Reduced dynamics of two oscillators collectively coupled to a thermal bath

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We study the reduced dynamics of a pair of nondegenerate oscillators coupled collectively to a thermal bath. The model is related to the trilinear boson model where the idler mode is promoted to a field. Due to nonlinear coupling, the Markovian master equation for the pair of oscillators admits non-Gaussian equilibrium states, where the modes distribute according to the Bose-Einstein statistics. These states are metastable before the nonlinear coupling is taken over by linear coupling between the individual oscillators and the field. The Gibbs state for the individual modes lies in the subspace with infinite occupation quantum number. We present the time evolution of a few states to illustrate the behaviors of the system.

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I. INTRODUCTION

We consider a composite system of two nondegenerate oscillators coupled collectively to a bath, i.e., the change in the occupation quantum number of one oscillator mode due to the environmental influence induces a corresponding change in the other mode. In the studies of environmental influence on a pair of oscillators, the oscillators are usually coupled separately to the bath through linear interactions [1]. In contrast, our main focus is on the three-body interactions between the oscillators and the field modes of the bath.

The system is related to the trilinear boson model in quantum optical systems, where it is used to describe the process of parametric amplification and frequency conversion [2–5]. The two oscillators then play the roles of the pump mode and the signal mode, respectively, which are coupled to the idler or vibrational mode of a nonlinear medium in which nonlinear interactions are assumed to be dominant. By promoting the idler mode to a field and assuming that the coupling is weak, we can employ the standard open quantum system approach [6,7] to study the damping of the system of oscillators in a thermal bath of the field.

As far as we know, the reduced dynamics of this model has not been discussed before for the general sectors in the nonintegrable region. We find that the Markovian master equation as a time independent eigenvalue problem can be solved analytically. Due to the collective coupling between the oscillators and the bath, the reduced dynamics exhibits the behaviors of finite-level systems [8–10] under the cascade process, even though we are dealing with a continuous variable system.

The Hamiltonian of the system has the same formal structure as the Lee model for the bosonic systems [11,12], and the one-particle sector of its oscillators' subsystem is equivalent to the Friedrichs model [13]. The Friedrichs-Lee model was originally devised to study the effect of perturbation on the spectra in the Hilbert space [13], the mathematical structure of renormalizable quantum field theory [11,14], and later on to study the nonintegrable systems where resonance states emerge [15–17].

The reduced dynamics for this type of interaction shows a few unique features. It gives rise to a family of non-Gaussian equilibrium states confined to their respective irreducible subspaces, whereas the Gibbs state of the individual oscillators is recovered in the subspace with unrestricted occupation quantum number. The oscillator modes in the equilibrium eigenstates distribute according to the Bose-Einstein statistics [18]. These states are metastable before the nonlinear coupling is taken over by linear coupling between the individual oscillators and the field.

In our discussion, we first present the Hamiltonian of the system in Sec. II. The Markovian master equation of the reduced system and its bosonic representation are then presented in Sec. III. We then solve for the equilibrium states in Sec. IV, and study the time evolution of some states in Sec. V.

II. THE HAMILTONIAN

We consider a system of two oscillators and a field in onedimensional space, labeled by 1,2, and k, respectively. The free Hamiltonian is

$$H_0 = \omega_1 a_1^{\dagger} a_1 + \omega_2 a_2^{\dagger} a_2 + \sum_k \omega_k a_k^{\dagger} a_k, \tag{1}$$

where we use the units $\hbar=c=1$, and ω_1,ω_2 are the natural frequencies of the respective oscillators. We assume that $\omega_1>\omega_2$, and the field is massless, with the dispersion relation $\omega_k=|k|$. The creation and annihilation operators a_i^\dagger and a_i obey the commutation relations $[a_i,a_j^\dagger]=\delta_{i,j},\,i,j=1,2,k$. We normalize the field in a box with length Ω , so that $k=2\pi i/\Omega$, with $i=1,2,3,\ldots$ The limit $\Omega\to\infty$ will be taken eventually, but we continue to use the discrete notation in the expressions below.

We assume that the oscillators are coupled collectively to the field through the interaction

$$V = \lambda \sum_{k} \frac{v(\omega_k)}{\sqrt{\Omega/2\pi}} (L_+ a_k + L_- a_k^{\dagger}), \tag{2}$$

in analogy to the linear coupling model between an oscillator (labeled by a,a^{\dagger}) and the field, i.e., $a^{\dagger}a_k + aa_k^{\dagger}$. A derivation of this interaction can be found in Ref. [4] in the context of the trilinear boson model, where it is used to describe parametric

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amplification and frequency conversion in quantum optical systems [2–5]. The L_{\pm} are the ladder operators of the SU(2) algebra

$$L_{+} = a_{1}^{\dagger} a_{2}, \quad L_{-} = a_{1} a_{2}^{\dagger}.$$
 (3)

They raise and lower the r-quantum number of the composite system, respectively [see Eqs. (15)–(17b)]. In Eq. (2), λ is a dimensionless coupling constant, and $v(\omega_k)$ is a form factor that contains a high frequency cutoff to regularize the interactions.

The total Hamiltonian of the system $H = H_0 + V$, when written explicitly in terms of the individual mode, has a form similar to the Lee model for the bosonic system [11,12]. Its one-particle sector of the oscillators' subsystem is equivalent to the Friedrichs model [13].

The system possesses two independent constants of motion,

$$N = a_1^{\dagger} a_1 + a_2^{\dagger} a_2, \quad N_{1k} = a_1^{\dagger} a_1 + \sum_k a_k^{\dagger} a_k.$$
 (4)

N remains a constant of motion of the reduced dynamics when the field modes are traced out. In the unstable regime, the system develops a resonance at the frequency

$$\omega_0 = \omega_1 - \omega_2,\tag{5}$$

where the 1-oscillator turns unstable and decays into the 2-oscillator and the field [15]. Complex poles corresponding to the unstable oscillator then arise in the complex energy plane.

From another point of view, the a_i 's can be regarded as the normal modes of two degenerate oscillators coupled through the SU(2) coupling interactions [19]. The system Hamiltonian H is then unitarily related to a family of Hamiltonians by SU(2) transformation. The details are presented in Appendix A. Therefore, the reduced dynamics discussed below is also applicable to these systems. It is then interesting to note that experimentally [20] it had been shown that spatial wave patterns generated by three-dimensional coherent waves obtained through the longitudinal and transverse coupling of laser modes in a cavity [21] are related to the eigenstates of the system with SU(2) coupling interactions. It may then be possible to realize the spatial profiles of a mixture of eigenstates of the system in the future.

We could gain further insights into the conditions under which the model is applicable by comparing the Hamiltonian to anharmonic interactions [22]. However, since this comparison is outside the main line of our discussion, we present it in Appendix B.

III. MARKOVIAN MASTER EQUATION: BOSONIC REPRESENTATION OF SU(2)

A. Markovian master equation

The reduced dynamics of the two oscillators' subsystem immersed in a thermal bath can be derived by tracing out the field degrees of freedom with standard methods [6,7], or through the complex spectral representation [23]. In the derivation, we assume that the oscillators and the fields are initially factorizable, and we use the weak coupling limit or, equivalently, the $\lambda^2 t$ approximation [22,24] or the Born-Markov approximation [6,7,25].

Since the structure of the interaction Hamiltonian is similar to the linear coupling model between a single oscillator and a field [7], the master equation acquires the Kossakowski-Lindblad (KL) form [26,27], with a modified unitary part. The reduced dynamics is therefore completely positive too [26,27]. The reduced density operator of the two oscillators, \hat{f} , evolves according to the equation $\partial \hat{f}/\partial t = -K \hat{f}$, where

$$K = K_0 + K_d \tag{6}$$

can be decomposed into a unitary part

$$K_0 \hat{f} = i[H_0', \hat{f}],$$
 (7)

$$H_0' = (\omega_1 - \delta\omega_1)a_1^{\dagger}a_1 + (\omega_2 - \delta\omega_2)a_2^{\dagger}a_2 - \delta\omega_0'a_1^{\dagger}a_1a_2^{\dagger}a_2$$

= $(\omega_0 - \delta\omega_0)L_0 - \frac{1}{2}(\omega_0' - \delta\omega_0')N + \delta\omega_0'(L_0^2 - \frac{1}{4}N^2),$
(8)

and a dissipative part

$$K_{d}\hat{f} = -\frac{1}{2}\gamma\bar{n}_{0}(2L_{+}\hat{f}L_{-} - L_{-}L_{+}\hat{f} - \hat{f}L_{-}L_{+})$$
$$-\frac{1}{2}\gamma(\bar{n}_{0} + 1)(2L_{-}\hat{f}L_{+} - L_{+}L_{-}\hat{f} - \hat{f}L_{+}L_{-}).$$
(9)

In the first equality of Eq. (8), we have presented H'_0 in terms of the creation and annihilation operators to exhibit the frequency renormalization $\delta \omega_i$ to the individual oscillator. The explicit expressions of the coefficients are

$$\delta\omega_0' \equiv \delta\omega_1 + \delta\omega_2,\tag{10a}$$

$$\delta\omega_0 \equiv \delta\omega_1 - \delta\omega_2,\tag{10b}$$

$$\delta\omega_1 \equiv \frac{\lambda^2}{\Omega/2\pi} \sum_k P \frac{|v(\omega_k)|^2}{\omega_k - \omega_0} (\bar{n}_k + 1), \qquad (10c)$$

$$\delta\omega_2 \equiv -\frac{\lambda^2}{\Omega/2\pi} \sum_{l} P \frac{|v(\omega_k)|^2}{\omega_k - \omega_0} \bar{n}_k, \qquad (10d)$$

$$\gamma = 2\pi \lambda^2 |v(\omega_0)|^2, \tag{10e}$$

where ω_0 is the resonant frequency (5) and γ is the decay constant. We note that the natural frequencies of the oscillators are renormalized with opposite signs [compare Eq. (10c) with (10d)]. In scattering problems, the average number of field modes is zero $\bar{n}_k = 0$. In this case, we find that only the frequency of the 1-oscillator is renormalized, consistent with the discussion in Ref. [11].

We assume that the field modes are in thermal equilibrium satisfying the Bose-Einstein distribution $\bar{n}_k = 1/[\exp(\omega_k \beta) - 1]$, where $\beta = 1/(k_B T)$. For the resonant mode, we label its occupation number by

$$\bar{n}_0 \equiv \frac{1}{e^{\omega_0 \beta} - 1}.\tag{11}$$

B. SU(2) bosonic representation

The generators of the SU(2) group in terms of the bosonic representation [28] are

$$L_0 = (a_1^{\dagger} a_1 - a_2^{\dagger} a_2)/2, \tag{12a}$$

$$L_1 = (L_+ + L_-)/2, \quad L_2 = (L_+ - L_-)/2i.$$
 (12b)

They obey the commutation relations $[L_i, L_j] = i\epsilon_{ijk}L_k$. The Casimir operator of the SU(2) is [29]

$$L^{2} \equiv \frac{1}{2}(L_{+}L_{-} + L_{-}L_{+}) + L_{0}^{2} = L_{+}L_{-} + L_{0}(L_{0} - 1),$$
(13)

which commutes with the L_i 's, $[L^2, L_i] = 0$. The total occupation number of both oscillators as denoted by N (4) remains a constant of motion of the reduced dynamics. It commutes with the generator of the SU(2) group $[N, L_i] = 0$, $i = 0, \pm$. The quantum number N will be used to label the irreducible representation [see Eq. (15)].

We make use of the occupation number basis

$$|n_1, n_2\rangle = \frac{(a_1^{\dagger})^{n_1}}{\sqrt{n_1!}} \frac{(a_2^{\dagger})^{n_2}}{\sqrt{n_2!}} |0, 0\rangle,$$
 (14)

and denote a state in the irreducible subspace labeled by N as

$$|r\rangle_N \equiv |n_1, n_2\rangle,\tag{15}$$

where

$$r \equiv n_1 - n_2, \quad N \equiv n_1 + n_2$$
 (16)

are related to the eigenvalues of L_0 and L^2 through Eqs. (17c) and (17d), respectively. Using these labels we can establish the following relations:

$$L_{+}|r\rangle_{N} = \frac{1}{2}\sqrt{(N+r+2)(N-r)}|r+2\rangle_{N},$$
 (17a)

$$L_{-}|r\rangle_{N} = \frac{1}{2}\sqrt{(N+r)(N-r+2)}|r-2\rangle_{N},$$
 (17b)

$$L_0|r\rangle_N = \frac{1}{2}r|r\rangle_N,\tag{17c}$$

$$L^2|r\rangle_N = \frac{1}{4}N(N+2)|r\rangle_N,\tag{17d}$$

$$N|r\rangle_N = N|r\rangle_N. \tag{17e}$$

Whenever a 1-oscillator is created, a 2-oscillator is annihilated, and vice versa. Consequently, the index r changes in step of ± 2 under L_{\pm} . There is a total number of N+1 substates in each irreducible subspace, and r ranges from $-N, -N+2, \ldots, N-2, N$. The highest and lowest states are $|N\rangle_N$ and $|-N\rangle_N$, annihilated by the raising and lowering operators $L_{\pm}|\pm N\rangle_N=0$, respectively. The state $|r\rangle_N$ has energy

$$E = \omega_1 n_1 + \omega_2 n_2 = \frac{1}{2} (N \omega_0' + r \omega_0)$$
 (18)

(see Fig. 1 for a plot of the energy levels). They are nondegenerate if the ratio ω_1/ω_2 is not a rational number.

We denote the basis in the Liouville space by

$$f_{r;\tilde{r}}^{(N,\tilde{N})} \equiv |r\rangle_{N\tilde{N}}\langle \tilde{r}| = |n_1, n_2\rangle \langle m_1, m_2|.$$
 (19)

We find that using this notation is more convenient for our later discussion since it is more compact and it manifests the fact that N and \tilde{N} are constants of motion of the reduced dynamics. Since the L_{\pm} operators come in pairs in K_d [Eq. (9)], $Kf_{r,\tilde{r}}^{(N,\tilde{N})}$ is a linear combination of $f_{r;\tilde{r}}^{(N,\tilde{N})}$, $f_{r\pm 2;\tilde{r}\pm 2}^{(N,\tilde{N})}$. Consequently, the quantity

$$v \equiv r - \tilde{r} \tag{20}$$

is a constant of motion under K, and the basis states in each $f^{(N,\tilde{N})}$ subspace are connected to the others with the same ν value only, under the reduced dynamics.

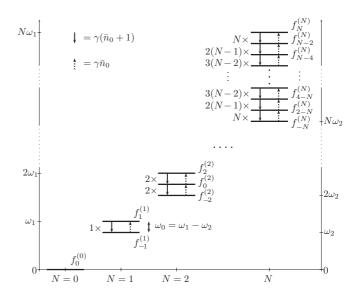


FIG. 1. Energy level diagram and the rate of transitions between different levels in the $f^{(N)}$ subspace (between diagonal elements) [see Eq. (21) for the definition of $f_r^{(N)}$]. The transition rates between two levels are proportional to the numerical constants on the left of the levels. Transitions between different irreducible $f^{(N)}$ subspaces are forbidden by the SU(2) symmetry of the reduced dynamics.

It is interesting to note that the generator of the time evolution K is invariant under a rotation along the L_0 axis, as shown in Appendix D. Furthermore, the SU(2) generalized coherent states [30–32] reside in the $f^{(N,N)}$ subspace (see Appendix E for details).

IV. EQUILIBRIUM STATES

In this section, we obtain the set of equilibrium states for the irreducible subspaces. These states are non-Gaussian states in the coordinate space, as opposed to the Gaussian Gibbs state for an oscillator in a thermal bath. We will also show that the oscillator modes in these equilibrium states distribute according to the Bose-Einstein statistics, and the Gibbs states of the individual oscillator are recovered when the occupation quantum numbers for these modes are not restricted.

We first note that the equilibrium states have nonzero trace and can be decomposed into the diagonal basis state labeled by

$$f_r^{(N)} \equiv f_{r;r}^{(N,N)},$$
 (21)

whenever $\tilde{N} = N$ and $\tilde{r} = r$. We can then write the equilibrium state as a sum of the diagonal basis states:

$$f_{\text{eq}}^{(N)} = \frac{1}{Z_N} \sum_{n=0}^{N} p_{N-2n} f_{N-2n}^{(N)}, \tag{22}$$

where Z_N is a normalization constant and p_{N-2n} are coefficients to be obtained below. Note that since each $f_r^{(N)} = |n_1\rangle\langle n_1| \otimes |n_2\rangle\langle n_2|$ is separable, $f_{\rm eq}^{(N)}$ is a separable state.

The action of K on these states is

$$Kf_r^{(N)} = u_r f_{r+2}^{(N)} + v_r f_r^{(N)} + w_r f_{r-2}^{(N)},$$
 (23)

where

$$u_r = -\frac{\gamma}{4}\bar{n}_0[(N+1)^2 - (r+1)^2],$$
 (24a)

$$v_r = \frac{\gamma}{4}(2\bar{n}_0 + 1)[(N+1)^2 - r^2 - 1] + \frac{\gamma}{2}r,$$
 (24b)

$$w_r = -\frac{\gamma}{4}(\bar{n}_0 + 1)[(N+1)^2 - (r-1)^2]$$
 (24c)

by using Eqs. (17a)-(17e). The transition rates between different energy levels within the same $f^{(N)}$ subspace are summarized in Fig. 1. They can be read off from Eqs. (23)-(24c) directly to give

$$f_r^{(N)} \to f_{r+2}^{(N)} = \begin{cases} 0, & \text{if } r = N \\ u_r, & \text{if } r \neq N \end{cases}$$
 (25a)

$$\rightarrow f_r^{(N)} = v_r, \quad \text{all} \quad r$$
 (25b)

The maximum rate is proportional to the numerical factor N(N+2)/4 for a transition between the $f_{\pm 2}^{(N)}$ and the $f_0^{(N)}$ levels for even N, whereas the maximum rate is proportional to $(N+1)^2/4$ for a transition between the $f_1^{(N)}$ and the $f_{-1}^{(N)}$ levels for odd N [7].

Since the transitions between the energy levels in the $f^{(N)}$ subspace are very similar to the Dicke model for N two-level systems collectively coupled to a field [33,34], superradiance [33] may also arise in this model. However, the energy levels for the two models are different (compare Fig. 1 in this paper to Fig. 1 in Ref. [33]). This is because a symmetrization of the underlying states is required for N two-level systems but not for a pair of nondegenerate oscillators.

The equilibrium condition $Kf_{eq}^{(N)} = 0$ can be written as a matrix equation

$$\mathbf{K} \cdot \mathbf{f}_{\text{eq}}^{(N)} = 0, \tag{26}$$

where

is a tridiagonal square matrix of dimension N + 1, and the eigenvector is a column matrix

$$\mathbf{f}_{eq}^{(N)} = \frac{1}{Z_N} \begin{pmatrix} p_N \\ p_{N-2} \\ p_{N-4} \\ \vdots \\ p_{4-N} \\ p_{2-N} \\ p_{N-N} \end{pmatrix}. \tag{28}$$

The solution to the coefficients is

$$p_{N-2n_2} = \bar{n}_0^{N-n_2} (1 + \bar{n}_0)^{n_2}, \quad n_2 = 0, 1, 2, \dots, N$$
 (29)

which can be checked to satisfy Eq. (26) easily. By requiring $Tr(f_{eq}^{(N)}) = 1$, we obtain the normalization constant

$$Z_N \equiv \sum_{n_2=0,1,2,\dots}^{N} p_{N-2n_2} = (1+\bar{n}_0)^{N+1} - \bar{n}_0^{N+1}$$
 (30)

by using the relation $(1 - r)(1 + r + r^2 + \dots + r^N) = 1 - r^{N+1}$.

Furthermore, the equilibrium occupation probability p_{2n_1-N} can be cast into the form (for fixed N)

$$p_{N-2n_2} = [e^{\beta\omega_2}(1+\bar{n}_0)]^N e^{-\beta E}$$

$$\propto e^{-n_1\beta\omega_1} e^{-n_2\beta\omega_2},$$
(31a)

$$\propto e^{-n_1\beta\omega_1}e^{-n_2\beta\omega_2},$$
 (31b)

where E is the energy of the configuration (18). Equation (31a) shows that p_{N-2n_2} satisfies the canonical distribution for fixed N, whereas Eq. (31b) shows that the system obeys the Bose-Einstein statistics with two energy modes [18], where the 1and 2-oscillator states play the roles similar to the excited and ground states in a two-level system, respectively.

A special case occurs at zero T or, equivalently, when the field mode has zero occupation number $\bar{n}_0 = 0$. Damping to the system can then be caused only by the spontaneous emission of a field quantum [3], accompanied by the lowering and raising of the 1- and 2-oscillator's occupation quantum numbers, respectively. In this situation, the lowest state in each subspace, such as $f_{-N}^{(N)}$ in the diagonal subspaces, and $f_{-N;-\tilde{N}}^{(N,\tilde{N})}$ for $N \neq \tilde{N}$ in the off-diagonal subspaces, are annihilated by K_d . Hence, the subspace spanned by these states does not experience decoherence. Therefore, they should be included in the expression of the equilibrium state for T=0,

$$f'_{\text{eq}} = \sum_{N=0,1,2,\dots} c_N f_{-N}^{(N)} + \sum_{N,\tilde{N}=0,1,2,\dots}^{N \neq \tilde{N}} \left(c'_{N\tilde{N}} f_{-N,-\tilde{N}}^{(N,\tilde{N})} + \text{H.c.} \right),$$
(32)

where c_N are real coefficients subjected to the normalization condition $\sum_{N} c_{N} = 1$, and $c'_{N\tilde{N}}$ are complex coefficients constrained by the positivity condition of f'_{eq} . f'_{eq} is a separable state by inspection. Indeed, since

$$f_{-N,-\tilde{N}}^{(N,\tilde{N})} = |0,N\rangle\langle 0,\tilde{N}| = |0\rangle\langle 0| \otimes |N\rangle\langle \tilde{N}|, \tag{33}$$

we can factor out the overall 1-oscillator state $|0\rangle\langle 0|$ from the right-hand side of Eq. (32), so that $f_{\rm eq}'=|0\rangle\langle 0|\otimes \rho^{(2)}$, where $\rho^{(2)}$ is the density matrix of the 2-oscillator. Hence, it is clear that f'_{eq} is a separable state.

It is interesting to compare the reduced dynamics of the 2-oscillators system to that of a single oscillator coupled to a bath. In the case of a single oscillator, the reduced dynamics too has the Kossakowski-Lindblad form [26,27], although it has a different set of L'_i operators, $L'_0 = a^{\dagger}a + 1/2$, $L'_+ = a^{\dagger}$, and $L'_{-}=a$, where a^{\dagger} , a are the creation and annihilation operators of the oscillator. When written in terms of superoperators, the reduced dynamics has the SU(1,1) symmetry [35]. This reduced dynamics connects all the number bases in this system, and the equilibrium state is the Gibbs state.

For the system we consider, if we trace out the 2-oscillator state from Eq. (22), we find that the 1-oscillator equilibrium state becomes

$$f_{\text{eq},1}^{(N)} = \frac{(1+\bar{n}_0)^N}{Z_N} \sum_{n_1=0,1,2}^{N} e^{-n_1\beta\omega_0} |n_1; n_1\rangle . \tag{34}$$

When *N* is unrestricted, we recover the Gibbs state. Tracing out the 1-oscillator state will produce a similar expression of $f_{\text{eq},2}^{(N)}$ as in Eq. (34), except $|n_1; n_1\rangle$ is replaced by $|N - n_2; N - n_2\rangle$, whereas the rest of the coefficients remain unchanged.

V. TIME EVOLUTION OF STATES

To study the time evolution of the reduced system, we first introduce the interaction picture for the reduced dynamics. We denote the density state in this picture as

$$\tilde{\rho} \equiv e^{K_0 t} \rho = e^{i H_0' t} \rho e^{-i H_0' t}.$$
 (35)

By expanding $\tilde{\rho} = \sum_i \tilde{c}_i f_i$ and $\rho = \sum_i c_i f_i$ in terms of the time-independent basis $f_i \equiv f_{r;\tilde{r}}^{(N,\tilde{N})}$, the coefficient \tilde{c}_i acquires a phase due to the action of $\exp(i K_0 t)$ on f_i ,

$$\tilde{c}_i = c_i \exp(i\theta_i t), \tag{36}$$

in which

$$\theta_{i} = \frac{1}{2}(\omega'_{0} - \delta\omega'_{0})(N - \tilde{N}) + \frac{1}{2}(\omega_{0} - \delta\omega_{0})(r - \tilde{r}) - \frac{1}{4}\delta\omega'_{0}(N^{2} - \tilde{N}^{2}) + \frac{1}{4}\delta\omega'_{0}(r^{2} - \tilde{r}^{2}).$$
(37)

Since the phase angle vanishes for the coefficients associated to the probability elements $f_r^{(N)}$, we have $\tilde{c}_i = c_i$. In this situation, we drop the tilde sign on the coefficients to simplify the notation.

In the interaction picture, the equation of motion becomes

$$\frac{\partial}{\partial t}\tilde{\rho} = -\tilde{K}_d\tilde{\rho},\tag{38}$$

where the effect of \tilde{K}_d on the basis state is

$$\tilde{K}_{d} f_{i} = -\frac{1}{2} \gamma \bar{n}_{0} (2e^{i\delta\omega'_{0}\nu t} L_{+} f_{i} L_{-} - L_{-} L_{+} f_{i} - f_{i} L_{-} L_{+})
- \frac{1}{2} \gamma (\bar{n}_{0} + 1) (2e^{-i\delta\omega'_{0}\nu t} L_{-} f_{i} L_{+} - L_{+} L_{-} f_{i}
- f_{i} L_{+} L_{-}),$$
(39)

in which ν is already defined in Eq. (20) (see Appendix F for details). The extra phase factors in front of two of the terms $L_+f_iL_-$ and $L_-f_iL_+$ are due to the action of L_0^2 in H_0' of Eq. (8). For basis states that lie in the probability subspace, we have $\nu=0$. In this case, \tilde{K}_df_i reduces to K_df_i .

have $\nu=0$. In this case, $\tilde{K}_d f_i$ reduces to $K_d f_i$. By noting that the $f^{(N,\tilde{N})}$ and $f^{(\tilde{N},N)}$ subspaces are related by the following relations, $f^{(\tilde{N},N)}_{\tilde{r};r}=[f^{(N,\tilde{N})}_{r;\tilde{r}}]^{\dagger},\ K_0 f^{(\tilde{N},N)}_{\tilde{r};r}=[K_0 f^{(N,\tilde{N})}_{r;\tilde{r}}]^{\dagger}$, and $\tilde{K}_d f^{(\tilde{N},N)}_{\tilde{r};r}=[\tilde{K}_d f^{(N,\tilde{N})}_{r;\tilde{r}}]^{\dagger}$, we obtain

$$Kf_{\tilde{r}r}^{(\tilde{N},N)} = \left[Kf_{r,\tilde{r}}^{(N,\tilde{N})}\right]^{\dagger}.$$
 (40)

Hence, we can deduce the action of K on one subspace from the other. We will next illustrate the general features of the time evolution of the system with a few examples.

A. States in lower subspaces

We begin by studying the time evolution of a state dwelling in the subspaces up to $N, \tilde{N} = 1$. The density matrix is

$$\tilde{\rho}(t) = d(t)f_0^{(0)} + \left[\tilde{g}(t)f_{0;1}^{(0,1)} + \tilde{h}(t)f_{0;-1}^{(0,1)} + \text{H.c.}\right] + a(t)f_1^{(1)} + b(t)f_{-1}^{(-1)} + \left[\tilde{c}(t)f_{1;0}^{(1,0)} + \text{H.c.}\right], \quad (41)$$

where the coefficients are subjected to the normalization condition d(t) + a(t) + b(t) = 1 and the positivity conditions of $\tilde{\rho}$. The effect $\tilde{K}_d f_i$ can be worked out using Eqs. (39), (17a), and (17b). By equating the coefficients associated to the same basis state f_i on both sides of Eq. (38), we find that the coefficients evolve as

$$\dot{d} = 0, (42a)$$

$$\dot{a} = -\gamma (1 + \bar{n}_0)a + \gamma \bar{n}_0 b, \tag{42b}$$

$$\dot{b} = \gamma (1 + \bar{n}_0)a - \gamma \bar{n}_0 b, \tag{42c}$$

$$\dot{\tilde{c}} = -\gamma \left(\bar{n}_0 + \frac{1}{2}\right)\tilde{c},\tag{42d}$$

$$\dot{\tilde{g}} = -\frac{1}{2}\gamma(\bar{n}_0 + 1)\tilde{g},\tag{42e}$$

$$\dot{\tilde{h}} = -\frac{1}{2}\gamma \bar{n}_0 \tilde{h},\tag{42f}$$

where we have omitted the time dependence on the coefficients for simplicity.

We first make a few observations. We find that only the coefficients under the same $f^{(N,\tilde{N})}$ subspace are connected. We also find that the coefficients for the $f^{(1,1)}$ subspace, namely, a,b,\tilde{c} , evolve in exactly the same way as the amplitude damping channel for qubits [8]. In general, it can be shown that the $f^{(N,N)}$ subspace evolves similar to the N-level system under the cascade process with a single decay constant in vacuum γ ,

$$|N\rangle \leftrightarrow |N-1\rangle \leftrightarrow \cdots \leftrightarrow |1\rangle \leftrightarrow |0\rangle,$$
 (43)

as depicted in Fig. 1. For instance, the dynamics in the $f^{(2,2)}$ subspace behaves similar to the three-level system [10] under the cascade process.

The solutions to Eqs. (42a)–(42f) are

$$d(t) = d_0, (44a)$$

$$a(t) = a_0 e^{-\gamma(1+2\bar{n}_0)t} + \frac{\bar{n}_0(1-d_0)}{1+2\bar{n}_0} [1 - e^{-\gamma(1+2\bar{n}_0)t}], \quad (44b)$$

$$b(t) = 1 - d_0 - a(t), (44c)$$

$$\tilde{c}(t) = \tilde{c}_0 e^{-(2\bar{n}_0 + 1)\gamma t/2},$$
(44d)

$$\tilde{g}(t) = \tilde{g}_0 e^{-\gamma(\bar{n}_0 + 1)t/2},\tag{44e}$$

$$\tilde{h}(t) = \tilde{h}_0 e^{-\gamma \bar{n}_0 t/2},\tag{44f}$$

where d_0 denotes the value of d at t = 0, and etc. From these expressions, we learn that the diagonal components eventually settle down at some equilibrium values, whereas all the off-diagonal components vanish asymptotically. In the general $f^{(N,N)}$ subspace, we find that the diagonal coefficients contain the time exponential factor $\exp[-N\gamma(2\bar{n}_0 + 1)t]$, hence they evolve towards the equilibrium value more rapidly, whereas all the off-diagonal coefficients vanish asymptotically. However, as already shown in Sec. IV, for the special case of zero temperature, the lowest off-diagonal coefficient in each

subspace does not experience decoherence, as is clear from the expression of $\tilde{h}(t)$ in Eq. (44f), which is independent of time when $\bar{n}_0 = 0$.

B. States involving infinite number of subspaces

We now investigate the time evolution of states involving the general subspaces. In principle, the initial states can be decomposed into basis states in the various $f^{(N,\tilde{N})}$ subspaces, and the time evolution can be analyzed subsequently. Since this is a tedious and not an illuminating process, we will discuss the general features of the time evolution by comparing the initial and the equilibrium states with two examples. To simplify the expressions, we will make use of the dimensionless position coordinate

$$x_i \equiv \sqrt{\frac{m_i \omega_i}{\hbar}} \, q_i, \tag{45}$$

where m_i is the mass of the *i* oscillator and q_i is the ordinary position coordinate with the dimension of length.

(1) Assume that the 1-oscillator is initially a superposition of two Gaussian states centered at $x_1 = \pm a$, respectively [36],

$$\phi_1(x_1) = N_1 e^{-(x_1 - a)^2} + N_2 e^{-(x_1 + a)^2} = \sum_{n=0}^{\infty} c_n \langle x_1 | n \rangle, \quad (46)$$

where N_1, N_2 are the normalization constants that give the relative height between the two Gaussians. In Eq. (46), we decompose ϕ_1 in terms of the harmonic oscillator wave function $\langle x|n\rangle = H_n(x) \exp(-x^2/2)/\sqrt{2^n n! \sqrt{\pi}}$, with the expansion coefficient $c_n = \langle n|\phi_1\rangle$, and $H_n(x)$ is the Hermite polynomial. As an example, we choose a=2 and $N_1/N_2=2$. A little calculation shows that c_2 is the dominant term

$$\{c_0, c_1, c_2, \ldots\} = \{0.34, 0.22, 0.78, 0.23, 0.41, 0.05, \ldots\}.$$

The density matrix $\rho_{1,\text{ini}}^{(1)}(x_1,\tilde{x}_1) = \langle x_1|\phi_1\rangle\langle\phi_1|\tilde{x}_1\rangle$ in the coordinate space is plotted in Fig. 2(a).

Consider an initially uncorrelated composite system of two oscillators $\Phi_{\rm ini}^{(1)}$ with the 2-oscillator initially in the ground state $\rho_{2,\rm ini}^{(1)} = |0\rangle\langle 0|$:

$$\Phi_{\text{ini}}^{(1)} \equiv \rho_{1,\text{ini}}^{(1)} \otimes \rho_{2,\text{ini}}^{(1)} = \sum_{n,m=0}^{\infty} c_n c_m^* |n,0\rangle \langle m,0|$$

$$= \sum_{n,m=0}^{\infty} c_n c_m^* f_{n,m}^{(n,m)}.$$
(48)

The spatial profile of $\rho_{2,\text{ini}}^{(1)}(x_1,\tilde{x}_2)$ is plotted in Fig. 2(e). Under the reduced dynamics, the off-diagonal coefficients undergo exponential decay, and $\Phi_{\text{ini}}^{(1)}$ eventually evolves into the equilibrium state

$$\Phi_{\text{eq}}^{(1)} = \sum_{N=0}^{\infty} |c_N|^2 f_{\text{eq}}^{(N)}$$
 (49a)

$$= \sum_{N=0}^{\infty} |c_N|^2 \sum_{n_2=0}^{N} \frac{p_{N-2n_2}}{Z_N} |N - n_2, n_2\rangle \langle N - n_2, n_2| \quad (49b)$$

[cf. Eq. (22) for the expression of $f_{eq}^{(N)}$].

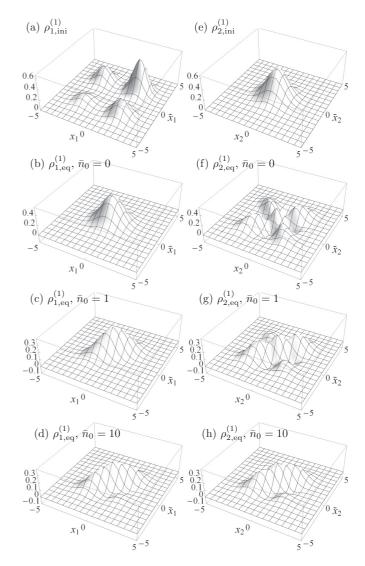


FIG. 2. Initial and equilibrium configurations of the individual modes of example (1) for several temperatures, with parameters a=2 and $N_1/N_2=2$ in Eq. (46).

In the special case of zero temperature, we have $\bar{n}_0 = 0$. The 1-oscillator then settles down to the ground state $\rho_{1,\rm eq}^{(1)} \equiv {\rm Tr}_2 \Phi_{\rm eq}^{(1)} = |0\rangle\langle 0|$ with a Gaussian profile as depicted in Fig. 2(b). Notice that the off-diagonal peaks in Fig. 2(a) have decohered away, much like its single oscillator counterpart in a thermal bath [36]. The information carried by the 1-oscillator in the c_i s is inherited by the 2-oscillator to some extent, as can be seen in the density matrix of the 2-oscillator,

$$\rho_{2,\text{eq}}^{(1)} \equiv \text{Tr}_1 \Phi_{\text{eq}}^{(1)} = \sum_{N=0}^{\infty} |c_N|^2 |N\rangle \langle N|, \tag{50}$$

in which only the highest energy level in each of the 2-oscillator subspaces is occupied. Since c_2 is the dominant term, $\rho_{2,\text{eq}}^{(1)} \approx |c_2|^2 |2\rangle\langle 2|$. The plot in the coordinate space then gives a characteristic three-peak profile of the wave function $\langle x|2\rangle$ along the diagonal [see Fig. 2(f)].

When temperature increases, the populations of the 1- and 2-oscillator start to distribute accordingly among different

levels in the $f^{(N)}$ subspace. The reduced states are then approximately given by

$$\rho_{1,\text{eq}}^{(1)} \approx \frac{|c_2|^2}{Z_2} \left(p_2^{(2)} |2\rangle\langle 2| + p_0^{(2)} |1\rangle\langle 1| + p_{-2}^{(2)} |0\rangle\langle 0| \right), \quad (51)$$

$$\rho_{2,\text{eq}}^{(1)} \approx \frac{|c_2|^2}{Z_2} \left(p_{-2}^{(2)} |2\rangle\langle 2| + p_0^{(2)} |1\rangle\langle 1| + p_2^{(2)} |0\rangle\langle 0| \right). \tag{52}$$

As a result, the three-peak profile is smoothed out, as shown in Figs. 2(c) and 2(g) for $\bar{n}_0=1$. By comparing Eq. (51) with (52), we notice that the order of the level populations between the 1- and 2-oscillator, i.e., $p_2^{(2)}$, $p_0^{(2)}$, and $p_{-2}^{(2)}$, is reversed. This feature recurs in the other subspaces as well. For large temperature, $p_{i,\text{eq}}^{(N)} \approx \bar{n}_0^N$ for all i. Consequently, we obtain a uniform distribution among all the levels. In this limit, $\rho_{1,\text{eq}}^{(1)} \approx \rho_{2,\text{eq}}^{(1)}$, as can be seen by comparing Figs. 2(d) and 2(h) for $\bar{n}_0=10$.

(2) We next consider an initially entangled state

$$\left|\Phi_{\text{ini}}^{(2)}\right\rangle = \sum_{n=0}^{\infty} c_n |n, n\rangle. \tag{53}$$

For a comparison with the first example, we choose c_n to be the same as those in Eq. (47). In this example, both the oscillators are initially in the same state,

$$\rho_{1,\text{ini}}^{(2)} = \rho_{2,\text{ini}}^{(2)} = \sum_{n=0}^{\infty} |c_n|^2 |n\rangle \langle n| \approx |c_2|^2 |2\rangle \langle 2|.$$
 (54)

The density matrix of the system $\Phi_{ini}^{(2)} \equiv |\Phi_{ini}^{(2)}\rangle\langle\Phi_{ini}^{(2)}|$ evolves into the equilibrium state

$$\Phi_{\text{eq}}^{(2)} = \sum_{N=0}^{\infty} |c_N|^2 f_{\text{eq}}^{(2N)},\tag{55}$$

which has been shown to be separable in Sec. IV. Hence, the entanglement between the pair of oscillators is lost eventually. In fact, any initial entanglement in the reduced system vanishes asymptotically in view of the separability of the equilibrium state $f_{eq}^{(N)}$ [Eq. (22)].

A comparison of Eq. (55) with (49a) also shows that the choice of the superpositions $|n,n\rangle$ in Eq. (53) has resulted in the exclusion of the odd $f^{(2N+1)}$ subspace from the equilibrium state. This is a consequence of the fact that the set of odd number subspaces are absent from the initial state. Therefore, by specifically preparing the initial state, some subspaces could be excluded from the equilibrium state. In this model, states with different initial conditions may evolve into different classes of equilibrium states.

VI. CONCLUSION

We have considered a system of two oscillators collectively coupled to a field through a three-body interaction. The model is applicable within a time frame in which nonlinear interaction is dominant. The two oscillator modes become effectively coupled as a result of their collective interaction with the field. Consequently, the reduced dynamics possesses the SU(2) symmetry that is common to finite-level systems, which leads to non-Gaussian equilibrium states for the collective modes. These are metastable states until linear interactions between the individual oscillators and the field take over and drive

them to new equilibrium states. The results suggest that new forms of equilibrium states could emerge when subsystems are collectively coupled to the environment under different symmetry of the reduced dynamics.

It is interesting to further explore the implications and manifestations of the results in other systems, such as in the orbital motion of two-electron quantum dots [37] and light-phonon systems [38], for example, the photosynthetic systems [39]. The existence of metastable states may prevent the oscillators from thermalizing too rapidly with the bath. We leave these interesting investigations to future works.

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APPENDIX A: OSCILLATORS COUPLED BY SU(2) COUPLING INTERACTIONS

Using the operator [19,20]

$$U = \exp(-i\phi L_0) \exp(-i\theta L_2), \tag{A1}$$

where L_i are the generators of the SU(2) group (12a) and (12b), the a_i 's are related to a corresponding set of b_i operators by

$$\begin{pmatrix} a_1 \\ a_2 \end{pmatrix} = U \begin{pmatrix} b_1 \\ b_2 \end{pmatrix}$$

$$U^{\dagger} = \begin{pmatrix} e^{i\phi/2} \cos(\theta/2) & e^{-i\phi/2} \sin(\theta/2) \\ -e^{i\phi/2} \sin(\theta/2) & e^{-i\phi/2} \cos(\theta/2) \end{pmatrix} \begin{pmatrix} b_1 \\ b_2 \end{pmatrix}. \quad (A2)$$

The family of Hamiltonian related to the oscillators' subsystem is then given by

$$H_{12} = \omega_1 a_1^{\dagger} a_1 + \omega_2 a_2^{\dagger} a_2$$

$$= \frac{\omega_0'}{2} (b_1^{\dagger} b_1 + b_2^{\dagger} b_2) + \omega_0 \cos \theta L_0$$

$$+ \omega_0 \sin \theta (\cos \phi L_1 + \sin \phi L_2), \tag{A3}$$

with the corresponding interaction

$$V \sim a_1^{\dagger} a_2 a_k + a_1 a_2^{\dagger} a_k^{\dagger}$$

$$= [-\sin\theta L_0 + \cos\theta(\cos\phi L_1 + \sin\phi L_2)](a_k + a_k^{\dagger})$$

$$-i(\sin\phi L_1 - \cos\phi L_2)(a_k - a_k^{\dagger}), \tag{A4}$$

where a_1, a_2 in L_i [Eqs. (12a) and (12b)] are replaced by b_1, b_2 , respectively. The second equality is obtained by substituting Eq. (A2) into the first line of Eq. (A4). Note that N [Eq. (4)] also commutes with U. Hence, the total occupation quantum number of the oscillators remains the same under the change of basis. Note also that the vacuum state does not change under U, i.e., $|0,0\rangle' = U^{\dagger}|0,0\rangle = |0,0\rangle$.

APPENDIX B: COMPARISON WITH ANHARMONIC INTERACTIONS

The anharmonic interactions couple the position operators of the three species. Using dimensionless position coordinate defined in Eq. (45), the position operator \hat{x}_i is related to the a_i by $\hat{x}_i = (a_i + a_i^{\dagger})/\sqrt{2}$. The anharmonic interactions then take the form

$$V' = \lambda \sum_{k} \frac{v(\omega_k)}{\sqrt{\Omega/2\pi}} (a_1 + a_1^{\dagger})(a_2 + a_2^{\dagger})(a_k + a_k^{\dagger}). \quad (B1)$$

In this situation, we find that aside from the ω_0 resonant mode, there is a second resonant mode that occurs at a higher frequency ω'_0 compared to ω_0 ,

$$\omega_0' = \omega_1 + \omega_2. \tag{B2}$$

This mode is initiated by the interaction terms $a_1a_2a_k^{\dagger} + \text{H.c.}$, where H.c. denotes Hermitian conjugate. If the occupation number of the ω_0' mode in the field is small enough, then this mode is not excited. This could be achieved if the temperature of the bath, T, is low enough so that the condition

$$\omega_0 < k_B T \ll \omega_0' \tag{B3}$$

is satisfied, where k_B is the Boltzmann constant. Another way to achieve this is to formally impose a high frequency cutoff ω_c to the form factor $v(\omega_k)$ in the interaction (B1) so that

$$k_B T < \omega_c \ll \omega_0'$$
 (B4)

is satisfied. Both conditions are separately consistent with the $\lambda^2 t$ approximation [22,24], or the Born-Markov approximation [7,25], used to derive the Markovian master equation of the system. Condition (B3) implies that ω_1,ω_2 should be of the same order since $\omega_1 - \omega_2 < k_B T$ for small T, whereas condition (B4) implies that they should not be too small since we require $\omega_c \ll \omega_1 + \omega_2$ for a large cutoff frequency. In the latter situation, since ω_1 and ω_2 are not small, only a few lower energy modes will be excited in the dynamics (see Fig. 1 for the energy levels of the system).

Aside from the resonant modes, there are two other virtual modes that involve the field quanta with negative frequencies, i.e., $\omega_v' = -(\omega_1 + \omega_2)$ and $\omega_v = -(\omega_1 - \omega_2) < 0$. They are initiated by the interaction terms $a_1^{\dagger} a_2^{\dagger} a_k^{\dagger} + \text{H.c.}$ and $a_1 a_2^{\dagger} a_k + \text{H.c.}$ of Eq. (B1), respectively. The ω_v' mode is a fast rotating mode so that its contribution to the reduced dynamics averages to zero in the relaxation time scale $\tau_R \sim 1/\gamma$, where the $\lambda^2 t$ approximation is valid. This is the rotating-wave approximation [7,40] usually implemented in the Markovian limit.

The virtual mode ω_v does not contribute to the reduced dynamics on the level of the $\lambda^2 t$ approximation with the collision operator $\psi_2^{\vec{\nu}}$ defined in Eq. (C1). A similar example is provided by Ref. [23] for the reduced dynamics of a single oscillator coupled to a field, in which the interactions contain a virtual transition mode. The effect of a possible extension of the collision operator $\psi_2^{\vec{\nu}}$ [Eq. (C1)] to include the contribution of the virtual transitions is presented in the next Appendix.

It is also interesting to note that if instead of assuming $\omega_1 > \omega_2$ in our discussion so far, the opposite situation $\omega_2 > \omega_1$ would interchange the role played by the ω_0 and the ω_0 modes,

i.e., now ω_0 becomes a virtual mode, whereas ω_v becomes a resonant mode. However, the ω_0' and ω_v' modes are not affected by this change.

In summary, in comparison to the full anharmonic interactions, the higher frequency resonant mode ω_0' is not included in the interaction Hamiltonian of the system we consider here. The contribution of this mode is negligible compared to the ω_0 mode at low temperature, or when a frequency cutoff is imposed to the form factor below ω_0' . Moreover, out of the two virtual modes we have discussed, the effect of the ω_v' mode is effectively dropped under the rotating-wave approximation, whereas the ω_v mode does not affect the dissipative part of the reduced dynamics using the usual definition of the collision operator.

APPENDIX C: EXTENSION OF ψ_2

Using the usual definition of the collision operator up to the second order in the coupling constant [23]

$$\psi_{2}^{\vec{\nu}} = P^{\vec{\nu}} \psi_{2} P^{\vec{\nu}}$$

$$\equiv P^{\vec{\nu}} \mathcal{L}_{0} P^{\vec{\nu}} - \lambda^{2} P^{\vec{\nu}} \mathcal{L}_{V} Q^{\vec{\nu}} \frac{1}{\mathcal{L}_{0} - \vec{\omega} \cdot \vec{\nu} - i\epsilon} Q^{\vec{\nu}} \mathcal{L}_{V} P^{\vec{\nu}},$$
(C2)

the virtual mode ω_v does not contribute to the collision operator on the level of the $\lambda^2 t$ approximation. However, if we incorporate the transitions between different $P^{\vec{v}}$ subspaces into the definition of the collision operator (C1), the virtual mode ω_v may contribute to the dissipative part of the reduced dynamics as follows [41].

We define the collision operator by $\psi_2' \equiv P\psi_2 P$, where $P \equiv \sum_{\vec{v}} P^{\vec{v}}$. Figure 3 shows an example of a possible transition of ψ_2' that involves different $P^{\vec{v}}$ subspaces. As indicated in the figure, the interaction vertex on the right is due to the virtual transition $a_1^{\dagger}a_2a_k^{\dagger}$, whereas the vertex on the left is due to a real transition mode in the original model (2). A standard calculation shows that Fig. 3 contributes a term $(a_1^{\dagger}a_2)^2\hat{f}$ to the dissipative part of the reduced dynamics, in addition to the contributions by the real transitions. When taking into account all these additional transitions, the terms $L_+\hat{f}L_+, L_+^2\hat{f}, \hat{f}L_+^2$ and their Hermitian conjugates are added to the dissipative operator, resulting in a not completely positive reduced dynamics [42].

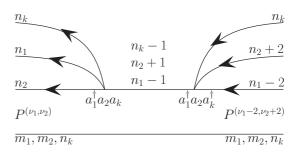


FIG. 3. $P^{(\nu_1,\nu_2)}\psi_2 P^{(\nu_1-2,\nu_2+2)}$ transition. It is one of the diagrams giving the contribution to the collision operator ψ_2 when virtual mode ω_v is considered. The operators on the right vertex cause the virtual transition.

We note that this extension to the definition of the collision operator does not affect the reduced dynamics of the original model (2) since all the possible transitions in this model involve the same $P^{\vec{v}}$ subspace. We already learned that in this situation they lead to a completely positive reduced dynamics [26,27] with the Kossakowski-Lindblad form (9).

APPENDIX D: ROTATIONAL SYMMETRY UNDER L_0

In this Appendix, we show that K is invariant under a rotation along the L_0 axis. Indeed, we find that

$$L'_{+} \equiv e^{i\theta L_0} L_{\pm} e^{-i\theta L_0} = e^{\pm i\theta} L_{\pm} \tag{D1}$$

by using the commutation relations $[L_+, L_-] = 2L_0$ and $[L_0, L_{\pm}] = \pm L_{\pm}$. Since L_+ and L_- appear pairwise in K [cf. Eqs. (6)–(9)], the phase in Eq. (D1) cancels out. Added by the fact that L_0 commutes with N, we conclude that K is invariant under the rotation.

APPENDIX E: SU(2) GENERALIZED COHERENT STATES RESIDE IN $f^{(N,N)}$ SUBSPACE

The SU(2) generalized coherent states [30–32]

$$|\tau\rangle_N = (1+|\tau|^2)^{-\frac{N}{2}} \sum_{n_1=0}^N \sqrt{\binom{N}{n_1}} \tau^{n_1} |n_1, N-n_1\rangle$$
 (E1)

reside in the corresponding $f^{(N,N)}$ subspace of the system, where $\tau = \tan(\theta/2) \exp(-i\phi)$, in which θ, ϕ are the parameters that label the coherent states. In terms of the $f^{(N,N)}$ basis, we

have

$$|\tau\rangle_N\langle\tau| = \sum_{n,m}^N c_{n,m}^{(N,N)} f_{N-2n;N-2m}^{(N,N)},$$
 (E2a)

$$c_{n,m}^{(N,N)} = \frac{1}{2^N} (\sin 2\theta)^N \sqrt{\binom{N}{n}} \binom{N}{m} e^{i\phi(n-m)}. \quad (E2b)$$

APPENDIX F: K_d IN THE INTERACTION PICTURE

In this Appendix, we calculate the effect of \tilde{K}_d on the f_i basis. Since L_0 and N commute, we have

$$e^{iH_0't} = e^{i(\omega_0' - \delta\omega_0')Nt/2 - i\delta\omega_0'N^2t/4} e^{i\delta\omega_0'L_0^2t} e^{i(\omega_0 - \delta\omega_0)L_0t}.$$
 (F1)

Using Eq. (D1), we find that

$$e^{i(\omega_0 - \delta\omega_0)L_0 t} L_+ e^{-i(\omega_0 - \delta\omega_0)L_0 t} = e^{\pm i(\omega_0 - \delta\omega_0) t} L_+.$$
 (F2)

Therefore, $\tilde{K}_d = e^{i\delta\omega_0'L_0^2t} K_d e^{-i\delta\omega_0'L_0^2t}$ since L_+ and L_- appear pairwise in K_d , and N commutes with L_0, L_\pm . Furthermore, using

$$[L_0^2, L_{\pm}] = \pm (2L_0 \mp 1)L_{\pm} = \pm L_{\pm}(2L_0 \pm 1),$$
 (F3)

we can show that

$$\tilde{L}_{\pm} \equiv e^{i\delta\omega_0' L_0^2 t} L_{\pm} e^{-i\delta\omega_0' L_0^2 t} = L_{\pm} e^{\pm i\delta\omega_0' (2L_0 \pm 1)t}$$
 (F4)

$$=e^{\pm i\delta\omega_0'(2L_0\mp1)t}L_{\pm},\tag{F5}$$

where $\tilde{L}_{\pm}^{\dagger} = \tilde{L}_{\mp}$. Using the relation (17c), we finally obtain Eq. (39). Therefore, \tilde{K}_d differs from K_d by a phase factor in two of the terms. For basis vectors that lie in the probability subspace, $\nu \equiv r - \tilde{r} = 0$, and we get $\tilde{K}_d = K_d$.

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