

# Entanglement Growth from Entangled States: A Unified Perspective on Entanglement Generation and Transport

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Studies of entanglement dynamics in quantum many-body systems have focused largely on initial product states. Here, we investigate the far richer dynamics from initial entangled states, uncovering universal patterns across diverse systems ranging from many-body localization (MBL) to random quantum circuits. Our central finding is that the growth of entanglement entropy can exhibit a counterintuitive nonmonotonic dependence on the initial entanglement in many nonergodic systems, peaking for moderately entangled initial states. To understand this phenomenon, we introduce a conceptual framework that decomposes entanglement growth into two mechanisms: “build” and “move.” The “build” mechanism creates new entanglement, while the “move” mechanism redistributes preexisting entanglement throughout the system. Specifically, we demonstrate that MBL dynamics are “move-dominated,” exhibiting a quantitative agreement with a random SWAP circuit that serves as a model of pure “move” dynamics by uniformly distributing preexisting entanglement. This implies that MBL acts as a redistributor of a hidden entanglement reservoir quantified by the bipartition-averaged entropy. This “build-move” framework offers a unified perspective for classifying diverse physical dynamics, deepening our understanding of entanglement propagation and information processing in quantum many-body systems.

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*Introduction*—Entanglement is a cornerstone of modern physics [1–3] and specifically a key diagnostic for non-equilibrium quantum dynamics [4–10]. The standard approach investigates the time evolution of entanglement entropy (EE) from an initial product state, revealing a rich landscape of dynamical behaviors. In thermalizing systems and generic random quantum circuits (RQCs), entanglement grows rapidly, often linearly in time [11–26]. By contrast, many-body localized (MBL) systems [27–63] exhibit a much slower, logarithmic growth [64–72], while integrable systems saturate to nonthermal values [73–77] and non-interacting Anderson localized (AL) systems show severely suppressed entanglement growth [78–90]. However, this fruitful paradigm inherently conflates two distinct aspects of entanglement dynamics: the initial generation of entangled resources and the subsequent transport of this entanglement. The observed half-chain entanglement entropy (HCEE) growth is thus always a composite effect, leaving the underlying pattern of entanglement propagation obscured.

To deconstruct this process and isolate the intrinsic entanglement handling capabilities of a given system, we shift the paradigm to study the evolution of initial states that already possess volume-law entanglement, which we term

partially thermalized states. These states are thus prepared to contain a rich, preexisting reservoir of entanglement, providing a nontrivial substrate for the subsequent dynamics. This paradigm shift allows us to move from the question “How fast is entanglement created?” to a more profound one: “How does a system’s inherent dynamics process and redistribute preexisting entanglement?” This setup facilitates a deep probe into the intricate interplay between a system’s fundamental character and an existing, complex entanglement structure.

In this Letter, we address these questions and uncover a remarkable phenomenon: in MBL dynamics, the HCEE growth exhibits a nonmonotonic behavior with its initial value, profoundly differing from thermalized systems [see Fig. 1(d)]. To rationalize the observed behaviors, we introduce a novel conceptual framework that decomposes HCEE growth into two primary mechanisms: “build” for creating new entanglement and “move” for redistributing preexisting entanglement, as schematically shown in Figs. 1(a) and 1(b), respectively. We demonstrate that this framework not only quantitatively explains our results by identifying MBL as a “move-dominated” dynamics but also provides a unified perspective for classifying diverse quantum dynamics.

*Model and methods*—We study the dynamics in a one-dimensional spin-1/2 chain of even length  $L$  using exact diagonalization. The central quantity of interest is the

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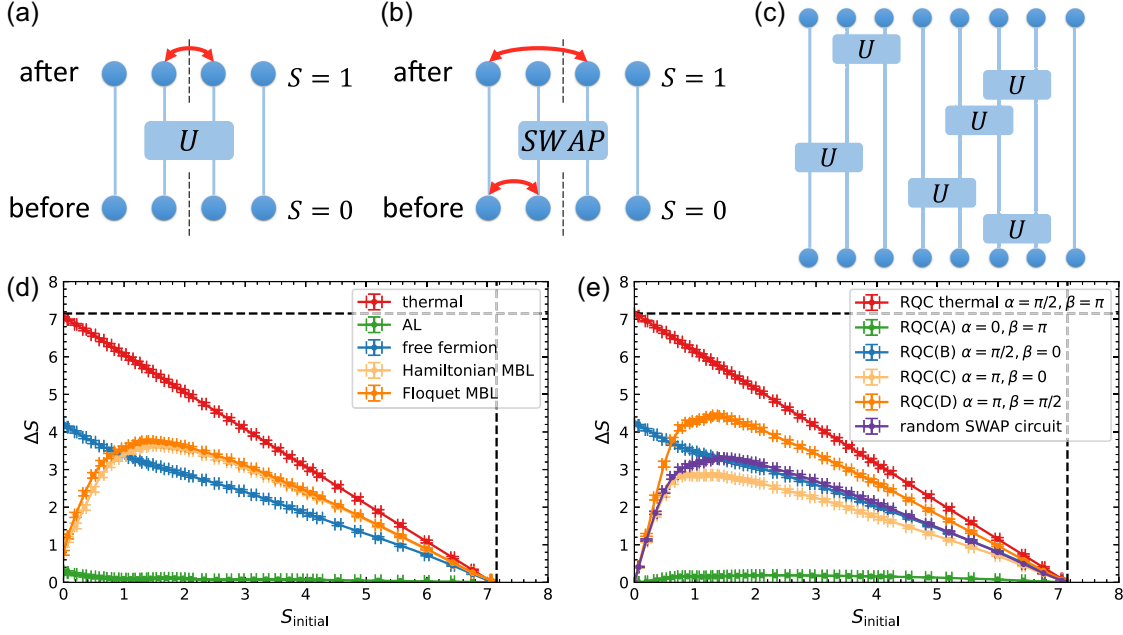


FIG. 1. “Build-move” framework and the increase of HCEE in various dynamics. (a)–(c) Schematic illustrations of “build” mechanism (a), “move” mechanism (b), and RQC composed of two-qubit gates  $U$  (c). Blue circles represent spins, and pairs of spins linked by red double arrows indicate Bell pairs. (d),(e) Classification of dynamics based on the HCEE growth:  $\Delta S = S_{\text{sat}} - S_{\text{initial}}$  as a function of the initial HCEE,  $S_{\text{initial}}$ . All data points are for system size  $L = 16$ .  $S_{\text{initial}}$  is controlled by varying the preparation time  $\tau$ . (d) Results for Hamiltonian-based dynamics. (e) Results for RQC-based dynamics. Note the striking similarity between the MBL curve in (d) and the Random SWAP curve in (e). The two-qubit gates  $U$  are defined in Eq. (2), with parameters in various classes detailed in Table I. For the calculation details of the data in all figures in the main text, please refer to Supplemental Material, Sec. I [91].

HCEE measured in bits:  $S = -\text{Tr}(\hat{\rho}_{\text{HC}} \log_2 \hat{\rho}_{\text{HC}})$ , defined as the von Neumann entropy of the first  $L/2$  spins whose reduced density operator is  $\hat{\rho}_{\text{HC}}$ . The various physical processes in our study, from initial state preparation to subsequent dynamics, are mainly based on the one-dimensional disordered XXZ Hamiltonian with open boundary conditions:

$$\hat{H} = \sum_{i=1}^{L-1} \left( \hat{S}_i^x \hat{S}_{i+1}^x + \hat{S}_i^y \hat{S}_{i+1}^y + J_z \hat{S}_i^z \hat{S}_{i+1}^z \right) + \sum_{i=1}^L h_i \hat{S}_i^z. \quad (1)$$

Here,  $\hat{S}_i^\alpha$  ( $\alpha = x, y, z$ ) are the spin-1/2 operators at site  $i$ , and we set the anisotropy  $J_z = 0.5$ . The on-site fields  $h_i$  are independent random variables drawn uniformly from the interval  $[-W, W]$ . This Hamiltonian has a  $U(1)$  symmetry which conserves the total magnetization in the  $z$  direction,  $\hat{S}_{\text{tot}}^z = \sum_{i=1}^L \hat{S}_i^z$ . All calculations are performed in the half-filling sector  $\hat{S}_{\text{tot}}^z = 0$ . The system is in the thermal and MBL phases at  $W = 0.5$  and  $W = 5.0$ , respectively [29] [see Supplemental Material (SM), Sec. II [91] for details].

Our primary focus is on the evolution from initial states with varying degrees of entanglement. We prepare initial states  $|\psi(\tau)\rangle$  by evolving a product state  $|\psi_0\rangle$  under the thermal Hamiltonian  $\hat{H}(W = 0.5)$  for a duration  $\tau$ .  $|\psi_0\rangle$  is randomly chosen from the computational basis states within the half-filling sector. The evolution time  $\tau$  directly

controls the initial entanglement and the level of thermalization as  $|\psi(\tau)\rangle = e^{-i\tau\hat{H}(W=0.5)}|\psi_0\rangle$ . Subsequently, we examine the dynamics of these prepared states under the following  $U(1)$ -symmetric dynamical protocols: (1) Thermal dynamics:  $|\psi(\tau)\rangle$  evolves under the thermal Hamiltonian  $\hat{H}(W = 0.5)$ . (2) Hamiltonian MBL dynamics:  $|\psi(\tau)\rangle$  evolves under the MBL Hamiltonian  $\hat{H}(W = 5.0)$ . (3) Free fermion dynamics: evolution under the free fermion Hamiltonian  $\hat{H}_{\text{XY}}$  with  $J_z = 0$  and  $W = 0$  in Eq. (1). (4) Floquet MBL dynamics: a periodic drive is applied using the Floquet operator  $\hat{F} = e^{-iT_0\hat{H}_0}e^{-iT_1\hat{H}_{\text{XY}}}$  [57]. Here,  $\hat{H}_0 = \sum_{i=1}^{L-1} \hat{S}_i^z \hat{S}_{i+1}^z + \sum_{i=1}^L h_i \hat{S}_i^z$ , with disorder strength  $W = 5.0$ ,  $T_0 = 1.0$ , and  $T_1 = 0.4$ . (5) AL dynamics:  $|\psi(\tau)\rangle$  evolves under the noninteracting AL Hamiltonian, with  $J_z = 0$  and  $W = 5.0$  in Eq. (1). (6) RQC protocols: we employ RQCs as a paradigmatic model for understanding universal aspects of entanglement dynamics [15,17,22,23,102–114]. Our implementation consists of circuits where, at each step, a single two-qubit gate  $U$  is applied to a randomly selected adjacent pair of spins ( $i, i + 1$ ) from the  $L - 1$  bonds [Fig. 1(c)]. The two-qubit unitary is given by

$$\hat{U}_{i,i+1} = e^{-i\alpha(\hat{S}_i^x \hat{S}_{i+1}^x + \hat{S}_i^y \hat{S}_{i+1}^y)} e^{-i\beta \hat{S}_i^z \hat{S}_{i+1}^z}. \quad (2)$$

**Results**—To analyze the entanglement dynamics under different protocols, we first introduce a conceptual

framework that decomposes the growth of HCEE into two fundamental mechanisms: “build” and “move.” As shown in Fig. 1(a), the “build” mechanism creates new entanglement directly. For instance, a two-qubit unitary gate  $U$  across the central bond can generate a Bell pair from a product state, increasing the HCEE from 0 to 1. The “move” mechanism, depicted in Fig. 1(b), involves the redistribution of preexisting entanglement. Here, an entangled pair within one subsystem is transported across the central bond, for instance, via a SWAP gate, also raising HCEE from 0 to 1. While both mechanisms often occur simultaneously in a generic interacting system, their proportional contributions are highly dependent on the specific dynamics.

To isolate the “move” mechanism, we employ the random SWAP circuit as an idealized model, which can be achieved up to a global phase by setting  $\alpha = \beta = \pi$  in our RQC protocols. The key insight is that a SWAP operation merely exchanges the state of two spins. This allows us to equivalently view the action of the random SWAP circuit on a fixed bipartition as an evolution of the bipartition itself across a static state. This evolution generates transitions between the  $N = \binom{L}{L/2}/2$  distinct bipartitions, establishing an ergodic Markov chain with a uniform stationary distribution (see the proof in SM, Sec. III [91]). Consequently, at sufficiently large circuit depth, every bipartition becomes equally probable. This directly implies that the steady-state HCEE under the random SWAP circuit corresponds precisely to the initial state’s *bipartition-averaged entanglement entropy* (BAEE)—the entanglement entropy averaged across all  $N$  possible equal-size bipartitions:  $\bar{S} = (1/N) \sum_{A \in \mathcal{P}} S_A$ , where  $\mathcal{P}$  is the set consisting of subsystems corresponding to all  $N$  possible bipartitions and  $S_A$  represents the entropy of the subsystem  $A$  on  $\mathcal{P}$ . The BAEE provides a more robust measure of the state’s overall entanglement, invariant under SWAP dynamics, which only redistributes existing entanglement. This model thus establishes a baseline for move-driven dynamics, enabling quantitative comparisons with more complex protocols. For more analysis about the BAEE, see SM, Sec. VI [91].

For RQC protocols of  $\alpha, \beta \in [0, 2\pi]$ , the specific points  $(\alpha, \beta) = (0, 0), (0, 2\pi), (2\pi, 0),$  and  $(2\pi, 2\pi)$  are trivial as the operator  $\hat{U}_{i,i+1}$  can be decomposed into a tensor product of single-qubit gates. Generic choices of  $(\alpha, \beta)$  are expected to lead to full thermalization, where the HCEE saturates near the Haar measure average value in the half-filling sector [115–119], which is marked by dashed lines in Figs. 1(d), 1(e), 2, and 3. Nonthermalization behavior occurs only for a small set of special parameter values, which we classify into five categories based on the distinct saturation values of their HCEE. Within each category, different parameter choices for  $(\alpha, \beta)$  lead to dynamics with almost the same saturating HCEE value. The random SWAP circuit mentioned above is one of these five

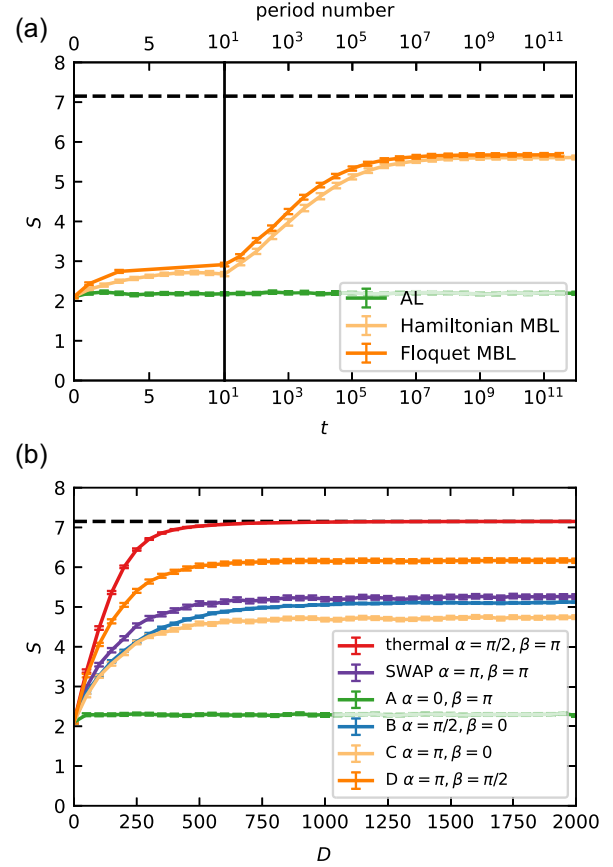


FIG. 2. Time evolution of HCEE  $S$  with initial state  $|\psi(\tau)\rangle$  for  $L = 16$  and  $\tau = 4.5$ . (a) Evolution under localized dynamics. The lower axis shows time  $t$  for AL and Hamiltonian MBL evolutions, and the upper axis shows the period number for the Floquet MBL dynamics. To visualize the dynamics across multiple timescales, the horizontal axis uses a hybrid scale of linear time for the initial evolution ( $t, \text{period number} < 10$ ) and logarithmic for long times. (b) Evolution under RQCs dynamics. The horizontal axis represents the circuit depth. The specific  $(\alpha, \beta)$  values shown in the legend are the parameters of the quantum gates selected from their respective categories.

categories. The parameter setting for these categories is summarized in Table I.

Figure 2(a) displays the localized dynamics for a long time starting from  $|\psi(\tau)\rangle$ , where the evolution under different nonthermal dynamics reveals distinct behaviors. For both the Hamiltonian MBL and the Floquet MBL protocols, the HCEE exhibits a characteristic slow, logarithmic growth in time. It eventually reaches a saturation value that, while significantly higher than the initial HCEE, remains well below the Haar measure average value in the half-filling sector. In sharp contrast, the dynamics governed by the noninteracting AL Hamiltonian [Fig. 2(a)] and the RQC(A) protocol [Fig. 2(b)] both show almost no further increase in HCEE, among which the former confirms that interactions are crucial for moving the entanglement.

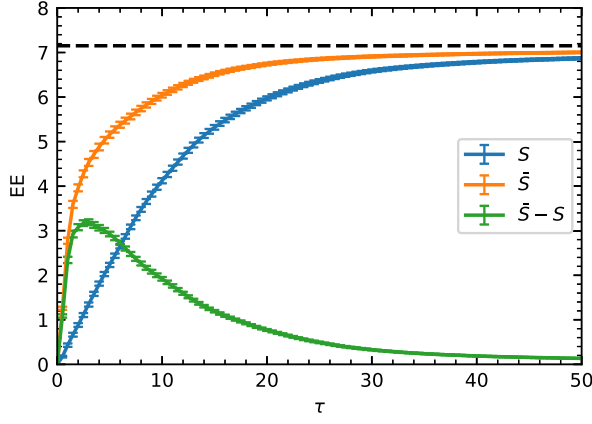


FIG. 3. Evolution of BAEE  $\bar{S}$ , HCEE  $S$ , and their difference under thermal Hamiltonian  $\hat{H}(W = 0.5)$  for  $L = 16$ , starting from the product state  $|\psi_0\rangle$ . The data point corresponding to the maximum value of  $\bar{S} - S$  is at  $\tau = 3.0$  with HCEE  $S \approx 1.34$ , roughly coinciding with the optimal  $S_{\text{initial}}$  for the peak  $\Delta S$  of move-dominated dynamics in Figs. 1(d) and 1(e).

The other nonthermal RQC protocols, RQC(B)–(D), all facilitate HCEE growth, but with distinct steady values.

To systematically quantify the relationship between initial entanglement and subsequent entanglement growth, we analyze the total increase of HCEE,  $\Delta S = S_{\text{sat}} - S_{\text{initial}}$ , as a function of  $S_{\text{initial}}$ . The results, plotted in Figs. 1(d) and 1(e), reveal that the diverse dynamical protocols can be classified into three classes based on their monotonicity, which reflects the capability to generate and transport entanglement. This classification not only is supported by the macroscopic behavior of  $\Delta S$  but also is consistent with the microscopic action of the gate for the RQC dynamics.

*Move-dominated dynamics*—The first class of protocols is defined by a predominant “move” mechanism, whose fundamental role is to redistribute existing entanglement. As depicted in Fig. 1(d), the dynamics of MBL clearly exhibit a nonmonotonic dependence of  $\Delta S$  on  $S_{\text{initial}}$ . This behavior, which differs significantly from that in thermal or AL dynamics, is a key finding of this Letter. For nearly product states ( $S_{\text{initial}} \approx 0$ ),  $\Delta S$  is negligible due to the minimal entanglement available for transport. It then increases to a peak for moderately entangled states before decreasing again as  $S_{\text{sat}}$  approaches the Haar measure average value. This distinct nonmonotonic behavior serves

as a hallmark of the move-dominated mechanism. More importantly, we find the quantitative agreement of the results from MBL and random SWAP circuits, although the time scales for entanglement saturation are distinct for these dynamics. This agreement is highly nontrivial: it implies that, despite the complex Hamiltonian origin of MBL, its effective entanglement dynamics approximately reduces to a simple shuffling of bipartitions. This class also encompasses RQC(C) and (D), as illustrated in Fig. 1(e). This is directly explained by their gate-level action, which functions similarly to a SWAP operator. For example, when an RQC(C) or (D) gate acts on the first two spins of the state  $|\uparrow\uparrow\downarrow\rangle + |\uparrow\downarrow\uparrow\rangle$ , the transformation yields  $e^{-i\beta/4}|\uparrow\uparrow\downarrow\rangle - ie^{i\beta/4}|\downarrow\uparrow\uparrow\rangle$ . This effectively converts entanglement from between the last two spins to between the first and third spins, thereby realizing the “move” operation at the microscopic level.

*Build and move hybrid dynamics*—The second class of dynamics exhibits a hybrid “build-and-move” character. In conventional thermal systems, a significant  $\Delta S$  emerges even from initial product states ( $S_{\text{initial}} = 0$ ), signifying the “build” mechanism that generates entanglement from scratch [Fig. 1(d)]. This entanglement increase,  $\Delta S$ , diminishes as  $S_{\text{initial}}$  rises. Similarly, free fermion systems also reach the equilibrium with weaker “build” capabilities due to extensive conserved quantities, resulting in a smaller  $\Delta S$ . Consistent with this, RQCs with general  $\alpha \in (0, \pi) \cup (\pi, 2\pi)$ , including both thermal RQC and RQC(B), also show the same monotonically decreasing trend of  $\Delta S$ , as shown in Fig. 1(e). The ability of these RQCs to effectively generate entanglement from product states—for instance, transforming  $|\uparrow\downarrow\rangle$  into the entangled state  $e^{i\beta/4}[\cos(\alpha/2)|\uparrow\downarrow\rangle - i\sin(\alpha/2)|\downarrow\uparrow\rangle]$ , further corroborates the pervasive “build” mechanism within this class. Specifically, RQC(B) with  $\beta = 0$  mimics free fermion dynamics. Moreover, with increasing  $S_{\text{initial}}$ , the “move” aspect for these systems also becomes a prominent contributor to the overall entanglement growth.

*Entanglement-inert dynamics*—Finally, the third class is essentially inert with respect to HCEE growth. Both AL and RQC(A) fall into this class, producing negligible  $\Delta S$  across the entire range of  $S_{\text{initial}}$  [see Figs. 1(d) and 1(e)]. This inert behavior stems from the lack of essential ingredients for either “build” or “move” contribution. For AL, the noninteracting and localized character of the Hamiltonian prevents entanglement from forming or

TABLE I. Classification of Hamiltonian-based and RQC-based dynamics on the behavior of  $\Delta S$  with  $S_{\text{initial}}$ .

Hamiltonian-based dynamics	RQC-based dynamics	Tendency of $\Delta S$ with $S_{\text{initial}}$
Thermalization	$\alpha \in (0, \pi) \cup (\pi, 2\pi), \beta \in (0, 2\pi)$	Monotonic decay
MBL	$\alpha = \pi, \beta = \pi$ (SWAP), $\{0, 2\pi\}$ (class C), other values (class D)	Nonmonotonic peak
Free fermion	$\alpha \in (0, \pi) \cup (\pi, 2\pi), \beta = \{0, 2\pi\}$ (class B)	Monotonic decay
AL	$\alpha = \{0, 2\pi\}, \beta \in (0, 2\pi)$ (class A)	Negligible

spreading. For RQC(A), the underlying gate is diagonal in the computational basis, which can only affect relative phases, thereby rendering it incapable of generating or transporting entanglement.

*Discussion and conclusion*—In this Letter, we systematically investigate entanglement dynamics across diverse quantum systems, moving beyond the conventional focus on product states to explore initial states that already possess entanglement. Our central finding is a universal three-class classification of entanglement growth. Most notably, we uncover a nonmonotonic dependence of HCEE increase on initial entanglement in MBL dynamics (we also considered the quantum sun model [120], more consistent results are detailed in SM [91]) and specific RQCs—a behavior qualitatively distinct from thermal dynamics. To fundamentally account for these novel results, we introduce the “build-move” framework. This framework offers a universal lens through which to understand complex entanglement growth by classifying dynamics based on their capacity to generate new entanglement (“build”) and effectively transport preexisting entanglement (“move”).

Our Letter yields another key insight by distinguishing between the entanglement of a single, fixed bipartition (HCEE) and the overall inherent entanglement in the quantum state, which we quantify using the BAEE. Figure 3 illustrates that, even in a generic thermalized system, the BAEE grows considerably faster than the HCEE during early dynamics. This initial divergence quantifies the rapid buildup of entanglement throughout the system at a local level, effectively establishing an accumulating intrasubsystem entanglement reservoir. This reservoir is central to understanding the initial increase of  $\Delta S$  with  $S_{\text{initial}}$  in move-dominated dynamics: a larger reservoir provides more entanglement potential to be “moved” across the central bond. Our results demonstrate that MBL dynamics effectively unlock this potential, converting the hidden reservoir  $\bar{S} - S$  into accessible half-chain entropy via the “move” mechanism. This picture is supported by a remarkable quantitative consistency: the value of  $S_{\text{initial}}$  where the entropy growth  $\Delta S$  peaks for MBL dynamics [Fig. 1(d)] aligns with the maximum difference between BAEE and HCEE. This alignment provides compelling evidence for the dominant “move” mechanism in MBL entanglement dynamics, fundamentally recasting MBL as a global redistributor of a “hidden” entanglement reservoir.

This perspective underscores the limitation of using a single bipartition’s EE to fully capture a system’s entanglement structure and highlights the importance of multipartition entanglement measures, a topic of growing interest [121–123]. Our Letter provides a clear dynamical context for this multipartition viewpoint. Moreover, the distinct dynamical signatures predicted by our framework present clear targets for experimental verification on near-term quantum simulation platforms. Finally, our results

have practical significance for numerical approaches: the saturation entanglement of MBL dynamics, which is reached only after exponentially long timescales, can be effectively approximated by the BAEE or random SWAP circuits. This provides a basis for utilizing efficient proxies to estimate steady-state properties while bypassing the prohibitive computational cost of simulating the full dynamical evolution.

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*Data availability*—The data that support the findings of this article are openly available [124].

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