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Mølmer–Sørensen entangling gate for cavity QED systems

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Abstract
The Mølmer–Sørensen gate is a state-of-the-art entangling gate in ion trap quantum computing where the gate fidelity can exceed 99%. Here we propose an analogous implementation in the setting of cavity QED. The cavity photon mode acts as the bosonic degree of freedom in the gate in contrast to that played by the phonon mode in ion traps. This is made possible by utilising cavity assisted Raman transitions interconnecting the logical qubit states embedded in a four-level energy structure, making the ‘anti-Jaynes–Cummings’ term available under the rotating-wave approximation. We identify practical sources of infidelity and discuss their effects on the gate performance. Our proposal not only demonstrates an alternative entangling gate scheme but also sheds new light on the relationship between ion traps and cavity QED, in the sense that many techniques developed in the former are transferable to the latter through our framework.

Keywords: quantum gate, cavity QED, quantum computing

(Some figures may appear in colour only in the online journal)

1. Introduction

Currently trapped atomic ions are among the most successful platforms for quantum information processing (QIP). A number of quantum algorithms [1–3], entanglement of up to 14 ions [4] and quantum simulation of spin systems [5, 6] have been demonstrated, to name a few. Many of these achievements rely on the realization of high fidelity two-qubit entangling gates known as Mølmer–Sørensen (MS) [7] or the geometric phase gate [8]. These gates exploit a collective phonon mode shared by the ions to mediate a state-dependent force and induce a quantum phase conditioned on the collective atomic states. Notable characteristics of this gate scheme are (1) individual addressing of the ions is not required, (2) the time evolution is cyclic such that the electronic and phonon degrees of freedom become disentangled at certain times and (3) the scheme is insensitive to the ions’ initial motional state [9]. Due to these favorable features, the gate can achieve a fidelity in excess of 99% [10, 11].

On the other hand, cavity QED is a paradigm where stationary quantum emitters (e.g. single atoms) interact with quantized radiation fields. It serves as a versatile platform for studies in quantum optics and for quantum information. Namely cavity QED systems are regarded as a vital building block in the development of quantum networks [12]. In the quantum network architecture, each network node is required to be a quantum register capable of multi-qubit quantum logic operations. Even though the recent experimental progress makes it possible to couple multiple qubits to a single optical cavity [13–16], entangling gate operations within a single cavity QED system have not been demonstrated despite a number of theoretical proposals [17–20].

In this article we propose an alternative implementation of a quantum entangling gate for cavity QED systems, which has a direct correspondence to the MS gate in ion traps. It is well known that trapped ions and cavity QED systems share similar physical compositions, i.e. effective spins coupled with a quantized bosonic mode [21]. The prime difference is that in ion traps both Jaynes–Cummings (JC) and anti-Jaynes–Cummings (AIC) Hamiltonians are naturally available by addressing the red and blue-sideband transitions of the ions, respectively [22], whereas in cavity QED normally only the JC Hamiltonian is available. However, this restriction can...
be lifted by utilizing two cavity-assisted Raman transitions interconnecting qubit states embedded in a four-level energy structure. This was first discovered in conjunction with the realization of the Dicke model [23, 24] and later used for studies of the Rabi model [25], but it has never been discussed in terms of a quantum logic gate as per our knowledge.

Even though our proposal is directly inspired by the MS-gate, there are notable differences from the ion-trap implementation. Firstly cavity QED systems are essentially a single-mode system as opposed to the inherent multiple mechanical modes in a string of ions. Therefore, our scheme is free from the issues in ion traps such as off-resonant excitations to irrelevant mechanical modes and spectral congestion in the mode structure with the increasing number of ions. Along the same line, there is no Lamb–Dicke parameter in our scheme, which means there is no compromise between the spatial localization of the qubits and the gate speed. Finally, the optical mode of a cavity can be regarded as being at zero temperature without the need of cooling. Hence, our scheme does not suffer from the heating of the bosonic mode as it is often problematic in ion trap QIPs. However, optical cavities normally have non-negligible field decay rates.

In the following we refer to the individual stationary qubits in the cavity as ‘atoms’ for the sake of convenience. However, in an actual implementation they do not need to be single atomic particles. Indeed they could be e.g. molecules, nitrogen-vacancy centers in diamond or artificial atoms such as semiconductor quantum dots, as long as they have the required energy structure and transitions addressable by a cavity and external laser fields (see figure 1). Therefore, we expect that our proposal is relevant to a broad class of physical systems where direct interaction between the qubits is difficult to attain, but they can be indirectly coupled to each other via an optical cavity field.

The article is structured as follows: in section 2 we introduce the MS Hamiltonian. Section 3 is devoted to the discussion of the cavity-induced Stark shift. The influence of the decay channels on the gate performance is presented in section 4. Finally, in section 4 we summarize our main conclusions.

2. Derivation of the Hamiltonian

We consider an ensemble of $N$ atoms coupled to a single cavity mode. All the atoms possess an identical four-level energy structure, as shown in figure 1. The ground states $|g\rangle$ and $|e\rangle$ form a qubit whereas the excited states $|r_1\rangle$ and $|r_2\rangle$ mediate the coupling between the qubit states via cavity-assisted Raman transitions. The cavity frequency $\omega_c$ is near resonant with the transitions $|g\rangle \leftrightarrow |r_1\rangle$ and $|e\rangle \leftrightarrow |r_2\rangle$ with corresponding detunings $\Delta_{C_1}$ and $\Delta_{C_2}$. We assume that the coupling to the cavity mode is uniform among the atoms and at the same strength of a vacuum Rabi frequency $2g$ for both $|g\rangle \leftrightarrow |r_1\rangle$ and $|e\rangle \leftrightarrow |r_2\rangle$. (The latter condition is not essential and can be relaxed.) In addition, the atoms are externally driven by two laser fields. These two lasers off-resonantly drive the transitions $|e\rangle \leftrightarrow |r_1\rangle$ and $|g\rangle \leftrightarrow |r_2\rangle$ with Rabi frequencies $\Omega_1$ and $\Omega_2$, and detunings $\Delta_{L_1}$ and $\Delta_{L_2}$, respectively. In this way a pair of Raman transitions are constructed to couple the qubit states. Each of them has a cavity-induced transition on one arm and a laser-induced transition on the other. The two-photon detunings $\delta_1$ and $\delta_2$ are given by

$$\delta_1 = \Delta_{C_1} - \Delta_{L_1},$$

$$\delta_2 = \Delta_{C_2} - \Delta_{L_2}.$$ (2)

The Hamiltonian for the total atom–cavity system is composed of the bare energy $H_0$ and the interaction Hamiltonian $H_i$, i.e. $H = H_0 + H_i$. $H_0$ is (assuming $\hbar = 1$)

$$H_0 = \omega_c a_1^t a_1 + \sum_{i=1}^{N} (\omega_i |g_i^{(0)}\rangle \langle g_i^{(0)}| + \omega_i |e_i^{(0)}\rangle \langle e_i^{(0)}| + \omega_i |r_i^{(0)}\rangle \langle r_i^{(0)}| + \omega_i |s_i^{(0)}\rangle \langle s_i^{(0)}|).$$ (3)

Here, $a$ is the annihilation operator of the cavity photon, $\omega_c$ and $|\xi^{(0)}\rangle (\xi = g, e, r_1, r_2)$ are the energy of the atomic level and the vector to the corresponding eigenstate of the $i$th atom, respectively. On the other hand, $H_i$ is given by a sum of the interaction Hamiltonians for the individual atoms:

$$H_i = \sum_{j=1}^{N} \left[ \frac{\Omega_1}{2} e^{-i\omega_1 t} |r_i^{(0)}\rangle \langle e^{(0)}| + \text{H.c.} \right]$$

$$+ \frac{\Omega_2}{2} e^{-i\omega_2 t} |s_i^{(0)}\rangle \langle r_i^{(0)}| + \text{H.c.}$$

$$+ g(a |r_i^{(0)}\rangle \langle g^{(0)}| + a^\dagger |s_i^{(0)}\rangle \langle e^{(0)}| + \text{H.c.}).$$ (4)

Here, $\omega_1$ and $\omega_2$ are the optical frequencies of the driving lasers and H.c. indicates the Hermitian conjugate of the
preceding term inside the same bracket. In the interaction picture, with respect to the operator,  
\[ H_I = \omega_i a^\dagger a + \sum_{i=1}^{N} (\omega_{g_i} |g^{(i)}\rangle \langle g^{(i)}|) \]

\[ + \omega_i |e^{(i)}\rangle \langle e^{(i)}| + (\omega_{r_1} + \Delta_1)|r_1^{(i)}\rangle \langle r_1^{(i)}| \]

\[ + (\omega_{r_2} + \Delta_2)|r_2^{(i)}\rangle \langle r_2^{(i)}|, \]

(5)

with  
\[ \Delta_1 = \frac{\Delta_G + \Delta_L}{2}, \]

\[ \Delta_2 = \frac{\Delta_G + \Delta_L}{2}, \]

the new Hamiltonian,  
\[ H' = H_I H U_t^\dagger + iU_t\frac{dh}{dt}, \]

(6)

results in  
\[ H'_{eff} = U_2 U_{eff} U_t^\dagger - H_2 \]

\[ = \sum_{i=1}^{N} \left[ \frac{\omega_i}{2} \left( \frac{1}{\Delta_1} + \frac{1}{\Delta_2} \right) a^\dagger a \langle |e^{(i)}\rangle \langle e^{(i)}| \right. \]

\[ - |g^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_i}{2\Delta_1} (e^{i\delta_i} a |e^{(i)}\rangle) \langle e^{(i)}| + \text{H.c.} \]

\[ + g \langle e^{i\delta_i} a |r_1^{(i)}\rangle \langle g^{(i)}| \]

\[ + e^{-i\delta_i} a^\dagger |r_2^{(i)}\rangle \langle g^{(i)}| + \text{H.c.} \right], \]

(8)

Assuming \( |\Delta_{1,2}| \gg |\omega_i|, |\Omega_{1,2}|, |\delta_{1,2}| \), the excited states \( |r_1\rangle, |r_2\rangle \) can be adiabatically eliminated and we obtain an effective Hamiltonian:  
\[ H_{eff} = \frac{\omega_i}{2} \left( \frac{1}{\Delta_1} + \frac{1}{\Delta_2} \right) a^\dagger a \]

\[ + \sum_{i=1}^{N} \left\{ \frac{\Omega_i^2}{4\Delta_2} |g^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_i^2}{4\Delta_1} |e^{(i)}\rangle \langle e^{(i)}| \right\} \]

\[ + \frac{\omega_i}{2} \left( \frac{1}{\Delta_1} - \frac{1}{\Delta_2} \right) a^\dagger a \langle |e^{(i)}\rangle \langle e^{(i)}| - |g^{(i)}\rangle \langle g^{(i)}| \rangle \]

\[ + \frac{g\Omega_i}{2\Delta_1} (e^{-i\delta_i} a |e^{(i)}\rangle) \langle g^{(i)}| + \text{H.c.} \]

\[ + \frac{g\Omega_i}{2\Delta_2} (e^{-i\delta_i} a |g^{(i)}\rangle) \langle e^{(i)}| + \text{H.c.} \right\}. \]

(9)

The first two terms inside the sum over the atom index \( i \) correspond to the ac Stark shifts caused by the driving lasers. These terms, in addition to the first term in (9), shifts the bare energy eigenfrequencies with constant offsets. Thus, they can be removed from the equation by moving to another interaction picture with respect to  
\[ H_2 = \frac{\omega_i}{2} \left( \frac{1}{\Delta_1} + \frac{1}{\Delta_2} \right) a^\dagger a \]

\[ + \sum_{i=1}^{N} \left\{ \frac{\Omega_i^2}{4\Delta_2} |g^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_i^2}{4\Delta_1} |e^{(i)}\rangle \langle e^{(i)}| \right\} \]

\[ = \Delta^{(i)} a^\dagger a \]

\[ + \sum_i \left\{ \Delta^{(i)}_2 |g^{(i)}\rangle \langle g^{(i)}| + \Delta^{(i)}_1 |e^{(i)}\rangle \langle e^{(i)}| \right\}, \]

(10)
\( \alpha(t) \) draws a circle with a radius of \( \frac{g_{\text{eff}}}{|\tau|} \) in phase space, as seen in figure 2, and it returns to the origin after every \( \tau = \frac{2\pi}{|\tau|} \).

Therefore, at \( t = n\tau (n = 0, 1, 2, \ldots) \), \( \alpha(t) = 0 \) and \( U_{\text{MS}} \) becomes a propagator involving only the atomic degrees of freedom:

\[
U_{\text{MS}}(t = n\tau) = e^{i\delta S_{I}},
\]  
(28)

where \( \beta_\alpha = 2n\pi \left( \frac{2\pi}{|\tau|} \right)^2 \text{sign}(\delta) \) can be expressed by the area \( A \) enclosed by \( \alpha(t) \) as \( \beta_\alpha = 2n \text{ sign}(\delta) A \), and hence is called geometric phase. By choosing the two-photon detuning such that

\[
|\delta| = 2\sqrt{m} g_{\text{eff}} \quad (m = 1, 2, 3, \ldots),
\]  
(29)

the geometric phase becomes \( \beta_\alpha = \text{sign}(\delta) \frac{\pi}{2} \) at \( t = m\tau \) and \( U_{\text{MS}} \) can be used to generate maximally entangled states [7].

We define the gate time

\[
t_{\text{gate}} = m\tau = \frac{\pi|\delta|}{2g_{\text{eff}}}
\]  
(30)

In particular, when \( N = 2 \), it accomplishes the following transformations of the two-qubit basis states, which are equivalent to the controlled-not gate up to single-qubit rotations:

\[
|\phi_1\rangle \equiv |gg\rangle \rightarrow |\Phi_1^{(\pm)}\rangle = \frac{1}{\sqrt{2}}(|gg\rangle \pm i|ee\rangle),
\]  
(31)

\[
|\phi_2\rangle \equiv |ge\rangle \rightarrow |\Phi_2^{(\pm)}\rangle = \frac{1}{\sqrt{2}}(|ge\rangle \pm i|eg\rangle),
\]  
(32)

\[
|\phi_3\rangle \equiv |eg\rangle \rightarrow |\Phi_3^{(\pm)}\rangle = \frac{1}{\sqrt{2}}(|eg\rangle \pm i|ge\rangle),
\]  
(33)

\[
|\phi_4\rangle \equiv |ee\rangle \rightarrow |\Phi_4^{(\pm)}\rangle = \frac{1}{\sqrt{2}}(|ee\rangle \pm i|gg\rangle).
\]  
(34)

Here the plus and minus signs correspond to the sign of \( \delta \).

So far we have neglected the first term in (19) which represents the differential ac Stark shift induced by the cavity field:

\[
H_{\text{AS}} = \chi a^\dagger a S_{C}.
\]  
(35)

This Hamiltonian can cause a deviation from the ideal time evolution of \( H_{\text{MS}} \). By setting \( \Delta_1 = \Delta_2 \), in addition to the conditions (17) and (18), which in turn means \( \Omega_1 = \Omega_2 \), \( \chi \) vanishes. However, in general, this additional condition may not be satisfied since it imposes a constraint \( \omega_1 - \omega_2 \approx \omega_{\text{r}} - \omega_{\text{e}} \) on the energy structure if \( |\delta| \ll |\Delta_1|, |\Delta_2| \). When this is not the case, the caused deviation may not be negligible depending on the magnitude of \( \chi \). Another possible deviation could arise from dissipative processes such as the cavity field decay and atomic spontaneous emissions from the excited states.

In the following sections, we treat these two kinds of imperfections—the effect of \( \chi \) and that of dissipative processes—in the case of two atoms.

### 3. Effect of the cavity-induced ac Stark shift

When \( \chi \neq 0 \), inclusion of \( H_{\text{AS}} \) results in deviations from the ideal entangling gate operations (34). Let us denote the propagator of Hamiltonian (19) by \( V(t) \). Then the wave function at time \( t \) is given by

\[
|\Psi(t)\rangle = V(t)|\Psi(0)\rangle,
\]  
(36)

where we assume that the initial state \( |\Psi(0)\rangle = |\phi_1\rangle |n\rangle \) (\( i = 1, 2, 3, 4 \)), that is a tensor product of one of the logical basis states of two qubits (see (34)) and a photon number state \( |n\rangle \) with an arbitrary \( n \). The fidelity of the state \( |\Psi(t)\rangle \), with respect to the ideal gate output is

\[
F_{\text{g,eff}}(t) = \langle \Phi_i^{(\pm)} | \text{tr}_{\text{photon}}(V(t)) |\Phi_i^{(\pm)} \rangle = \langle \Phi_i^{(\pm)} | \text{tr}_{\text{photon}}(V(t)) |\phi_i\rangle \langle \phi_i | \phi_i \rangle \cong |n\rangle \langle n| V^+(t) |\Phi_i^{(\pm)} \rangle.
\]  
(37)

Here \( \text{tr}_{\text{photon}} \) is a partial trace over the photon degree of freedom and \( |\Phi_i^{(\pm)} \rangle \) is one of the atomic states for the ideal gate operation shown in (34), chosen accordingly to the initial state \( |\phi_i\rangle \). The subscripts \( i \) and \( n \) represents the initial atomic state and photon number, respectively.

Since \( H_{\text{AS}} \) is proportional to the photon number operator \( a^\dagger a \), the perturbation caused by this Hamiltonian is expected to increase with the number of photons in the cavity. Figure 3(a) shows time evolutions of \( F_{\text{g,eff}}(t) \), i.e. state fidelity to \( |\Phi_i^{(\pm)} \rangle \) when the initial state is prepared in \( |\phi_i\rangle |n\rangle \). Here we set \( \delta = 2g_{\text{eff}} > 0 \). When \( \chi = 0 \), the fidelity reaches the maximum value of unity at \( t_{\text{eff}} t = \pi \) (black dashed curve) and this behavior does not depend on the initial cavity state. However, when \( \chi \neq 0 \) the fidelity shows strong dependence on the number of photons in the initial cavity state. In particular, we see that there is an acute fidelity drop for cavity initial states \( |n\rangle \) with \( n > 0 \). By preparing the initial cavity state in the vacuum state, the maximum fidelity remains close to unity even though the effect through dynamically generated photons is still present as a small drop of the fidelity and a time shift of the peak. Therefore, in order to obtain the best gate performance for a given \( \chi \), the cavity state has to be prepared close to the vacuum state. However, since thermal excitations are negligible in the optical domain at the room temperature \( (n) \sim 10^{-20} \), the cavity state remains essentially in vacuum unless we intentionally drive the cavity field. Hence, this restriction to the initial cavity state causes little
problem for cavity QED systems in the optical domain. This is in stark contrast with ion traps where active cooling of phonon modes is always required.

In addition to the dependence on the initial cavity states, $H_{\text{AS}}$ also leads to a dependence of the gate performance on the initial atomic states. Figure 3(b) shows the fidelities for different atomic initial states to their target states. As can be seen, the initial states $|\phi_0\rangle$ and $|\phi_4\rangle$ are more sensitive to the perturbation than $|\phi_1\rangle$ and $|\phi_2\rangle$ when the cavity mode is prepared in vacuum.

These dependencies of the gate performance on the initial photonic and atomic states when $\chi \approx 0$ can be illustrated by explicitly considering an approximation of the state fidelity for small $\chi$. In doing so, the main difficulty arises from the fact that $H_{\text{MS}}$ and $H_{\text{AS}}$ do not commute with each other. In [25], the authors presented a steady-state analysis of the same Hamiltonian (19). However, here we are interested in explicit time evolutions of states under the Hamiltonian. We compute them by moving into a suitable interaction picture, defined by the transformation

$$|\Psi(t)\rangle = U_{\text{as}}(t)|\Psi(0)\rangle,$$

where $U_{\text{as}}(t) = \exp(-i\delta \sigma^z t)$. $U_{\text{as}}(t) = \exp(iH_{\text{MS}}t)$, and

$$H_{\text{MS}} = U_{\text{as}}(t)H_{\text{MS}}U_{\text{as}}^\dagger(t) = \delta \sigma^z a + g_{\text{eff}}(a + a^\dagger)S_z.$$

In this picture, the evolution of $|\Psi(t)\rangle$ is given by the time-ordered exponential of the Hamiltonian

$$H_{\text{as}}(t) = e^{iH_{\text{MS}}}H_{\text{AS}}e^{-iH_{\text{MS}}}.$$

Our strategy will be the perturbative expansion of this propagator in powers of $H_{\text{AS}}$. Assuming the initial state $|\phi_i\rangle$ is given in (37) as we did not trace out the photonic degree of freedom in (41). However, as long as the additional excitation of photons due to $H_{\text{AS}}$ is small for small $\chi$, $|\phi_i\rangle$ is a good approximation of $F_{\text{as}}(t)$. Consequently, the state overlap $\eta_{\text{as}}(t)$ can be expanded in a perturbative series as (cf. A.13):

$$\eta_{\text{as}}(t) = \eta_{\text{as}}^{(0)}(t) + \eta_{\text{as}}^{(1)}(t) + \eta_{\text{as}}^{(2)}(t) + \cdots,$$

with

$$\eta_{\text{as}}^{(0)}(t) = \left\langle \Phi_i | a^\dagger n U_{\text{as}}(t) U_{\text{as}}^\dagger(t) | \phi_i \right\rangle |n\rangle,$$

$$\eta_{\text{as}}^{(1)}(t) = -i \left\langle \Phi_i | a^\dagger n U_{\text{as}}(t) U_{\text{as}}^\dagger(t) \int_0^t H_{\text{as}}(t')dt' | \phi_i \right\rangle |n\rangle,$$

$$\eta_{\text{as}}^{(2)}(t) = -i^2 \left\langle \Phi_i | a^\dagger n U_{\text{as}}(t) U_{\text{as}}^\dagger(t) \int_0^t \int_0^{t'} H_{\text{as}}(t'')dt'' | \phi_i \right\rangle |n\rangle.$$

$$\eta_{\text{as}}^{(k)}(t)$$ is on the order of $\mathcal{O}(\chi^k)$ and $\eta_{\text{as}}^{(0)}(t)$ is the state overlap for the ideal MS-gate without the cavity-induced ac Stark shift. Using the energy eigenstates of $H_{\text{MS}}$ (see A.4)-(A.7) for their specific definition), the initial states $|\phi_i\rangle |n\rangle$ can be written as

$$|\phi_{i,n}\rangle = \frac{1}{\sqrt{2}} \sum_m d_m \chi^m |0, m\rangle + \frac{1}{2} \sum_m d_m (\chi^m - 1, m\rangle.$$

Even though $H_{\text{MS}} \propto a^\dagger a$ alone preserves the number of the cavity photons, due to the fact that $H_{\text{AS}}$ and $H_{\text{MS}}$ do not commute, the commutators between them arising in the propagator $V(t)$ causes a change of the cavity photons in addition to that caused by the MS-gate process.
\[ |\phi_{2,3}\rangle |n\rangle = \frac{1}{2} \sum_{m} d_{mn} (-\alpha) |1, m\rangle \\
\pm \frac{1}{\sqrt{2}} |S = 0, S_z = 0\rangle |n\rangle \\
- \frac{1}{2} \sum_{m} d_{mn} (\alpha) |1, m\rangle, \tag{47} \]

where we have introduced the abbreviated state notation

\[ |j, n\rangle = |S = 1, S_z = j\rangle |n, j\rangle, \quad j = -1, 0, 1 \tag{48} \]

along with the definitions \( d_{mn} (\pm \alpha) = \langle m | D (\pm \alpha) | n \rangle \). The plus and minus signs in front of the second term in (46) differentiate \( |\phi_{2}\rangle \) and \( |\phi_{3}\rangle \), respectively, and similarly \( |\phi_{1}\rangle \) (plus) and \( |\phi_{0}\rangle \) (minus) in (47). Likewise the target states \(|\Phi_{1.4}\rangle |n\rangle\) are

\[ |\Phi_{1.4}\rangle |n\rangle = \frac{e^{i\frac{\pi}{4}}}{2} \sum_{m} d_{mn} (-\alpha) |1, m\rangle \\
\pm \frac{1}{\sqrt{2}} |S = 0, S_z = 0\rangle |n\rangle \\
- \frac{1}{2} \sum_{m} d_{mn} (\alpha) |1, m\rangle. \tag{49} \]

\[ |\Phi_{2.3}\rangle |n\rangle = \frac{e^{i\frac{\pi}{4}}}{2} \sum_{m} d_{mn} (-\alpha) |1, m\rangle \\
\pm \frac{1}{\sqrt{2}} |S = 0, S_z = 0\rangle |n\rangle \\
- \frac{1}{2} \sum_{m} d_{mn} (\alpha) |1, m\rangle. \tag{50} \]

In terms of these energy eigenstates, the ideal MS-gate without the cavity ac Stark shift can be understood solely by the evolution of quantum phases in the interaction picture: during the ideal MS gate operation, each term in the form of the evolution of quantum phases in the interaction picture:

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In order to assess the relative significance of \( \eta_{i,0}^{(2)} \) for \( i = 1, 2, 3, 4 \) with respect to each other, we consider the independent functions on \( \alpha \). Since \( d_{mn} (\alpha) = O (\alpha^{m-n}) \), \( d_{mn} (\alpha) d_{n0} (\alpha) d_{m0} (\alpha) Y_{mn}(t) \), and \( d_{n0} (\alpha) d_{m0} (\alpha) d_{n0} (\alpha) Y_{mn}(t) \), we have

\[ \eta_{1,0}^{(2)} = \eta_{2,0}^{(2)} = \eta_{3,0}^{(2)} = \text{O}(\alpha^{2}), \tag{59} \]

\[ \eta_{1,0}^{(2)} = \eta_{2,0}^{(2)} = \text{O}(\alpha^{2}), \tag{60} \]

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\pm \frac{1}{\sqrt{2}} |S = 0, S_z = 0\rangle |n\rangle \\
- \frac{1}{2} \sum_{m} d_{mn} (\alpha) |1, m\rangle. \tag{50} \]

In order to assess the relative significance of \( \eta_{i,0}^{(2)} \) for \( i = 1, 2, 3, 4 \) with respect to each other, we consider their independent functions on \( \alpha \). Since \( d_{mn} (\alpha) = O (\alpha^{m-n}) \), \( d_{mn} (\alpha) d_{n0} (\alpha) d_{m0} (\alpha) Y_{mn}(t) \), and \( d_{n0} (\alpha) d_{m0} (\alpha) d_{n0} (\alpha) Y_{mn}(t) \), we have

\[ \eta_{1,0}^{(2)} = \eta_{2,0}^{(2)} = \eta_{3,0}^{(2)} = \text{O}(\alpha^{2}), \tag{59} \]

\[ \eta_{1,0}^{(2)} = \eta_{2,0}^{(2)} = \text{O}(\alpha^{2}), \tag{60} \]

The difference between (59) and (60) originates from the factors \( (1 \pm (-1)^i) \) in (54) and (55). Since \( \alpha = g_{00} |< \beta| 0 \rangle \), this difference indicates that the second order correction is smaller for the initial states \(|\phi_{1/2}\rangle |0\rangle\) than \(|\phi_{2/3}\rangle |0\rangle\), which is also consistent with figure 3(b). Note that in figure 3(b) there is a small difference between the initial states \(|\phi_{1/2}\rangle |0\rangle\) (blue) and \(|\phi_{1/2}\rangle |0\rangle\) (red) whereas calculations based on (54) and (56) and even higher order terms in (42) predict that they should exactly coincide. This is due to the partial trace of the photonic degree of freedom carried out in (37) which is missing in the calculation of the state overlap (41).

Figure 4 shows attainable maximal state fidelities as a function of \( \chi \) for different initial states. For the blue and red solid lines we numerically calculated max(\( \langle F_{1,0}(t) \rangle \) and
max\((F_{2,0}(t))\), respectively, by solving the Schrödinger equations where the maximum is taken over time \(t\) for a given \(\chi\). On the other hand the dashed lines are approximated fidelities using (42) up to the second order. That is, \(\max (|\gamma_{2,0}^{(0)}(t) + \gamma_{2,0}^{(2)}(t)|^2)\) for the blue dashed line and \(\max (|\gamma_{1,0}^{(0)}(t) + \gamma_{1,0}^{(2)}(t)|^2)\) for the red dashed line. They reproduce the behaviors of the exact solutions, namely the distinction between different initial states, well up to \(\chi/g_{\text{eff}} \sim 0.2\) with an error in fidelity less than \(10^{-3}\), confirming the validity of our perturbative approach.

As the state fidelities to the target states differ among different initial states in the presence of \(H_{\text{AS}}\), the performance of the gate as a whole is quantified by the fidelity averaged over all possible initial states, called average gate fidelity. According to [27], the average gate fidelity, that we refer to as \(\bar{F}(t)\), is calculated using the following formula:

\[
\bar{F}(t) = \frac{1}{80} \sum_{i,j} \text{tr}(U_{\text{MS}}(t) (\sigma_i \otimes \sigma_j) U_{\text{MS}}^\dagger(t) \mathcal{E}(\sigma_i \otimes \sigma_j)) + 16,
\]

where \(\sigma_{ij}\) is one of the Pauli matrices or identity matrix and the sum runs through all possible combinations of those. \(\mathcal{E}\) is a map between the atomic density operators to represent the time evolution governed by \(H_{\text{eff}}\):

\[
\mathcal{E}(\rho) = \text{tr}_{\text{photons}}(V(t)(\rho \otimes |0\rangle\langle 0|) V^\dagger(t)).
\]

Taking the maximum of (61) in terms of time \(t\), figure 5 shows the attainable average gate fidelities as a function of \(\chi\) for different two-photon detunings \(\delta\). In the limit of \(\delta/g_{\text{eff}} \gg 1\), excitation of photons during the gate operation is increasingly suppressed as the radius of the trajectory \((= g_{\text{eff}}/\delta)\) in phase space reduces to zero. As a consequence, the perturbation by \(H_{\text{AS}}\) proportional to \(a^\dagger a\) becomes negligible such that the average gate fidelity remains close to unity with a large value of \(\delta\) (see the dash-dotted yellow trace in figure 5). However, this general trend does not necessarily apply to relatively small values of \(\delta\) due to their oscillatory nature, as seen in figure 5. For example the average gate fidelity for \(\delta/g_{\text{eff}} = 4\) is smaller than that for \(\delta/g_{\text{eff}} = 2\) up to \(\chi/g_{\text{eff}} \sim 0.8\).

4. Cavity and atomic decay

A deviation from the ideal gate operation could arise from dissipative processes such as the cavity decay and atomic spontaneous emissions. In this section, the effects of these processes are studied while assuming \(\chi = 0\) for brevity.

The optical fields in cavity QED systems suffer inevitable scattering/absorption or transmission losses at the cavity mirrors. In order to incorporate these losses, we use a master equation for the time evolution:

\[
\frac{d\rho}{dt} = -i[H_{\text{eff}}, \rho] + \kappa(2a\rho a^\dagger - a^\dagger a\rho - \rho a^\dagger a),
\]

where \(\rho\) is the density operator for the total system including the atomic and photonic degrees of freedom and \(\kappa\) is the amplitude dissipation rate of the cavity field. Since the field dissipation does not depend on the atomic states, there is no difference between the average gate fidelity and the fidelity for a specific initial atomic state. Figure 6 shows attainable...
maximum fidelities for $\kappa = 0$. As can be seen, the detrimental effect of the field dissipation can be mitigated by increasing the two-photon detuning $\delta$ with expense of increasing $t_{\text{gate}}$. This can be understood as follows: in the regime where the photonic excitation is very small, the probability $P_d$ that a photon loss occurs per unit time is proportional to the mean photon number $\sim O(g_{\text{eff}}^2/\delta^2)$, which decreases quadratically with $\delta$. On the other hand $t_{\text{gate}}$ only scales linearly with $\delta$ (see (30)). Therefore, the probability that a photon loss occurs within the gate time is $P_d t_{\text{gate}} \sim O(1/\delta)$, and hence increasing $\delta$ improves the achievable gate fidelity.

The detunings required for a gate fidelity $\ge 0.99$ are $\delta/g_{\text{eff}} \approx 20, 100$ and $900$ for $\kappa/g_{\text{eff}} = 0.1, 1.0$ and $10$, respectively (see figure 6). Hence, in principle the gate can be implemented with a high fidelity even in the regime where $g_{\text{eff}} < \kappa$. However, since large detunings $\delta$ means long gate times $t_{\text{gate}}$, increasing $\delta$ arbitrarily may not be an option if the system suffers from other decoherence mechanisms, such as the decay of atomic coherence, that cannot be mitigated by increasing $\delta$.

By adiabatically eliminating the excited states, the atomic spontaneous emissions from the excited states $|e\rangle$ and $|g\rangle$ as in the following master equation [28]:

$$\frac{\text{d}\rho}{\text{d}t} = -i[H_{\text{eff}}, \rho] + \sum_i (\mathcal{L}(C^{(i)}_1, \rho) + \mathcal{L}(C^{(i)}_{1e}, \rho) + \mathcal{L}(C^{(i)}_{2g}, \rho) + \mathcal{L}(C^{(i)}_{2e}, \rho)), \quad (64)$$

where

$$\mathcal{L}(C, \rho) \equiv 2C\rho C^\dagger - C^\dagger C\rho - \rho C^\dagger C,$$

$$C^{(i)}_{1g} \equiv \sqrt{\gamma_{1g}} \left( \frac{g}{\Delta_1} a|g\rangle\langle e| + \frac{\Omega_1}{2\Delta_1} |g\rangle\langle e| \right), \quad (65)$$

$$C^{(i)}_{1e} \equiv \sqrt{\gamma_{1e}} \left( \frac{g}{\Delta_1} a|e\rangle\langle e| + \frac{\Omega_1}{2\Delta_1} |e\rangle\langle e| \right), \quad (66)$$

$$C^{(i)}_{2g} \equiv \sqrt{\gamma_{2g}} \left( \frac{g}{\Delta_2} a|g\rangle\langle e| + \frac{\Omega_2}{2\Delta_2} |g\rangle\langle e| \right), \quad (67)$$

$$C^{(i)}_{2e} \equiv \sqrt{\gamma_{2e}} \left( \frac{g}{\Delta_2} a|e\rangle\langle e| + \frac{\Omega_2}{2\Delta_2} |e\rangle\langle e| \right), \quad (68)$$

and $\gamma_{1j}$ ($j = 1, 2, \xi = g, e$) is the spontaneous decay rate associated with the $|r\rangle \rightarrow |\xi\rangle$ transition. These decay rates and their branching ratios (e.g. $\gamma_{1e}/\gamma_{1g}$) differ to a large extent among different physical systems depending on their specific energy structures. Here we first consider the simplest case where $\gamma_{1g} = \gamma_{1e} = \gamma_{2g} = \gamma_{2e}$ ($\equiv \gamma$) holds. In addition we assume that $\Delta_1 = \Delta_2$ ($\equiv \Delta$) and $\Omega_1 = \Omega_2$ ($\equiv \Omega$) hence $\chi = 0$ is satisfied. Figure 7 shows numerical calculations of the average gate fidelity using this setting for different values of $\Delta$ and $\delta$.

If the contributions from the terms proportional to the photon annihilation operator $a$ in the collapse operators (66)-(69) are negligible due to the small photon excitation, the effective spontaneous decay rate $\gamma_{\text{eff}}$ is given by

$$\gamma_{\text{eff}} \approx \frac{\gamma_0^2}{4\Delta}. \quad (70)$$

![Figure 7](attachment:image.png)

**Figure 7.** Average gate fidelity as a function of $\gamma$ normalized by $g$. Here we set $\Omega = g$, $\Delta/g = 100$ (red, upper three lines) and $10$ (blue, lower three lines). $\delta/g_{\text{eff}} = 2$ (solid), $50$ (dashed) and $200$ (dotted).

**Table 1.** Parameter sets used in the MS-gate simulations for $^{87}$Rb atoms. All the numbers are in units of $2\pi$ MHz.

<table>
<thead>
<tr>
<th>$g$</th>
<th>$\kappa$</th>
<th>$\Delta_1$</th>
<th>$\Delta_2$</th>
<th>$\Omega_1$</th>
<th>$\Omega_2$</th>
<th>$\delta$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>60</td>
<td>1.5</td>
<td>10000</td>
<td>3980</td>
<td>−50</td>
<td>−17.6</td>
</tr>
<tr>
<td>2</td>
<td>200</td>
<td>0.1</td>
<td>20000</td>
<td>13977</td>
<td>−50</td>
<td>−33.9</td>
</tr>
</tbody>
</table>

Therefore, the probability that a spontaneous emission occurs during the gate time is

$$P_{\text{spont}} t_{\text{gate}} = \frac{\pi^2 \delta}{8g^2} = \frac{\pi \sqrt{m} \gamma \Omega}{4g\Delta}. \quad (71)$$

This probability needs to be sufficiently small in order for the atomic decay to be negligible for the gate fidelity. In that regard one can see that increasing $\delta$ is rather detrimental than beneficial, which can also be seen in figure 7.

In the general case where the cavity and atomic decays are both simultaneously present in the system, one needs to carefully choose experimental parameters depending on the relative magnitudes of $\kappa$ and $\gamma$ to minimize their total effect on the gate performance. As an example of such practical systems, we now consider single neutral $^{87}$Rb atoms coupled to a high finesse optical cavity. Note that this particular system was theoretically and experimentally studied in [25] and [24], respectively. See appendix B for the details of modeling this system. First we assume a state-of-the-art conventional Fabry–Perot cavity with a length of 50 $\mu$m and finesse of 10$^6$. This results in $g/2\pi = 60$ MHz and $\kappa/2\pi = 1.5$ MHz for a single $^{87}$Rb atom. With the other parameters listed in set 1 of table 1, we obtain $g_{\text{eff}} /2\pi = 123$ kHz and $\gamma /2\pi = −240$ kHz, and the maximum average gate fidelity of $F = 0.844$ at $t = 260 \mu$s (see figure 8). In order to further improve the gate fidelity, one can use an ultra-high Q microsphere/toroidal cavity where a greater atom–cavity coupling is expected. Here we assume $(g, \kappa, \gamma) /2\pi = (200, 0.1) \text{ MHz}$ [25] together with the parameters listed in set 2 of table 1. Then we obtain $g_{\text{eff}} /2\pi = 204$ kHz, $\gamma /2\pi = −331$ kHz and $F = 0.986$ at $t = 98 \mu$s (figure 8).
parameters in table 1. Set 1 (blue solid) and set 2 (red dashed).

5. Conclusion

In this article we have proposed a novel entangling quantum logic gate for general cavity QED systems. This scheme is inspired by the MS gate in ion traps whose well-established physics enables us to have a clear picture of the basic gate dynamics and the effects of the imperfections. We have analysed and evaluated possible adverse effects caused by the cavity-induced ac Stark shift, cavity field decay and atomic spontaneous emissions. Using the $^{87}$Rb system as a practical example, we have demonstrated that a high fidelity gate operation is possible with sufficiently large atom–cavity coupling. However, we believe that our scheme is applicable to a broader class of physical systems, not limited to atomic cavity QED, due to the fact that cavity interactions are ubiquitous in many different physical systems. In particular, nitrogen-vacancy centres in diamond and semiconductor quantum dots may serve as good qubit candidates in solid state, due to the rapidly improving optical control of their electronic quantum states [29].

In ion traps, there have been various proposals for extension of the MS-gate and entangling operations using a state-dependent force in general [30–32]. Using the framework that we presented in this article, potentially these schemes can be also translated to cavity QED and used to further improve the gate fidelity. Furthermore, we would like to point out that the realization of the AJC Hamiltonian in cavity QED means that not only the MS-gate, but many techniques developed in the context of ion trapping are transferable to optical cavity QED systems. Such examples can include engineered spin–spin interactions [33] and quantum state preparation of the bosonic modes [34, 35]. Moreover, the open-system nature of cavity QED could add another interesting aspect to such realizations through the external driving and extraction of the optical field.

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Appendix A. Perturbative expansion of the propagator in an interaction picture

Starting from Hamiltonian (19) and moving on to an interaction picture defined by a unitary operator

$$U_I(t) = e^{-i\theta a^\dagger a},$$

we get a new system Hamiltonian

$$H_I = \chi a^\dagger a S_x + \delta a^\dagger + g(a + a^\dagger) S_z.$$  

We define a new Hamiltonian

$$H_{MS}^{\text{eff}} = \delta a^\dagger a + g_{\text{eff}}(a + a^\dagger) S_z,$$

such that $H_I = H_{AS} + H_{MS}^{\text{eff}}$. It can be easily shown that the following states form a complete set of the energy eigenstates of $H_{MS}^{\text{eff}}$ for $N = 2$.

$$|S = 1, S_z = 1\rangle |n, \alpha \rangle,$$

$$|S = 1, S_z = 0\rangle |n\rangle,$$

$$|S = 1, S_z = -1\rangle |n, -\alpha \rangle,$$

$$|S = 0, S_z = 0\rangle |n\rangle,$$

where $|n, \alpha \rangle = D(\alpha)|n\rangle$ and $D(\alpha)$ is the displacement operator with an amplitude $\alpha = -g/\delta$. Their energy eigenvalues are a function of $S_z$ and $n$ but not of $S$ and are given by

$$E_n = \delta(n - (j\alpha)^2),$$

for $|S_z = j\rangle |n, j\alpha \rangle$ ($j = 0, \pm 1$).

Now we further move on to a second interaction picture with

$$U_{II}(t) = e^{i\theta a^\dagger a},$$

that leads to a new system Hamiltonian

$$H_{II}(t) = i\theta a^\dagger a H_{AS} e^{-i\theta a^\dagger a}.$$  

A wave function in this second interaction picture $|\Psi_{II}(t)\rangle$ is related to the one in the original picture $|\Psi(t)\rangle$ by

$$|\Psi_{II}(t)\rangle = U_{II}(t)|\Psi(t)\rangle.$$  

On the other hand, the time evolution of $|\Psi(t)\rangle$ is described by a propagator $V_{II}(t)$.

$$|\Psi_{II}(t)\rangle = V_{II}(t)|\Psi(0)\rangle$$

$$V_{II}(t) = 1 - i \int_0^t \int_0^t H_{II}(t') dt'' + (-i)^2 \int_0^t \int_0^t \int_0^t H_{II}(t') H_{II}(t'') dt'dt'' + \cdots,$$

where we have used a Dyson series expansion for $V_{II}(t)$ as $H_{II}(t)$ is not commutative at different times. Note also that the initial state $|\Psi(0)\rangle$ is identical among all the pictures. From (38) and (A.12), we get

$$|\Psi(t)\rangle = U_I^\dagger(t) U_{II}^\dagger(t) V_{II}(t)|\Psi(0)\rangle.$$  

Figure 8. Average gate fidelities as a function of time for two sets of parameters in table 1. Set 1 (blue solid) and set 2 (red dashed).
From (36)
\[ V(t) = U(t)^* U(t) V(t). \] (A.15)

Since \( H_0(t) |S = 0, S_z = 0\rangle = 0 \) and \([H_0(t), S^2] = 0\), non-zero matrix elements of \( H_0(t) \) are limited in the \( S = 1 \) manifold. In other words \( |S = 0, S_z = 0\rangle |n\rangle \) with an arbitrary photon number \( n \) is an eigenstate of the perturbation Hamiltonian \( H_{AS} \) with an eigenvalue of 0. Therefore, this state is not affected by the perturbation and the gate works for this state even in the presence of \( H_{AS} \).

From now on we only consider the matrix elements of \( H_0(t) \) in the \( S = 1 \) manifold and we use the notation introduced in (48). The matrix elements of \( H_0(t) \) are calculated as follows:
\[ \langle 1, m | H_0(t) | 1, n \rangle = 0, \] (A.16)
\[ \langle -1, m | H_0(t) | 1, n \rangle = 0, \] (A.17)
\[ \langle 0, m | H_0(t) | 1, n \rangle = \frac{m \chi e^{i \Delta_0 t}}{\sqrt{2}} \langle m | D(\alpha) | n \rangle = \chi \Lambda_{m0}(t; \alpha), \] (A.18)
\[ \langle 0, m | H_0(t) | -1, n \rangle = \frac{m \chi e^{i \Delta_0 t}}{\sqrt{2}} \langle m | D(-\alpha) | n \rangle = \chi \Lambda_{m0}(t; -\alpha). \] (A.19)

Here we have defined
\[ \Lambda_{m0}(t; \alpha) = \frac{m \chi e^{i \Delta_0 t}}{\sqrt{2}} \langle m | D(\alpha) | n \rangle. \] (A.20)

With these matrix elements \( H_0(t) \) can be expressed as
\[ H_0(t) = \chi \sum_{m,n} (\Lambda_{m0}(t; \alpha) |0, m\rangle \langle 1, n| + \Lambda_{m0}(t; -\alpha) |0, m\rangle \langle -1, n|) + \text{H.c.} \] (A.21)

Likewise
\[ H_0(t)H_0(t') = \chi^2 \sum_{l,m,n} (\Lambda_{ml}(t; \alpha) \Lambda_{m0}(t'; \alpha) + \Lambda_{ml}(t; -\alpha) \Lambda_{m0}(t'; -\alpha)) \]
\[ \times |0, m\rangle \langle 0, n| \]
\[ + \chi^2 \sum_{j,j'=\pm 1} \sum_{l,m,n} \Lambda_{lm}(t; j\alpha) \Lambda_{m0}(t'; j'\alpha) \]
\[ \times |j, m\rangle \langle j', n|. \] (A.22)

Higher order integrands in (A.13) can be as well calculated straightforwardly.

**Appendix B. A model for \( ^{87}\text{Rb} \) atoms coupled to an optical cavity**

We employ the \( D_1 \) transition between \( 5^2S_{1/2} \) and \( 5^2P_{1/2} \) states of single \( ^{87}\text{Rb} \) atoms. Among the eight Zeeman sublevels in the \( 5^2S_{1/2} \) manifold, we pick up \( |F = 2, m_F = -2\rangle \) and \( |F = 1, m_F = -1\rangle \) as our qubit states \(|g\rangle \) and \(|e\rangle \), respectively. These qubit states can be coupled to each other via the upper \( 5^2P_{1/2} \) states. Due to the relatively small hyperfine splitting in the \( 5^2P_{1/2} \) manifold (812 MHz), we take not only \( |F' = 1, m_F = -1\rangle \) (=|\( r_2 \rangle \)) but \( |F' = 2, m_F = -1\rangle \) into account as a mediating upper level for one of the cavity-assisted Raman transitions (see figure B1). The expressions for the effective parameters are obtained as follows [25]:
\[ \Omega_{\text{eff}}^{(1)} = \frac{\mu \Omega}{\sqrt{6} \Delta_{1}}, \] (B.1)
\[ g_{\text{eff}}^{(2)} = \frac{g \Omega_2}{2\sqrt{6}} \left( \frac{1}{\Delta_1} + \frac{1}{\Delta_2 + \omega_{12}} \right), \quad (B.2) \]

\[ \chi = g^2 \left( \frac{1}{4(\Delta_2 + \omega_{12})} + \frac{1}{12\Delta_2} - \frac{1}{3\Delta_1} \right). \quad (B.3) \]

Here \( \omega_{12}/2\pi = 812 \text{ MHz} \) and \( g_{\text{eff}}^{(1)} \) and \( g_{\text{eff}}^{(2)} \) are the effective coupling strengths of the Raman transitions corresponding to the JC and AJC terms, respectively. Note that the above expressions are modified from (20) and (21) due to the Clebsch–Gordan coefficients and the off-resonant coupling to \( |r_z^1\rangle \). The condition \( g_{\text{eff}}^{(1)} = g_{\text{eff}}^{(2)} \) imposes a constraint for \( \Omega_1 \) and \( \Omega_2 \).

Each upper state can decay to the ground qubit states at a rate proportional to the Clebsch–Gordan coefficient for the relevant transition. In addition they can also decay to other Zeeman sublevels in the \( S_{1/2}^\pm \) manifold, which are not shown in figure B1, effectively bringing the system out of the qubit subspace. In order to incorporate such decays into the model, we introduce a virtual auxiliary level \( |w\rangle \). For example the decay rates to \( |u\rangle \) from \( |r_1\rangle \) is equal to the total sum of all the decay rates from \( |r_1\rangle \) to the ground states except for the ones for \( |r_1\rangle \to |g\rangle \) and \( |r_1\rangle \to |e\rangle \). The same applies to the decays from \( |r_2\rangle \) and \( |r_2^\prime\rangle \) to \( |u\rangle \). In this way the decays to \( |u\rangle \) embody all the decays to the outside of the qubit subspace. Note that in the real system it is possible for the atomic population outside the qubit subspace to be pumped back to the subspace again. Here we ignore such processes and the population in \( |u\rangle \) only accumulates in the simulation.

In the end we have nine different Lindblad operators per atom in the form of (65) with the following collapse operators (here \( 2\gamma = 2\pi \cdot 5.75 \text{ MHz} \)):

\[ C^{(i)}_{2g} = \frac{\sqrt{\pi}}{6} \left( \frac{g}{2\sqrt{3} \Delta_1} a |g^{(i)}\rangle \langle e^{(i)}| + \frac{\Omega_1}{2\sqrt{2} \Delta_1} |g^{(i)}\rangle \langle e^{(i)}| \right), \quad (B.4) \]

\[ C^{(i)}_{2e} = \frac{\sqrt{\pi}}{2} \left( \frac{g}{2\sqrt{3} \Delta_1} a |e^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_1}{2\sqrt{2} \Delta_1} |e^{(i)}\rangle \langle g^{(i)}| \right), \quad (B.5) \]

\[ C^{(i)}_{2g} = \frac{\sqrt{\pi}}{2} \left( \frac{g}{2\sqrt{3} \Delta_2} a |g^{(i)}\rangle \langle e^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |g^{(i)}\rangle \langle e^{(i)}| \right), \quad (B.6) \]

\[ C^{(i)}_{2e} = \frac{1}{2} \frac{\sqrt{\pi}}{3} \left( \frac{g}{2\sqrt{3} \Delta_2} a |e^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |e^{(i)}\rangle \langle g^{(i)}| \right), \quad (B.7) \]

\[ C^{(i)}_{2u} = \frac{\sqrt{\pi}}{6} \left( \frac{g}{2(\Delta_2 + \omega_{12})} a |u^{(i)}\rangle \langle e^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |u^{(i)}\rangle \langle e^{(i)}| \right), \quad (B.8) \]

\[ C^{(i)}_{2g} = \frac{\sqrt{\pi}}{2} \left( \frac{g}{2(\Delta_2 + \omega_{12})} a |e^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |e^{(i)}\rangle \langle g^{(i)}| \right), \quad (B.9) \]

\[ C^{(i)}_{2u} = \frac{\sqrt{\pi}}{6} \left( \frac{g}{2\sqrt{3} \Delta_1} a |u^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_1}{2\sqrt{2} \Delta_1} |u^{(i)}\rangle \langle g^{(i)}| \right), \quad (B.10) \]

\[ C^{(i)}_{2g} = \frac{\sqrt{\pi}}{12} \left( \frac{g}{2\sqrt{3} \Delta_2} a |g^{(i)}\rangle \langle e^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |g^{(i)}\rangle \langle e^{(i)}| \right), \quad (B.11) \]

\[ C^{(i)}_{2e} = \frac{\sqrt{\pi}}{12} \left( \frac{g}{2(\Delta_2 + \omega_{12})} a |e^{(i)}\rangle \langle g^{(i)}| + \frac{\Omega_2}{2\sqrt{2} \Delta_2} |e^{(i)}\rangle \langle g^{(i)}| \right). \quad (B.12) \]

References