## Speed limit, dissipation bound, and dissipation-time trade-off in thermal relaxation processes

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We investigate bounds on speed, nonadiabatic entropy production, and the trade-off relation between them for classical stochastic processes with time-independent transition rates. Our results show that the time required to evolve from an initial to a desired target state is bounded from below by the information-theoretical  $\infty$ -Rényi divergence between these states, divided by the total rate. Furthermore, we conjecture and provide extensive numerical evidence for an information-theoretical bound on the nonadiabatic entropy production and a dissipation-time trade-off relation that outperforms previous bounds in some cases.

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Introduction. Optimal control of a system's evolution from an initial to a desired target state is a crucial task [1–4], with close ties to the optimal transport problem [5,6]. The definition of "optimal" varies depending on the specific cost function employed, which may include time, energy consumption, dissipation, error, robustness, or trade-offs between them. In addition to optimal control protocols, non-modelspecific fundamental bounds on the cost functions are of significant interest [7]. For instance, in quantum systems, rapid state transformations are usually desirable, thereby motivating extensive investigations of the so-called "quantum speed limit" [8–10]. For a quantum system with a timeindependent Hamiltonian *H*, the time it needs to evolve from the initial state  $\rho^{(i)}$  to the final state  $\rho^{(f)}$  is bounded from below by ( $\hbar = 1$ )

$$\tau \ge \max\left\{\frac{\mathcal{L}(\rho^{(i)}, \rho^{(f)})}{\Delta H}, \frac{2\mathcal{L}^2(\rho^{(i)}, \rho^{(f)})}{\pi \langle H \rangle}\right\},\tag{1}$$

where  $\mathcal{L}(\rho^{(i)}, \rho^{(f)})$  is the Bures angle that quantifies the distance between the end points, and  $\Delta H$  and  $\langle H \rangle$  are variance and average of the Hamiltonian, respectively [11,12].

Recent developments have extended the concept of speed limits to classical stochastic processes, where entropy production plays a crucial role [13–18]. For Markovian stochastic processes with given initial and final probability distributions  $p^{(i)}$  and  $p^{(f)}$ , speed limits can be expressed as a trade-off between entropy production  $\Sigma$  and time duration  $\tau$  given by  $(k_{\rm B} = 1)$ 

$$\Sigma \geqslant f_{\boldsymbol{p}^{(i)} \to \boldsymbol{p}^{(f)}}(\mathcal{R}\tau), \tag{2}$$

where f is a monotonically decreasing function of  $\mathcal{R}\tau$ . The subscript  $p^{(i)} \rightarrow p^{(f)}$  denotes the dependence of the function on the end points, and  $\mathcal{R}$  quantifies the system's timescale. This inequality encompasses the following three important cases in the literature.

In generic processes entropy production can be split into adiabatic and nonadiabatic (Hatano-Sasa [19]) contributions [20]. The former persists even when the probability distribution equals the instantaneous steady state, while the latter arises from deviations from this state. It was reported that nonadiabatic entropy production  $\Sigma_{na}$  is lower-bounded by (referred to as activity bound) [15,16]

$$\Sigma \geqslant \Sigma_{\text{na}} \geqslant \frac{\mathcal{L}^2(\boldsymbol{p}^{(i)}, \boldsymbol{p}^{(f)})}{2\mathcal{A}\tau},$$
(3)

where  $\mathcal{L}(\boldsymbol{p}^{(i)}, \boldsymbol{p}^{(f)})$  represents the total variation distance between end points and  $\mathcal{A}$  denotes time-averaged dynamical activity [21].

In the quasistatic limit with  $\tau \to \infty$ , the time-averaged dynamical activity becomes time independent, resulting in an inequality  $\Sigma_{na} \ge \mathcal{O}(1/\tau)$ . This observation bears resemblance to the result for slow but finite-time Markovian stochastic processes [22–26], where the entropy production bound is given by  $\Sigma \ge \mathcal{L}_T^2/\tau$ . Here  $\mathcal{L}_T$  denotes the thermodynamic length [27–44], which already encodes the information about the timescale derived from the transition rates.

For relaxation processes with a time-independent transition rate matrix satisfying the detailed balance condition, the general dissipation-time trade-off relation Eq. (2) is superseded by a simpler  $\tau$ -independent bound [45]

$$\Sigma \geqslant D_1(\boldsymbol{p}^{(1)} \| \boldsymbol{p}^{(1)}), \tag{4}$$

where  $D_1(\mathbf{p}^{(i)} \| \mathbf{p}^{(f)})$  is the 1-Rényi divergence [46,47], also known as Kullback-Leibler divergence or relative entropy [48]. Since the bound is independent of time duration, the trade-off between dissipation and evolution time is, in fact, concealed. In the context of these kinds of simple but important processes, i.e., relaxation processes with time-independent transition rates where the detailed balance condition is not necessarily met [49], three natural questions arise. First, given the similarity between a quantum process with a time-independent Hamiltonian and a thermodynamic process with a time-independent transition matrix, one may wonder whether there exists a speed limit analogous to Eq. (1), where the bound is given by the distance between the end points divided by a timescale constant [50]. Second, it is tempting to see whether Eq. (4) also holds for the nonadiabatic entropy production. Third, one may

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FIG. 1. (a)  $\Sigma_{na}(\tau)/D_1(p^{(i)}, p^{(f)})$  represented by black dots and  $\Sigma_{na}(\tau/2)/D_1(p^{(i)}, p^{(f)})$  represented by gray circles. (b) Nonadiabatic entropy production  $\Sigma_{na}$  vs corresponding two-state entropy production  $\Sigma_{\Omega_i}$ . The inset depicts a schematic of the pseudo-coarse-graining procedure. (c) Nonadiabatic entropy production  $\Sigma_{na}$  vs dimensionless time duration  $W\tau$ , with the end points fixed. The inset presents the two-state curves for different partitioning schemes, and the horizontal and vertical dashed lines represent the asymptotes. In both (a) and (b), there are 10<sup>5</sup> data points, with each generated as follows. An initial three-state probability distribution is sampled randomly, together with a transition rate matrix with positive entries drawn uniformly from [0,1]. An evolution time  $\tau$  is also drawn uniformly from [0,1]; then the state at  $t = \tau/2$  and the final state are calculated. Using Eq. (7), we compute the nonadiabatic entropy production  $\Sigma_{na}(\tau/2)$  and  $\Sigma_{na}(\tau)$ . We then randomly partition the states into two disjoint nonempty sets and use Eqs. (8) and (7) to calculate  $\Sigma_{\Omega_i}$ . In (c) there are 40 000 data points, each corresponding to a transition from  $p^{(i)} = [0.5, 0.3, 0.2]^T$  to  $p^{(f)} = [0.6, 0.22, 0.18]^T$  with different W and  $\tau$  obtained as follows: We randomly select a transition rate matrix W and an evolution time  $\tau$  as initial guesses and minimize the total-variation distance between  $\exp(W\tau)p^{(i)}$  and  $p^{(f)}$  with the error threshold set to  $10^{-6}$ . The optimal values of W and  $\tau$  are then used to calculate and plot the nonadiabatic entropy production as a function of  $W\tau$ .

inquire the trade-off relation in the form of Eq. (2), if any, between dissipation and time in these relaxation processes. Specifically, it is desirable to obtain a suitable timescale constant.

In this Letter we answer these three questions and show that the relevant timescale constant is the total rate, i.e., the sum of all positive transition rates. We also demonstrate with examples that our dissipation-time trade-off relation outperforms previous bounds.

Speed limit. Consider a stochastic Markov jump process with finite N states. The dynamics of the probability distribution  $\boldsymbol{p} = [p_1, p_2, \dots, p_N]^T$  is described by a Pauli master equation [49]

$$\frac{\mathrm{d}p_m}{\mathrm{d}t} = \sum_n W_{mn} p_n,\tag{5}$$

where  $p_m$  is the probability of state m,  $W_{mn}(m \neq n)$  is the time-independent transition rate from state n to m, and  $W_{mm} = -\sum_{n,n\neq m} W_{nm}$ . For later use, we define the *total rate* as

$$\mathcal{W} = \sum_{m \neq n} W_{mn} = -\mathrm{Tr}\mathbf{W}.$$
 (6)

In experiments, the transition rate matrix and its trace can be inferred from trajectory data [51]. Provided that the Markov chain is ergodic, a steady-state distribution  $p^{(ss)}$  is expected, satisfying  $\sum_{n} W_{mn} p_n^{(ss)} = 0$ . Steady states can be divided into two categories depending on whether they meet the detailed balance condition,  $W_{mn} p_n^{(ss)} = W_{nm} p_m^{(ss)}$ , which is not assumed throughout this Letter. The total entropy production  $\Sigma$  consists of an adiabatic contribution and a nonadiabatic one, both of which are non-negative. Given an initial state  $p^{(i)}$  and a target state  $p^{(f)}$ , the nonadiabatic entropy production is given by [15,52]

$$\Sigma_{\rm na} = D_1(\boldsymbol{p}^{(\rm i)} \| \boldsymbol{p}^{(\rm ss)}) - D_1(\boldsymbol{p}^{(\rm f)} \| \boldsymbol{p}^{(\rm ss)}), \tag{7}$$

where  $D_1(\boldsymbol{p} || \boldsymbol{q}) = \sum_n p_n \ln(p_n/q_n)$  is the 1-Rényi divergence between the two probability distributions [46–48]. For arbitrary dynamics,  $\Sigma \ge \Sigma_{na}$ , and the equality is attained when the detailed balance condition is satisfied. Notably, the detailed balance condition is always fulfilled in two-state systems.

The fixed end points impose constraints on the transition rate matrix W but generally do not uniquely determine it. The choice of W affects the magnitude of entropy  $\Sigma$  and  $\Sigma_{na}$ produced by the stochastic process connecting the same end points over time  $\tau$ . For two-state systems, the steady state  $p^{(ss)}$ is uniquely determined, with  $p_1^{(ss)}$  expressible in terms of  $W\tau$ as [53]

$$p_1^{(\rm ss)} = p_1^{(\rm i)} + \frac{p_1^{(\rm f)} - p_1^{(\rm i)}}{1 - e^{-W\tau}},$$
(8)

with  $p_2^{(ss)} = 1 - p_1^{(ss)}$ .

By substituting Eq. (8) into (7), it can be observed that, given the end points, the nonadiabatic entropy production in a two-state system is a monotonically decreasing function of  $W\tau$ , see the inset of Fig. 1(c). In other words, the  $\Sigma$  versus  $W\tau$  curve indicates a trade-off between entropy production and time duration and is bounded by a vertical and a horizontal asymptote. The horizontal asymptote signifies that the minimum of the entropy production is reached when  $W\tau \to \infty$ . In this limit,  $p^{(ss)} = p^{(f)}$ , so Eq. (7) implies that the minimum entropy production is given by (4).

It is unphysical for the population  $p_1^{ss}$  to be 0 or 1, as this would require one of the transition rates to vanish. Combining Eq. (8) with this constraint ( $0 < p_1^{(ss)} < 1$ ), we obtain a speed

limit represented by the vertical asymptote, i.e., a lower bound for the evolution time  $\tau$ :

$$\tau > \frac{\max\left\{\ln\left(p_1^{(i)}/p_1^{(f)}\right), \ln\left(p_2^{(i)}/p_2^{(f)}\right)\right\}}{\mathcal{W}}.$$
 (9)

The numerator coincides with the  $\infty$ -Rényi divergence, defined as  $D_{\infty}(\boldsymbol{p}||\boldsymbol{q}) = \max_{n} \{\ln p_{n}/q_{n}\}$  [46,47]. Specifically, the lower bound on evolution time is given by the quotient of the  $\infty$ -Rényi divergence between the two end points and the total rate of transitions  $\mathcal{W}$ .

This lower bound reflects an information-theoretical limit on speed in two-state systems; hence, it is relevant to examine whether this bound can be extended to generic *N*-state systems. The answer is affirmative. We state our main result here while deferring the proof to the end of the Letter: the time it needs to evolve from the initial state  $p^{(i)}$  to the final state  $p^{(f)}$ is bounded from below by

$$\tau > \frac{D_{\infty}(\boldsymbol{p}^{(i)} \| \boldsymbol{p}^{(f)})}{\mathcal{W}},\tag{10}$$

where  $D_{\infty}(p^{(i)}||p^{(f)})$  is the  $\infty$ -Rényi divergence between the distributions  $p^{(i)}$  and  $p^{(f)}$ , and W is the total rate. Remarkably, this lower bound on evolution time resembles the distance-based formulation in Eq. (1), where a constant rate characterizes the timescale. It is noteworthy that the  $\infty$ -Rényi divergence plays a significant role in the resource theory of thermodynamics [47,54,55], and its application to speed limits highlights its versatility and significance in various physical contexts. If  $p^{(f)} = p^{(ss)}$ , the time it needs to reach the steady state is infinite, so this bound is trivially satisfied. When the final state  $p^{(f)}$  is not of full rank, both the  $\infty$ -Rényi divergence and the bound of time diverge. This is a manifestation of the third law of thermodynamics, which states that nonfull rank states cannot be attained in finite time [56].

Information-theoretical bound on nonadiabatic entropy production. Building upon the information-theoretical bound on entropy production described in Ref. [45] for systems with detailed balance, we extend our findings to scenarios without detailed balance. Our second result introduces a conjecture (referred to as divergence bound):

$$\Sigma_{\mathrm{na}} \geqslant D_1(\boldsymbol{p}^{(\mathrm{i})} \| \boldsymbol{p}^{(\mathrm{f})}). \tag{11}$$

This conjecture, presented as an extension of Eq. (4), suggests a lower bound on the nonadiabatic entropy production given by the 1-Rényi divergence between the end points. Figure 1(a) provides numerical evidence for three-state systems to support it, and the code used to generate numerical verification for other numbers of states can be found in the Supplemental Material [53]. In fact, we confirm a tighter bound,  $\Sigma_{na}(\tau/2) \ge D_1(p^{(i)}||p^{(f)})$ , where  $\Sigma_{na}(\tau/2)$  represents the nonadiabatic entropy production during the time interval  $[0, \tau/2]$ , which strengthens our conjecture. Our findings have implications for the excess entropy production  $\Sigma_{ex}$ , which is defined by a variational principle and is always greater than  $\Sigma_{na}$  [57,58]. As a consequence, we observe the following relationships:

$$\Sigma_{\text{ex}}(\tau) \geqslant \Sigma_{\text{ex}}(\tau/2) \geqslant D_1(\boldsymbol{p}^{(i)} \| \boldsymbol{p}^{(f)}), \quad (12)$$

as reported in Refs. [57,58].

*Trade-off between dissipation and time.* Let us state the third main result: for *N*-state dynamics evolving from  $p^{(i)}$  to  $p^{(f)}$  during  $\tau$  with a total rate W, the conjecture is that the entropy production  $\Sigma$  is bounded from below by (referred to as trade-off bound)

$$\Sigma_{\mathrm{na}} \geqslant \Sigma_{\Omega_i}(\mathcal{W}\tau) \quad i = 1, 2, \dots, S_N^{(2)}. \tag{13}$$

Here,  $\Sigma_{\Omega_i}(\mathcal{W}\tau)$  is the entropy production of two-state dynamics obtained from a pseudo-coarse-graining procedure as follows: partition the set of N states into two nonempty sets where the first set is denoted by  $\Omega_i$ , then we have the two end points  $\boldsymbol{P}_{\Omega_i}^{(i)} = [\sum_{n \in \Omega_i} p_n^{(i)}, 1 - \sum_{n \in \Omega_i} p_n^{(i)}]^T$  to  $\boldsymbol{P}_{\Omega_i}^{(f)} = [\sum_{n \in \Omega_i} p_n^{(f)}, 1 - \sum_{n \in \Omega_i} p_n^{(f)}]^T$ . We then use Eq. (8) to find the steady-state distribution, where  $\mathcal{W}$  and  $\tau$  of the original N-state dynamics are used. The entropy production of the two-state dynamics is obtained by substituting the steadystate distribution into (7). The inset of Fig. 1(b) gives a schematic of the pseudo-coarse-graining procedure for three states. As shown in the upper panel, the system evolves from  $\boldsymbol{p}^{(i)} = [0.5, 0.17, 0.33]^{T}$  to  $\bar{\boldsymbol{p}}^{(f)} = [0.1, 0.4, 0.5]^{T}$ , visually depicted by bars of varying colors and lengths. There are three partitioning schemes, and the lower panel shows one in which the second and third state are grouped together, resulting in  $\boldsymbol{P}_{\Omega_1}^{(i)} = [0.5, 0.5]^{\mathrm{T}}$  and  $\boldsymbol{P}_{\Omega_1}^{(i)} = [0.1, 0.9]^{\mathrm{T}}$ . Given N states, we have  $S_N^{(2)}$  different ways of partitioning the N states into two nonempty sets, where  $S_N^{(2)}$  is the Stirling number of the second kind [59]. Explicitly,  $S_N^{(2)} = 2^{N-1} - 1$ . For example, four states can be partitioned in  $S_4^{(2)} = 7$  ways, where  $\Omega_i = \{1\}, \{2\}, \{3\}, \{4\}, \{1, 2\}, \{1, 3\}, \{1, 4\}$ , with the elements in each set being the indices of states. Thus there are  $S_N^{(2)}$  different two-state bounds in total, jointly bounding  $\Sigma_{na}$ . As discussed above, these two-state bounds are monotonically decreasing function of  $W\tau$ , which has the form of Eq. (2), with  $\mathcal{W}$  playing the role of  $\mathcal{R}$ . We stress that the two-state dynamics cannot be obtained from the standard coarse-graining procedure [60,61]: the population  $P_{\Omega}(t)$  is in general not equal to  $\sum_{n\in\Omega} p_n(t)$  except at the two end points; the transition rate matrix for each two-state dynamics is time independent, and its entries are not simple linear combinations of the original transition rates. In the following we consider the three-state case as an example, and numerical evidence to support Eq. (13) for other numbers of states can be generated using the code in the Supplemental Material [53].

As a direct verification of the inequality (13), Fig. 1(b) shows a plot of the nonadiabatic entropy production  $\Sigma_{na}$  versus the corresponding two-state  $\Sigma_{\Omega_i}$ . All data points lie above the diagonal, confirming the new bound given by (13). In the long- $\tau$  limit, the entropy production of the two-state dynamics is  $D_1(\boldsymbol{P}_{\Omega_i}^{(i)} || \boldsymbol{P}_{\Omega_i}^{(f)})$ . By applying the theorem that refinement cannot decrease divergence [62], which is essentially the log-sum inequality, this quantity is not greater than  $D_1(\boldsymbol{p}^{(i)} || \boldsymbol{p}^{(f)})$  and also  $\Sigma_{na}$ . This leads to two questions: (1) Can the bound ever be saturated, as it is not evident from Fig. 1(a)? (2) Is the new bound ever tighter than the divergence bound and the activity bound, or is it otherwise redundant?

There are at least two simple cases in which the bound is (nearly) saturated. By examining the condition for equality



FIG. 2. Nonadiabatic entropy production  $\Sigma_{na}$ , corresponding divergence bound (4), activity bound (3), and the trade-off bound (13) as a function of the dimensionless time Wt. There are three ways to partition a set with three states: one of the two nonempty sets is  $\Omega_i = \{1\}, \{2\}, \text{ and } \{3\}, \text{ respectively.}$ 

in the log-sum inequality, it can be observed that as long as a subset  $\Omega$  of the states such that  $p_m^{(i)}/p_m^{(ss)}$  are equal for all  $m \in \Omega$ , the bound is saturated in the long- $\tau$  limit. Another scenario in which the bound is nearly saturated throughout the entire process is when all the N - 2 states' initial populations and the transition rates between them are vanishingly small, effectively reducing the dynamics to only two states. This case also partially addresses the second question, as the divergence bound is not generally saturated. The superiority of the new bound over the previous bounds is also evident in nontrivial cases, as will be seen in Figs. 1(c) and Fig. 2. We will also prove that there always exists a parameter range in which the trade-off bound outperforms the divergence bound for any given pair of end points.

Figure 1(c) displays a representative scenario where the end points are given and fixed, while also demonstrating universal behavior. Equations (4), (11), and (13) jointly bound  $\Sigma_{na}$ , as represented by the shaded area. The trade-off between dissipation and evolution time quantified by Eq. (13) is clearly exemplified in this figure. It demonstrates that, when the end points are fixed, faster state transformations are accompanied by larger amounts of dissipation. As shown in the inset, each pseudo-coarse-grained two-state dynamics also has a speed limit, and the log-sum inequality implies that it is not greater than the speed limit of the N-state dynamics. The vertical asymptote of the two-state curve for  $\Omega_3$  is exactly the speed limit for three states, and this is not a coincidence. For any given pair of end points, there must be at least one two-state curve whose vertical asymptote coincides with the genuine speed limit for N states. Thus there must exist a parameter range in which the trade-off bound outperforms the divergence bound.

Figure 2 shows the entropy production  $\Sigma$  and the corresponding divergence bound, activity bound, and the new bound as a function of the dimensionless time Wt for a representative case whose initial distribution and transition rate matrix (in arbitrary units) are

given by

$$\boldsymbol{p}^{(i)} = \begin{bmatrix} 1/3 & 1/3 & 1/3 \end{bmatrix}^{\mathrm{T}},$$
$$\boldsymbol{W} = \begin{bmatrix} -1.60 & 0.20 & 0.10 \\ 0.70 & -0.80 & 0.70 \\ 0.90 & 0.60 & -0.80 \end{bmatrix}.$$
(14)

At each time instant t, we calculate the instantaneous distribution  $p(t) = \exp(Wt)p^{(i)}$  and calculate the divergence bound and the new bound using corresponding equations by replacing  $p^{(f)}$  therein with p(t). The time-averaged dynamical activity is calculated using Eq. (8) in Ref. [15], with the upper bound of the integral set to t. Hence the time derivative of the divergence bound, the activity bound, and the new bound cannot be regarded as entropy production rate and are not guaranteed to be non-negative. The divergence bound is initially loose but saturates as  $t \to \infty$ , as expected. The activity bound has a much better performance than the divergence bound in the beginning but gradually loses its advantage. The activity bound even decreases as the steady state is approached, showing the expected  $1/\tau$  asymptotic behavior as previously mentioned. The new bound with  $\Omega_1 = \{1\}$  has a performance that is almost all the time better than both the divergence bound and the activity bound. As discussed above, in the long-time limit the divergence bound's performance should be the best as it saturates. For this special case, the tradeoff bound is almost as good, as can be numerically verified. The steady-state distribution is  $p^{(ss)} = [0.086, 0.467, 0.447]^{T}$ , and the divergence bound is 0.2405. The trade-off bound in the long-time limit is given by the divergence between  $[1/3, 2/3]^{T}$  and  $[0.086, 0.914]^{T}$ , which is 0.2403.

Proof of the speed limit. Let us prove that Eq. (10) gives a speed limit for relaxation processes in generic *N*-state systems, where the denominator is still the total rate  $\mathcal{W} \equiv -\text{Tr} \boldsymbol{W}$ , i.e., the sum of all the positive rates. First of all, let us divide  $[0, \tau]$  into  $\mathcal{K} \gg 1$  intervals so a time sequence  $t_0 = 0 < \cdots < t_k < \cdots < t_{\mathcal{K}} = \tau$  is obtained. Consider an arbitrarily selected infinitesimal interval  $[t_k, t_{k+1}]$ , then the master equation (5) gives

$$\delta p_m^{(k)} = (t_{k+1} - t_k) \sum_n W_{mn} p_n^{(k)}, \qquad (15)$$

where  $\delta p_m^{(k)} = p_m^{(k+1)} - p_m^{(k)}$ . Without loss of generality, we assume the state l satisfies  $\ln (p_l^{(k)}/p_l^{(k+1)}) = \max_n \{ \ln (p_n^{(k)}/p_n^{(k+1)}) \}$ . Therefore

$$t_{k+1} - t_k = \frac{\delta p_l^{(k)}}{W_{ll} p_l^{(k)} + \sum_{n \neq l} W_{ln} p_n^{(k)}}.$$
 (16)

If the sum of all positive rates is fixed, it is not hard to see that if  $W_{ll} = -W$  with all the other transition rates vanishing,  $t_{k+1} - t_k$  reaches the minimum,

$$t_{k+1} - t_k \ge \frac{\delta p_l^{(k)}}{(-\mathcal{W})p_l^{(k)}} \approx \frac{1}{\mathcal{W}} \ln \frac{p_l^{(k)}}{p_l^{(k+1)}}$$
$$= \frac{1}{\mathcal{W}} \max_{1 \le n \le N} \left\{ \ln \frac{p_n^{(k)}}{p_n^{(k+1)}} \right\}, \tag{17}$$

where  $\approx$  is exact to first order in  $\delta p_l^{(k)}/p_l^{(k)}$ . Technically speaking, the equality is not achievable in physical processes because  $W_{ll} = -W$  is impractical. Summing over k results in

$$\mathcal{W}\tau > \sum_{k=0}^{\mathcal{K}-1} \max_{1 \le n \le N} \left\{ \ln \frac{p_n^{(k)}}{p_n^{(k+1)}} \right\}.$$
 (18)

The inequality that the maximum of sum is at most the sum of maxima gives

$$\sum_{k=0}^{K-1} \max_{1 \le n \le N} \left\{ \ln \frac{p_n^{(k)}}{p_n^{(k+1)}} \right\} > \max_{1 \le n \le N} \left\{ \sum_{k=0}^{K-1} \ln \frac{p_n^{(k)}}{p_n^{(k+1)}} \right\}$$
(19)
$$= \max\{ \ln(p_n^{(i)}/p_n^{(f)}) \}.$$

Combining with (18) completes the proof.

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Conclusion. We present a speed limit, a bound on the nonadiabatic entropy production and a trade-off relation between dissipation and evolution time for time-independent relaxation processes. Our result in Eq. (10) resembles the quantum speed limit (1) and indicates that the minimum transition time from an initial to a target state is constrained by the ratio of their  $\infty$ -Rényi divergence to the total rate. Equation (11) gives a lower bound on the nonadiabatic entropy production in terms of 1-Rényi divergence. Furthermore, Eq. (13), implicitly in the form of Eq. (2), reveals a new trade-off between dissipation and time that surpasses the divergence bound for certain parameters. Given successful quantum extensions of both divergence and activity bounds [6,63–65], we anticipate that our results can also be generalized to quantum settings.

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