# Dynamics of entanglement entropy and entanglement spectrum crossing a quantum phase transition 

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#### Abstract

We study the time evolution of entanglement entropy and entanglement spectrum in a finite-size system which crosses a quantum phase transition at different speeds. We focus on the transverse-field Ising model with a time-dependent magnetic field, which is linearly tuned on a time scale $\tau$. The time evolution of the entanglement entropy displays different regimes depending on the value of $\tau$, showing also oscillations which depend on the instantaneous energy spectrum. The entanglement spectrum is characterized by a rich dynamics where multiple crossings take place with a gap-dependent frequency. Moreover, we investigate the Kibble-Zurek scaling of the entanglement entropy and Schmidt gap.


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## I. INTRODUCTION

In recent times, there have been considerable experimental and consequent theoretical advances in the study of the dynamics of closed quantum many-body systems (for a review of both the experimental and theoretical aspects, see Ref. [1]). In this work we will deal with the problem of studying the time evolution of a closed quantum many-body system at $T=0$, when it is driven from one phase to another by allowing the coupling constants in the Hamiltonian to change in time. Typical evolution protocols are the so-called slow quenches, where the velocities of the variations of the couplings are finite, or, on the contrary, sudden quenches, when such couplings are instantaneously varied to a different value and then left constant. Here we will discuss the physical behavior of an integrable system, namely the Ising chain in a transverse field [2-4], in the whole range of velocities according to which we let the magnetic field to vary.

One-dimensional problems may be tackled with generally powerful numerical methods such as t-DMRG [5-7] or TEBD [8]. It is well known that the final efficiency of such methods is related to the amount of entanglement of the considered state [9], a quantity which is expected to diverge when getting closer to a phase transition. However, at least in the static case, the behavior of entanglement (and more specifically of entanglement entropy) has a universal character so that it can be used as an estimator of quantum correlations [10] and to detect as well as to classify quantum phase transitions also in fully interacting models [11-16].

Thus, it is natural to ask whether the dynamical behavior of a closed quantum system, especially when crossing a phase transition, can be described by looking at the dynamics of entanglement entropy and entanglement spectrum, a topic on which there are only a few general results [17-20].

The aim of this work is to investigate this question in a paradigmatic example: the Ising chain in a time-dependent transverse field, a problem which allows for an exact solution at any instant of time [2-4]. The plan of the work is the following. In Sec. II we define the notion of entanglement entropy and entanglement spectrum and present the model and its phases. In Sec. III we describe the dynamics when letting the system go from the paramagnetic to the ferromagnetic phase
by controlling the speed with which we change the magnetic field. We will examine the adiabatic regime, the sudden-quench situation, and the cases with intermediate speeds. Then we will see how these results are related to the so-called Kibble-Zurek mechanism [21,22] in its quantum version [23], by looking both at the scaling of entanglement entropy and the so-called Schmidt gap $[24,25]$ in the entanglement spectrum. In Sec. IV we perform a similar analysis when the system evolves from the ferromagnetic to the paramagnetic phase. We end the work with conclusions and outlooks in Sec. V, and with three appendices where we have reported technical details of the calculations.

## II. THE MODEL

In this work, we are interested in the time evolution of bipartite quantities, such as the entanglement entropy and the entanglement spectrum, which are defined in the following way [26]. Starting from a one-dimensional chain of $L$ sites, we consider a subsystem $A$ containing $\ell<L$ adjacent sites, $\bar{A}$ being its complement. The reduced density matrix is obtained from the pure density matrix of the ground state of the whole chain, $\rho=|\mathrm{GS}\rangle\langle\mathrm{GS}|$, as

$$
\begin{equation*}
\rho_{A}=\operatorname{Tr}_{\bar{A}} \rho \tag{1}
\end{equation*}
$$

The so-called entanglement spectrum is the set $\left\{\lambda_{n}\right\}$ of the eigenvalues of $\rho_{A}$; the entanglement entropy is defined as

$$
\begin{equation*}
S=-\operatorname{Tr}_{A} \rho_{A} \log _{2} \rho_{A} \tag{2}
\end{equation*}
$$

and computed as

$$
\begin{equation*}
S=-\sum_{n} \lambda_{n} \log _{2} \lambda_{n} \tag{3}
\end{equation*}
$$

In what follows we take as subsystem the half chain; i.e., $\ell=L / 2$. We checked that the main findings of this work are not changed qualitatively if we take a different $\ell$.

In the following, we consider the Ising model in a transverse field [2-4]:

$$
\begin{equation*}
H=-\frac{J}{2} \sum_{j=1}^{L}\left[\sigma_{j}^{z} \sigma_{j+1}^{z}+h \sigma_{j}^{x}\right] \tag{4}
\end{equation*}
$$

where $L$ is the system size and we have periodic boundary conditions (PBCs); $\vec{\sigma}$ are the Pauli matrices, and $h=h(t)$ is a time-dependent magnetic field (moreover, we assume $\hbar=J=1$, so that the energies are measured in units of $J)$. As recalled in Appendix A, the model is exactly solvable by a sequence of Jordan-Wigner-Fourier-Bogolyubov transformations; the eigenenergies and the corresponding eigenstates are completely known. Remarkably, the spectrum divides into two sectors, labeled by the quantum number $\alpha \equiv \prod_{j=1}^{L} \sigma_{j}^{x}= \pm 1$; at finite size, the ground state always belongs to the $\alpha=1$ sector [27], which is the one we will deal with in this work. Moreover, the model is one of the prototypical playgrounds for quantum phase transitions [3]. Indeed, varying $h$, the Hamiltonian in Eq. (4) displays (in the thermodynamic limit) a quantum critical point at $h=1$, which separates the paramagnetic $(h>1)$ and the ferromagnetic ( $0 \leqslant h<1$ ) phases [3] (the negative part of the phase line is the mirror reflection of the positive one, because of the $\mathbb{Z}_{2}$ symmetry under the canonical transformation $\sigma_{j}^{x} \rightarrow-\sigma_{j}^{x}$ [4]). The low-energy physics of such a quantum critical point is described by a conformal field theory [28] of central charge $c=1 / 2$. Its correlation length and dynamic critical exponents are given by $v=z=1$ [29].

We make the Hamiltonian in Eq. (4) explicitly time dependent by letting $h=h(t)$ change linearly in time, from an initial value $h_{i}$ to a final one $h_{f}$ :

$$
\begin{equation*}
h(t)=h_{i}+\operatorname{sgn}\left(h_{f}-h_{i}\right) \frac{t}{\tau}, \tag{5}
\end{equation*}
$$

where $\tau$ is the time scale of the ramping and $t \in\left[0, t_{f}\right]$, with $t_{f}=\left|h_{f}-h_{i}\right| \tau$ (time is measured in units of $1 / J$ ). The dynamics of the model is also exactly accessible [30], as we recall in Appendix B (see also Ref. [31]). As we shall see, very different dynamic behaviors are expected for different values of $\tau$.

## III. PARAMAGNET TO FERROMAGNET

In this section we study the ramping from the paramagnetic sector of the phase diagram $\left(h_{i}>1\right)$ to the ferromagnetic one ( $h_{f}<1$ ). This is the setting for the study of the Kibble-Zurek mechanism in the one-dimensional quantum Ising model [23].

## A. Initial structure of the entanglement spectrum

Let us start by studying the initial condition for bipartite quantities. To this aim it is useful to first understand the limit $h_{i} \rightarrow \infty$.

In this case the ground state of the system at $t=0$ looks like

$$
\begin{equation*}
|0\rangle \equiv \prod_{j=1}^{L}|\rightarrow\rangle_{j} \tag{6}
\end{equation*}
$$

where we denote with $|\rightarrow\rangle_{j},|\leftarrow\rangle_{j}$ the state with $S_{j}^{x}= \pm \hbar / 2$, respectively. Of course, this is not the exact ground state for finite $h_{i} \gg 1$, but, at first order in perturbation theory, it is easy
to show that the latter is given by

$$
\begin{equation*}
|\mathrm{GS}\rangle=N\left[|0\rangle+\frac{1}{4 h} \sum_{j=1}^{L}|j, j+1\rangle\right] \tag{7}
\end{equation*}
$$

with $|j, j+1\rangle$ being the state with two spin flips at sites $j$, $j+1$ :

$$
\begin{equation*}
|j, j+1\rangle \equiv|\leftarrow\rangle_{j}|\leftarrow\rangle_{j+1} \prod_{\substack{k=1 \\ k \neq j, j+1}}^{L}|\rightarrow\rangle_{k} \tag{8}
\end{equation*}
$$

$N \equiv\left(1+\frac{L}{16 h^{2}}\right)^{-\frac{1}{4}}$ is the normalization coefficient (this usually neglected normalization factor is necessary to obtain a good agreement with numerical results).

The zero-temperature density matrix of the system is given by

$$
\begin{equation*}
\rho=|\mathrm{GS}\rangle\langle\mathrm{GS}| \tag{9}
\end{equation*}
$$

and the reduced density matrix $\rho_{A}=\operatorname{Tr}_{\bar{A}} \rho$ of the half chain $A=\{1, \cdots, L / 2\}$ (we will always choose this bipartition) is seen to take the form

$$
\rho_{A}=\left(|0\rangle_{A}, \quad|2 p\rangle_{A}, \quad|1\rangle_{A}, \quad|L / 2\rangle_{A}\right) \mathbb{R}_{A}\left(\begin{array}{l}
{ }_{A}\langle 0|  \tag{10}\\
A\langle 2 p| \\
A\langle 1| \\
A\langle L / 2|
\end{array}\right)
$$

$|0\rangle_{A}$ being the paramagnetic state relative to subsystem $A$. Also,

$$
\begin{align*}
|2 p\rangle_{A} & \equiv\left(\frac{L}{2}-1\right)^{-\frac{1}{2}} \sum_{j=1}^{\frac{L}{2}-1}|j, j+1\rangle \\
|1\rangle_{A} & \equiv|\leftarrow\rangle_{1} \prod_{j=2}^{\frac{L}{2}}|\rightarrow\rangle_{j}  \tag{11}\\
|L / 2\rangle_{A} & \equiv|\leftarrow\rangle_{L / 2} \prod_{j=1}^{\frac{L}{2}-1}|\rightarrow\rangle_{j}
\end{align*}
$$

and

$$
\mathbb{R}_{A} \equiv N^{2}\left(\begin{array}{cccc}
1+\frac{\frac{L}{2}-1}{16 h^{2}} & \frac{\sqrt{\frac{L}{2}-1}}{4 h} & 0 & 0  \tag{12}\\
\frac{\sqrt{\frac{L}{2}-1}}{4 h} & \frac{\frac{L}{2}-1}{16 h^{2}} & 0 & 0 \\
0 & 0 & \frac{1}{16 h^{2}} & 0 \\
0 & 0 & 0 & \frac{1}{16 h^{2}}
\end{array}\right)
$$

The form of $\mathbb{R}_{A}$ shows that $|1\rangle_{A}$ and $|L / 2\rangle_{A}$ are true eigenstates of $\rho_{A}$; diagonalizing the remaining block it can be seen that, for large enough $h$, the remaining two eigenstates are superpositions of $|0\rangle_{A}$ and $|2 p\rangle_{A}$, one in which the paramagnet dominates and the other in which $|2 p\rangle_{A}$ dominates. A numerical analysis shows that the largest eigenvalue is associated with the first one, while the smallest with the second; the ones associated with the single flipped states $|1\rangle_{A}$ and $|L / 2\rangle_{A}$ are of course degenerate, and occupy the second and third position in magnitude.


FIG. 1. (Color online) Dynamics of the entanglement entropy for $L=50, h_{i}=1.4, h_{f}=0.4$. Main panel: $S(t)$ for different values of $\tau$ (dashed vertical line: location of the critical point). Inset: $\tilde{S}(t)$ as defined in Eq. (14) for the same values of $\tau$ as in the main panel.

## B. General dynamical features

The aim of this section is to show that just few eigenvalues contribute to the entanglement dynamics of the system. To see this, we first compute the entanglement entropy for a half-chain bipartition, as explained in Appendix C [see Eq. (C5)]. The results for $L=50, h_{i}=1.4, h_{f}=0.4$, and different values of $\tau$ are shown in the main panel of Fig. 1.

Then, we compute the first four eigenvalues of the reduced density matrix of $A$ and consider their sum

$$
\begin{equation*}
W_{4}(t)=\sum_{n=1}^{4} \lambda_{n}(t) \tag{13}
\end{equation*}
$$

which is shown in Fig. 2.
We notice that $W_{4}$ is very close to unity for fast rampings ( $\tau \lesssim 1$ ) and, away from the critical value of $h$, for nearly


FIG. 2. (Color online) Sum of the first four eigenvalues of the reduced density matrix $W_{4}$ [see Eq. (13)] for $L=50, h_{i}=1.4, h_{f}=$ 0.4 , and different values of $\tau$.
adiabatic rampings ( $\tau \gtrsim 500$ ). In all other cases the weight of the first four eigenvalues is always at least $\sim 0.97$; moreover, this seems to be quite $L$ independent (not shown). Therefore, in the following, we will always consider just the first four eigenvalues of the reduced density matrix. By means of the latter eigenvalues, we can compute a "partial" entanglement entropy:

$$
\begin{equation*}
\tilde{S}(t)=-\sum_{n=1}^{4} \lambda_{n}(t) \log _{2} \lambda_{n}(t) . \tag{14}
\end{equation*}
$$

We show this quantity in the inset of Fig. 1, where we see that, apart from a small quantitative discrepancy, the qualitative behavior is indeed the same as the one of the true entanglement entropy $S(t)$.

In Secs. III C-IIIE we will study the dynamics of the entanglement spectrum in detail (Fig. 3). The entanglement spectrum at the initial value $h_{i}$ is very close to the one described in Sec. III A (large-field case). As we shall see, it displays, as well as entanglement entropy, a different dynamical qualitative behavior depending on the value of $\tau$. Unless explicitly stated, we choose $L=50$ (postponing the discussion of size effects to Sec. III F) and show our results for a ramping from $h_{i}=1.4$ to $h_{f}=0.4$. We choose these values of the initial and final magnetic field in order to restrict the range of integration of the ODEs, Eq. (B6).

## C. Adiabatic and sudden regimes

We begin by considering very large values of $\tau$, i.e., a quasiadiabatic quench; see for example the curve at $\tau=500$ of Fig. 1 and panel (a) of Fig. 3.

We observe that during the evolution the entanglement entropy and the entanglement spectrum closely follow the static values, i.e., those obtained from the ground state of the system at each value of $h(t)$, the only difference being represented by some small oscillations, that will be discussed in Sec. IIIE. This behavior is expected from the adiabatic theorem [32] and as a consequence of the finite size of the system. Indeed the energy gap closes as a function of the inverse size, remaining nonzero for any finite $L$, so that in this case it is always possible to reach the adiabatic limit provided $\tau$ is large enough (see also Sec. III F). More precisely, as shown by Cincio et al. in Ref. [20], the probability of having an adiabatic evolution at size $L$ is given by $P(\tau)=$ $1-\exp \left(-2 \pi^{3} \tau / L^{2}\right)$, so that the maximum rate $(\sim 1 / \tau)$ at which the evolution is adiabatic decays as $1 / L^{2}$.

We then consider the opposite regime, with very small values of $\tau$, i.e., very fast quenches: we show this situation in Fig. 1 (curve with $\tau=0.1$ of the main panel) and Fig. 3(b). The entanglement entropy and the entanglement spectrum do not evolve at all, as expected from the adiabatic theorem, independently of the size of the system.

## D. Fast sweeps

We consider now rampings that are slower than sudden ones, but much faster than adiabatic ones; we call them fast sweeps, and, for our system sizes, they correspond to $\tau=$ $1-20$. For the sake of clarity, for both the entanglement entropy


FIG. 3. (Color online) Dynamics of the entanglement spectrum for $L=50, h_{i}=1.4$, and $h_{f}=0.4$. Black dashed, red dot-dot-dashed, green solid, and blue dash-dotted lines correspond to the dynamical first, second, third, and fourth eigenvalue of the reduced density matrix of the half chain, respectively. Different panels refer to (a): $\tau=500$; (b): $\tau=0.1$; (c): $\tau=1$; (d): $\tau=8$; (e): $\tau=30$; (f): $\tau=100$. In panels (b), (c), and (d) the red and green lines are overlaid. The cyan lines in panel (a) show the ground-state values of the first four eigenvalues. The inset in panel (d) is a zoom of the crossing point.
and the entanglement spectrum it is useful first to consider the faster regime $\tau \sim 1$ and then slower rampings $\tau \sim 10-20$.

Starting from faster rampings (see curves with $\tau=1$ and 5 in the main panel of Fig. 1), the entanglement entropy increases linearly in the region close to the phase transition: this behavior can be related to the results of Calabrese and Cardy [18] relative to a sudden quench to a conformal critical point. In their case, the entanglement entropy is predicted to grow (at least in the first part of its evolution) linearly, with a slope related to the central charge of the underlying conformal field theory. Even if our case is different from the cited one because of the finite ramping speed, we can try to apply this picture. Indeed, close to the critical point, the correlation length and relaxation time are large, so that the system behaves as critical for a finite interval of $h$.

The behavior of the entanglement spectrum is of course related to the one of the entanglement entropy and is shown in panel (c) of Fig. 3. In this regime of $\tau$, the first eigenvalue decreases, while the remaining three increase: this results in a growth of the entanglement entropy that we observe [26]. Remarkably, the second and third eigenvalues of the reduced density matrix remain degenerate even during this kind of evolution: indeed, these eigenvalues correspond, at $t=0$, to the eigenstates $|1\rangle$ and $|L / 2\rangle$ (see Sec. III A), and the time evolution, as shown by a perturbative analysis (that we are not going to report), does not break this degeneracy, at least for these values of $\tau$.

The second regime is encountered by further increasing $\tau$ (see for example curves with $\tau=8,10$, and 30 in the main panel of Fig. 1). In such cases, the entanglement entropy still presents a linear-growth region, which does not last to the end of the sweep, ending in an oscillatory region, in which the entanglement entropy alternates between maxima and minima, with variable frequency. This behavior has already
been observed in a thermodynamic-limit study of the dynamics of entanglement entropy [19], and has been ascribed to the fact that the system ends up, after passing the critical point, in a superposition of excited states of the instantaneous Hamiltonian. In particular, the oscillation frequency has been predicted to scale as

$$
\begin{equation*}
\omega(t) \approx \Delta(t) \tag{15}
\end{equation*}
$$

$\Delta(t)$ being the energy gap of the instantaneous Hamiltonian, given by Eq. (4). To verify this prediction in our finite-size system, we evolve the ground state of $H(t=0)$ according to the protocol

$$
h(t)= \begin{cases}h_{i}-\frac{t}{\tau}, & 0 \leqslant t \leqslant\left(h_{i}-h_{f}\right) \tau  \tag{16}\\ h_{f}, & t>\left(h_{i}-h_{f}\right) \tau\end{cases}
$$

i.e., the usual ramping of Eq. (5) followed, in the end, by an evolution according to the final Hamiltonian. The result is that in the second part of the evolution the entropy oscillates with a constant frequency [Fig. 4(a)], and such an oscillation is superimposed on another one, much smaller in amplitude and period (plus an increasing power-law trend). To determine the period of the first one, we fit the right part of the curve in Fig. 4(a) by means of the seven-parameter formula $y=a_{0}+a_{1} x^{a_{2}}+a_{3} \cos \left(\frac{2 \pi x}{a_{4}}+a_{5}\right) / x^{a_{6}}$, that turns out to be a very suitable fitting equation (apart from the subdominant oscillatory behavior); $a_{4}$ is directly the period of the oscillation. After getting from the fits the values of the periods for several values of $h_{f}$, we can plot them as a function of $\Delta(t)$ (the gap is computed from the exact solution; see Appendix A). The results are shown in Fig. 4(b): the behavior of the oscillation period as a function of the gap is compatible with a $1 / x$ law. This check confirms, as we shall see in Sec. III F, that the


FIG. 4. (Color online) Panel (a): evolution of the entanglement entropy according to the protocol (16) for $\tau=10$. Panel (b): oscillation period as a function of the energy gap ( $h_{i}=1.4, h_{f} \in$ [ $0.31,0.4]$ ); the fit is performed by means of the formula $y=$ $a_{0}+a_{1} / x$. Panel (c): time at which the eigenvalues of the reduced density matrix cross as a function of $\tau$, for $\tau \in[8,17]$. Black dots: numerical data; red line: fitting formula $y=a_{0}+a_{1} / x^{a_{2}}$, giving $a_{0}=0.428193$ (critical point: 0.4). Panel (d): same as in (b), but with $h_{i}=1.5, h_{f} \in[0.41,0.5]$, and $\tau=30$.
physics in this regime is the same as in the thermodynamic limit.

We now investigate the behavior of the entanglement spectrum in this regime. As shown in Fig. 3(d), the decreasing of the first eigenvalue and the growth of the remaining ones continue until they cross, all at the same point. Moreover, this crossing structure recurs also for later times in an almost periodic pattern (not shown). This behavior is very peculiar, and we shall investigate it in detail. First of all, it must be noticed that the crossings correspond, as expected, to the maxima of entanglement entropy and that this oscillatory behavior starts only after the system has crossed the critical point. This fact is easily confirmed by plotting the crossing time $t_{c r}$ as a function of $\tau$. The result is shown in Fig. 4(c): the data can be fitted by a power law, showing that, for $\tau \rightarrow \infty$, the crossing point converges with good precision to the critical point (strictly speaking, we could not take the limit $\tau \rightarrow \infty$, since, for larger $\tau$, the behavior of the system tends to become adiabatic; however, this extrapolation shows that the oscillations, also present for larger $\tau$, always have the same nature; see Sec. IIIE). We have also verified that the crossing time $t_{c r}$ does not depend on the size of the system at fixed $\tau$ (not shown): this fact represents further evidence of the fact that the physics, for these values of $\tau$, coincides with the thermodynamic-limit one.

By magnifying the crossing region [see inset of Fig. 3(d)], it becomes manifest that the fourfold crossing is actually a crossing between the first and the fourth eigenvalue, while the second and the third continue evolving parallel to each other.

A question which might arise is, what are the signatures of the crossing of the eigenvalues of the reduced density matrix on observable quantities? As an example, we can get an insight


FIG. 5. (Color online) Main panel: time evolution of the expectation value of the density of left spins on the first four eigenstates of the reduced density matrix [cf. Eq. (17)] for a ramping of the magnetic field from $h_{i}=6$ to $h_{f}=-1$ with $\tau=1$ and $L=18$ sites. Panel (a): zoom of the main panel around the crossing point, with the usual color code. Panel (b): corresponding dynamics of the entanglement spectrum in the same time interval of panel (a).
in this direction by computing the expectation value of the density of left spins on each of the eigenstates, defined by

$$
\begin{equation*}
\rho_{x}(j, t) \equiv \frac{1}{\ell} \sum_{i=1}^{\ell}\langle j(t)| \frac{1}{2}\left(1-\sigma_{i}^{x}\right)|j(t)\rangle, \tag{17}
\end{equation*}
$$

$|j(t)\rangle$ being the $j$ th eigenstate of the reduced density matrix at time $t(j=1,2,3,4)$. Since the latter states are many-body objects, in order to compute $\rho_{x}$ we use exact diagonalization, with the time evolution performed via the time-dependent Lanczos algorithm [33] (in a recent study [34], Zamora and collaborators were able to compute similar quantities in the free-fermions approach). Our results are summarized in Fig. 5 for a quench with $h_{i}=6$ and $h_{f}=-1$ and $L=18$ sites. The choice of this interval of $h$, much larger than the ones considered up to now, is due to the small size of the system. However, the crossing of the eigenvalues $\lambda_{i}$ has the same structure as in the inset of Fig. 3(d).

First of all we notice from the main panel that for all the four eigenstates the density $\rho_{x}$ increases during the time evolution. This is expected, since the operator $\sigma_{j}^{z} \sigma_{j+1}^{z}$ rules the evolution of the system, by creating pairs of left spins. Almost at the end of the evolution the densities approach each other, in a way which is magnified in Fig. 5(a). The densities corresponding to the first and the fourth eigenvalue, i.e., $\rho_{x}(1, t)$ and $\rho_{x}(4, t)$, are first exchanged at a time $t \sim 6.81$, in correspondence with the crossing of $\lambda_{1}$ and $\lambda_{4}$, and then cross each other a $t \sim 6.86$. The important information which we can extract from Fig. 5(a) is that the main contribution to the density of flips in the half chain [which is the physical meaning of $\rho_{x}(1, t)$ ] changes the profile of its time evolution in correspondence with the crossings of the eigenvalues of the reduced density matrix. In particular, if we follow the black symbols in Fig. 5(a), the density $\rho_{x}(1, t)$ increases more slowly after the crossing.


FIG. 6. (Color online) Panel (a): dynamics of the entanglement spectrum for a ramping from $h_{i}=6.0$ to $h_{f}=0$, with $\tau=10$ and $L=18$. Panel (b): expectation value of the density of left spins.

## E. Slow sweeps

The last regime is observed for $\tau \gtrsim 20$; a typical situation is shown in Figs. 3(e) and 3(f). As Fig. 3(e) shows, the second and the third eigenvalues begin to separate, and then the crossing of the first and the fourth begins to become an avoided crossing; for larger values of $\tau$, as shown in Fig. 3(f), this separation continues and the dynamical structure of the spectrum gets closer to the static one, i.e., the one of Fig. 3(a). In all cases, the crossings take now place between the first and second, third and fourth eigenvalues; remarkably, they take place at the same times for the first and the second couple. On the other hand, the entanglement entropy, as shown in the main panel of Fig. 1 (curves with $\tau=100$ ), at the beginning of the evolution is practically coincident with the static one, and at a certain point begins to grow; however, it begins to oscillate around a value that is smaller than the ones of Sec. IIID, and that decreases as $\tau$ increases. This behavior of entanglement spectrum and entanglement entropy can be ascribed to the approaching of the adiabatic regime. However, as already observed in Sec. III C, the oscillation studied in Sec. IIID survives as a sign of nonadiabaticity, but this time between the first and the second two eigenvalues. Even in this case, performing the same analysis as in Sec. IIID, we can show that the period of such oscillation at the instant $t$ decreases as a function of the inverse gap of $H(t)$ [see Fig. 4(d)].

By taking a small system with $\tau$ in this slow regime, we can proceed as in Sec. IIID to compute the expectation value of the density of left spins [see Eq. (17)]; we show our results in Fig. 6. In panel (a) the dynamics of the eigenvalues is shown to be analogous to that of Fig. 3(f). In panel (b) of Fig. 6 we first observe a crossing of the density of left spins corresponding to the first and third eigenstate, immediately followed by a crossing of the second and the fourth and another involving the first and the fourth eigenstate. When the crossings of the $\lambda_{j}(t)$
take place, the densities of left spins exchange and, at different times, cross each other, until at the end of the evolution we observe two pairs of self-avoiding levels. Even in this case, the density of left spins is the only quantity displaying crossings in correspondence of the crossing of the eigenvalues (up to a nonsynchronization of the exchanges of the eigenvalues and the crossings themselves).

## F. Kibble-Zurek physics

In this section, we discuss the Kibble-Zurek scaling [21,22] of two quantities, i.e., the already considered entanglement entropy and the Schmidt gap [24,25], i.e., the difference between the two largest eigenvalues in the entanglement spectrum. A discussion of this mechanism for the $X Y$ model may be found in Refs. [23], [35], [36], and [37].

In its original formulation, the Kibble-Zurek mechanism is able, on the basis of extremely simple approximations, to predict the scaling of the number of topological defects produced after the dynamical transition of a critical point. The key assumption underlying the mechanism is that the evolution can be divided, for suitable ramping velocities, into three parts: a first adiabatic one, where the wave function of the system coincides with the ground state of $H(t)$; a second impulsive one, where the wave function of the system is practically frozen, due to the large relaxation time close to the critical point; a third adiabatic one, as the system is driven away from the critical point [23]. This division takes the name of adiabatic-impulse-adiabatic approximation [38]. What plays a role in this kind of mechanism is the correlation length $\hat{\xi}$ at the times of passage between the different regimes, that can be seen to scale, for a linear quench of inverse velocity $\tau$, as [22]

$$
\begin{equation*}
\hat{\xi} \approx \tau^{\frac{v}{1+z v}} \tag{18}
\end{equation*}
$$

$v$ and $z$ being the critical exponents of the crossed quantum critical point [29].

## 1. Entanglement entropy

Any quantity that is directly related to the correlation length is suitable to a Kibble-Zurek analysis. In particular, close to a conformal critical point of conformal charge $c$, the entanglement entropy has been shown by Calabese and Cardy to diverge, in an infinite system, as [39]

$$
\begin{equation*}
S=\frac{c}{6} \log _{2} \xi+\text { const. } \tag{19}
\end{equation*}
$$

In particular, we remark that, because of the infinite size of the system, subsystem $A$ possesses just one effective boundary. The entanglement entropy after the quench is therefore easily seen to scale as [19]

$$
\begin{equation*}
S=\frac{c v}{6(1+z v)} \log _{2} \tau+\text { const. } \tag{20}
\end{equation*}
$$

The prefactor of the logarithm is $1 / 24$, since in the Ising case $v=z=1$ and $c=1 / 2$. This clearly holds in the thermodynamic limit, where the gap is strictly closed at the critical point. In our case, at finite size, we expect some deviations from the Kibble-Zurek behavior for large $\tau$. We plot the results we obtain in Fig. 7: as expected, we observe a progressive breakdown of the Kibble-Zurek prediction lowering $L$ (see also the inset of Fig. 7).


FIG. 7. (Color online) Main panel: entanglement entropy at the final instant of the evolution for $\tau \in[1,200]$ at different system sizes ( $L=10-100$, from bottom to top). Dashed maroon line: $y=\frac{1}{12} \log _{2} x+$ const. Inset: entanglement entropy at the final instant of the evolution for $\tau=100$ as a function of the system size $L$.

A few other remarks are in order: First, Eq. (19) has to be modified, since, because of its finite size, subsystem $A$ possesses two boundaries; therefore, Eq. (19) is modified by doubling the prefactor of the logarithm [39] (see also Ref. [20]). Moreover, it is evident that the logarithmic behavior expected from the Kibble-Zurek mechanism is superimposed on an oscillating behavior, as already observed in Ref. [19]: it is clearly a reflex of the oscillating structure of the entanglement entropy as a function of time, studied in Secs. III D and III E. Third, we observe that for small values of $\tau$, the curves at different sizes are practically coincident. This coincidence is lost for larger values of $\tau$ 's, depending on $L$ : the velocities at which this coincidence is observed are the ones at which the physics is practically the one of the thermodynamic limit. For example, at $L=50$, the physics is practically the thermodynamic limit one up to $\tau \approx 15$.

Finally, we note that, remarkably, the $\tau$ 's that correspond to the passage from the fast to the slow regime (the $\tau$ 's for which the crossing between the first and the fourth eigenvalue of the reduced density matrix begin to disappear) correspond to the breakdown of the Kibble-Zurek, or, equivalently, thermodynamic-limit physics. This fact could be verified by a direct thermodynamic-limit investigation (as, e.g., in Ref. [19]), and could represent, in principle, a very simple tool to check the equivalence between finite-size and thermodynamic-limit physics.

## 2. Schmidt gap

As already mentioned above, the Schmidt gap $\Delta_{S}$ is defined as the difference between the two highest eigenvalues of the reduced density matrix. It has been very recently shown [24] to be related to the correlation length, i.e.,

$$
\begin{equation*}
\Delta_{S} \approx \xi^{-z} \tag{21}
\end{equation*}
$$

and therefore its Kibble-Zurek scaling is

$$
\begin{equation*}
\Delta_{S} \approx \tau^{-\frac{z v}{1+z v}} \tag{22}
\end{equation*}
$$



FIG. 8. (Color online) Main panel: Schmidt gap at the final instant of the evolution for $\tau \in[15,150]$ at different system sizes ( $L=80-150$, from bottom to top). Black line: $y \approx x^{-1 / 2}$. Inset: Schmidt gap at the final instant of the evolution for $\tau=100$ as a function of the system size $L$.

In our case the exponent of $\tau$ in Eq. (22) is $-1 / 2$.
In Fig. 8 we present the data for the scaling of the Schmidt gap at the end of the ramping as a function of $\tau$. At fixed $L$ the shape of each curve shows cusps as a function of $\tau$ (which correspond to crossings of the first two eigenvalues of the reduced density matrix) superimposed to an overall power-law decay. First we comment on these non-analyticities and then discuss the scaling with $\tau$. We suggest a possible relation of the zeros of the Schmidt gap with the dynamical phase transitions first discussed by Heyl et al. [40]. These transitions manifest themselves as periodic non-analyticities of the free-energy density in the thermodynamic limit. Although in the latter paper only the case of a sudden quench is explicitly considered, in the same reference a connection is done with the non-analytic behavior of the Loschmidt echo [19]. In that work they consider the case of finite-velocity ramping followed by


FIG. 9. (Color online) Main panel: dynamics of the entanglement entropy $S(t)$ for $L=50, h_{i}=0.5, h_{f}=1.5$, and different values of $\tau$ (dashed vertical line: location of the critical point). Inset, black dots: period of oscillation of the entropy as a function of the gap at $h=h_{f}\left(h_{i}=0.5, h_{f}=1.41,1.42, \ldots, 1.5\right)$ with $\tau=50$. Red line: fit with $y=a_{0}+a_{1} / x$.


FIG. 10. (Color online) Dynamics of the entanglement spectrum with $L=50, h_{i}=0.5$, and $h_{f}=1.5$. Red/blue/green/brown line: dynamical first/second/third/fourth eigenvalue of the reduced density matrix of the half chain. Panels (a)-(f): $\tau=500,0.1,5,30,40,100$. The cyan lines in panel (a) show the ground-state values of the first four eigenvalues. In panels (b) and (c) black and red, green and blue data are almost coincident.
an evolution with constant magnetic field [19], i.e., a protocol similar to Eq. (16). We have verified that applying the same protocol, during the evolution at fixed magnetic field, the Schmidt gap as a function of time shows periodic cusps, with the period depending on $\tau$, in analogy with the Loschmidt echo [19]. In the case of a sudden quench, cusps of the Schmidt gap are also found in Ref. [41] by Torlai et al., who also suggest a relation with the dynamical phase transitions. We then suggest that the non-analyticities of the Schmidt gap at $t_{f}=\tau\left|h_{f}-h_{i}\right|$ for different $\tau$ shown in Fig. 8 may be interpreted as an early-time indication of dynamical phase transition.

Concerning the power-law scaling as a function of $\tau$, as expected from the behavior of the entanglement entropy, for each $\tau$ the Schmidt gap tends to converge to a finite value increasing the size $L$, as shown in the inset of Fig. 8. We expect


FIG. 11. (Color online) Schmidt gap at the final instant of the evolution for $\tau \in[1,200]$ at different system sizes for a ramping from $h_{i}=1.5$ to $h_{f}=0.5$.
the scaling function of Eq. (22) to be compatible with the numerical results. What we actually find is something more, i.e., that Eq. (22) almost perfectly interpolates the maxima of the curves at large system sizes.

## IV. FERROMAGNET TO PARAMAGNET

In this section, we perform the same analysis discussed in Sec. III, but now for the transition from the ferromagnetic to the paramagnetic region. This transition is much less studied in literature than the previous one (but see Ref. [35]), since it is not related to the Kibble-Zurek mechanism; however, it is still interesting to consider it in the present work.

## A. Initial structure of the entanglement spectrum

Let us consider now the Hamiltonian in Eq. (4), for simplicity, with $h=0$. The ground space takes the form span $\{\mid$ up $\rangle, \mid$ down $\rangle\}$, with $\mid$ up $\rangle \equiv \prod_{j=1}^{L}|\uparrow\rangle_{j}$ and $\mid$ down $\rangle \equiv$ $\prod_{j=1}^{L}|\downarrow\rangle_{j}$. At finite size, the ground state always belongs to the $\alpha=1$ sector (see Appendix A): therefore, since $\sigma^{x}|\uparrow\rangle=|\downarrow\rangle$ and vice versa, it is easily seen to be (up to a phase)

$$
\begin{equation*}
\left.\left.|\mathrm{GS},+\rangle \equiv \frac{1}{\sqrt{2}}(\mid \text { up }\rangle+\mid \text { down }\right\rangle\right) \tag{23}
\end{equation*}
$$

The zero-temperature density matrix of the system is therefore

$$
\begin{equation*}
\rho=|\mathrm{GS},+\rangle\langle\mathrm{GS},+| \tag{24}
\end{equation*}
$$

and the correspondent reduced density matrix is, for general $A$,

$$
\left.\left.\rho_{A}=\frac{1}{2}(\mid \text { up }\rangle_{A}, \quad \mid \text { down }\right\rangle_{A}\right) \mathbb{I}_{2}\left(\begin{array}{l}
{ }_{A}\langle\text { up }|  \tag{25}\\
A
\end{array} \text { down } \mid\right) .
$$

The initial entanglement structure is much simpler than in the other case, just two equally weighted eigenstates playing a role.

## B. Dynamics of entanglement entropy and entanglement spectrum

We work with $L=50$ and set the initial value of the magnetic field to $h_{i}=0.5$. The reason of this choice is that the entanglement spectrum is practically identical to the predicted one at $h=0$, with just two equal eigenvalues different from zero, but we can take the advantage of solving the ODEs to shorter times at fixed $\tau$. The final value of the magnetic field is set to $h_{f}=1.5$.

Similarly to what found in Sec. III, we observe the two limiting regimes, the adiabatic and the sudden one, respectively for large and small values of $\tau$ (see Figs. 9 and 10) and, at intermediate values of $\tau$ we find the regimes described in Sec. III. Remarkably, as already mentioned, oscillations in the entanglement entropy are present (see Fig. 9) and, similarly to what we discussed in Sec. III, their origin can be traced to the partially excited nature of the wave function after passing the critical point. Indeed, we can let the system evolve with protocol of Eq. (16) and study the period of the oscillations of the entropy when $t>\left(h_{f}-h_{i}\right) \tau$ as a function of the dynamical gap at the final value of the magnetic field $h=h_{f}$. Analogously to the paramagnetic to ferromagnetic quench, we find that the period scales inversely proportional to the gap, as we show in the inset of Fig. 9.

The four largest eigenvalues of the entanglement spectrum are degenerate in pairs before crossing the critical point [see all the panels of Fig. 10 for $\left.\left|h(t)-h_{i}\right|<0.5\right]$. After crossing the transition, the structure of the eigenvalues changes, so that only the second and the third eigenvalues are degenerate in the adiabatic limit, as we show in Fig. 10(a). For slow rampings, these eigenvalues cross each other with a regular pattern analogous to that of Figs. 3(a) and 3(f). The Schmidt gap is finite in this limit, apart from cusps originating from the oscillation of $\lambda_{1}$ and $\lambda_{2}$, as shown in Fig. 11 for $\tau \gtrsim 30$. On the contrary, if the ramping of the magnetic field is fast, the two pairs of eigenvalues remain almost degenerate, so that the Schmidt gap is several orders of magnitude smaller than in the adiabatic limit, as shown in Fig. 11 for $\tau \lesssim 10$.

We conclude this section with two comments: first, going from the sudden to the adiabatic limit, the Schmidt gap increases sharply (see Fig. 11 for $\tau \sim 10-40$ ); second, comparing Figs. 10 and 11 at $L=50, \tau=30$, it emerges that the crossing of the first and the second eigenvalue, giving rise to the cusp in the Schmidt gap, is isolated, because it corresponds to a narrow region of $\tau$ where the degeneracy of $\lambda_{1}$ and $\lambda_{2}$ is only slightly lifted, but for slower quenches these crossings disappear.

## V. CONCLUSIONS

In this work we have examined the dynamical evolution of the quantum Ising chain in a transverse magnetic field by looking at the entanglement entropy and entanglement spectrum, in particular in the case of a ramping from the paramagnetic to the ferromagnetic phase, and vice versa. We made the Hamiltonian time-dependent by letting the magnetic
field vary linearly in time with a varying time scale $\tau$, obtaining three qualitatively different regimes: an adiabatic one (large $\tau$ ) when the system evolves according the instantaneous ground state; a sudden quench (small $\tau$ ) when the system is essentially frozen to its initial state; and an intermediate one, where complicated behaviors occur. In particular, the most interesting feature we observe is the arising, in this regime, of dynamical multiple crossings of the first Schmidt eigenvalues: this effect is partially understood by means of an analysis of the fine structure of the corresponding eigenvectors, even if its general explanation, together with its observation in different models, is still missing. However, the physics of the dynamical evolution is well understood by looking at the behavior in time of the entanglement spectrum, starting from which one can study both universal quantities (scaling exponents) and physical phenomena, such as the Kibble-Zurek mechanism, that may manifest during the evolution.

We may conclude that entanglement entropy and entanglement spectrum seem to be, for the dynamical evolution as in the static case, a powerful tool to investigate the physics of a closed quantum many-body system crossing a phase transition at $T=0$. We have explicitly used this technique to study a paradigmatic exactly integrable system such as the quantum Ising chain in a transverse magnetic field, but investigation is under way to examine different situations where we either break integrability and/or introduce disorder.

We recently learned that two independent studies of the dynamics of the entanglement spectrum in one-dimensional models have appeared [34,41].

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## APPENDIX A: EXACT SOLUTION OF THE ISING MODEL

In this appendix we show how to diagonalize the Hamiltonian in Eq. (4). We follow Ref. [4] quite closely.

By defining the raising and lowering operators $\sigma_{j}^{ \pm} \equiv\left(\sigma_{j}^{z} \mp\right.$ $\left.i \sigma_{j}^{y}\right) / 2$, Eq. (4) reads
$H=-\frac{1}{2} \sum_{j=1}^{L}\left[\left(\sigma_{j}^{+} \sigma_{j+1}^{+}+\sigma_{j}^{+} \sigma_{j+1}^{-}+\right.\right.$H.c. $\left.)+2 h \sigma_{j}^{+} \sigma_{j}^{-}\right]+\frac{L h}{2}$.

Performing a Jordan-Wigner transformation by means of

$$
\begin{align*}
c_{j} & \equiv \prod_{k=1}^{j-1}\left(2 \sigma_{k}^{+} \sigma_{k}^{-}-1\right) \sigma_{j}^{+} \\
c_{j}^{\dagger} & \equiv \sigma_{j}^{-} \prod_{k=1}^{j-1}\left(2 \sigma_{k}^{+} \sigma_{k}^{-}-1\right) \tag{A2}
\end{align*}
$$

allows us to rewrite the Hamiltonian Eq. (A1) in fermionic form:

$$
\begin{align*}
H= & -\frac{1}{2} \sum_{j=1}^{L-1}\left[c_{j+1}^{\dagger} c_{j}+c_{j+1} c_{j}+\text { H.c. }\right] \\
& +\frac{\alpha}{2}\left[c_{1}^{\dagger} c_{L}+c_{1} c_{L}+\text { H.c. }\right]+h \sum_{j=1}^{L} c_{j}^{\dagger} c_{j}-\frac{L h}{2} \tag{A3}
\end{align*}
$$

where $\alpha \equiv \prod_{j=1}^{L}\left(1-2 c_{j}^{\dagger} c_{j}\right)=\prod_{j=1}^{L} \sigma_{j}^{x}$. It is easy to show that $\alpha$ commutes with $H$, and therefore it is a constant of motion; moreover, $\alpha^{2}=1$, so that $\alpha= \pm 1$. As is manifest from its definition, the case $\alpha= \pm 1$ corresponds to the case in which in the chain an even/odd number of down spins is present, and, in fermionic language, to a chain with antiperiodic/periodic boundary conditions (APBCs/PBCs) and an even/odd number of fermions. We choose to work in the sector of even parity in the number fermions, i.e., $\alpha=1$, being, at finite size, the ground state of the model always in this sector. One ends up with

$$
\begin{equation*}
H=-\frac{1}{2} \sum_{j=1}^{L}\left[\left(c_{j+1}^{\dagger} c_{j}+c_{j+1} c_{j}+\text { H.c. }\right)-2 h c_{j}^{\dagger} c_{j}\right]-\frac{L h}{2} \tag{A4}
\end{equation*}
$$

with fermions satisfying APBCs. The diagonalization proceeds by means of a Fourier transform

$$
\begin{equation*}
c_{j} \equiv \frac{e^{i \pi / 4}}{\sqrt{L}} \sum_{m=0}^{L-1} e^{i p_{m} j} d_{m} \tag{A5}
\end{equation*}
$$

with $p_{m} \equiv 2 \pi(m+1 / 2) / L$, in order to automatically implement the APBCs. With some algebra, it is possible to show that the Hamiltonian takes the form

$$
\begin{equation*}
H=\frac{1}{2} \sum_{m=0}^{L-1}\left(d_{m}^{\dagger}, d_{L-m-1}\right) M_{m}\binom{d_{m}}{d_{L-m-1}^{\dagger}} \tag{A6}
\end{equation*}
$$

with

$$
M_{m} \equiv\left(\begin{array}{cc}
A_{m} & -B_{m}  \tag{A7}\\
-B_{m} & -A_{m}
\end{array}\right)
$$

and

$$
\begin{equation*}
A_{m} \equiv h-\cos p_{m}, \quad B_{m} \equiv \sin p_{m} \tag{A8}
\end{equation*}
$$

that, remarkably, satisfy $A_{L-m-1}=A_{m}, B_{L-m-1}=-B_{m}$; i.e., the Hamiltonian decouples into the sum of $L$ noninteracting modes, each one independently diagonalizable.

The last step of the procedure consists of a Bogolyubov transformation, which puts each $M_{m}$ in diagonal form. The eigenvalues of each $M_{m}$ are given by the two values $\pm E_{m}$, with

$$
\begin{equation*}
E_{m}=\sqrt{A_{m}^{2}+B_{m}^{2}} \tag{A9}
\end{equation*}
$$

and the orthogonal transformation $U_{m}$ making $M_{m}$ diagonal, i.e., giving $U_{m}^{\dagger} M_{m} U_{m}=\operatorname{diag}\left(E_{m},-E_{m}\right)$, is given by

$$
U_{m} \equiv\left(\begin{array}{cc}
u_{m} & v_{m}  \tag{A10}\\
-v_{m} & u_{m}
\end{array}\right)
$$

where

$$
\begin{equation*}
u_{m}=\frac{-(-1)^{m} \frac{A_{m}+E_{m}}{B_{m}}}{\sqrt{1+\left(\frac{A_{m}+E_{m}}{B_{m}}\right)^{2}}}, v_{m}=\frac{-(-1)^{m}}{\sqrt{1+\left(\frac{A_{m}+E_{m}}{B_{m}}\right)^{2}}} \tag{A11}
\end{equation*}
$$

satisfying $u_{L-m-1}=u_{m}, v_{L-m-1}=-v_{m}$. The diagonalizing operators are

$$
\begin{equation*}
\binom{b_{m}}{b_{L-m-1}^{\dagger}} \equiv U_{m}\binom{d_{m}}{d_{L-m-1}^{\dagger}} \tag{A12}
\end{equation*}
$$

and the orthogonality of $U_{m}$ ensures their fermionic nature. The Hamiltonian takes, by means of the inverse of Eq. (A12), the final form

$$
\begin{equation*}
H=\sum_{m=0}^{L-1} E_{m}\left(b_{m}^{\dagger} b_{m}-\frac{1}{2}\right) \tag{A13}
\end{equation*}
$$

and its ground state is, for $\alpha=1$, the vacuum state $|\mathrm{GS}\rangle$ such that $b_{m}|\mathrm{GS}\rangle=0$. Excited states, in the APBC sector, are obtained by applying couples of Bogolyubov creation operators on $|\mathrm{GS}\rangle$.

## APPENDIX B: DYNAMICS OF THE ISING MODEL

In this appendix, we show how to describe the dynamics of a state according to the Hamiltonian in Eq. (4). We follow the procedure of Ref. [42].

The time evolution of the system in Eq. (4) is described by the Heisenberg equation for the $c$ operators:

$$
\begin{equation*}
i \frac{d}{d t} c_{j, H}(t)=\left[c_{j, H}(t), H_{j, H}(t)\right] \tag{B1}
\end{equation*}
$$

which can be rewritten as

$$
\begin{equation*}
i \frac{d}{d t} c_{j, H}(t)=\sum_{k=1}^{L}\left[A_{j k}(t) c_{k, H}(t)+B_{j k}(t) c_{k, H}^{\dagger}(t)\right] \tag{B2}
\end{equation*}
$$

with

$$
\begin{align*}
A_{j k}(t) & \equiv h(t) \delta_{j k}-\frac{1}{2}\left(\delta_{j, k+1}+\delta_{j+1, k}-\delta_{j 1} \delta_{k L}-\delta_{j L} \delta_{k 1}\right) \\
B_{j k}(t) & \equiv-\frac{1}{2}\left(\delta_{j+1, k}-\delta_{j, k+1}+\delta_{j 1} \delta_{k L}-\delta_{j L} \delta_{k 1}\right) \tag{B3}
\end{align*}
$$

In order to solve such an equation, we make the following ansatz, known as the time-dependent Bogolyubov transformation:

$$
\begin{equation*}
c_{j, H}(t) \equiv \sum_{m=0}^{L-1}\left[u_{j m}(t) b_{m}+v_{j m}^{*}(t) b_{m}^{\dagger}\right] \tag{B4}
\end{equation*}
$$

with the initial conditions $u_{j m}(0)=u_{j m}$ and $v_{j m}(0)=v_{j m}$ given by the exact solution:

$$
\begin{align*}
& u_{j m} \equiv \frac{1}{\sqrt{L}} e^{i\left(p_{m} j+\frac{\pi}{4}\right)} u_{m}, \\
& v_{j m} \equiv \frac{1}{\sqrt{L}} e^{i\left(p_{m} j+\frac{\pi}{4}\right)} v_{m} \tag{B5}
\end{align*}
$$

By putting the ansatz of Eq. (B4) in the Heisenberg equation, we come to the set of linear coupled ODEs

$$
\begin{align*}
i \frac{d}{d t} u_{j m}(t) & =\sum_{k=1}^{L}\left[A_{j k}(t) u_{k m}(t)+B_{j k}(t) v_{k m}(t)\right]  \tag{B6}\\
-i \frac{d}{d t} v_{j m}(t) & =\sum_{k=1}^{L}\left[B_{j k}(t) u_{k m}(t)+A_{j k}(t) v_{k m}(t)\right] .
\end{align*}
$$

## APPENDIX C: BIPARTITE QUANTITIES IN FREE FERMIONIC SYSTEMS

In this appendix we review computing the entanglement entropy and the entanglement spectrum for free fermionic systems.

As is known from recent literature [43,44] (see also Refs. [45,46]), for fermionic biquadratic (static) Hamiltonians the density matrix can be obtained from correlation functions. In order to evaluate the time evolution of the entanglement entropy and spectrum we need a step forward, which is the introduction of Majorana fermions:

$$
\begin{align*}
& \bar{c}_{2 m-1}=c_{m}^{\dagger}+c_{m}  \tag{C1}\\
& \bar{c}_{2 m}=i\left(c_{m}^{\dagger}-c_{m}\right) \tag{C2}
\end{align*}
$$

which satisfy anticommutation rules $\left\{\bar{c}_{r}, \bar{c}_{s}\right\}=2 \delta_{r s}$. The correlation matrix of the Majorana fermions has the form

$$
\begin{equation*}
\left\langle\bar{c}_{r} \bar{c}_{s}\right\rangle=\delta_{r, s}+i \Gamma_{r s} \tag{C3}
\end{equation*}
$$

where $r, s=1, \ldots, 2 \ell$. The matrix $\Gamma_{r s}$ is antisymmetric and its eigenvalues are purely imaginary $\pm i v_{r}, r=1, \ell$. It can be shown that this matrix describes a set of uncorrelated (true) fermions $\left\{a_{m}\right\}$ satisfying

$$
\begin{equation*}
\left\langle a_{m} a_{n}\right\rangle=0, \quad\left\langle a_{m}^{\dagger} a_{n}\right\rangle=\delta_{m n} \frac{1+v_{n}}{2} \tag{C4}
\end{equation*}
$$

Each of the $\ell$ blocks is then in the state $\rho_{j}=p_{j} a_{j}^{\dagger}|0\rangle\langle 0| a_{j}+$ $\left(1-p_{j}\right)|0\rangle\langle 0|$, with $p_{j}=\left(1+v_{j}\right) / 2$ so that the entropy is the sum of the single-particle entropies, thus yielding for the reduced $\ell$-site system

$$
\begin{equation*}
S(\ell)=\sum_{j=1}^{l} H_{2}\left(\frac{1+v_{j}}{2}\right) \tag{C5}
\end{equation*}
$$

where $H_{2}(x) \equiv-x \log _{2} x-(1-x) \log _{2}(1-x)$. The eigenvalues $\lambda_{j}, \quad j=1, \ldots, 2^{\ell}$ of the reduced density matrix can in principle be found by taking properly chosen products of either $p_{j}$ or $\left(1-p_{j}\right)$, with $j=1, \ldots, \ell$ [47].

The procedure described above works equally well for the time-dependent case, provided that the Majorana fermions are constructed using the time-evolved true fermions $c_{i, H}(t)$. In this way we can obtain the time-dependent entropy $S(\ell, t)$ and entanglement spectrum $\lambda_{i}(\ell, t)$.
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