A new route toward orthogonality

Andrea Valdés-Hernández* and Francisco J Sevilla

Instituto de Física, Universidad Nacional Autónoma de México, Apdo. Postal 20-364, 01000, Ciudad de México, México

E-mail: andreavh@fisica.unam.mx and fjsevilla@fisica.unam.mx

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Abstract
We revisit the problem of determining conditions under which a pure state, that evolves under an arbitrary unitary transformation, reaches an orthogonal state in a finite amount of the transformation parameter. Simple geometric considerations disclose the existence of a fundamental limit for the minimal amount required, providing, in particular, an intuitive hint of the Mandelstam–Tamm bound. The geometric considerations leads us to focus on a particular, yet relevant, family of states that evolve toward orthogonality. Several dynamical features are discussed, which include the (relative) entropy production during transformation, and special attention is paid to multipartite systems of N bosons that are allowed to tunnel between two sites. The effects of the tunneling in the amount of transformation required for the system to attain an orthogonal state are revealed, and the relation between the latter, the tunneling intensity and the mode-entanglement is explored.

Keywords: quantum speed limit, entanglement, Bose Hubbard Hamiltonian

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

1. Introduction

The reliable manipulation of quantum states to reach particular target states lies at the core of quantum information science [1, 2]. Particularly, the determination of the amount of transformation required to take a specific state to an orthogonal—distinguishable—one, acquires relevance in quantum information tasks, and has received attention under different contexts, mainly in relation with the so-called quantum speed limit, in which the minimal time required to evolve between two states separated by a given ‘distance’ is considered [3–14]. Investigations regarding the amount of transformation to take a state into an orthogonal one, are also significant in the broader theoretical and experimental analysis of the dynamic evolution of

*Author to whom any correspondence should be addressed.
quantum states, in the establishment of fundamental physical limitations (usually expressed as uncertainty relations), as well as in other applications of quantum physics [15]. Moreover, ongoing theoretical analysis on this topic are still revealing new aspects on its connection with several quantum features [16–21].

In this paper we revisit the problem of determining the amount of transformation required for an initial pure state to reach an orthogonal one under an arbitrary unitary transformation. Our analysis allows to unveil, clarify, and identify different novel aspects related with the dynamics toward orthogonality, and adds to preceding works in several ways. In particular, in section 2, we present a fresh discussion—based on associating the overlap between two states with the superposition of two-dimensional rotating vectors—that brings out the notion of ‘speed limit’ in a most natural geometrical way, and particularly hints out on the Mandelstam–Tamm bound in an intuitive fashion. Further, the amount of evolution (and its lower bound) toward orthogonality is determined for a particular, yet important, family of initial states. This amount defines a characteristic scale for which the evolution leads to ‘clock’ models [22], and the dynamics between the ‘tickings’ is analyzed from an entropic perspective.

In section 3 special attention is paid to multipartite systems of $N$ bosons allowed to tunnel between two sites (or rather $N$ qubits in an effective Hilbert space of dimension $N + 1$). The ‘slowest’ and ‘fastest’ states in reaching an orthogonal state are identified. In the non-tunneling regime, a simple relation between the amount of transformation needed to attain orthogonality and the states’ mode-entanglement is obtained. The effects of the transition between sites are also explored, showing that the tunneling favors a ‘faster’ evolution toward orthogonality, and that variations of the tunneling intensity gives rise to a rich dynamics of the entanglement between modes (as opposed to the extensively studied case of entanglement between particles). Section 4 presents a summary and some final remarks.

Our results enrich the current comprehension on the dynamics of quantum systems, and in particular of multipartite entangled systems which are of extended theoretical and experimental interest [23].

2. The geometrical shortcut to orthogonality

2.1. The geometrical picture

We consider a pure-state vector in a Hilbert space $\mathcal{H}$ that evolves under the unitary transformation

$$\hat{U}_\gamma(\gamma) = e^{-i\hat{g} \gamma},$$

with $\hat{g}$ an Hermitian operator, and $\gamma$ a real and continuous parameter that defines the $\gamma$-translation symmetry transformation. The initial (normalized) state is expanded according to

$$|\psi_0\rangle = \sum_{k=1}^{N} \sqrt{r_k} e^{i\varphi_k} |g_k\rangle,$$

so the evolved state reads

$$|\psi_\gamma\rangle = \hat{U}_\gamma(\gamma) |\psi_0\rangle = \sum_{k=1}^{N} \sqrt{r_k} e^{i(\varphi_k - g_k \gamma)} |g_k\rangle,$$

where $\hat{g} | g_k \rangle = g_k | g_k \rangle$, $\varphi_k \in [0, 2\pi]$, and the coefficients $r_k \in [0, 1]$ satisfy the normalization condition

$$\sum_{k=1}^{N} r_k = 1,$$
thus \( \{ r_k \} \) conforms the probability distribution from which the average \( \langle F(\hat{g}) \rangle = \sum_{k=1}^{N} r_k F(\hat{g}_k) \) is computed. This distribution has associated the information (Shannon) entropy \( S_{\text{SH}}[\{ r_k \}] = -\sum_{k=1}^{N} r_k \) which measures the lack of information about the state \( |\psi_\gamma \rangle \) relative to the basis \( \{ |g_k \rangle \} \). Its minimum value \( (S_{\text{SH}} = 0) \) is attained when there is maximal information, corresponding to \( r_k = g_k \), whereas its maximum value \( (S_{\text{SH}} = 1) \) occurs when the available information is minimal, so the distribution is \( r_k = 1/N, \) in line with the equal a priori probability principle.

One quantity of interest that characterizes the evolution of pure states is the overlap

\[
\langle \psi_\gamma | \psi_0 \rangle = \sum_{k=1}^{N} r_k e^{i\hat{g}_k \gamma} = \langle e^{i\hat{g} \gamma} \rangle,
\]

which is a \( \gamma \)-dependent complex number that measures how similar is the transformed state \( |\psi_\gamma \rangle \) with respect to the original one, \( |\psi_0 \rangle \). Equation (4) vanishes provided \( |\psi_\gamma \rangle \) is orthogonal to (distinguishable from) \( |\psi_0 \rangle \). A relative entropy between such states can be defined in terms of (4) as \( S_{\text{rel}} \{ \{ r_k \}, \gamma \} = 1 - \langle \langle \psi_\gamma | \psi_0 \rangle \rangle^2 \), with \( |\langle \psi_\gamma | \psi_0 \rangle |^2 \) known as the non-decay or survival probability whenever \( \gamma \) and \( \hat{g} \) are identified, respectively, with time and the Hamiltonian operator \( \hat{H} \) (we take \( h = 1 \)). The relative entropy acquires its maximum \( (S_{\text{rel}} = 1) \) and minimum \( (S_{\text{rel}} = 0) \) value when the transformed state is orthogonal and equal (up to a phase factor) to the original one, respectively.

Each term of the expansion in (4) can be mapped into the two-dimensional vector \( r_k(\gamma) = r_k(0) \cos \theta_k, \sin \theta_k \) lying inside the unitary circle \( (r_k \leq 1) \) with \( \theta_k = g_k \gamma \). Consequently \( \langle \psi_\gamma | \psi_0 \rangle \) is mapped into the vector \( \hat{R}(\gamma) = \sum_{k=1}^{N} r_k(\gamma) \), and thus the overlap vanishes if and only if \( \hat{R} = 0 \). This observation offers a simple geometrical prescription to guarantee that \( |\psi_0 \rangle \) evolves toward an orthogonal state, namely: sum up the vectors \( r_k(\gamma) \), each performing uniform circular motion with frequency \( g_k = d\theta_k/d\gamma \) along orbits of radius \( r_k \), and from this sum, determine the conditions under which the vector’s tips are distributed such that the centroid of the polygon they form coincides with the origin (this is the case for example, when the tips form a regular polygon centered at the origin). From this geometrical picture it follows immediately that:

(a) The number of vectors \( r_k \) equals \( N \), the number of states that contribute in the expansion (2).
(b) Each of the \( N \) phase oscillators is accommodated in one of the \( R \) different radii \( (1 \leq R \leq N) \), with \( R \) the number of different coefficients \( r_k \).
(c) The number of frequencies equals the number of different eigenvalues \( g_k \) appearing in (2).
(d) At least two vectors with different frequencies, i.e., two different eigenvalues \( g_k \), must contribute in (2).
(e) The value of \( \gamma \) at which this occurs, denoted as \( \tilde{\gamma} \), must be larger than or equal to a certain minimum value \( \tilde{\gamma} \), that is, \( \tilde{\gamma} \geq \tilde{\gamma} > 0 \).

From (d) it follows that a nonzero value of the \( (\gamma\text{-invariant}) \) dispersion \( \sigma_\gamma^2 = \langle \hat{g}^2 \rangle - \langle \hat{g} \rangle^2 \) is necessary for orthogonality to be attained. Since this quantity is equivalent to the variance of the eigenvalues \( \{ g_k \} \) computed with the distribution \( \{ r_k \} \), the latter plays a fundamental role
Figure 1. Left panel: Initially ($\gamma = 0$) all $r_k$ ($k = 1, 2, 3$) lie along the horizontal axis, and for $\gamma > 0$ they perform a counter-clockwise rotation with $\theta_k = g_k \gamma$. For $0 < \gamma < \tilde{\gamma}$ clearly $R(\gamma) = \sum_{k=1}^{3} r_k(\gamma)$ is not a vanishing vector. Right-panel: at $\gamma = \tilde{\gamma}$ the tips of the vectors form a polygon with its centroid at the origin, hence the vector $R(\tilde{\gamma})$ vanishes, indicating that the system has reached a state orthogonal to $|\psi_0\rangle$.

in attaining the orthogonal state. In particular, a distribution $\{r_k\}$ with nonzero $S_{5h}$ is required for orthogonality to be reached. On the other hand, (e) implies that if $\langle \psi_\gamma | \psi_0 \rangle$ vanishes, then a sort of speed limit—or $\gamma$-$\hat{g}$ uncertainty relation—exists, determined only by $\gamma_{\text{min}}$. When the unitary transformation (1) refers to the temporal evolution, that is, when $\hat{g}$ and $\gamma$ correspond, respectively, to a time-independent Hamiltonian and to time, these intuitive conditions become formally stated and quantified in the form of the Mandelstam–Tamm bound for the so-called quantum speed limit, or the time–energy uncertainty relations [4, 5].

Notice that according to the present geometrical argument, the existence of a speed limit originates in the fact that the dynamics of each term in the overlap (4) is performed (uniformly) on a circle, and requires no explicit quantum hypothesis. Interestingly, speed limits based on the dynamics on the circle [24] can also found, for example, in the emergence of collective synchronization of interacting phase oscillators [25]. These observations are pertinent in relation with recent works that point toward an ubiquitous nature of the limitations in the speed of evolution, by analyzing the existence of speed limits in the classical realm [26–28].

2.2. Lower bound for the ‘speed’ limit

The problem of determining the conditions on the distribution $\{r_k\}$ and the set $\{g_k\}$ for $|\psi_0\rangle$ to reach an orthogonal state at some finite $\tilde{\gamma}$ is an interesting one [14, 29], yet it represents a difficult task in the most general case. However, we can make progress either by assuming a particular distribution $\{r_k\}$ and determining the requirements on $\{g_k\}$, or the other way around. We shall adopt the first strategy and consider a distribution $\{r_k\}$ that fulfills the equal a priori probability principle. Thus, in what follows we will consider the family $I$ of initial states that are equally-weighted superpositions of eigenvectors of $\hat{g}$, i.e., the collection of states of the form

$$|\psi_0\rangle = \frac{1}{\sqrt{N}} \sum_{k=1}^{N} e^{i\phi_k} |g_k\rangle,$$

(5)
where we have chosen \( r_k = 1/N \), with finite \( N \geq 2 \) (the case \( N = 1 \) clearly never reaches an orthogonal state). This is a natural choice for the distribution \( \{ r_k \} \), assuming we have the minimal information, hence maximal Shannon entropy. We further assume that all the \( g_k \) appearing in the expansion (2) are different, and that (without loss of generality) the index \( k \) is such that \( g_k < g_{k+1} \). According to (b) above, this means that the tips of the \( N \) vectors lie on the circle of radius \( 1/N \), they rotate with different angular frequencies \( g_1 < \cdots < g_k < \cdots < g_N \), and after a transformation by \( \gamma \) units they point at the directions \( \theta_1 < \cdots < \theta_k < \cdots < \theta_N \), respectively.

From this geometrical description, it is clear that an orthogonal state is reached for the first time at \( \tilde{\gamma} \)—that is, \( \langle \psi_\tilde{\gamma} | \psi_0 \rangle = 0 \) or equivalently, \( R(\tilde{\gamma}) = 0 \)—whenever the tip vectors form (for the first time) the \( N \) vertices of a regular polygon inscribed in the circle of radius \( 1/N \), as illustrated in figure 2 for \( N = 5 \). In this case it holds that \( \Delta \theta = \theta_{k+1} - \theta_k = 2\pi/N \), meaning that

\[
\tilde{\gamma} = \frac{2\pi}{\Delta g} \frac{1}{N}, \tag{6a}
\]

where

\[
\Delta g = g_{k+1} - g_k \tag{6b}
\]

is independent of \( k \), which occurs only if the elements of the set \( \{ g_k \} \) are of the form

\[
g_k = (k - 1)\Delta g + g_1. \tag{7}
\]

With this, all the \( g \)-distributed moments, \( \langle g^n \rangle = \sum_{k=1}^{N} r_k g^n_k = N^{-1} \sum_{k=1}^{N} g^n_k \), (which are not all independent among themselves) are explicitly determined in terms of \( g_1, N \), and \( \Delta g \) only.

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1 The condition of the regular polygon is sufficient, although not necessary to attain orthogonality; however, for \( N = 2 \) it is a necessary and sufficient condition. Indeed, for \( N = 2 \) there are only two vectors that perform phase rotation, and for these to sum up a null vector both radii must necessarily be equal (i.e., \( r_1 = r_2 = 1/2 \)), and one vector opposite to the other forming a diameter of the circle (the ‘polygon’ in this case). Consequently, an equally weighted qubit state always reaches an orthogonal state at \( \tilde{\gamma} = \pi/(g_2 - g_1) \), as follows from equation (6a).
Thus, different, yet equivalent, expressions for $\tilde{\gamma}$ are possible when this is written in terms of any of the moments of $\hat{g}$. Of particular interest are those expressions for $\tilde{\gamma}$ that involve the first two moments [9, 30] since these play the most determinative role in the transformation (1) [9, 22].

In terms of $\langle \hat{g} \rangle - g_1$, the mean value of $\hat{g}$ as measured from the lowest eigenvalue appearing in equation (5), equation (6a) rewrites as

$$\tilde{\gamma} = \frac{\pi}{2(\langle \hat{g} \rangle - g_1)} \geq \frac{\pi}{2(\langle \hat{g} \rangle - g_1)}.$$  (9)

whereas in terms of the standard deviation $\sigma_{\hat{g}} = \sqrt{\langle (\hat{g} - \langle \hat{g} \rangle)^2 \rangle} = \sqrt{\langle \hat{g}^2 \rangle - \langle \hat{g} \rangle^2}$, $\gamma$ can be written as

$$\tilde{\gamma} = \frac{\pi}{2\sigma_{\hat{g}}} \sqrt{\frac{4(N^2 - 1)}{3N^2}} \geq \frac{\pi}{2\sigma_{\hat{g}}}. $$  (10)

The inequalities in equations (9) and (10) are, respectively, the extension to the $\gamma$-translation transformation of the Margolus–Levitin [7] and the Mandelstam–Tamm [5] bounds that determine the quantum speed limit. In particular, equation (10) leads to the inequality $\sigma_{\hat{g}} \tilde{\gamma} \geq \pi / 2$, which plays the role of a $\gamma$-$\hat{g}$ “uncertainty” relation [4]. Moreover, the quantity $s = 2\sigma_{\hat{g}} \tilde{\gamma}$ is independent of the particular generator $\hat{g}$ that defines the unitary transformation (1), and thus corresponds to the analogousof a geometric phase as discussed in reference [22].

The bounds indicated in equations (9) and (10) have been synthesized in the following expression for the lower bound of $\tilde{\gamma}$ [31, 32]

$$\tilde{\gamma}_{\text{min}} = \max \left\{ \frac{\pi}{2(\langle \hat{g} \rangle - g_1)}, \frac{\pi}{2\sigma_{\hat{g}}} \right\}. $$  (11)

The quantities inside the brackets are not necessarily independent of each other, and as shown by Levitin and Toffoli [9], the bound (11) can be attained only when $\langle \hat{g} \rangle - g_1 = \sigma_{\hat{g}}$, and is asymptotically attainable otherwise.

Now, the equalities in equations (9) and (10) lead to

$$\frac{\sigma_{\hat{g}}}{(\langle \hat{g} \rangle - g_1)} = \frac{1}{\sqrt{\frac{3}{N}}(N - 1)} \leq 1,$$

thus disclosing the relation between $\langle \hat{g} \rangle - g_1$ and $\sigma_{\hat{g}}$ (which is of course particular for the specific distribution $\{ r_k \}$ under consideration), and accordingly for the family of initial states $\mathcal{I}$ we have

$$\tilde{\gamma}_{\text{min}} = \frac{\pi}{2\sigma_{\hat{g}}}. $$  (13)

The quotient in equation (12) approaches $1/\sqrt{3} \approx 0.57735$ as $N \to \infty$ meaning that the fluctuations are not negligible even in the regime of large systems. On the other hand, the quotient in (12) saturates the inequality for $N = 2$, and it is only for this value of $N$ that
the Margolus–Levitin and the Mandelstam–Tamm bounds coincide and \( \tilde{\gamma} \) attains its minimum value, as follows from equation (10) (thus recovering the main statement established in theorem 1 of reference [9]). Therefore, for fixed \( \langle \tilde{g} \rangle - g_1 \) (or \( \sigma_k \)), a superposition involving two states only (an effective qubit), reaches an orthogonal state faster than any other superposition. In contrast, for the same \( \langle \tilde{g} \rangle - g_1 \) we have that \( \sigma_k < \langle \tilde{g} \rangle - g_1 \) for \( \mathcal{N} > 2 \), and thus \( \tilde{\gamma} \) asymptotically approaches the value \( (2/\sqrt{3})\gamma_{\min} \) with \( \mathcal{N} \to \infty \).

Throughout the paper, when we say that a given state evolves ‘faster’ than other one, we are tacitly considering the absolute amount of evolution \( \gamma \) (a universal parameter that characterizes the pace of change of all systems’ states on equal footing), so that a state \( |\psi_a\rangle \) is said to reach orthogonality faster than \( |\psi_b\rangle \) whenever \( \tilde{\gamma}_a < \tilde{\gamma}_b \). However, it should be bear in mind that other state-dependent scales can be considered, leading to different and relative concepts of speed of evolution that are also useful in characterizing the dynamics of the system. One example is the ratio \( \tilde{g} = \gamma / \gamma_{\min} \), which quantifies the amount of evolution relative to the minimal amount imposed by the system’s own dynamical limitations. In this case, \( |\psi_a\rangle \) is said to be faster than \( |\psi_b\rangle \) if \( \tilde{g}_a < \tilde{g}_b \), with

\[
\tilde{g} = \frac{\tilde{\gamma}}{\gamma_{\min}} \geq 1. \tag{14}
\]

Another notion of speed of evolution ensues by focusing on the quantity \( \gamma_{\min} \), according to which the faster state is that with the minimum quantum speed limit. Notice, however, that this figure of merit does not necessarily compare the amount of transformation required to actually reach orthogonality, because (as stated above) the bound \( \gamma_{\min} \) is in general not attainable.

Since \( \gamma_{\min} \) varies in general from state to state, it may occur that \( \tilde{\gamma}_a < \tilde{\gamma}_b \) but \( \tilde{g}_b < \tilde{g}_a \), so clearly the notion of ‘faster’ depends on the quantity employed to measure the speed of evolution. In particular, for the family of states considered above, equations (10) and (13) lead to

\[
\tilde{g} = \frac{2}{\sqrt{3}} \sqrt{\frac{\mathcal{N}^2 - 1}{\mathcal{N}^2}}, \tag{15}
\]

so \( \tilde{g} \) depends only on \( \mathcal{N} \), whereas \( \tilde{\gamma} \) depends on \( \mathcal{N} \) and also on the separation \( \Delta \tilde{g} \), as seen from equation (6a). Therefore, for example, two states \( |\psi_a\rangle \) and \( |\psi_b\rangle \) having the same number of terms in their expansion (5) are equally fast in terms of \( g \), yet in terms of the absolute parameter \( \gamma \) the state \( |\psi_a\rangle \) is faster provided \( \Delta g_b < \Delta g_a \), that is, provided the separation between its states is larger.

2.3. Evolution between orthogonal states

Besides determining the value for which the transformed state is orthogonal to the initial one, \( \tilde{\gamma} \) also defines a collection of \( \mathcal{N} \) mutually orthonormal states. First notice that for integers \( n, m = 0, 1, \ldots, \mathcal{N} - 1 \), the evolution operator (1) implies that

\[
\langle \psi_m | \psi_{m'} \rangle = \langle \psi_{(n-m)} | \psi_m \rangle = \sum_{k=1}^{\mathcal{N}} r_k e^{i k_1 (n-m) \gamma}, \tag{16}
\]

In the present case we have \( r_k = 1/\mathcal{N} \), \( g_k \) is given by equation (7), and \( \tilde{\gamma} \) is given by equation (6a). Gathering these results equation (16) becomes

\[
\langle \psi_m | \psi_{m'} \rangle = e^{i k_1 (n-m) \gamma} \frac{1}{\mathcal{N}} \sum_{l=0}^{\mathcal{N}-1} e^{i 2 \pi l (n-m) / \mathcal{N}} = e^{i k_1 (n-m) \gamma} \delta_{mm} = \delta_{mm}, \tag{17}
\]
Figure 3. The relative entropy $S_{\text{rel}} = 1 - |\langle \psi_\gamma | \psi_0 \rangle |^2$ as a function of $\gamma/\Gamma$, with $\Gamma = N\tilde{\gamma}$, for $N = 2$ (blue, dash-dotted line), 3 (red, dashed line) and 10 (pink, solid line). $S_{\text{rel}}$ attains its maximum value 1 each time an orthogonal state is reached, and becomes 0 only when the initial state is recovered, at $\gamma = N\tilde{\gamma}$. For sufficiently large $N$, the state $|\psi_\gamma \rangle$ is highly distinguishable from $|\psi_0 \rangle$ in the entire interval $\tilde{\gamma} \leq \gamma \leq (N-1)\tilde{\gamma}$.

meaning that $O = \{|\psi_0 \rangle, |\psi_{\tilde{\gamma}} \rangle, \ldots, |\psi_{(N-1)\tilde{\gamma}} \rangle \}$ is indeed a set of $N$ mutually orthonormal states, obtained by the successive application: once, twice, ..., $N-1$ times, of the operator $\hat{U}_g(\tilde{\gamma})$ to the initial state $|\psi_0 \rangle$. From equation (17) we have that the relative entropy $S_{\text{rel}}$ between any pair of these states acquires its maximum value 1.

Figure 3 shows the relative entropy between $|\psi_0 \rangle$ and $|\psi_\gamma \rangle$ for three different initial states, corresponding to $N = 2, 3$ and 10 (for comparison purposes $S_{\text{rel}}$ is plotted as a function of $\gamma/\Gamma$ with $\Gamma = N\tilde{\gamma}$). As seen in the figure, the relative entropy is produced in the passing from the initial state ($\gamma = 0$) to the first orthogonal state ($\gamma = \tilde{\gamma}$), when $S_{\text{rel}}$ attains its maximum value for the first time. Then $S_{\text{rel}}$ exhibits oscillations reaching its maximum value periodically, as $\gamma$ take the values $2\tilde{\gamma}, \ldots, (N-1)\tilde{\gamma}$, that is, each time the system visits an element of $O$. From that point on, the relative entropy is degraded as the state goes back to the initial state ($\gamma = N\tilde{\gamma}$), as expected by the recurrent nature of the dynamics. Remarkably, for $N > 2$, there is little variation of the relative entropy production when the system evolves between successive orthogonal states ($S_{\text{rel}}$ remains close to 1), meaning that $|\psi_\gamma \rangle$ remains 'far' from $|\psi_0 \rangle$ almost all time, a feature that gets conspicuous as $N$ gets larger.

The basis $O$ is encountered, for instance, in the theory of a one-dimensional locally periodic media [33], in which case the $n$th element of $O$ corresponds to the state of a particle in the $n$th cell of the system, the unitary transformation corresponds to the spatial translation, and $\tilde{\gamma}$ to the length of the unit cell. When $\hat{U}_g(\gamma)$ corresponds to the time-translation transformation, the states $\{|\psi_{m\gamma} \rangle \}$ coincide with the pointer positions of the simple model of the quantum clock proposed by Anandan and Aharonov in reference [22]. In our geometrical picture, that associates the condition $\langle \psi_\gamma | \psi_0 \rangle = 0$ with a regular polygon with $N$ (labeled) vertices, each application of $\hat{U}_g(\tilde{\gamma})$ amounts to perform a counterclockwise rotation that takes one vertex into the next one (in the time-translation case, takes the hands of the clock into each 'second' positions). Thus, at each rotation by the angle $\tilde{\gamma} \Delta g$, an element of the basis $\{|\psi_{m\gamma} \rangle \}$ is reached. Clearly the $N$th rotation brings back the vertices to their original position, meaning that the evolution is cyclic, with period
\[ \Gamma = N \tilde{\gamma} = \frac{2\pi}{\Delta g}, \]  

(18)
a quantity that has been identified as a characteristic time scale in the context of time evolution [34].

With the aid of equation (7), a direct calculation shows that indeed, up to a global \((g_1\)-dependent) phase factor, the following condition holds

\[ |\psi_\gamma\rangle = \hat{U}_g(\Gamma) |\psi_\gamma\rangle = \left[ \hat{U}_g(\tilde{\gamma}) \right]^N |\psi_\gamma\rangle, \]  

(19)
or equivalently,

\[ \hat{U}_g(\tilde{\gamma}) |\psi_0\rangle = \hat{U}_g(\tilde{\gamma} + \Gamma) |\psi_0\rangle. \]  

(20)
Consequently, the state \(|\psi_\gamma\rangle\) is invariant under the single-period transformation \(\hat{U}_\gamma\Gamma\). When \(\hat{U}_\gamma\) corresponds to the time-evolution operator, \(\hat{U}_\gamma\Gamma\) is known as the Floquet operator [35], and equation (20) is consistent with the quantum recurrence theorem [36].

Before ending this section it is important to make some remarks on the applicability of the previous analysis. Two basic assumptions regarding the initial state \(|\psi_0\rangle = \sum_{k=1}^{N} \sqrt{r_k} e^{i\phi_k} |g_k\rangle\) were made; the first one fixed \(r_k = 1/N\), and the second one imposed the (equally spacing) condition (6b) on the eigenvalues \(g_k\), namely that \(\Delta g = g_{k+1} - g_k\) is constant for all \(k \leq N - 1\). It should be stressed, however, that this is not a condition imposed on the spectrum of \(\tilde{g}\), but rather a restriction on the particular \(g\)-eigenstates that appear in the expansion of \(|\psi_0\rangle\). That is, the spectrum of \(\tilde{g}\) needs no to be an equally-spaced one (as, i.e., in the harmonic oscillator case), yet all the (equally probable) contributing eigenstates must comply with the restriction \(\Delta g = \text{constant}\). In its turn, it is clear that \(\{r_k\}\) is a uniform distribution only in the restricted subspace of eigenstates that satisfy condition (6b), though in the complete space of all \(g\)-eigenstates it is highly non-uniform.

3. An application: \(N\)-two-level systems

Two-level, or qubit systems, have proven to be useful in shedding light on some aspects of the quantum speed limit [10, 29, 37–39]. Here we will consider multi-qubit systems, and apply the above results to get further insight into the dynamics of systems composed of \(N\) bosons in two sites (which are equivalent to \(N\) qubits in a Hilbert space with an effective dimension equal to \(N + 1\)).

3.1. The generator of unitary transformations

When describing a system of \(N\) bosons in two modes, it is usual to resort to the standard annihilation (\(\hat{a}_i\)) and creation (\(\hat{a}_i^\dagger\)) operators of a particle occupying the level \(i = 0, 1\). These operators define the number operator \(\hat{n}_i = \hat{a}_i^\dagger \hat{a}_i\) describing the number of particles in the level \(i\), and satisfy the usual commutation relations

\[ [\hat{a}_i, \hat{a}_j^\dagger] = \delta_{ij}, \quad [\hat{a}_i, \hat{a}_j] = 0. \]  

(21)

\(^2\)For example, for an energy spectrum of the form \(E_k = E_1 k^2\), with \(k\) a positive integer (particle inside an impenetrable box), or \(E_k = E_1 k^{-1}\) (hydrogenoid atom), it is not difficult to choose a set of eigenstates that comply with condition (6b).
In terms of the \( \hat{a} \)'s operators, the generator \( \hat{g} \) of unitary transformations can be written as

\[
\hat{g} = \sum_{i,j=0,1} G_{ij} \hat{a}_i^\dagger \hat{a}_j + \sum_{i,j,k=0,1} G_{ijk} \hat{a}_i^\dagger \hat{a}_j \hat{a}_k + \cdots,
\]

where the \( G \)'s are real parameters. Here we will consider the simpler structure

\[
\hat{g} = G_0 \hat{n}_0 + G_1 \hat{n}_1 + G_{01} \left( \hat{a}_0^\dagger \hat{a}_1 + \hat{a}_1^\dagger \hat{a}_0 \right),
\]

which amounts to disregard many-body interactions and keep (to first order) one-body terms, and from which many well-known system models can be recognized. For example, if \( G_{01} = 0 \), \( G_0 = \omega_0 \) and \( G_1 = \omega_1 \), \( \hat{g} \) corresponds to the Hamiltonian of two independent harmonic oscillators of frequencies \( \omega_0 \) and \( \omega_1 \). If, instead, \( G_0 = G_1 = 0 \) and \( G_{01} \neq 0 \) is constant, \( \hat{g} \) can be identified with the part of the Bose–Hubbard Hamiltonian related with the tunneling of particles between neighboring sites (with \( |G_{01}| \) the intensity of the tunneling), or simply, the Hamiltonian that describes the transitions between the two levels considered.

Now, the Schwinger transformation \( |G| \) maps the bosonic two-mode system into a system of \( N \) elementary 1/2-spins (\( N \) qubits) with total angular momentum \( J \). This allows us to perform our analysis resorting to angular momentum operators (and therefore applicable to angular momentum systems), defined in terms of the creation and annihilation operators as follows

\[
\hat{J}_+ = \hat{J}_x + i \hat{J}_y = \hat{a}_0^\dagger \hat{a}_1, \quad \hat{J}_- = \hat{J}_x - i \hat{J}_y = \hat{a}_1^\dagger \hat{a}_0, \quad \hat{J}_z = \frac{1}{2} (\hat{n}_1 - \hat{n}_0),
\]

where

\[
\hat{J}_l = \frac{1}{2} \left( \hat{\sigma}_l \otimes I_{2^{N-1}} + I_2 \otimes \hat{\sigma}_l \otimes I_{2^{N-2}} + \cdots + I_{2^{N-1}} \otimes \hat{\sigma}_l \right),
\]

\( l = x, y, z \), \( I_d \) stands for the identity operator in a \( d \)-dimensional Hilbert space, and \( \hat{\sigma}_l \) denotes a Pauli operator, expressed in the basis \( \{ |0\rangle, |1\rangle \} \) such that \( \hat{\sigma}_z |0\rangle = |0\rangle \) and \( \hat{\sigma}_z |1\rangle = |1\rangle \). By use of the Schwinger’s transform \( (24a) \), \( (23) \) can be written as

\[
\hat{g} = \frac{1}{2} (G_0 + G_1) \hat{N} + (G_1 - G_0) \hat{J}_z + 2 G_{01} \hat{J}_x,
\]

with \( \hat{N} \) the total number operator \( \hat{N} = \hat{n}_0 + \hat{n}_1 \).

The eigenvectors of \( \hat{g} \), denoted as \( |g_n\rangle \) with \( n \) a quantum number that characterizes the state, correspond to a nondegenerate and equally-spaced spectrum, that is,

\[
\hat{g} |g_n\rangle = g_n |g_n\rangle, \quad \text{with} \quad g_n = A + na,
\]

where \( n = 0, 1, \ldots, N \) and

\[
A = \frac{N}{2} (G_0 + G_1 - a), \quad a = \sqrt{(G_1 - G_0)^2 + 4 G_{01}^2}.
\]

For \( G_{01} = 0 \), the common eigenvectors of \( \hat{J}_z \) and \( \hat{N} \), \( |N, n\rangle \), correspond to the two-mode Fock states \( |n_0\rangle \otimes |n_1\rangle = |N - n\rangle \otimes |n\rangle = |N - n\rangle |n\rangle \) that denote a state for which \( N - n \) particles dwell at level 0 and \( n \) at level 1 (here \( |n_i\rangle \) denotes an eigenvector of \( \hat{n}_i \), such that \( \hat{n}_i |n_i\rangle = n_i |n_i\rangle \)). Thus, (for \( G_{01} = 0 \)) we get

\[
\hat{g} |g_n\rangle = g_n |g_n\rangle, \quad |g_n\rangle = |N - n\rangle |n\rangle, \quad g_n = A^0 + na^0,
\]

with \( A^0 = G_0 N \) and \( a^0 = G_1 - G_0 \), where we have assumed that \( G_1 > G_0 \).
3.2. Reaching orthogonality

We have seen that the eigenvectors of the generator $\hat{g}$ given in (25) constitute a set of nondegenerate states whose eigenvalues satisfy $g_n < g_{n+1}$. Therefore, any superposition of the form

$$|\psi_0\rangle = \sum_{n=0}^{N} \sqrt{r_n} e^{i\phi_n} |g_n\rangle$$

is amenable to the analysis presented in section 2.2, provided the set $\{r_n\}$ is such that the states that contribute to the sum are all equally probable and correspond to a set of equally-spaced eigenvalues. Thus, we will assume that there are $N$ nonzero and equal coefficients $r_n$ (with $2 \leq N \leq N + 1$), so that

$$|\psi_0\rangle = \frac{1}{\sqrt{N}} \sum_{k=1}^{N} e^{i\phi_{nk}} |g_{nk}\rangle,$$

where $\{nk\} = \{n_1, n_2, \ldots, n_N\}$ is a subset of equidistant integers of the set $\{n\} = \{0, 1, \ldots, N\}$, i.e.,

$$\Delta g = g_{n_{k+1}} - g_{n_k} = (n_{k+1} - n_k)a = ma,$$

with $m$ a fixed integer that satisfies $1 \leq m \leq N$. This bound follows from the fact that $m$ is the difference between a pair of integers in the set $\{n\}$; however, it should be noticed that the specific value of $m$ is also restricted by the value of $N$ [see equation (33) below].

We can now resort to equation (6a) to conclude that the state (30) evolving under the transformation $e^{-i\hat{g}\gamma}$ with $\hat{g}$ given by (25), reaches an orthogonal state at

$$\tilde{\gamma} = \frac{2\pi}{N ma}.$$

Now, the (equally probable) $N$ states contributing to (30) allow us to depict the state (30) as an `$N$-teeth comb'—each tooth representing the state $|g_{nk}\rangle$ with eigenvalue $g_{nk}$—with teeth separation $ma$. The length of the comb is therefore $L = ma(N - 1)$. The largest comb corresponds to the expansions (29) that include the extremal states $|g_0\rangle$ and $|g_N\rangle$, and has the maximum length $L_{\text{max}} = Na$. In its turn, the shortest combs correspond to the expansions that involve only two adjacent (minimally-spaced) states, $|g_n\rangle$ and $|g_{n+1}\rangle$, thus fixing the minimum length as $L_{\text{min}} = a$ (see figure 4 below). From here it follows that

$$1 \leq m(N - 1) \leq N.$$  \hspace{1cm} (33)

In terms of the length and the teeth spacing, equation (32) rewrites as

$$\tilde{\gamma} = \frac{2\pi}{L + ma}.$$  \hspace{1cm} (34)

The maximal and minimal values of the length impose lower and upper bounds to $\tilde{\gamma}$ in terms of $N$ and the single parameter of the comb $m$, as follows

$$\tilde{\gamma}_l \leq \tilde{\gamma} \leq \tilde{\gamma}_s,$$  \hspace{1cm} (35)
Figure 4. Comb representation of four possible states of the form (30) for $N = 7$.

(a) The slowest state, for which $\tilde{\gamma} = \pi/a$. The comb’s parameters are $L = a$, $m = 1$.
(b) A state with three contributing $\hat{g}$-eigenstates, thus $\tilde{\gamma} = \pi/3a$. The comb’s parameters are $L = 4a$, $m = 2$.
(c) A state with seven contributing $\hat{g}$-eigenstates, and $\tilde{\gamma} = 2\pi/7a$. The comb’s parameters are $L = 6a$, $m = 1$.
(d) The fastest state, for which $\tilde{\gamma} = \pi/7a$. The comb’s parameters are $L = 7a$, $m = 7$. If the speed of evolution toward orthogonality is measured in terms of $\tilde{\gamma}$, (c) would correspond to the slowest state, followed by (b), and then (a) and (d) would be the (equally) fastest states.

where

$$
\tilde{\gamma}_l = \frac{2\pi}{(N + m)a} \quad \text{and} \quad \tilde{\gamma}_s = \frac{\pi}{a}
$$

(36)
correspond to the largest ($L_{\text{max}}, m = 1, 2, \ldots, N$) and the shortest ($L_{\text{min}}, m = 1$) comb, respectively. That is, larger combs reach orthogonality before shorter combs.

For fixed $N$, $\tilde{\gamma}_l$ attains its lowest value for $m = m_{\text{max}} = N$, whereas $\tilde{\gamma}_s$ admits only $m = m_{\text{min}} = 1$. In both cases ($L = Na$ with $m = N$, and $L = a$ with $m = 1$) we must have $\gamma = 2$, precisely the value of $\gamma$ for which the minimum value $\gamma_{\text{min}}$ is reached. Therefore

$$
\gamma_{\text{min}} \bigg|_{m=N} = \frac{\pi}{Na} \leq \tilde{\gamma}_l \leq \tilde{\gamma} \leq \tilde{\gamma}_s = \frac{\pi}{a} = \gamma_{\text{min}} \bigg|_{m=1}.
$$

(37)
From equation (37) it follows that the ‘fastest’ states are associated to the two-teeth, largest comb with the largest possible separation, i.e.,

$$
\left| \psi_0^{\text{fast}} \right\rangle = \frac{1}{\sqrt{2}} \left( \left| g_0 \right\rangle + e^{i\phi} \left| g_N \right\rangle \right),
$$

(38)
with $\phi$ an arbitrary phase. Analogously, the ‘slowest’ states are also represented by a two-teeth comb, but now with the minimal teeth separation, i.e., by any of the states

$$
\left| \psi_0^{\text{slow}} \right\rangle = \frac{1}{\sqrt{2}} \left( \left| g_{n_1} \right\rangle + e^{i\phi} \left| g_{n_1+1} \right\rangle \right), \quad g_{n_1+1} - g_{n_1} = a,
$$

(39)
with $n_1 = 0, 1, \ldots, N - 1$. Figure 4 shows four possible comb arrays for a system with $N = 7$. The gray (dotted) lines depict the eight possible teeth the comb may have. From top to
bottom, the figure represents four combs obtained by varying the number \( N \) of actual teeth and the separation \( m \), whose corresponding values of \( \tilde{\gamma} \) are in decreasing order. In particular, the first comb saturates the last inequality in (37) \( (\tilde{\gamma} = \pi/a) \) and therefore represents the slowest state \( |\psi_0^{\text{slow}}\rangle \), whereas the last comb saturates the left inequalities in (37) \( (\tilde{\gamma} = \pi/7a) \) hence represents the fastest state \( |\psi_0^{\text{fast}}\rangle \). Notice that in terms of the relative, dimensionless parameter \( \tilde{g} \), both states are equally fast, in line with the discussion below equation (15).

If the eigenvectors \( |g_n\rangle \) are of the form (28) (which means no tunneling between sites, or no mode-interaction), then \( |\psi_0^{\text{fast}}\rangle \) is a balanced superposition of the Fock states \( |N\rangle|0\rangle \) and \( |0\rangle|N\rangle \), which, when written in the computational basis, is identified with the \( N \)-qubit Greenberger–Horne–Zeilinger (GHZ) state (with a relative phase)

\[
|\psi_{\text{ghz}}\rangle = \frac{1}{\sqrt{2}} \left( |0\ldots0\rangle + e^{i\varphi} |1\ldots1\rangle \right). \tag{40}
\]

In its turn, in absence of tunneling, the state \( |g_{n_1}\rangle \) in (39) is the Fock state \( |g_{n_1}\rangle = |N - n_1\rangle|n_1\rangle \).

Taking, for example, \( n_1 = 0 \), the state \( |\psi_0^{\text{slow}}\rangle \) written in the computational basis reads

\[
|\psi_{\text{w}}\rangle = \frac{1}{\sqrt{2}} \left( |0\ldots0\rangle + e^{i\varphi} |W_N\rangle \right), \tag{41}
\]

with \( |W_N\rangle \) the \( N \)-qubit W state [41]

\[
|W_N\rangle = \frac{1}{\sqrt{N}} \left( |100\ldots0\rangle + |010\ldots0\rangle + |00\ldots01\rangle \right). \tag{42}
\]

The states \( |\psi_{\text{ghz}}\rangle \) and \( |\psi_{\text{w}}\rangle \) are thus extremal states that bound the value \( \tilde{\gamma} \) at which any state of the form (30), with \( |g_{n_k}\rangle \) a two-mode Fock state, will evolve toward an orthogonal state. That is, none of such states will attain an orthogonal state ‘faster’ than \( |\psi_{\text{ghz}}\rangle \) nor ‘slower’ than \( |\psi_{\text{w}}\rangle \).

### 3.3. Role of mode-entanglement in the non-tunneling regime

Interestingly, the states \( |\psi_{\text{ghz}}\rangle \) (40) and \( |\psi_{\text{w}}\rangle \) (41) are representative of different types of multiparticle entanglement between \( N \)-distinguishable qubits [41–44]. The connection between entanglement and the speed of quantum evolution goes back to the seminal work of Giovannetti et al [31, 32], that reveals that entanglement can ‘speed up’ the evolution toward an orthogonal state when compared to separable (nonentangled) states. Since then, several investigations exploring the effects of entanglement in the quantum speed limit problem have been advanced, for example, in [45–49].

In order to get further insight into the role of entanglement in the unitary evolution of composite systems toward orthogonality, instead of focusing on multiparticle entanglement between \( N \) parties, here we pay attention to the bipartite entanglement between the system’s modes (or sites), as has been considered experimentally in [23, 50]. Since our analysis involves pure states with only two modes (0 and 1), the mode-entanglement can be directly computed resorting to the concurrence [51]

\[
C_{AB} = C(|\psi\rangle_{AB}) = \sqrt{2 \left( 1 - \text{Tr} \rho_A^2 \right)} = \sqrt{2 \left( 1 - \text{Tr} \rho_B^2 \right)} \leq \sqrt{2(d-1)/d}, \tag{43}
\]
which measures the (bipartite) entanglement between subsystems $A$ and $B$ when the whole composite $A + B$ is in the state $|\psi\rangle = |\psi\rangle_{AB} \in \mathcal{H}_A \otimes \mathcal{H}_B$. Here $d = \min\{\dim \mathcal{H}_A, \dim \mathcal{H}_B\}$, and $\rho_{AB}$ stands for the reduced density matrix of subsystem $A(B)$, i.e., $\rho_{AB} = \text{Tr}_{B(A)} \{ |\psi\rangle \langle \psi| \}$.

In the case where no interaction between modes occur ($G_{01} = 0$), the eigenstates $|g_{nk}\rangle$ in (30) are the two-mode Fock states (28), and a straightforward calculation gives for the entanglement between the modes

$$C_{01} = \sqrt{\frac{2(N - 1)}{N}},$$

which increases with $N$ and attains its minimum value when $N = 2$, precisely when the Margolus–Levitin and Mandelstam–Tamm bounds are saturated. It thus follows that the extremal states $|\psi^{ghz}\rangle$ and $|\psi^{w}\rangle$ possess the same and the minimum value of entanglement between the modes. Comparison of equations (9) and (44) leads to an explicit relation between $\tilde{\gamma}$ and mode-entanglement, namely

$$\tilde{\gamma} = \frac{\pi}{2(g - g_1)} C_{01}^2,$$

meaning that for fixed $(g - g_1)$, the greater the amount of mode-entanglement the greater the ‘delay’ in reaching an orthogonal state. Thus, if we consider a set of equally-weighted and spaced superpositions of Fock states, all having the same $g$-(relative)-expectation value, the first states that will attain an orthogonal state will be those with less entanglement, and correspond (in order of arrival at orthogonality) to combs with 2, 3, . . . , and so on, number of teeth. When the evolution is a Hamiltonian one, this means that for a given mean energy, or for the same energetic resources, mode-entanglement tends to slow down the evolution toward orthogonality.

If instead of fixing the (relative) mean value of $\langle g \rangle$ we fix the teeth separation $ma$, then $\tilde{\gamma}$ and $C_{01}$ exhibit an inverse behavior with respect to $N$. The former is a decreasing function of $N$ [see equation (32)] whereas the latter increases with $N$. Therefore, for the same spacing $ma$, the mode-entanglement favors a ‘faster’ evolution to an orthogonal state.

3.4. The effects of the transition between sites: $G_{01} \neq 0$

Resorting to equation (27), we can compare the value of $\tilde{\gamma}$ when the generator $\hat{g}$ includes the mode-interaction terms ($G_{01} \neq 0$), denoted as $\tilde{\gamma}_{\text{int}}$, with its value in absence of interaction ($G_{01} = 0$), represented by $\tilde{\gamma}_{\text{no int}}$. This gives

$$\frac{\tilde{\gamma}_{\text{int}}}{\tilde{\gamma}_{\text{no int}}} = \left[ 1 + \frac{4G_{01}^2}{(G_1 - G_0)^2} \right]^{-1/2} \leq 1,$$

which shows that $\tilde{\gamma}$ decreases as $G_{01}$ increases, or rather, that the transition between sites favors a faster transit toward an orthogonal state, as depicted in figure 5 where the quotient in (46) is plotted as a function of $G_{01}/(G_1 - G_0)$ (logarithmic scale).

The tunneling between sites will also affect the relation between $C_{01}$ and $\tilde{\gamma}$ expressed in equation (45), where $G_{01} = 0$ was assumed. Thus, in what follows we shall discuss the effects of finite values of $G_{01}$ in the connection between mode-entanglement and $\tilde{\gamma}$. We focus our analysis in the case when the initial state is $|\psi^{\text{fast}}_0\rangle$, given explicitly in (38), for which $\tilde{\gamma} = \pi/Na$ with $a$ given in (27). Notice that contrary to the $\gamma$-independent expression (44), $C_{01}$ will in general depend on $\gamma$ when $G_{01} \neq 0$. In particular, it can be shown that the concurrence (43) for
the $N$-boson state $|\psi_{01}^{\text{fast}}\rangle$ is given by

$$C_{01}^{\text{fast}} = \left[ 2 \left( 1 - \frac{1}{4} \sum_{n=0}^{N} |c_{n,0} + c_{n,N} e^{i(\phi - \pi \bar{\gamma}/2)}|^4 \right) \right]^{1/2} \leq \sqrt{\frac{2N}{N+1}},$$

where $c_{n,0}$, $c_{n,N}$ are the coefficients in the decomposition of the states $|g_0\rangle$ and $|g_N\rangle$, that define (38), in the basis of the Fock states $|N-n\rangle |n\rangle$, i.e., $c_{n,0} = (\langle N-n | \langle n | |g_0\rangle)$ and $c_{n,N} = (\langle N-n | \langle n | |g_N\rangle)$. Such coefficients make $C_{01}^{\text{fast}}$ to depend also on $G_0$, $G_1$, and $\bar{\gamma}$. For the sake of simplicity we set $G_0 = 0$ and use $G_1$ as a scale of units for $g$, so that $C_{01}^{\text{fast}}$ depends only on $N$, $\bar{\gamma}$, and $G_0/G_1$.

In order to illustrate the effects of $G_0/G_1 > 0$ on the connection between $C_{01}^{\text{fast}}$ and $\bar{\gamma}$, we put $N = 2$ and $\phi = 0$ for the sake of simplicity. The concurrence in this case has a maximum value equal to $2/\sqrt{3} \approx 1.1547$, and is shown in figure 6 as a function of $G_0/G_1$, and $\bar{\gamma}$ [within the interval $[0,2]$, corresponding to a period of the dynamics as follows from equation (18)]. For $G_0 = 0$, $|g_0\rangle$ and $|g_2\rangle$ correspond, respectively, to the Fock states $|2\rangle |0\rangle$ and $|0\rangle |2\rangle$, and equation (47) reduces to $C_{01}^{\text{fast}} = 1$ for all $\bar{\gamma}$, in agreement with equation (44) (for $N = 2$), as expected. For small values of $G_0/G_1$, approximately for $G_0/G_1 \lesssim 0.3535$ (or $\log_{10}(G_0/G_1) \lesssim -0.4516$), $C_{01}^{\text{fast}}$ is a concave function of $\bar{\gamma}$ exhibiting a maximum of entanglement at $\bar{\gamma} = \bar{\gamma}$, and the state [see equation (26)]

$$|\psi_{01}^{\text{fast}}\rangle = \frac{e^{-i\bar{\gamma}7}}{\sqrt{2}} \left( |g_0\rangle - |g_2\rangle \right)$$

becomes maximally entangled, i.e. $C(|\psi_{01}^{\text{fast}}\rangle) = 2/\sqrt{3}$, at $G_0/G_1 \approx 0.3535$.

Interestingly, for $G_0/G_1 \gtrsim 0.3535$ the concurrence’s maximum bifurcates into two local maxima at $\bar{\gamma}_m$ and $2\bar{\gamma} - \bar{\gamma}_m$, with $0 < \bar{\gamma}_m < \bar{\gamma}$, while decreases monotonically with $G_0/G_1$ at $\bar{\gamma} = \bar{\gamma}$. Thus, for large enough amplitudes of the transition between modes, $G_{01}$, two states emerge, namely $|\psi_{01}^{\text{fast}}\rangle$ and $|\psi_{01}^{\text{fast}}\rangle$, that exhibit the corresponding maximum mode-entanglement (although slightly smaller than $2/\sqrt{3}$). In contrast, the mode-entanglement of the state $|\psi_{01}^{\text{fast}}\rangle$ vanishes as $G_{01}/G_1 \to \infty$. Indeed, a numerical analysis shows that in such a
Figure 6. Concurrence or mode-entanglement $C_{01}^{\text{fast}}$ considering the initial state $|\psi_{\text{fast}}^{0}\rangle$ for a system of two bosons as a function of the dimensionless parameter of the transformation, $\gamma/\tilde{\gamma}$, and (the logarithm of) the dimensionless transition coefficient between modes, $G_{01}/G_1$ (for $G_0 = 0$).

In the limit we have

$$|\psi_{\text{fast}}^{0}\rangle \to |1\rangle |1\rangle = -i |\Psi^+\rangle,$$

with $|1\rangle |1\rangle$ a Fock state, whose mode-entanglement clearly vanishes. Notice that in the computational basis this state corresponds to the two-qubit Bell state $|\Psi^+\rangle = (|01\rangle + |10\rangle)/\sqrt{2}$, with maximal qubit–qubit entanglement.

4. Summary and final remarks

We revisited the problem of determining conditions under which a pure state that evolves under an arbitrary unitary transformation reaches an orthogonal state in a finite amount of the transformation parameter. Simple geometric considerations disclose the existence of a ‘speed limit’, and the necessity of dispersive dynamical variables—specifically the variable associated to the generator of the evolution—for the initial state to evolve toward a distinguishable one. In particular, when the evolution is generated by the Hamiltonian operator, these geometric observations go in line with the celebrated Mandelstam–Tamm bound.

The geometrical picture presented here leads to consider the family of equally-weighted superpositions of, at least two, non-degenerate eigenvectors of the generator of the transformation, for which the evolution toward orthogonality is guaranteed. The distinguishable state is thus attained after ‘rotating’ the initial state by a characteristic ‘angle’ $\gamma_{\Delta g}$ [see equation (6a)], and the generalized expressions (valid for arbitrary transformations) of the Mandelstam–Tamm and the Margolus–Levitin bounds are obtained. Moreover, for each state in the family under consideration, the relation $\sigma_g \leq \langle g \rangle - g_1$ is satisfied, thus privileging the Mandelstam–Tamm bound over the Margolus–Levitin one.

Some features of the dynamics are disclosed and the entropy production is analyzed, showing that in the simple model of a quantum clock advanced in [22], the relative entropy becomes maximal with each ticking of the clock, and for large $N$ it does not change appreciable until the clock’s hands point to the initial hour, after $(N - 1)$ tickings. In terms of the distinguishability of the states, this means that for a superposition of a sufficiently large number ($\gtrsim 10$)...
of (equally-weighted and spaced) non-degenerate states, the evolved state will remain almost distinguishable from the original one during the entire period $\Gamma$.

Application of our generic analysis to multipartite systems allows, first, to contribute to the establishment and analysis of the relations between entanglement and the quantum speed limit, and second, to establish contact with physical systems that are customarily realized in the lab, such as bosons in optical lattices subject to a Bose–Hubbard Hamiltonian. In this respect, our results show that when no tunneling between sites exist, $\tilde{\gamma}$ is directly proportional to the squared of the entanglement between the modes ($C_{01}$), so for $\langle \hat{g} \rangle - g_1$ fixed, the higher the mode-entanglement is, the more amount of transformation is required to reach an orthogonal state. In contrast, for fixed $ma$, the higher the mode-entanglement is, the less amount of transformation is needed to attain a distinguishable state. Therefore, a high amount of mode-entanglement alone does not guarantee, nor prevents, a more ‘rapid’ transit toward orthogonality.

Now, when the bosons are allowed to tunnel between sites, a more complex relation between $\tilde{\gamma}$ and $C_{01}$ is obtained, and the dynamics of the mode-entanglement becomes richer. This is exemplified for the initial state (38), which corresponds to the superposition of eigenstates of the transformation generator with minimum and maximum eigenvalues. The observation of a more complex dynamics of $C_{01}$ is expected, since the transition between sites—encoded in the ‘mode-interaction’ part of the generator $\hat{g}$—typically affects the entanglement between modes.

Our discussion on the relation between the mode-entanglement and the amount of transformation required to reach orthogonality, complements previous analysis [45–49] regarding the role of particle entanglement on the quantum speed limit. As for the (multipartite) entanglement between the particles, here we found that in the nontunneling regime, the ‘slowest’ ($|\psi^w\rangle$) and the ‘fastest’ ($|\psi^{ghz}\rangle$) states correspond to paradigmatic states that exhibit multipartite entanglement [41]. Further investigations will certainly reveal the more subtle relation between the quantum speed limit and the different types of multipartite entanglement.

There exist other geometrical approaches that bring out new and tighter bounds on the speed of evolution, such as that advanced in [16], and some comments on the main differences with the present treatment are in place. Here we make use of a rather simple mathematical framework which takes advantage of the fact that the overlap between an initial pure state and its evolved one is a complex scalar, a property that endows it with a two-dimensional representation and ultimately allows us to map the overlap into a set of (rotating) vectors in $\mathbb{R}^2$. Instead, other works analyze the distinguishability between two states (either pure or mixed) by focusing directly in their mutual ‘angle’ (for instance the Bure or the Wigner–Yanase angle), which naturally leads to a description involving differential geometry on Riemannian manifolds, where a family of metrics based on the geodesic distance between states is used to determine tighter quantum speed limits, that add to the Mandelstam–Tam and Margolus–Levitin bounds. Considering if the approach presented here may be extended to recover those and other bounds is an interesting task that is left for future investigations, aimed at offering a simpler way of envisaging and constructing fundamental limitations in the evolution of dynamical systems.

Within this latter aim, another line of research that also deserves further analysis is related to the universality of the limitations on the speed of evolution, as suggested by works showing that the speed limit remains in the transition from the quantum to the classical regime [26–28]. As stated before, the method followed in this paper is based on the mapping of the probability amplitude between two pure states into a vector in $\mathbb{R}^2$, whose direct classical counterpart remains unclear. There are, however, some directions within the same spirit of the present work that can be explored. For instance, the limit of continuous spectrum could be analyzed to provide some insights into the semi-classical limit; also an interesting approach would be
to resort to the path functional description of the overlap. In a more ambitious scenario, some of the main ideas presented here and applicable to pure states only, may be extended to mixed states and then put into correspondence with probability distributions describing a classical system. All these open problems go beyond the scope of the present work, yet certainly are worthy of investigation.

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ORCID iDs

Andrea Valdés-Hernández  
https://orcid.org/0000-0003-3347-1637

Francisco J Sevilla  
https://orcid.org/0000-0003-0346-2967

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