L$. In this 1D cavity, the free-space electromagnetic Hamiltonian H_0 is given by

$$H_0 = \sum_{\beta} \omega_{\beta} \hat{a}_{\beta}^{\dagger} \hat{a}_{\beta},$$

with the commutator relationship $[\hat{a}_{\beta}, \hat{a}_{\beta'}^{\dagger}] = \delta_{\beta, \beta'}$. Now, we will convert the sum into an integral by the equivalence $(\Delta\beta \sum_{\beta}) \rightarrow (\int d\beta)$ to get

$$H_0 = \frac{L}{2\pi} \int d\beta \omega_{\beta} \hat{a}_{\beta}^{\dagger} \hat{a}_{\beta}.$$

The continuous mode operator \tilde{a}_{β} is related to the discrete mode \hat{a}_{β} by

$$\tilde{a}_{\beta} = \sqrt{\frac{L}{2\pi}} \hat{a}_{\beta},$$

which results in

$$H_0 = \int d\beta \omega_{\beta} \tilde{a}_{\beta}^{\dagger} \tilde{a}_{\beta}.$$

The commutator relationship $[\tilde{a}_{\beta}, \tilde{a}_{\beta'}^{\dagger}] = \frac{L}{2\pi} \delta_{\beta, \beta'}$ in the limit $L \rightarrow \infty$ becomes

$$[\tilde{a}_{\beta}, \tilde{a}_{\beta'}^{\dagger}] = \delta(\beta - \beta').$$

To see this result, define $f(\beta) = \frac{L}{2\pi} \delta_{\beta, 0}$. Integrating $f(\beta)$ will give

$$\int d\beta f(\beta) \rightarrow \frac{2\pi}{L} \sum_{\beta} f(\beta) = \frac{2\pi}{L} \frac{L}{2\pi} = 1.$$

As a result, the correct Hamiltonian in the continuum limit is

$$H_0 = \int d\beta \omega(\beta) \tilde{a}_{\beta}^{\dagger} \tilde{a}_{\beta},$$

with $[\tilde{a}_{\beta}, \tilde{a}_{\beta'}^{\dagger}] = \delta(\beta - \beta')$. It is then easy to show that

$$\mathbf{1} = \int d\beta |\beta\rangle \langle \beta|,$$

where $|\beta\rangle = \tilde{a}_{\beta}^{\dagger} |0\rangle$, since

$$\langle \gamma | \int d\beta |\beta\rangle \langle \beta | \zeta \rangle = \int d\beta \delta(\gamma - \beta) \delta(\beta - \zeta) = \delta(\gamma - \zeta).$$

In the discrete case

$$H_1 = \frac{1}{2} \tilde{\Omega} \sigma_z + \frac{V'}{\sqrt{L}} \sum_{\beta} (\sigma_+ \hat{a}_{\beta} + \hat{a}_{\beta}^{\dagger} \sigma_-),$$

where V' is the physical coupling constant. The factor $L^{-1/2}$ arises because the photon as created by \hat{a}_β^\dagger has a normalization constant $L^{-1/2}$. In the continuum case we get

$$\begin{aligned} H_1 &= \frac{1}{2}\tilde{\Omega}\sigma_z + \frac{V'}{\sqrt{L}}\frac{L}{2\pi}\sqrt{\frac{2\pi}{L}}\int d\beta(\sigma_+\tilde{a}_\beta + \tilde{a}_\beta^\dagger\sigma_-) \\ &= \frac{1}{2}\tilde{\Omega}\sigma_z + \frac{V'}{\sqrt{2\pi}}\int d\beta(\sigma_+\tilde{a}_\beta + \tilde{a}_\beta^\dagger\sigma_-). \end{aligned}$$

Thus, the coupling constants in the discrete (V') and the continuum (V) cases differ by a factor of $(2\pi)^{-1/2}$.

APPENDIX C: DERIVATION OF THE INPUT-OUTPUT FORMALISM

Here we provide a derivation of the input-output equations (9)–(11). This derivation closely follows [21,22]. Based on the Hamiltonian (2)–(3), and the definition $\tilde{V} \equiv V/\sqrt{v_g}$, the Heisenberg equations of motion for the operators are

$$i\frac{da_k}{dt} = ka_k + \tilde{V}\sigma_-, \quad (\text{C1})$$

$$i\frac{d\sigma_-}{dt} = \Omega\sigma_- - \tilde{V}\int dk\sigma_z a_k, \quad (\text{C2})$$

$$i\frac{d\sigma_z}{dt} = 2\tilde{V}\int dk(-a_k^\dagger\sigma_- + \sigma_+a_k). \quad (\text{C3})$$

After multiplying (C1) by the integration factor $\exp(ikt)$, we integrate it from an initial time $t_0 < t$ to get

$$a_k(t) = a_k(t_0)e^{-ik(t-t_0)} - i\tilde{V}\int_{t_0}^t dt'\sigma_-(t')e^{-ik(t-t')}. \quad (\text{C4})$$

We define the input operator as

$$a_{\text{in}}(t) = \frac{1}{\sqrt{2\pi}}\int dk a_k(t_0)e^{-ik(t-t_0)},$$

which satisfies the commutation relation

$$[a_{\text{in}}(t), a_{\text{in}}^\dagger(t')] = \delta(t - t').$$

We further introduce a field operator

$$\Phi(t) = \frac{1}{\sqrt{2\pi}}\int dk a_k(t)$$

and integrate (C4) with respect to k to get

$$\Phi(t) = a_{\text{in}}(t) - i\frac{\tilde{V}}{2}\sqrt{2\pi}\sigma_-(t) = a_{\text{in}}(t) - i\sqrt{\frac{1}{2\tau}}\sigma_-(t). \quad (\text{C5})$$

Here, notice that we integrate over half the delta function [21], which results in a factor of $\frac{1}{2}$ and τ is defined as

$$\frac{1}{\tau} \equiv \pi\tilde{V}^2. \quad (\text{C6})$$

Furthermore, plugging (C5) into (C2) and (C3) results in

$$\begin{aligned} \frac{d\sigma_-}{dt} &= i\sqrt{\frac{2}{\tau}}\sigma_z a_{\text{in}} - \frac{1}{\tau}\sigma_- - i\Omega\sigma_-, \\ \frac{dN}{dt} &= -i\sqrt{\frac{2}{\tau}}(\sigma_+ a_{\text{in}} - a_{\text{in}}^\dagger\sigma_-) - \frac{2}{\tau}N. \end{aligned}$$

Here $N = (\sigma_z + 1)/2$. Thus the spontaneous emission rate is $2/\tau$. We could have also directly calculated dN/dt from $d\sigma_-/dt$, since $N = \sigma_+\sigma_-$.

Similarly, we integrate (C1) up to a final time $t_1 > t$, and define an output operator

$$a_{\text{out}}(t) = \frac{1}{\sqrt{2\pi}}\int dk a_k(t_1)e^{-ik(t-t_1)},$$

which results in

$$\Phi(t) = a_{\text{out}}(t) + i\sqrt{\frac{1}{2\tau}}\sigma_-(t). \quad (\text{C7})$$

Combining (C5) and (C7), we finally obtain

$$a_{\text{out}}(t) = a_{\text{in}}(t) - i\sqrt{\frac{2}{\tau}}\sigma_-(t).$$

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