

Temperature effects on quantum non-Markovianity via collision models

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Quantum non-Markovianity represents memory during system dynamics, which is typically weakened by temperature. We study here the effects of environmental temperature on the non-Markovianity of an open quantum system by virtue of collision models. The environment is simulated by a chain of ancillary qubits that are prepared in thermal states with a finite temperature T . Two distinct non-Markovian mechanisms are considered via two types of collision models, one where the system S consecutively interacts with ancillas and a second where S collides only with an intermediate system S' , which in turn interacts with the ancillas. We show that in both models the relation between non-Markovianity and temperature is nonmonotonic. In particular, revivals of non-Markovianity may occur as temperature increases. We find that the physical reason behind this behavior can be revealed by examining a peculiar system-environment coherence exchange, leading to ancillary qubit coherence larger than system coherence, which triggers information backflow from the environment to the system. These results provide insights into the mechanisms underlying the counterintuitive phenomenon of temperature-enhanced quantum memory effects.

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I. INTRODUCTION

In most practical situations a quantum system is open, being coupled to an environment that induces decoherence and dissipation of the system's quantum properties [1]. The dynamics of an open quantum system is usually described with a Markov approximation through a family of completely positive trace-preserving reduced dynamical maps and a corresponding quantum master equation with a Lindblad generator [2,3]. In this case, the memoryless environment is assumed to be able to recover instantly from the interaction, which induces a monotonic one-way flow of information from the system to the environment. However, due to the increasing capability to manipulate quantum systems, in many scenarios the Markov approximation is no longer valid, leading to the occurrence of non-Markovian dynamics [4,5] and a backflow of information from the environment to the system. The non-Markovian dynamics not only embodies an important physical phenomenon linked to dynamical memory effects but also proves useful to enhance practical procedures, such as quantum-state engineering and quantum control [6–13].

Non-Markovianity has recently attracted considerable attention, particularly concerning the formulation of its quantitative measures [14–20], its experimental demonstration [21–25], and the exploration of its origin [26,27]. Nevertheless, the role of non-Markovianity for the assessment of

the properties of nonequilibrium quantum systems has so far remained little explored [28–32]. Non-Markovian dynamics can lead to a new type of entropy production term which is indispensable to recover the fluctuation relations for entropy [28]. In a bipartite system interacting dissipatively with a thermal reservoir in a cascaded model, the emerging non-Markovianity of one of the subsystems enables a heat flow with nonexponential time behavior [29]. By means of Landauer's principle, it has also been shown that memory effects are strategic in maintaining work extraction by erasure in realistic environments [30]. Moreover, non-Markovian dynamics can induce the breakdown of the validity of Landauer's principle [31,32].

An efficient tool that makes the study of quantum thermodynamics in the non-Markovian regime possible [29,31,32] is the *collision* model [33–56]. In the collision model, the environment is taken as a collection of N ancillas organized in a chain and the system of interest S interacts, or collides, at each time step with an ancilla. It has been shown that when the ancillas are initially uncorrelated and no correlations are created among them along the process, a Lindblad master equation can be derived [34,35]. By introducing either correlations in the initial state of the ancillas or interancilla collisions, one can then recover the dynamics of any indivisible, and thus non-Markovian, channel [36–39]. In other words, the non-Markovian dynamics can be achieved in the collision model when the system-environment interaction is mediated by the ancillary degrees of freedom. In analogy to the well-known situation where the non-Markovian dynamics of a system arises when it is coherently coupled to an auxiliary system in contact with a Markovian bath, a class of Lindblad-type master

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equations for a bipartite system has been also found through collision models such that the reduced master equation of the system of interest is derived exactly [40]. By constructing such composite collision models, one can simulate many known instances of quantum non-Markovian dynamics, such as the emission of an atom into a reservoir with a Lorentzian, or multi-Lorentzian, spectral density or a qubit subject to random telegraph noise [49].

Although it is generally believed that quantum memory effects are more important at low temperatures [57], the way temperature influences non-Markovianity depends on both quantum thermodynamics and open quantum system dynamics. For a qubit subject to a dephasing bath with an Ohmic class spectrum, there exists a temperature-dependent critical value of the Ohmicity parameter for the onset of non-Markovianity which increases for high temperatures [58]. For a qubit in contact with a critical Ising spin thermal bath it has been shown that the non-Markovianity decreases close to the critical point of the system in such a way that the higher the temperature, the higher the decrease [59]. Moreover, it is known that the non-Markovianity of a chromophore qubit in a super-Ohmic bath is reduced when the temperature increases [60]. However, temperature may also enhance the non-Markovianity in some situations. For an inhomogeneous bosonic finite-chain environment, temperature has been shown to be a crucial factor in determining the character of the evolution and for certain parameter values non-Markovianity can increase with the temperature [61]. In a spin-boson model made of a two-level system which is linearly coupled to an environment of harmonic oscillators, a nonmonotonic behavior of non-Markovianity as a function of temperature has been reported, with the system dynamics being strongly non-Markovian at low temperatures [62]. Another analysis, studying both entanglement and non-Markovianity measures to reveal how second-order weak-coupling master equations either overestimate or underestimate memory effects, suggests that non-Markovianity can be enriched by temperature [63].

The above results, limited to specific situations, already show how subtle the effect of temperature on quantum non-Markovianity can be during an open system dynamics. In particular, the occurrence of temperature-enhanced memory effects remains counterintuitive and requires further studies which can unveil the underlying mechanisms. In this work we address this issue by means of suitable collision models, which reveal themselves to be specially advantageous to unveil the role of environmental elements in ruling the temperature-dependent non-Markovian dynamics of the system. We consider two types of collision models with different non-Markovian mechanisms, finding that in both models the variation of non-Markovianity as a function of temperature is not monotonic and providing the possible physical reason behind this phenomenon.

II. MEASURE OF NON-MARKOVIANITY

The degree of non-Markovianity in a dynamical process can be quantified by different measures, such as the Breuer *et al.* measure based on the distinguishability between the evolutions of two different initial states of the system [14], the Lorenzo *et al.* measure based on the volume of accessible states of the

system [15], and the Rivas *et al.* measure [16] and the Hall *et al.* measure [20] based on the time behavior of the master equation decay rates. The trace distance between the evolutions of two different initial states $\rho_1(0)$ and $\rho_2(0)$ of an open system is one of the most employed quantifiers. Since a Markovian evolution can never increase the trace distance, when this happens it is a signature of non-Markovian dynamics of the system. Based on this concept, the non-Markovianity can be quantified by the Breuer *et al.* measure \mathcal{N} defined as [14]

$$\mathcal{N} = \max_{\rho_1(0), \rho_2(0)} \int_{\sigma > 0} \sigma[t, \rho_1(0), \rho_2(0)] dt, \quad (1)$$

where $\sigma[t, \rho_1(0), \rho_2(0)] = dD[\rho_1(t), \rho_2(t)]/dt$ is the rate of change of the trace distance given by

$$D[\rho_1(t), \rho_2(t)] = \frac{1}{2} \text{Tr}|\rho_1(t) - \rho_2(t)|, \quad (2)$$

with $|A| = \sqrt{A^\dagger A}$.

Since the dynamics of the system in the collision model is implemented via N equal discrete time steps, in the following the measure \mathcal{N} will be computed by substituting $\sigma[t, \rho_1(0), \rho_2(0)]dt$ with the difference $\Delta D[n] = D[\rho_{1,n}, \rho_{2,n}] - D[\rho_{1,n-1}, \rho_{2,n-1}]$ between the trace distances at steps n and $n-1$ and then summing up all the positive contributions, that is,

$$\mathcal{N} = \max_{\rho_1(0), \rho_2(0)} \sum_{n, \Delta D[n] > 0}^N \Delta D[n] \quad (n = 1, 2, \dots, N). \quad (3)$$

The value of the final collision step N is taken such as to cover all the oscillations of the trace distance during the evolution.

To evaluate the non-Markovianity \mathcal{N} , one then has to find a specific pair of optimal initial states to maximize the time derivative of the trace distance. In Ref. [64] it is proved that, for any non-Markovian quantum process of a qubit, the maximal backflow of information occurs for a pair of pure orthogonal initial states corresponding to antipodal points on the surface of the Bloch sphere. Through a numerical simulation, we find that for our model the optimal initial states alternate between $\{|0\rangle + |1\rangle\}/\sqrt{2}, \{|0\rangle - |1\rangle\}/\sqrt{2}$ and $\{|0\rangle, |1\rangle\}$ depending on the concrete parameters of the environments, which are consistent with the results in [11, 22, 48, 65].

III. NON-MARKOVIANITY IN THE DIRECT COLLISION MODEL

In the first model, illustrated in Fig. 1, the system qubit S directly interacts with the environment \mathcal{R} which comprises N identical qubits R_1, R_2, \dots, R_N . The system qubit and a generic environment qubit are described, respectively, by the Hamiltonians ($\hbar = 1$)

$$\hat{H}_S = \omega_S \hat{\sigma}_z^S / 2, \quad \hat{H}_R \equiv \hat{H}_{R_n} = \omega_R \hat{\sigma}_z^{R_n} / 2, \quad (4)$$

where $\hat{\sigma}_z^\mu = |1\rangle_\mu \langle 1| - |0\rangle_\mu \langle 0|$ is the Pauli operator and $\{|0\rangle_\mu, |1\rangle_\mu\}$ are the logical states of the qubit $\mu = S, R_n$ ($n = 1, 2, \dots, N$) with transition frequency ω_μ (hereafter, for simplicity, we take $\omega_{R_n} = \omega_R = \omega_S = \omega$). The system-bath coupling is assumed to be white noise (very large environment) so the system never collides twice with the same qubit [56]. As a consequence, at each collision step n the system S collides with a “fresh” R_n . Such a model can emulate, for a suitable

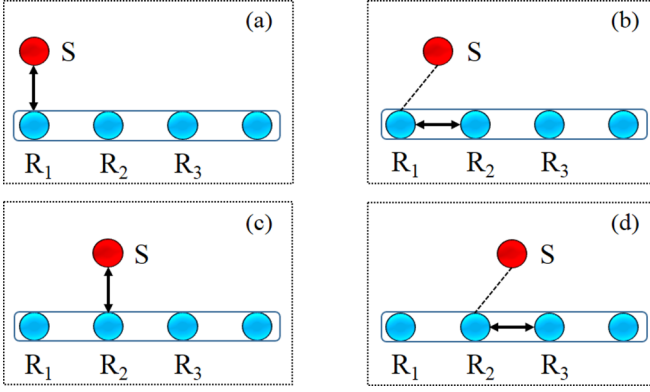


FIG. 1. Schematic diagram of the direct collision model. (a) The system S collides with the ancilla qubit R_1 and they become correlated, as denoted by the dashed line in (b). (b) The intracollision between R_1 and R_2 takes place and the tripartite correlation SR_1R_2 may be generated. (c) The ancilla R_1 is traced out and the process is iterated, namely, the collision of S - R_2 in (c) is followed by that of R_2 - R_3 in (d) and so on.

combination of parameters and interactions, an atom coupled to a lossy cavity [38].

Among the possible choices for the interaction between S and environment qubit R_n , here we focus on a Heisenberg-like coherent interaction described by the Hamiltonian

$$\hat{H}_{\text{int}} = g(\hat{\sigma}_x^S \otimes \hat{\sigma}_x^{R_n} + \hat{\sigma}_y^S \otimes \hat{\sigma}_y^{R_n} + \hat{\sigma}_z^S \otimes \hat{\sigma}_z^{R_n}), \quad (5)$$

where $\hat{\sigma}_j^\mu$ ($j = x, y, z$) is the Pauli operator, g denotes a coupling constant, and each collision is described by a unitary operator $\hat{U}_{S, R_n} = e^{-i\hat{H}_{\text{int}}\tau}$, τ being the collision time. By means of the equality

$$e^{i(\phi/2)(\hat{\sigma}_x \otimes \hat{\sigma}_x + \hat{\sigma}_y \otimes \hat{\sigma}_y + \hat{\sigma}_z \otimes \hat{\sigma}_z)} = e^{-i(\phi/2)}(\cos \phi \hat{\mathbb{I}} + i \sin \phi \hat{S}), \quad (6)$$

with $\hat{\mathbb{I}}$ the identity operator and \hat{S} the two-qubit SWAP operator with the action $|\psi_1\rangle \otimes |\psi_2\rangle \rightarrow |\psi_2\rangle \otimes |\psi_1\rangle$ for all $|\psi_1\rangle, |\psi_2\rangle \in \mathbb{C}^2$, the unitary time evolution operator can be written as

$$\hat{U}_{S R_n} = (\cos J) \hat{\mathbb{I}}_{S R_n} + i(\sin J) \hat{S}_{S R_n}, \quad (7)$$

where $J = 2g\tau$ is a dimensionless interaction strength between S and R_n which is supposed to be the same for any $n = 1, 2, \dots, N$. One can immediately see that $J = \pi/2$ induces a complete swap between the state of S and that of R_n . Thus, $0 < J < \pi/2$ means a partial swap conveying the intuitive idea that, at each collision, part of the information contained in the state of S is transferred into R_n . In the ordered basis $\{|00\rangle_{S R_n}, |01\rangle_{S R_n}, |10\rangle_{S R_n}, |11\rangle_{S R_n}\}$, $\hat{U}_{S R_n}$ reads

$$\hat{U}_{S R_n} = \begin{pmatrix} e^{iJ} & 0 & 0 & 0 \\ 0 & \cos J & i \sin J & 0 \\ 0 & i \sin J & \cos J & 0 \\ 0 & 0 & 0 & e^{iJ} \end{pmatrix}. \quad (8)$$

In the present model, the non-Markovian dynamics of the system is introduced via the interactions between two nearest-neighbor qubits R_n and R_{n+1} . Such interactions are described

by an operation similar to that of Eq. (8), namely,

$$\hat{V}_{R_n R_{n+1}} = \begin{pmatrix} e^{i\Omega} & 0 & 0 & 0 \\ 0 & \cos \Omega & i \sin \Omega & 0 \\ 0 & i \sin \Omega & \cos \Omega & 0 \\ 0 & 0 & 0 & e^{i\Omega} \end{pmatrix}, \quad (9)$$

where $0 \leq \Omega \leq \pi/2$ is the dimensionless R_n - R_{n+1} interaction strength which is taken to be the same for any n .

As illustrated in Fig. 1 exemplifying the first two steps of collisions, in each step we consider the ordered triplet (S, R_{n-1}, R_n) in such a way that after the collision between S and R_{n-1} via the unitary operation $\hat{U}_{S R_{n-1}}$, the system shifts by one site while R_{n-1} collides with R_n via $\hat{V}_{R_{n-1} R_n}$. Notice that the R_{n-1} - R_n collision occurs before the S - R_n collision so S and R_n are already correlated before they collide with each other. The three qubits after the two collisions can now all be correlated with the total state $\rho_{S R_{n-1} R_n}$ (the correlations are labeled by the dashed lines in Fig. 1). Then we trace out the qubit R_{n-1} giving rise to the reduced state $\rho_{S R_n}$ of S - R_n and proceed to the next step with the new ordered triplet (S, R_n, R_{n+1}) . Under the actions of $\hat{U}_{S R_n}$ of Eq. (8) and $\hat{V}_{R_n R_{n+1}}$ of Eq. (9), the total state of $S R_n R_{n+1}$ at the step n is obtained from the step $n-1$ as

$$\rho_{S R_n R_{n+1}} = \hat{V}_{R_n R_{n+1}} \hat{U}_{S R_n} (\rho_{S R_n} \otimes \rho_{R_{n+1}}) \hat{U}_{S R_n}^\dagger \hat{V}_{R_n R_{n+1}}^\dagger, \quad (10)$$

where $\rho_{R_{n+1}} \equiv \rho_R$ is the precollision state of the environmental qubit. Here, to reveal the effect of environmental temperature on the non-Markovianity, we assume that the environmental qubits are initially prepared in the same thermal states $\rho_R = e^{-\beta \hat{H}_R} / Z$ at temperature T_R , where $\beta = 1/k_B T_R$ (k_B being the Boltzmann constant) and Z is the partition function. In our analysis, we consider a dimensionless temperature T defined by $T \equiv k_B T_R / \hbar \omega_0$, where ω_0 is a reference frequency. We also take values of $J \ll \omega / \omega_0$ so as to have small collision times and a weak interaction between the system and the environment qubits.

In such a model, the system experiences a homogenization process and reaches asymptotically the very same state ρ_R [32]. The forward transfer of the lost information of the system S via intracollisions of environment qubits triggers dynamical memory effects of the system, so the non-Markovianity is closely related to the intracollision strength Ω . Figure 2(a) shows the dependence of non-Markovianity \mathcal{N} on Ω for different temperatures T of the environment. In both zero-temperature ($T = 0$) and thermal environments ($T > 0$), the non-Markovianity is activated when Ω exceeds a given threshold [see the inset of Fig. 2(a) for a more evident demonstration] and then monotonically increases with Ω . From this first analysis, it emerges that the thermal environment does not affect the monotonic relation between \mathcal{N} and Ω , while the thresholds of Ω triggering the non-Markovianity depend on the temperature. On the other hand, the variations of non-Markovianity \mathcal{N} with respect to the temperature T can be rich and nonmonotonic, as shown in Fig. 2(b). For relatively small values of Ω (e.g., $\Omega = 0.83$), the increase of T can enable the non-Markovianity which maintains nonzero values within a finite region of $T > 0$. For larger values of Ω , the system dynamics exhibits non-Markovian character already for a zero-temperature environment. In this case, the non-

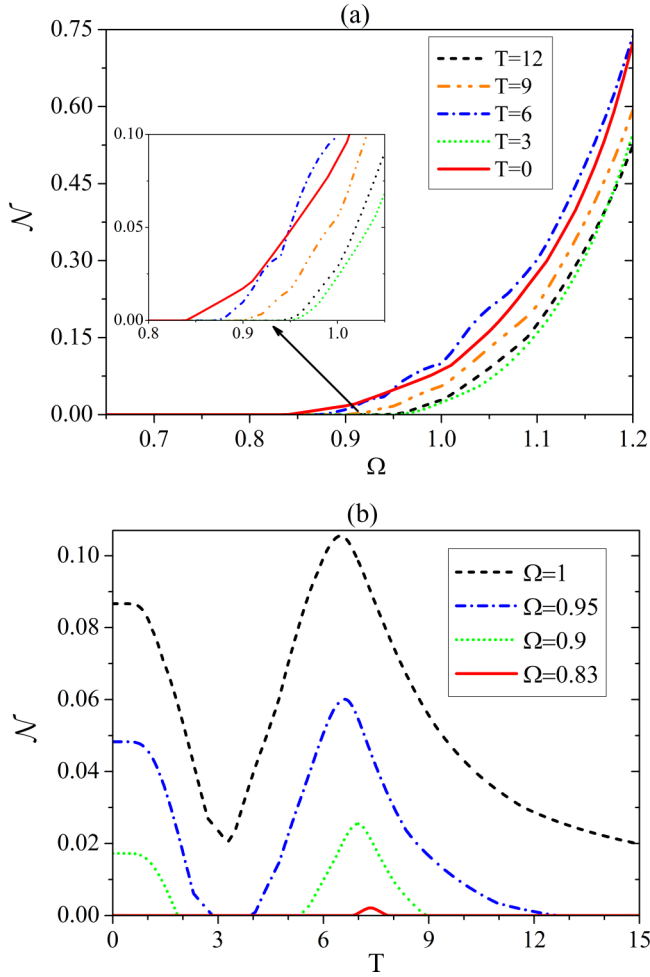


FIG. 2. (a) Non-Markovianity \mathcal{N} vs the collision strength Ω between the environment qubits for different temperatures T of the environment. (b) Non-Markovianity \mathcal{N} vs the temperature T for different Ω . In both plots, the remaining parameters are $\omega = 5\omega_0$ and $J = 0.3 = 0.06\omega/\omega_0$ (small collision time and weak interaction between system and environment qubits).

Markovianity approximately exhibits a plateau for small T and then experiences successive decreasing and increasing behaviors, eventually vanishing at high temperatures. For particular values of the environment qubits interaction strength (e.g., $\Omega = 0.9, 0.95$), when T increases we also observe that the non-Markovianity \mathcal{N} may vanish within a finite interval of T and then revive again. In other words, manipulations of the environment temperature T can induce successive transitions between non-Markovian and Markovian regimes for the system dynamics. A comprehensive picture for the dependence of \mathcal{N} on T and Ω is shown in Fig. 3, where we can see the non-Markovianity thresholds of Ω (identified by the dotted red line) for a given T and the crossovers between non-Markovian and Markovian regimes as T increases for a given Ω .

In order to shed some light on the temperature effects on the non-Markovianity, we examine the exchange of information in terms of coherence between the system and the environment. The initial state of the system is prepared as $|\psi\rangle_{S,0} = (|0\rangle_{S,0} + |1\rangle_{S,0})/\sqrt{2}$ and $C_{S,n} = |\langle 0|\rho_{S,n}|1\rangle| = |\langle 1|\rho_{S,n}|0\rangle|$ is defined

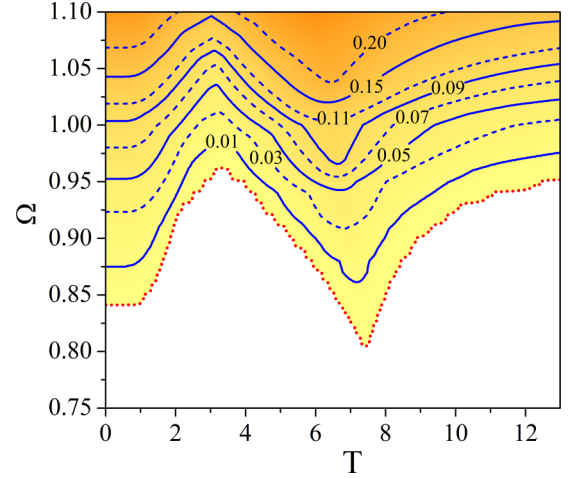


FIG. 3. Contour plot of the non-Markovianity \mathcal{N} for different T and Ω . The non-Markovian regime is colored while the Markovian regime is white. The blue lines are the contour lines of \mathcal{N} . The dotted red line is the curve of the thresholds of Ω triggering non-Markovian dynamics. The remaining parameters are $\omega = 5\omega_0$ and $J = 0.3 = 0.06\omega/\omega_0$ (small collision time and weak interaction between system and environment qubits).

as the coherence degree of the state $\rho_{S,n}$ of S after the n th collision, which is a bona fide quantifier of coherence, being the half of the so-called l_1 -norm measure within a resource theory [66,67]. Therefore, the initial coherence of S has the maximum value $C_{S,0} = 0.5$. The temporary growth of $C_{S,n}$ can serve as a witness for the onset of non-Markovian dynamics. Moreover, to assess the role of the environmental constituents, we consider the coherence $C_{R,n}$ of the environment qubit R_{n+1} transferred from S after the n th collision of S , R_n , and R_{n+1} . In Figs. 4(a)–4(f), we illustrate the evolution of $C_{S,n}$ and $C_{R,n}$ versus n for different temperatures with $\Omega = 0.95$, whose non-Markovian character is plotted in Fig. 2(b) (blue dot-dashed curve). An overall comparison of the panels in Figs. 4(a)–4(f), which indicate a temperature range from $T = 1$ to $T = 7$, verifies the fact that the initial increase of temperature speeds up the decay of the system coherence $C_{S,n}$. Therefore, on the one hand, the increase of temperature suppresses and eventually terminates the non-Markovianity, as can be seen in Figs. 4(a) and 4(b) for $T = 1$ and $T = 2$ and already confirmed in Fig. 2(b). On the other hand, the quick decay of $C_{S,n}$ can cause the coherence $C_{R,n}$ of the environment qubit R_{n+1} to approach [see Fig. 4(c)] and even to exceed [see Figs. 4(d)–4(f)] the coherence $C_{S,n}$ of the system. This behavior in turn induces the information backflow from the environment to the system, namely, a revival of the non-Markovian regime. In fact, $C_{R,n}$ overcomes $C_{S,n}$ in correspondence to the recovery of the non-Markovian character of the system dynamics from a Markovian one [compare Figs. 4(d)–4(f) and the blue dot-dashed curve of Fig. 2(b)]. In the high-temperature regime, the system coherence decays more quickly and the non-Markovian dynamics will cease if the intracollision strength Ω is not sufficiently large. For instance, from Fig. 2(b) one sees that at $T = 10$ the non-Markovianity vanishes if $\Omega = 0.9$, while it remains nonzero when $\Omega = 0.95$. In other words, to get non-Markovian dynamics (quantum memory effects) at high

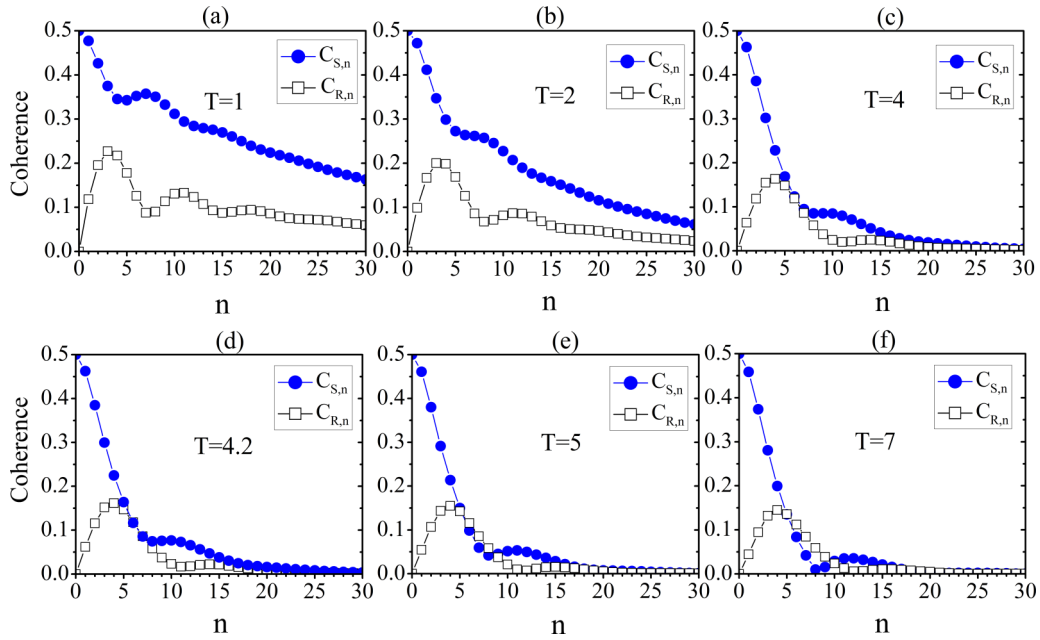


FIG. 4. Coherences $C_{S,n}$ of the system S and $C_{R,n}$ of the environment qubit R_{n+1} after the n th collision as a function of n . The relevant parameters are $\Omega = 0.95$, $\omega = 5\omega_0$, and $J = 0.3 = 0.06\omega/\omega_0$ (small collision time and weak interaction between system and environment qubits).

temperatures, one has to increase the interaction strength Ω between environmental ancillary qubits, which allows a more efficient transfer of environmental quantum coherence.

IV. NON-MARKOVIANITY IN THE INDIRECT COLLISION MODEL

We now consider a mechanism of non-Markovian dynamics based on another collision model, where the interaction of the system qubit S with the environment qubit R_n is mediated by an intermediate qubit S' , as depicted in Fig. 5. Such a scenario implies that the information contained in S is first transferred

to S' and then damped into \mathcal{R} via the collisions between S' and R_n . It is known that, in the absence of environmental intracollisions, this composite model can emulate (for short collision times and Jaynes-Cummings-type interactions) a two-level atom in a lossy cavity, S' playing the role of the cavity mode [49]. For straightforward extension, in the presence of environmental intracollisions, this model may represent a two-level atom in a reservoir with a photonic band gap [49,68,69].

We choose the Heisenberg-type coherent interaction between S and S' , with interaction strength $0 \leq \kappa \leq \pi/2$, represented by the unitary operator $\hat{U}_{SS'}$, analogous to that of

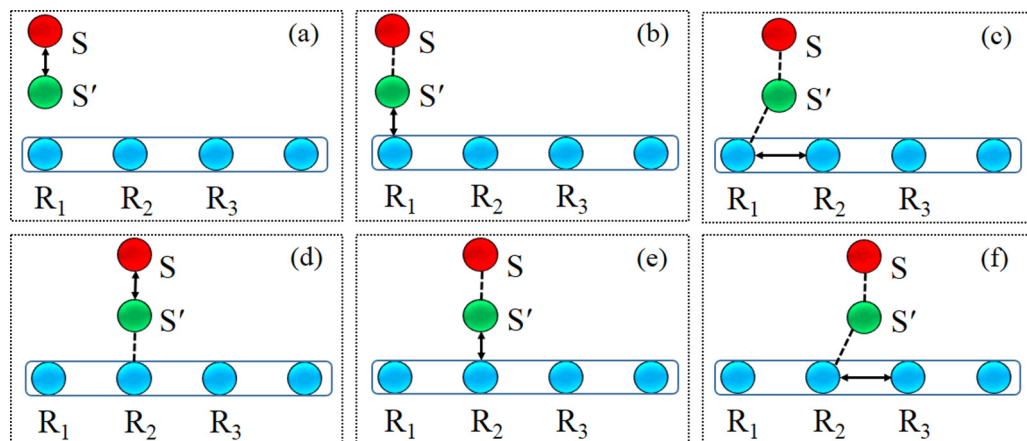


FIG. 5. Schematic diagram of the indirect collision model. (a) The system S collides with the intermediate qubit S' and they become correlated, as denoted by the dashed line in (b). (b) The qubit S' interacts with R_1 and the correlation among $SS'R_1$ is generated, as denoted by the dashed line in (c). (c) The intracollision between R_1 and R_2 takes place and the correlation of $SS'R_1R_2$ is then established. (d) The ancilla qubit R_1 is traced out and the process is iterated, namely, the collision of $S-S'$ in (d) is followed by the collision $S'-R_2$ in (e) and R_2-R_3 in (f) and so on.

Eq. (8), having the form

$$\hat{U}_{SS'} = \begin{pmatrix} e^{i\kappa} & 0 & 0 & 0 \\ 0 & \cos \kappa & i \sin \kappa & 0 \\ 0 & i \sin \kappa & \cos \kappa & 0 \\ 0 & 0 & 0 & e^{i\kappa} \end{pmatrix}. \quad (11)$$

The unitary operators $\hat{U}_{S'R_n}$ and $\hat{V}_{R_n R_{n+1}}$ representing, respectively, the S' - R_n interaction and the interaction between adjacent environment qubits are the same as Eqs. (8) and (9) with interaction strengths J and Ω .

As shown in Fig. 5, in each round of collisions we deal with four qubits (S, S', R_{n-1}, R_n) in such a way that, after the S - S' and S' - R_{n-1} collisions, the qubits S and S' shift by one site while R_{n-1} collides with R_n , which results in the correlated total state $\rho_{SS'R_{n-1}R_n}$ (the correlations are indicated by dashed lines). Then we trace out the qubit R_{n-1} obtaining the reduced state $\rho_{SS'R_n}$ and proceed to the next step with the new ordered group (S, S', R_n, R_{n+1}). As a consequence, the total state of $SS'R_n R_{n+1}$ at the n th collision is determined from the $(n-1)$ th collision as

$$\begin{aligned} \rho_{SS'R_n R_{n+1}} &= \hat{V}_{R_n R_{n+1}} \hat{U}_{S'R_n} \hat{U}_{SS'} (\rho_{SS'R_n} \otimes \rho_{R_{n+1}}) \\ &\times \hat{U}_{SS'}^\dagger \hat{U}_{S'R_n}^\dagger \hat{V}_{R_n R_{n+1}}^\dagger. \end{aligned} \quad (12)$$

The temperature effects are included in this model by considering the qubit S' and all the environmental qubits prepared in the same thermal states $\rho_R = e^{-\beta \hat{H}_R} / Z$ with temperature T .

A. Absence of collisions between environment qubits

In this section, we consider the non-Markovianity in the absence of collisions between environment qubits, i.e., for $\Omega = 0$. For this indirect collision model, the information of the system S is first transferred to the qubit S' via the coherent interaction and then dissipated to the environment through the collisions between S' and environment qubits. In this case, the intermediate qubit S' can have the role of a quantum memory leading to the non-Markovian dynamics even without collisions between environment qubits. The interaction strength κ between S and S' is then crucial in activating the non-Markovianity, as verified in Fig. 6, where \mathcal{N} increases with κ for a given J at a fixed temperature. Moreover, the non-Markovianity achieves a nonzero value only when κ is greater than a threshold and the larger the value of J , the larger the threshold of κ required to trigger the non-Markovian regime. From Fig. 6 one also observes that, for a given κ , the non-Markovianity \mathcal{N} decreases with J , which implies that a strong interaction between S' and the environment qubits weakens the non-Markovianity of the system S .

In Fig. 7, the effect of the temperature T on the non-Markovianity is taken into account for different values of κ . We notice that the non-Markovianity as a function of T is very sensitive to the value of κ , in that it can decrease even directly to zero or decrease to a minimum value and then slowly grow. Remarkably, non-Markovianity can persist at high temperatures provided the values of κ are sufficiently large.

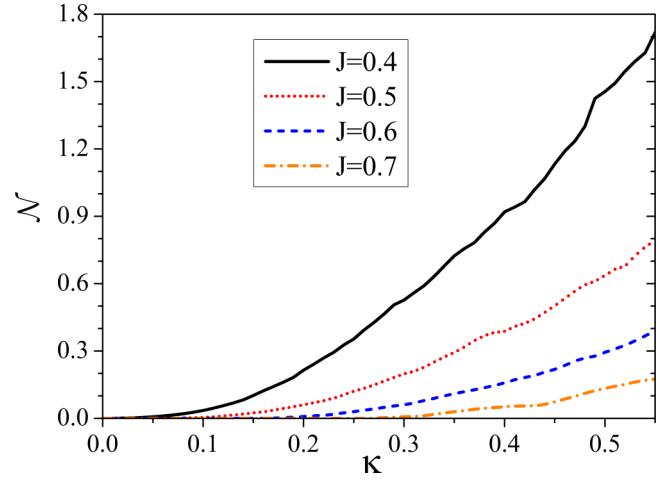


FIG. 6. Non-Markovianity \mathcal{N} versus κ (the interaction strength between S and S') for different values of J (the interaction strength between S' and environment qubits) in the absence of intra-interactions between environment qubits, that is, $\Omega = 0$. The remaining parameters are $T = 1$ and $\omega = 5\omega_0$.

B. Presence of collisions between environment qubits

Now we take the intracollisions between environment qubits R_n and R_{n+1} into account so that the two mechanisms of non-Markovian dynamics, namely, the S - S' interaction ruled by κ and the R_n - R_{n+1} interaction ruled by Ω , coexist in one and the same model. We first explore the role of Ω in enhancing the non-Markovianity at zero temperature. When the coupling between S and S' is weak with relatively small κ , we know from the preceding section that the dynamics of S is Markovian ($\mathcal{N} = 0$) if $\Omega = 0$. As shown in Fig. 8(a), by introducing the R_n - R_{n+1} interactions a threshold of Ω exists which triggers a non-Markovian regime. Such a threshold increases with κ , namely, the smaller the value of κ , the larger the threshold of Ω . The subsequent variations of \mathcal{N} with Ω are nonmonotonic. In particular, we find that the activated non-Markovianity \mathcal{N} can disappear within a finite interval of Ω and then reappear (e.g., for $\kappa = 0.04, 0.06$). When the coupling between S and

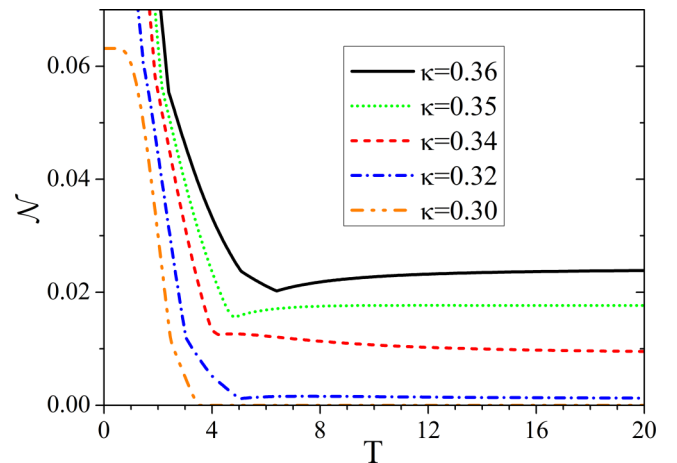


FIG. 7. Non-Markovianity \mathcal{N} as a function of T for different κ with $\Omega = 0$, $\omega = 5\omega_0$, and $J = 0.6 = 0.12\omega/\omega_0$.

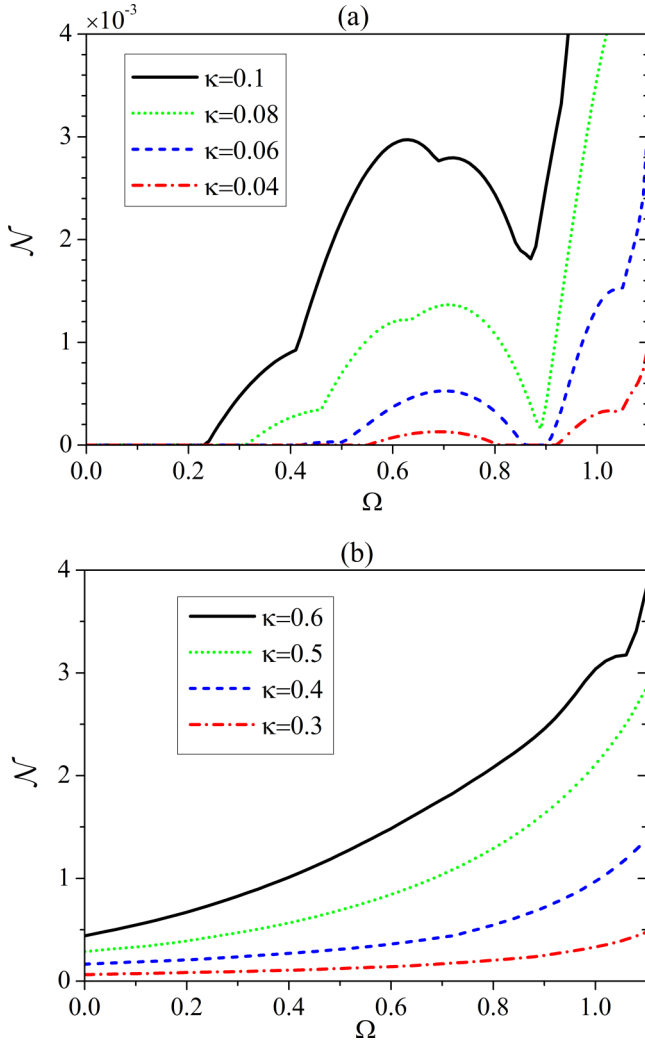


FIG. 8. Non-Markovianity \mathcal{N} versus Ω for different κ at $T = 0$ with $\omega = 5\omega_0$ and $J = 0.6 = 0.12\omega/\omega_0$.

S' is strong with larger κ , the dynamics may be already non-Markovian even for $\Omega = 0$, as seen in Fig. 7 at $T = 0$ and shown in more in detail in Fig. 8(b). In this case, the non-Markovianity can be further enriched by introducing the interactions between environmental qubits R_n and R_{n+1} .

The effects of the environment temperature on non-Markovianity are displayed in Fig. 9. In particular, \mathcal{N} exhibits a nonmonotonic variation with respect to the environmental qubit interaction strength Ω [see Fig. 9(a)], with a first descent and a successive ascent. Once again, we notice that the non-Markovianity can completely disappear for a finite range of Ω and then revive. The non-Markovianity \mathcal{N} as a function of T is then shown in Fig. 9(b), where we observe that the non-Markovianity is unavoidably weakened by increasing T from zero for all the given Ω , but it does not necessarily vanish for larger values of temperature. In fact, \mathcal{N} can collapse and revive (e.g., for $\Omega = 0.9$) and oscillate (e.g., for $\Omega = 1.0$). A comprehensive picture of the variation of \mathcal{N} as a function of both Ω and T , for fixed κ and J , is given in Fig. 10, where the above-detailed behaviors can be retrieved. The values of κ and J are such that the system dynamics is non-Markovian for $T = 0$ and $\Omega = 0$. Such a plot is useful to immediately see

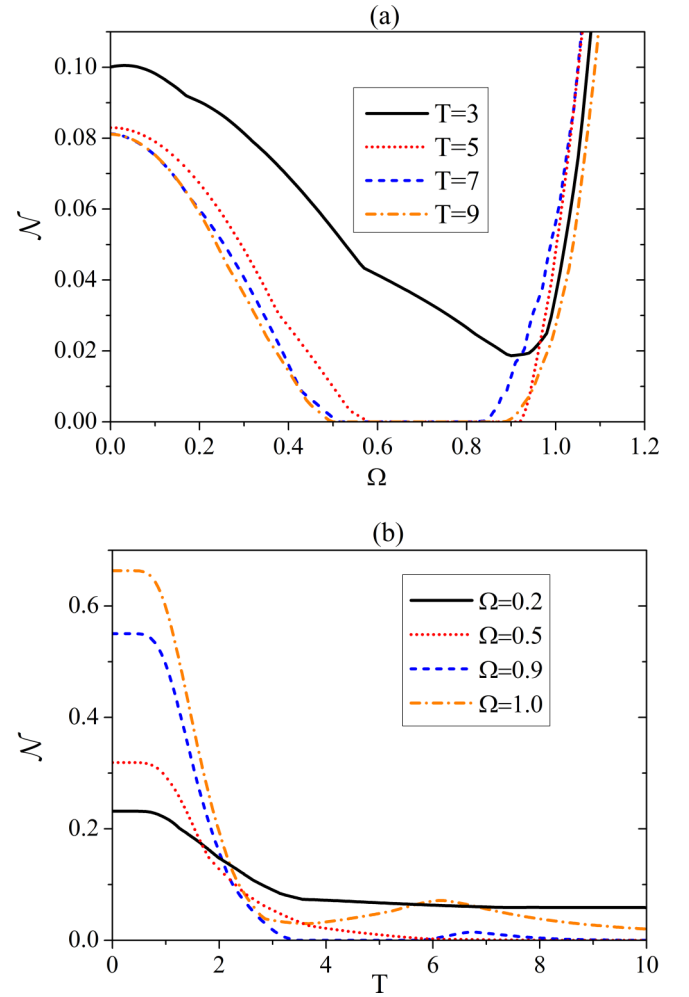


FIG. 9. (a) Non-Markovianity \mathcal{N} versus Ω for different T . (b) Non-Markovianity \mathcal{N} versus T for different Ω . The other parameters are $\omega = 5\omega_0$, $\kappa = 0.3$, and $J = 0.5 = 0.1\omega/\omega_0$.

how, in this composite indirect collision model, the temperature affects the system non-Markovianity in a different way from the case of the direct collision model treated in Sec. III. As a matter of fact, Fig. 10 shows that the temperature has a general detrimental effect on non-Markovianity, which can never overcome its value at $T = 0$ for the higher values of T , as instead happens for the direct collision model [see Fig. 2(b)]. However, a range of values of Ω exists for which a temperature threshold can be found which reactivate dynamical memory effects for the system lost at lower temperatures. In analogy with the direct collision model, such a feature is to be related to peculiar coherence exchanges from the system S to the environmental components.

V. CONCLUSION

We have studied the effects of temperature on the non-Markovian character of an open quantum system dynamics by means of two types of collision models which entail different mechanisms for the occurrence of non-Markovianity. In the first model, that is, the direct collision model, the system S consecutively interacts with a chain of environment qubits that

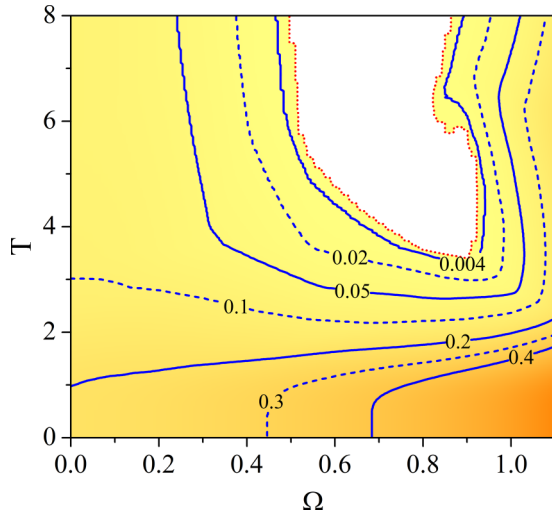


FIG. 10. Contour plot of the non-Markovianity \mathcal{N} for different T and Ω . The non-Markovian regime is colored while the Markovian regime is white. The alternate solid and dashed blue lines are the contour curves of \mathcal{N} . The dotted red line is the curve of the thresholds of Ω and T triggering non-Markovianity. The other parameters are $\omega = 5\omega_0$, $\kappa = 0.3 = 0.06\omega/\omega_0$, and $J = 0.5 = 0.1\omega/\omega_0$.

are prepared in the same thermal states at temperature T and the non-Markovianity \mathcal{N} is induced by the intracollisions of environment qubits. As expected, the non-Markovian dynamics can be triggered when the intracollision strength is greater than a temperature-dependent threshold. In striking contrast to the usual understanding of the effect of the temperature on the non-Markovianity [58–61,63], we have found that the behavior of \mathcal{N} as a function of T is nonmonotonic, exhibiting a process of reduction and enhancement when temperature increases. In particular, we have shown that the non-Markovianity can vanish within a finite interval of T and then reappear when T increases. We have given a possible interpretation of this counterintuitive revival of dynamical memory effects by resorting to the exchanges of coherence between the system and environment qubits. In fact, although the temperature can accelerate the decay of coherence of the system and suppress

the non-Markovianity until certain values, in the regime of high temperature this quick decay of system coherence can cause the coherence transferred to environment qubits to exceed that of the system. This mechanism in turn induces a backflow of information from the environment to the system and thus non-Markovian dynamics.

In the second model, that is, the indirect collision model, the system S indirectly interacts with the environment qubits through collisions with an intermediate qubit S' . In this case, S' serves as the memory for the transferred information from S towards the environment, representing a distinct non-Markovian mechanism. Without intracollisions between environment qubits, the non-Markovian dynamics for the system can still arise provided the strength of the S - S' interaction is sufficiently large. Moreover, the nonmonotonic relation between the non-Markovianity measure \mathcal{N} and T is once again observed. When the environmental intracollisions are taken into account, the two mentioned non-Markovian mechanisms coexist in the same model. In this case we have found that the presence of interactions between environmental qubits enriches non-Markovianity. The temperature now has the general effect of reducing the degree of non-Markovianity with respect to its value at zero temperature. However, once again, the non-Markovianity of the system can exhibit revivals as a function of the temperature.

Our findings within collision models are confirmed by some realistic composite quantum systems which exhibit a nonmonotonic relation between non-Markovianity and temperature [61,62]. More in general, our results contribute towards the capability of engineering suitable environments with optimal temperature conditions to exploit dynamical memory effects of an open quantum system, which is strategic for noisy intermediate-scale quantum information processing [70].

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