The initial-state dependence of the quantum speed limit

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Abstract

The generic bound of quantum speed limit time (the minimal evolution time) for a qubit system interacting with a structural environment is investigated. We define a new bound for the quantum speed limits. It is shown that the non-Markovianity and the population of the excited state can fail to signal the quantum evolution acceleration, but the initial-state dependence is an important factor. In particular, we find that different quantum speed limits could produce contradictory predictions on the quantum evolution acceleration.

Keywords: quantum speed limit bound, Margolus–Levitin bound, open system

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

1. Introduction

The maximal dynamic speed of a quantum system is a fundamental concept in many areas of quantum physics, such as quantum communication [1, 2], quantum metrology [3, 4], optimal control [5], etc. The quantum speed limit (QSL), which is defined as the minimal evolution time between two states was first introduced by Anandan–Aharonov using the Fubini-Study metric [6]. Later, the unified QSL is given for closed systems subject to the unitary evolution, by combining the conclusion of the variance of the energy, i.e., Mandelstam–Tamm bound (MT-QSL) [6–10], and the average of the energy, i.e., Margolus–Levitin bound (ML-QSL) [11–13], which is rigorously described by \( t_{QSL} = \max \{ \pi \hbar / (2 \Delta E), \pi \hbar / (2E) \} \). The QSL bound for the driven system [14–17], and for the mixed quantum state evolution [18–20] have also been investigated. In reality, the system inevitably interacts with its environment, so theories of open quantum systems are usually employed [21]. Recently, the QSL for an open

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system (nonunitary evolution) has attracted intense interest, such as characterizing QSL for an open system using quantum Fisher information [22], relative purity [23], Bures angle [24], and so on [20, 25–29]. Some works [22, 23] have also applied this to non-Markovian dynamics. In particular, in [24], the authors extended the ML-QSL from a closed system to an open system using the technology of the von Neumann trace inequality [30, 31] and operator norm [32], and showed that the non-Markovian effects [33–37] can speed up the quantum evolution and the ML-QSL is tight for the open system. In [28], the authors showed that the mechanism of the acceleration in open systems is determined not only by the non-Markovianity but also by the population of excited states during the quantum evolution.

In this paper, we find that the initial-state dependence is also a key factor to the acceleration of quantum evolution. In particular, the predications for the acceleration of quantum evolution based on different QSLs could be quite inconsistent. Here we first define a tight bound for the quantum speed limit without using the von Neumann trace inequality (we denote this by NI-QSL). Then we use the ML-QSL and the NI-QSL to investigate the damped Jaynes–Cumming (J–C) model with a superposition state as the initial state and the dephasing model. It is found that neither the non-Markovianity nor the population of the initial excited state is competent for the acceleration of quantum evolution, but the different initial states bring great effects. In addition, the ML-QSL and the NI-QSL sometimes produce contradictory predictions on the evolution acceleration, even though they are consistent in some particular cases. This implies that the QSL could need us for further consideration and investigation. This paper is organized as follows. In section 2, we give a brief introduction about the ML-QSL and define the NI-QSL. In sections 3 and 4, we study the damped J–C model and the dephasing model, respectively. Some discussions and conclusions are drawn at the end.

2. The quantum speed limit

In [24], the authors derived the ML-type quantum speed limit bound for a pure state $|\psi_0\rangle$ based on the von Neumann trace inequality and operator norm in the open system through a geometric approach. The ‘distance’ between the initial state $\rho_0 = |\psi_0\rangle\langle\psi_0|$ and its target state $\rho_t$ is measured by the Bures angle, which reads

$$\mathcal{L}(\rho_0, \rho_t) = \arccos \left( \sqrt{\langle \psi_0 | \rho_t | \psi_0 \rangle} \right).$$

(1)

The generalized time-dependent nonunitary equation can be expressed as $\dot{\rho}_t = L_t(\rho_t)$. The time derivative on the Bures angle leads to the relation

$$2 \cos (\mathcal{L}) \sin (\mathcal{L}) \mathcal{L} \leq \left| \langle \psi_0 | \rho_t | \psi_0 \rangle \right| \leq \left| \text{tr} \left\{ L_t(\rho_t) \rho_0 \right\} \right|.$$  

(2)

Based on the von Neumann trace inequality, the right-hand side of equation (2) can arrive at

$$\left| \text{tr} \left\{ L_t(\rho_t) \rho_0 \right\} \right| \leq \sum_{i=1}^{n} \sigma_{1,i} \sigma_{2,i}$$

(3)

with $\sigma_{1,i}$ and $\sigma_{2,i}$ being the singular values of matrix $L_t(\rho_t)$ and $\rho_0$, in decreasing order. By integrating equation (2) over time, one can obtain the ML-QSL for the nonunitary generator as

$$t_{\text{QSL}}^{\text{ML}} \geq \frac{\sin^2 \left[ \mathcal{L}(\rho_0, \rho_t) \right]}{\Lambda_t^{\text{op}}},$$

(4)
where \( \Lambda^{op} = (1/t) \int_0^t \| L_t(\rho_t) \|_{op} \) with \( \| \cdot \|_{op} \) being the operator norm (the maximum singular value) of the matrix.

However, if the von Neumann trace inequality equation (3) and operator norm are not used, then we can directly obtain, through integrating equation (2) over time,

\[
\int_0^t \tau \rho \tau d\tau \geq \tau t L(1) d(t),
\]

which we call the NI-QSL to distinguish it from the ML-QSL. The prominent feature of equation (5) is that it provides a tighter bound than the ML-QSL given by equation (4). Our main conclusion will be obtained by comparing the two types of QSL for the dynamics of open quantum systems.

3. The quantum speed limit time for the damped Jaynes–Cumming model

The first model we will consider includes a single qubit interacting with a vacuum reservoir [21], which is also called the damped J–C model. The whole Hamiltonian of the system and reservoir is \( H = \omega_0 \sigma_x + \sum_k \omega_k b_k^\dagger b_k + \sum_k (g_k \sigma_x b_k + g_k^* \sigma_x^+ b_k) \). The dynamics of the system can be described by

\[
L_t(\rho_t) = \tau_t \left( \sigma_- \rho \sigma_+ - \frac{1}{2} \sigma_+ \rho \sigma_- - \frac{1}{2} \sigma_- \rho \sigma_+ \right).
\]

The structure of environment is assumed as the Lorentzian form:

\[
J(\omega) = \sum_k |g_k|^2 \delta(\omega_0 - \omega_k) = \frac{\gamma_0}{2\pi} \frac{\lambda^2}{(\omega_0 - \omega)^2 + \lambda^2},
\]

where \( \lambda \) is the spectral width of the reservoir and \( \gamma_0 \) is the decay of the system. The Markovian and non-Markovian regimes can be distinguished by the relation of parameters \( \gamma_0 \) and \( \lambda \). In the Markovian regime, we have \( \gamma_0 < \lambda/2 \) and in the non-Markovian regime, we have \( \gamma_0 > \lambda/2 \) [33–37]. If the initial state of the system is a superposition state

\[
|\psi\rangle = a e^{i\theta} |1\rangle + \sqrt{1 - a^2} |0\rangle,
\]

then the final reduced state after the evolution will be [21]

\[
\rho_t = \begin{pmatrix}
\frac{a^2 |q(t)|^2}{\alpha \sqrt{1 - a^2}} & \frac{\alpha \sqrt{1 - a^2} e^{i\theta} q(t)}{\alpha \sqrt{1 - a^2}} \\
\frac{\alpha \sqrt{1 - a^2} e^{-i\theta} q(t)}{\alpha \sqrt{1 - a^2}} & \frac{1 - a^2 |q(t)|^2}{\alpha \sqrt{1 - a^2}}
\end{pmatrix},
\]

where \( q(t) \) is determined by the integro-differential equation \( q(t) = \int_0^t d\tau (t - \tau) q(\tau) \) with the correlation kernel related to the spectral density of the reservoir as \( f(t - \tau) = \int d\omega F(\omega) e^{i(\omega t - \omega \tau)} \). By using the Laplace transformation and its inverse transformation, \( q(t) \) can be given by

\[
q(t) = e^{-\frac{\tilde{d}}{2}} \left[ \cosh \left( \frac{\tilde{d}t}{2} \right) + \frac{\lambda}{d} \sinh \left( \frac{\tilde{d}t}{2} \right) \right],
\]

with \( d = \sqrt{\lambda^2 - 2\lambda \gamma_0} \). The time dependent decay rate \( \gamma_t \) in the equation (6) is given by

\[
\gamma_t = -\text{Im} \left( \frac{q(t)}{q(0)} \right).
\]

After some calculations, we can obtain the ML-QSL for the initial state \( |\psi\rangle \).
\[ \tau(q(t)) = -\left(1 - 2\alpha^2\right)q(t) \]

From equation (11), one can find that the ML-QSL \( t_{QSL}^{ML} \) is related to the \( \alpha, q(t) \) and \( \dot{q}(t) \). The variation of ML-QSL \( t_{QSL}^{ML} \) with the parameter \( \gamma_0 \) and \( \alpha \) is plotted in panel (a) of figure 1. In figure 1, the actual evolution time is chosen as \( t = 1 \) and \( \lambda = 15 \). \( t_{QSL}^{ML} \) is the symmetry about the population parameter \( \alpha = 0 \), so we just study the part \( \alpha > 0 \). It is obviously shown that for \( \gamma_0 < \lambda/2 \), the quantum speed limit time is below the actual evolution time 1. Namely, the quantum evolution displays the acceleration in the Markovian regime. This is quite different from that the acceleration that only appears in the non-Markovian regime due to non-

![Figure 1](image_url)

**Figure 1.** The quantum speed limit time for the initial state given by equation (8). Panel (a) is the ML-QSL \( t_{QSL}^{ML} \), and panel (b) is the NI-QSL \( t_{QSL}^{NI} \), where the parameter \( \lambda = 15 \) and the actual evolution time is \( t = 1 \).
Markovianity, which was discussed $\alpha = 1$ in [24]. One could imagine that the non-Markovianity is not the only reason. The population of the excited state is also a factor for the such an acceleration as that given in [28], which is determined by

$$t_{\text{QSL}} = \frac{t}{2N(t)} + \frac{1}{1 - |\psi(t)|^2},$$

(12)

where $N(\cdot)$ is the degree of non-Markovianity for the dynamics defined as the total backflow of information [34] and $|\psi(t)|^2$ is the population of the excited state, one should notice that in the Markovian regime, $N(\cdot) = 0$, it will directly eliminate the role of the population $|\psi(t)|^2$ in equation (12). Thus, one possible reason for the acceleration subject to the ML-QSL is the dependence of the initial state. The details of the non-Markovianity are given in appendix A.

If we use the NI-QSL to characterize the quantum speed limit, then we have to calculate the NI-QSL for the initial state $|\psi\rangle = (|1\rangle + |0\rangle)/\sqrt{2}$. Here $\lambda = 15$. The dashed line is the actual evolution time $t = 1$, the dash-dotted line stands for the ML-QSL $t_{\text{QSL}}^{\text{ML}}$ and the solid line represents the NI-QSL $t_{\text{QSL}}^\text{NI}$.

Figure 2. The quantum speed limit time for the initial state $|\psi\rangle = (|1\rangle + |0\rangle)/\sqrt{2}$. Here $\lambda = 15$. The dashed line is the actual evolution time $t = 1$, the dash-dotted line stands for the ML-QSL $t_{\text{QSL}}^\text{ML}$ and the solid line represents the NI-QSL $t_{\text{QSL}}^\text{NI}$.

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If we use the NI-QSL to characterize the quantum speed limit, then we have to calculate the NI-QSL for the initial state $|\psi\rangle$ as

$$t_{\text{QSL}}^\text{NI} = \frac{t (1 - q(t)) \left[ 1 - \left(1 - 2\alpha^2\right)q(t) \right]}{\int_0^t \left[ 1 - q(\tau) - \left(1 - 2q(\tau)^2\right)\alpha^2\right]q(\tau) \, d\tau}.$$

(13)

The variation of ML-QSL $t_{\text{QSL}}^\text{ML}$ with the parameter $\gamma_0$ and $\alpha$ is plotted in panel (b) of figure 1. One can find that the quantum speed limit time is fixed to the actual evolution time in the Markovian regime, while the evolution acceleration is demonstrated in the non-Markovian regime, i.e., $\gamma_0 > \lambda/2$. In order to give a more intuitive illustration, the QSLs $t_{\text{QSL}}^\text{ML}$ and $t_{\text{QSL}}^\text{NI}$ for the system with initial state $|\psi\rangle = (|1\rangle + |0\rangle)/\sqrt{2}$ are plotted in figure 2. In figure 2, the parameters are chosen to be the same as those in figure 1. The initial-state dependence can also be found from figure 1 by different $\alpha$.

By comparing the quantum speed limit $t_{\text{QSL}}^\text{ML}$ and $t_{\text{QSL}}^\text{NI}$, one can find that $t_{\text{QSL}}^\text{ML} = t_{\text{QSL}}^\text{NI}$ for the initial state $|1\rangle$ (i.e., $\alpha = 1$). However, if $\alpha \neq 1$, then $t_{\text{QSL}}^\text{ML}$ and $t_{\text{QSL}}^\text{NI}$ usually demonstrates different behaviors, which has been analyzed previously. Thus, the ML-QSL and the NI-QSL produce contradictory predictions on the evolution acceleration.
4. Quantum speed limit for the dephasing model

In the following, we will consider another exactly solvable model: a two-level system coupling with a harmonic oscillator reservoir, which is also called the dephasing model [21]. In the Schrödinger picture, the total Hamiltonian is taken to be
\[ H = \frac{\omega}{2} \sigma_z + \sum_k \omega_k b_k^\dagger b_k + \sum_k \sigma_x (g_k b_k^\dagger + g_k^* b_k) \]
The evolution operator of the system is
\[ \rho_{\sigma} = \frac{\gamma}{2} (\sigma_+ \rho_{\sigma} \sigma_- - \sigma_- \rho_{\sigma} \sigma_+) \]
The initial state is
\[ \ket{\phi} = \beta e^{i\theta} \ket{1} + \sqrt{1 - \beta^2} \ket{0} \]
and the dynamics of the reduced system is expressed as [21]
\[ \rho_s = \left( \begin{array}{cc} \beta^2 & \beta \sqrt{1 - \beta^2} e^{-i\gamma(t)} \\ \beta \sqrt{1 - \beta^2} e^{-i\gamma(t)} & 1 - \beta^2 \end{array} \right) \]
By taking the continuum limit of the bath mode and introducing the spectrum \( J(\omega) \) of the environment, we can find the dephasing factor \( \gamma(t) \) given by
\[ \gamma(t) = \int_0^\infty \omega J(\omega) \coth \left( \frac{\omega}{2k_B T} \right) \frac{1 - \cos(\omega t)}{\omega^2} \]
The spectrum of the environment is chosen as the Ohmic-like spectrum with soft cutoff [38, 39]
\[ J(\omega) = \frac{\omega}{\omega_c^s - 1} \exp \left( -\frac{\omega}{\omega_c} \right) \]
where \( \omega_c \) is the cutoff frequency, \( \eta \) is the dimensionless coupling constant and the parameter \( s > 0 \). For simplicity, we will assume that the cutoff frequency \( \omega_c \) is 1. \( s \) determines the property of the environment, such as the sub-Ohmic reservoir for \( s < 1 \), the Ohmic reservoir for \( s = 1 \) and the super-Ohmic reservoir for \( s > 1 \). Under the condition \( T = 0, \gamma > 0 \) and \( s > 0 \), the dephasing factor \( \gamma(t) \) can be obtained [40] by
\[ \gamma(t) = \eta \left[ 1 - \cos \left[ (s - 1) \arctan \left( \frac{\omega_t}{\omega_c} \right) \right] \right] \]
where \( \Gamma(\cdot) \) is the Euler gamma function. Thus, the ML-QSL for the state in equation (14) can be given by
\[ t_{QSL}^{ML} = \frac{2\beta}{\gamma} \sqrt{1 - \beta^2} \left( 1 - e^{-\gamma(t)} \right) \]
From equation (19), one can find that the ML-QSL \( t_{QSL}^{ML} \) depends on the dephasing rate \( \gamma'(t) \), which is the time derivative of the dephasing factor \( \gamma(t) \) and is determined by
\[ \gamma'(t) = \int_0^\infty \omega J(\omega) \sin \left( \omega t \right) \frac{d\omega}{\omega} \]
From equation (19), one can find that, if \( \gamma'(t) > 0 \), then the integral in denominator of equation (19) can be given analytically as \( 1 - e^{-\gamma(t)} \). This will obviously lead to that \( t_{QSL}^{ML} = 2\beta \sqrt{1 - \beta^2} \). In other words, the quantum speed limit time is fixed for the given initial state. It is irrelevant to the property of the environment. However,
in the region of \( \gamma'(<t) \), usually the denominator of equation (19) is hard to give explicitly. So it is important to tell the sign of \( \gamma'(<t) \) in order to give the ML-QSL. For the Ohmic-type spectral, the dephasing rate \( \gamma'(<t) \) can be given analytically as

\[
\gamma'(<t) = \eta \left( 1 + t^2 \right)^{-\nu/2} \Gamma (s) \sin (s \arctan (t)).
\]
One can find that $\gamma'(t)$ is always positive if $\sin \left( s \arctan \left( t \right) \right) > 0$, i.e., $s$ and $t$ satisfy

$$k\pi < s \arctan \left( t \right) < (2k + 1)\pi,$$

otherwise, $\gamma'(t)$ is negative. A vivid illustration of equation (21) is provided in figure 3. The green region corresponds to the positive $\gamma'(t)$, and the other blank region stands for the negative $\gamma'(t)$.

The ML-QSL $t_{\text{QSL}}^{\text{ML}}$ versus $s > 2$ and $\eta$ is plotted in figure 4, where the lower layer corresponds to $2\beta\sqrt{1 - \beta^2} = 0.5$ and the upper layer corresponds to $\beta = \frac{1}{\sqrt{2}}$. It can be found that the ML-QSL for $\beta = \frac{1}{\sqrt{2}}$ serves the tightest ML-QSL for all possible $\beta$. Here we choose the actual evolution time as $t = 3$. Since $\arctan(t) \in [0, \pi/2)$, if $s \leq 2$, $\gamma'(t)$ will always be positive, which can be seen from figure 3. In this case, the ML-QSL will be a constant and independent of $s$ and $\eta$, for example, $t_{\text{QSL}}^{\text{ML}} = 1.5$ for the lower layer and $t_{\text{QSL}}^{\text{ML}} = 3$ is just the actual evolution time for the upper layer. So the part for $s < 2$ is not shown in figure 4.

Compared with figure 3, one can find that the ML-SQL keeps invariant until $s \approx 2.5$. For all $s$ the ML-SQL given by the lower layer shows the acceleration of the quantum evolution, but no acceleration is shown by the upper layer for $s \ll 2.5$. In both layers, the population of the excited state is not changed due to the dephasing dynamics. Therefore, the acceleration with different degrees could be understood by the non-Markovianity. In [39], it is shown that for $s > 2$ the dynamics for $t \to \infty$ enter the non-Markovian regime. In fact, if $t$ is finite, then one can easily find that the negative $\gamma'(t)$ will lead to the non-Markovian dynamics (see appendix B for details). In other words, figure 3 just shows the bounds of the Markovian region and the non-Markovian region. However, the two layers in figure 4 corresponding to the different initial states imply the opposite properties of the acceleration, especially for $2 < s \ll 2.5$. Based on figure 5, the non-Markovianity could effectively signal the change of quantum evolution velocity for $\beta = \frac{1}{\sqrt{2}}$, but it fails for the lower layer. This shows the strong dependence of the initial state. Similarly, the NI-QSL can be given by

$$t_{\text{QSL}}^{\text{NI}}(s, \eta) = \frac{1}{\int_0^t \left| e^{-\gamma(t)'(t)} \right| \, dt}.$$ (23)

By comparing equation (19) with (23), one can immediately find that the NI-QSL $t_{\text{QSL}}^{\text{NI}}$ does not depend on the population of the excited state. It is interesting that in this particular model, the NI-QSL does not depend on the initial state either. The ML-QSL $t_{\text{QSL}}^{\text{ML}}$ is connected with the NI-QSL $t_{\text{QSL}}^{\text{NI}}$ by a factor $2\beta\sqrt{1 - \beta^2}$. Thus, the NI-QSL versus $s$ and $\eta$ is just consistent with the ML-QSL for $\beta = \frac{1}{\sqrt{2}}$, i.e., the upper layer in figure 4. Since the opposite prediction on the evolution acceleration has been analyzed by the two layers, it is also implied that the NI-QSL and the ML-QSL demonstrate contradictory predictions.

5. Discussions and conclusion

One may notice that the ML-QSL $t_{\text{QSL}}^{\text{ML}}$ is zero for some special states, such as the ground state $|0\rangle$ for damped J–C model, and the excited state $|1\rangle$ or ground state $|0\rangle$ for the dephasing model. However, the ‘distance’ (Bures angle) between the initial state and some other target state is also zero. So it is not difficult to understand the zero evolution time. In addition, the NO-QSL is obtained without the von Neumann trace inequality and operator norm used, so it is obviously tighter than (at least, consistent with) the previous bounds.
In summary, we have studied the ML-QSL and the NI-QSL of the damped J–C model and the dephasing model, and have found the importance of the initial-state dependence for the quantum evolution acceleration. To some extent, this is consistent with the derivation and conclusion of the MT-QSL in open system dynamics, which is evaluated with respect to the initial state [23]. We also find that the predictions on the evolution acceleration based on different QSLs could produce contradictory conclusions. This implies that the QSL could deserve our further consideration. We hope that our work could deepen the understanding of the quantum speed limit on the dynamics of open systems.

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Appendix A. The non-Markovianity of damped J–C model

Following [34], the non-Markovianity measure of the damped J–C model we employed is based on the trace distance $D(\rho_{p_1}, \rho_{p_2}) = \frac{1}{2} \text{tr} | \Phi_{p_1} - \Phi_{p_2} |$. The change rate of the trace distance is $\partial_t D(\rho_{p_1}, \Phi_{p_2})$. The positive change rate stands for the flow of information from the environment back to the system. The non-Markovianity of the quantum process $\Phi(t)$ can be given by

$$N(\Phi) = \max_{\rho_{p_1}, \rho_{p_2}} \int_0^\lambda D(\rho_{p_1}, \rho_{p_2}).$$

(A.1)

Figure A1. The non-Markovianity versus $\gamma_0$. The parameter $\lambda = 15$. The non-Markovianity measure is normalized to unity.

In summary, we have studied the ML-QSL and the NI-QSL of the damped J–C model and the dephasing model, and have found the importance of the initial-state dependence for the quantum evolution acceleration. To some extent, this is consistent with the derivation and conclusion of the MT-QSL in open system dynamics, which is evaluated with respect to the initial state [23]. We also find that the predictions on the evolution acceleration based on different QSLs could produce contradictory conclusions. This implies that the QSL could deserve our further consideration. We hope that our work could deepen the understanding of the quantum speed limit on the dynamics of open systems.

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$$N(\Phi) = \max_{\rho_{p_1}, \rho_{p_2}} \int_0^\lambda D(\rho_{p_1}, \rho_{p_2}).$$

(A.1)

It is hard to obtain the optimal initial state pair ($\rho_1$, $\rho_2$) for a general process. However, for the damped J–C model, it proves that the excited state $|1\rangle$ and the ground state $|0\rangle$ are the optimal state pair, so $\partial_t D(\rho_1, \rho_2)$ can be given by a simple expression $\partial_t D(\rho_1(t), \rho_2(t)) = 2 |q(t)| \cdot |q(t)|$ with $|q(t)|$ being the time derivative of $|q(t)|$. Thus one can easily obtain the relation between the non-Markvianity and the parameter $\gamma_0$, which is plotted in figure A1. By comparing with figures 1 and 2, one can find that the quantum evolution can be accelerated in the non-Markovian regime based on NI-QSL. However, based on the ML-QSL, the quantum evolution is accelerated not only in the non-Markovian regime but also in the Markovian regime.
Appendix B. The non-Markovianity of the dephasing model

The memory effects (non-Markovianity) for the dephasing model associated with the quantum channel capacity \( \Phi(\Phi) \) can be defined by [39, 41]

\[
N = \int_{\Phi(Q)>0} dt \partial_t \Phi(\Phi),
\]

where quantum channel capacity \( Q(t) = 1 - H_2(\frac{1+e^{-\eta t}}{2}) \) with \( H_2(\cdot) \) being the binary Shannon entropy. Due to equation (B.1), the non-Markovian regime is determined by \( \partial_t \Phi(\Phi) > 0 \). It is easy to find that \( \partial_t \Phi(\Phi) > 0 \) means \( \gamma'(t) < 0 \), based on which we plot the Markovian and non-Markovian regions in figure 3. The non-Markovianity versus the
parameter $s$ and $t$ is plotted in figure B1(a). Comparing with figure 3, one can find that the blank regime in the figure 3 corresponds to the non-Markovian regime. In figure B1(b), we plot the non-Markovianity versus the Ohmic parameter $s$ and coupling constant $\eta$ when the evolution time $t = 3$. Comparing with the figure 4, one can easily find that the quantum evolution is accelerated within the non-Markovian regime based on the ML-QSL for $\beta = 1/\sqrt{2}$ (the upper layer in figure 4). However, for the lower layer in figure 4 ($2\beta\sqrt{1 - \beta^2} = 0.5$), the quantum evolution can also be accelerated in the Markovian regime. This implies the initial-state dependence.

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