Simulating a measurement-induced phase transition for trapped-ion circuits

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The rise of programmable quantum devices has motivated the exploration of circuit models which could realize novel physics. A promising candidate is a class of hybrid circuits, where entangling unitary dynamics compete with disentangling measurements. Novel phase transitions between different entanglement regimes have been identified in their dynamical states, with universal properties hinting at unexplored critical phenomena. Trapped-ion hardware is a leading contender for the experimental realization of such physics, which requires not only traditional two-qubit entangling gates, but also a constant rate of local measurements accurately addressed throughout the circuit. Recent progress in engineering high-precision optical addressing of individual ions makes preparing a constant rate of measurements throughout a unitary circuit feasible. Using tensor network simulations, we show that the resulting class of hybrid circuits, prepared with native gates, exhibits a volume-law to area-law transition. Our simulations are able to characterize the critical exponents using circuit sizes with tens of qubits and thousands of gates. We argue that this transition should be robust against additional sources of experimental noise expected in modern trapped-ion hardware and will rather be limited by statistical requirements on postselection. Our work highlights the powerful role that tensor network simulations can play in advancing the theoretical and experimental frontiers of critical phenomena.

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I. INTRODUCTION

The study of quantum criticality is experiencing a sudden transformation originating in the idea of measurementinduced phase transitions (MPTs) [1-3]. These exist in many-body qubit wave functions, where entangling unitary dynamics compete with nonunitary measurements that act to disentangle the state. Such phase transitions can be driven by a finite rate of measurement throughout the circuit, or by the choice of measurement bases [4,5] or the strength of weak measurements [6-9]. Numerical studies of a variety of random quantum circuits have demonstrated that such phase transitions share broad features of conventional conformal field theories (CFTs) [2,4,10-14], with some quantities depending on details of the circuit, e.g., the interaction range between the qubits [15,16]. The observation of conventional CFT features, such as a critical scaling in the entanglement entropy characterized by universal exponents, was surprising and distinguishes the observed phase transition from a classical crossover. While special cases have been related to well-known statistical or spin models [7,14,17,18], percolation theories [19,20], or minimal-cut pictures [21], a general comprehensive theoretical framework is still absent and no rigorous connection to CFT has yet been made. This leaves

MPTs are known to exist in models, and significant debates exist as to whether they can be realized in real experimental systems. The most immediate candidate for engineering MPTs is a highly controlled qubit system capable of programming a *hybrid* quantum circuit [1–3]. These circuits are composed of two ingredients: entangling unitary dynamics, and disentangling nonunitary measurements. In most current experimental quantum devices, significant effort is being expended to produce high-quality two-qubit gates suitable for producing entanglement. At sufficiently large circuit depths, these can induce a *volume-law* entanglement scaling in the underlying quantum state. Commonly studied circuits to generate volume-law wave functions are built out of two-qubit unitaries drawn from the Clifford group or the random Haar measure [26–29].

The second ingredient typically used to induce MPTs in hybrid circuits are repeated local measurements applied at a constant rate p throughout the circuit. These measurements act to disentangle the wave function and for large p will reduce the entanglement to an *area law* scaling. It is the intermediate

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open the prospect that MPTs may be offering a glimpse at a deeper, as yet undiscovered set of critical points, and possibly a new class of universal phenomena. Further observations of MPTs in higher-dimensional systems [22] or as purification phase transitions [23] and their relation to quantum error correction [24,25] demonstrate the broad consequences such critical points can have in many fields of physics.

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regime of small but finite p where conditions are ripe for MPTs to exist.

The experimental realization of a quantum system harboring an MPT requires accurate programmable manipulation of individual qubits. Experimental platforms based on lasercooled trapped ions achieve such a high degree of control, where highly pure quantum states of tens of ions are routinely prepared [30,31]. In a trapped-ion computer, the entangling dynamics of a random quantum circuit can be achieved with a combination of native two-qubit entangling Mølmer-Sørensen (MS) gates and random single-qubit rotations. In addition, recent experiments employing programmable and high-precision single-ion optical addressing technologies [32,33] demonstrate the feasibility of a constant rate of measurements throughout the circuit.

In this article, we propose a class of hybrid circuit which is natural to implement using this set of native trapped-ion gates. We focus on non-Clifford circuits and study the behavior of the entanglement scaling using state-of-the-art tensor network algorithms, capable of simulating tens of qubits acted on by thousands of gates. We find a class of circuit that exhibits a phase transition in entanglement under a number of plausible experimental conditions. We investigate the critical point, which displays universal hallmarks of an MPT, and calculate the critical measurement density and the universal scaling exponent of the correlation length. We argue that this transition should be robust against sources of experimental noise expected in modern trapped-ion hardware, such as incoherent gate errors, state preparation and measurement (SPAM) errors, and optical cross-talk between the target and neighboring ions. We further demonstrate that bringing the qubit into an outcome-independent state after each measurement does not influence the MPT. This enables an efficient experimental protocol and mitigates potential leakage of ions into atomic states outside the qubit Hilbert space during projective measurements [34]. Beyond this robustness, near-term experiments will likely be limited by statistical requirements on postselection [7,12,21]. Due to the inherent random behavior of measurement outcomes, replicating the exact state after every run of the circuit becomes the main experimental challenge, especially for large system sizes and higher measurement rates. We conclude with a discussion of some recent proposals which may be able to bypass this bottleneck, such as detecting the MPT in a local order parameter via coupling a reference qubit to the system [5,7,12,35] or considering unitary evolutions in space-time duals [36].

II. HYBRID CIRCUIT DESIGN

We consider *N* trapped atomic ions, each with two internal energy levels behaving as an effective qubit system. In current experimental hardware, these two qubit states have been shown to be long-lived and highly coherent [37], with accurate initialization through optical pumping and qubit readout via fluorescence imaging. The qubits are manipulated coherently [30] using external laser fields, capable of applying phasesensitive single- and two-qubit quantum gates mediated by collective vibrational (phonon) modes.

We focus on modeling the dynamics generated by a hybrid circuit consisting of a mixture of entangling gates and disen-



FIG. 1. One time cycle realization of a quantum circuit setup to probe measurement-induced phase transitions, with vertical lines indicating the N = 6 evolved qubits. Squares denote single-qubit unitary rotation gates $R(\theta, \varphi)$, rectangles correspond to two-qubit Mølmer-Sørensen gates $M(\Theta)$, and circles show single-qubit projective measurements in the Z-basis, which are injected uniformly in space and time at a rate p. The time cycle consists of an even (upper, U_e) and an odd (lower, U_o) unitary layer, indicated in gray, each followed by a measurement layer. In the even unitary layer the outer two qubits are not affected by the Mølmer-Sørensen gates due to open boundary conditions.

tangling measurements, acting on a pure initial state $|\Psi_0\rangle = |0\rangle^{\otimes N}$. To begin we examine randomized quantum gates that are native to the hardware, for their ability to induce strong entanglement in the wave function. In trapped-ion hardware, the native entangling gate—the MS gate—is equivalent to a rotation around the *XX*-axis by an angle Θ , with unitary propagator acting on qubits *j* and *k* [38]:

$$M_{j,k}(\Theta) = \begin{bmatrix} \cos(\Theta) & 0 & 0 & -i\sin(\Theta) \\ 0 & \cos(\Theta) & -i\sin(\Theta) & 0 \\ 0 & -i\sin(\Theta) & \cos(\Theta) & 0 \\ -i\sin(\Theta) & 0 & 0 & \cos(\Theta) \end{bmatrix}.$$
(1)

There are multiple ways of introducing randomness in the circuit, such as the angles of each MS gate, the rotation axes, or the gate connectivity, for example. Here we study quantum circuits constructed in a brick-layer fashion (as illustrated in Fig. 1), with each layer consisting of a set of MS gates between neighboring qubits (alternating the bonds between layers), supplemented by a layer of random single-qubit rotations. Those circuits have a straight-forward implementation in trapped-ion experiments. For the MS gates, we fix the angle to $\Theta = \pi/4$ uniformly throughout the circuit, generating maximal entanglement between the pairs at each application. Note that we restrict connectivity to nearest neighbors only, despite the capability of trapped-ion hardware to execute longrange gates. As will become clear in the next section, this choice allows us to leverage powerful tensor-network algorithms for simulating the circuit dynamics.

The native random single-qubit gates for trapped-ion hardware are rotations around arbitrary axes in the *X*-*Y*-plane, with propagator acting on qubit j [38]:

$$R_{j}(\theta,\varphi) = \begin{bmatrix} \cos(\theta/2) & -ie^{-i\varphi}\sin(\theta/2) \\ -ie^{i\varphi}\sin(\theta/2) & \cos(\theta/2) \end{bmatrix}.$$
 (2)

As demonstrated in [39] for superconducting qubits, this random ensemble of rotations can be configured to rapidly generate entanglement when combined with deterministic entangling gates. Inspired by this work, we fix the angle θ in the single-qubit rotations to $\theta = \pi/2$ and randomly choose one of three angles $\varphi \in \{0, \pi/4, \pi/2\}$. These gates describe $\pi/2$ -rotations around the X-axis ($\varphi = 0$), around the Y-axis $(\varphi = \pi/2)$, and around the axis between X and Y $(\varphi = \pi/4)$. We note that, though the MS gate and the rotations around the X- and Y-axis are Clifford gates, the single-qubit rotation with $(\theta = \pi/2, \varphi = \pi/4)$ defines a gate beyond the Clifford group. This choice makes it so that our circuit can not be simulated efficiently (i.e., in polynomial time) with a classical computer. However, the local configuration of gates is amenable to tensor network simulations, which will form the core of our numerical strategy below.

The combined application of single- and two-qubit gates with fixed angle choices can be summarized in a unitary operator:

$$U_i(\varphi_i, \varphi_{i+1}) = R_i(\pi/2, \varphi_i)R_{i+1}(\pi/2, \varphi_{i+1})M_{i,i+1}(\pi/4).$$
 (3)

We thus build the brick-layer construction in the quantum circuit by alternately applying operators U° and U^{e} , with

$$U^{o} = \prod_{i \text{ odd}} U_{i}(\varphi_{i}, \varphi_{i+1}), \qquad (4)$$

$$U^{e} = \prod_{i \text{ even}} U_{i}(\varphi_{i}, \varphi_{i+1}), \qquad (5)$$

and refer to odd and even unitary layers, respectively. The random angles φ_i , φ_{i+1} are chosen independently for each unitary operator $U_i(\varphi_i, \varphi_{i+1})$.

By repeated application of U° and U^{e} , we create a circuit with D layers of gates, each evolving a pure quantum state $|\Psi\rangle$ as

$$|\Psi\rangle \to U^k |\Psi\rangle, \ k \in \{e, o\}.$$
 (6)

We further define a time cycle as the combined application of odd and even unitary layers, so that a periodic behavior is induced in time and the circuit consists of T = D/2 time cycles [11], where one such cycle is depicted in Fig. 1 for a system with N = 6 qubits.

For large times t, the pure quantum state evolving under the dynamics in Eq. (6) reaches a constant entanglement entropy, which shows a volume-law behavior as we will demonstrate numerically in the next section. In order to induce a phase transition into area-law entanglement at long times, we inject measurements uniformly through the circuit (i.e., after each single MS layer) at a fixed rate p (see Fig. 1). Specifically, at any space-time location, we perform a local projective measurement $P_i^{\pm} = (1_i \pm \sigma_i^z)/2$ into the computational basis with a probability p, associated to the σ^z Pauli matrix. On average, pND measurements are performed in the circuit, where each measurement projects the wave function according to the measurement outcome $\sigma_i^z = \pm 1$:

$$|\Psi\rangle \to \frac{P_i^{\pm}|\Psi\rangle}{\|P_i^{\pm}|\Psi\rangle\|}.$$
(7)

Independent of the results, these projections disentangle qubit i from the rest of the system, and stand in contrast to the entangling unitary gates.

III. TENSOR NETWORK IMPLEMENTATION

The hybrid circuit that we propose in the last section is not amenable to any asymptotically efficient classical simulation scheme [i.e., scaling poly(N) in time]. This is in contrast to Clifford circuits, for which efficient simulation is ensured through the Gottesman-Knill theorem [40-42], or quantum automaton circuits, which are tractable despite their high entanglement [43,44]. In order to simulate the long-time dynamics of our model circuit, we employ tensor-network methods for approximate time evolution. Despite not being poly(N) scalable in general for the physics we explore, we will demonstrate below that these techniques allow us to reach circuit sizes sufficiently large to resolve the universal properties of the MPT critical point. Because of the design choice of circuit topology, we leverage efficient and well-refined algorithms based on matrix product states (MPSs) [45-47], a class of parametric wave functions that has revolutionized large-scale simulations on strongly correlated quantum matter in one (and quasi-one) dimension [48–51].

An MPS is a classical representation of a many-body wave function $|\Psi\rangle$, obtained by decomposing $|\Psi\rangle$ into a set of local tensors $A_{i}^{\sigma_{j}}$:

$$|\Psi\rangle = \sum_{\sigma} \prod_{j} A_{j}^{\sigma_{j}} |\sigma_{1}, \dots, \sigma_{N}\rangle .$$
 (8)

Each rank-3 tensor (rank-2 at boundaries) has a physical index σ_j and two auxiliary indices with dimension \mathcal{D}_j . The total number of parameters scales as $O(N\mathcal{D}^2)$, where $\mathcal{D} = \max{\{\mathcal{D}_j\}}$ is the *bond dimension* of the MPS. In the limiting case of $\mathcal{D} = 1$ one recovers mean-field theory, while larger bond dimensions allow the MPS to capture increasingly large amounts of quantum correlations. This, however, comes at a cost that may become exponential in the number of qubits.

We evolve the initial product state using a procedure analogous to the time-evolving block decimation algorithm [52,53]. We examine the MPS wave function after *T* time cycles for different measurement rates *p*. At each application of a gate, the contraction is approximated by only keeping a reduced number of states (i.e., the bond dimension) following singular value decomposition, or by discarding all singular values below a fixed accuracy threshold ε . Because of the orthogonality property of MPSs, this approximation remains controlled, in that the local truncation error is equal to the global error in approximating the wave function [50,54]. For p > 0, independent measurements are realized by sampling the measurement outcome from the local one-body reduced density matrix, and projecting the MPS accordingly.

The numerical experiments shown in the following sections are performed using state-of-the-art Julia implementations of tensor network algorithms for quantum circuit simulations. We specifically use PastaQ.jl [55], a toolkit of tensor-network algorithms for simulation and benchmarking of quantum hardware, based on the ITensors.jl package for efficient tensor manipulation [56]. We tested our tensor network



FIG. 2. (a) The maximum bond dimension \mathcal{D} observed in the long-time states of 500 random circuits as a function of error rate p for tensor network simulations with a truncation cutoff of $\varepsilon = 10^{-10}$. Different system sizes N are considered, where for N = 28 no data exist for $p \leq 0.09$ due to exceeding computation times for large bond dimensions. (b) Illustration of the entanglement phase diagram. A measurement-induced phase transition separates the two phases at the critical measurement rate p_{c} ; see Sec. IV.

simulations for small circuit sizes against exact time evolution using the Yao.jl package [57].

Before we turn to numerical calculations of physical properties in the next section, we first examine the simulation complexity associated with the tensor network method. A natural measure of complexity is the bond dimension discussed above, which determines the total number of parameters in an MPS wave function. Physically, the bond dimension indicates the amount of entanglement in the quantum state. For the separable state $|\Psi_0\rangle$ at time t = 0, the bond dimension required is $\mathcal{D} = 1$. During the random circuit dynamics, the bond dimension of the simulation grows to capture the spreading of correlations at long distances associated with increasing entanglement. Because MPS wave functions possess exponentially decaying correlations, \mathcal{D} may in fact grow very rapidly in time, hindering simulation in the low-*p* regime.

We examine the computational cost required to reach quantum states deep in the volume-law phase (i.e., low p). In Fig. 2 we show the maximum bond dimension reached in the long-time states of 500 random circuits for different rates p and increasing system size N, obtained with a fixed truncation cutoff of $\varepsilon = 10^{-10}$. For large measurement rates we find convergence to a constant value of \mathcal{D} , which does not depend on the system size. This matches our expectations for an area-law entangling phase for large p. In contrast, for small measurement rates we observe increasing bond dimensions with increasing N. This is consistent with volume-law entanglement, and becomes more pronounced as p decreases. From these first observations, we can make a preliminary conclusion that a critical measurement rate p_c may exist between $p \approx 0.17$ and $p \approx 0.2$, below which the dependence on the system size starts to form, as indicated in Fig. 2(b).

IV. SIMULATION RESULTS

In order to characterize the putative phase transition, we focus on the entanglement entropy, the most relevant physical indicator of an MPT. We focus on generalized α -Rényi entanglement entropies

$$S_{\alpha}(A) = \frac{1}{1-\alpha} \ln\left(\mathrm{Tr}\big[\rho_{A}^{\alpha}\big]\big),\tag{9}$$

evaluated on subsystem A after a bipartition of the state $|\Psi\rangle$ into subsystems A and \bar{A} , with reduced density matrix

$$\rho_A = \mathrm{Tr}_{\bar{A}} |\Psi\rangle \langle \Psi|. \tag{10}$$

In Eq. (9) the limit $\alpha = 1$ recovers the familiar von-Neumann entanglement entropy. The second Rényi entropy, $\alpha = 2$, is most amenable to experimental measurement protocols [58,59]. In the tensor network simulations, the full family of Rényi entropies is accessible from a Schmidt decomposition of the MPS at any given bond.

A. Entanglement scaling

In this section we detail our numerical results for the entanglement entropy of our hybrid circuit shown in Fig. 1, where we choose even numbers of qubits N and open boundary conditions in accordance with numerical and experimental constraints. Thus, the first and last qubits in the one-dimensional chain are not affected by MS gates in every second layer of unitaries (see Fig. 1). There are three sources of randomness that play a role in the dynamics: the choice of single-qubit rotations, the location of projective measurements, and the measurement outcomes [11]. We carry out independent simulations and average over these three realizations of disorder. Specifically, we study entanglement entropies of half-chain bipartitions, so that the two subsystems contain equal numbers of qubits, $|A| = N_A = N_{\bar{A}} = N/2$. For convenience we write $S_{\alpha}(N, p, t)$, which refers to the halfchain entanglement entropy for a system of N qubits after ttime cycles in a circuit with measurement rate p.

We start with considering the evolution of the half-chain entanglement entropy under time cycles of the circuit for different system sizes N and measurement rates p. Figure 3 shows $S_1(N, p, t)$ and $S_2(N, p, t)$ as functions of time cycles t, evaluated after each even unitary layer. We consider two measurement rates p = 0.1 and p = 0.2 for different system sizes N, averaging over 500 circuit runs. All simulations show a clear convergence to a fixed entanglement entropy for $t \gtrsim N$ time cycles, confirming a circuit depth of T = 2N to study the constant long-time entanglement behavior.

While for p = 0.2 the entanglement entropy saturates at similar values for all system sizes, a strong growth with increasing N is observed for p = 0.1. This dependence on the subsystem size indicates volume-law entanglement for small measurement rates, while its absence for larger measurement rates is a sign for convergence to area-law entanglement at late times. Figure 3 hence suggests a phase transition between volume-law and area-law entanglement at a critical measurement rate $0.1 \le p_c \le 0.2$.

As already observed in Fig. 2, the simulations in the volume-law phase require large bond dimensions and hence excessive computation times. This limits our numerical



FIG. 3. The evolution of the first and second half-chain entanglement Rényi entropies under the quantum circuit are plotted as functions of time cycles t for different system sizes N and measurement rates p = 0.1 (left) and p = 0.2 (right). Convergence to a constant value for large t indicates a saturated long-time entanglement entropy. The dependence on N for p = 0.1 for both measures indicates a volume-law entanglement, while the right column shows area-law entanglement since the entropies converge to similar values independent of the system size. All data points result from averaging 500 circuit runs, except for sizes N = 30 and N = 36 with measurement rate p = 0.1, where only 85 and 19 circuit runs are averaged, respectively, as the large entanglement entropy leads to excessive computation times. Shaded areas show the standard mean error, which is barely visible when averaging 500 runs. Entanglement entropies for subsystems with an odd number $N_A = N/2$ of spins are smaller than those for subsystems with even N_A due to open boundary conditions [60,61].

studies with p = 0.1 in Fig. 3 to smaller system sizes, so that for N = 30 and N = 36 only 85 and 19 circuit runs are averaged, respectively, explaining stronger fluctuations. Further, simulations are not feasible for N > 36.

As an upper bound to the second-order Rényi entropy [3], the von Neumann entropy takes larger values, while both entanglement measures show similar dependencies on system size and measurement rate. This confirms that both entanglement measures can be used to detect the entanglement phase transition, which is also expected to happen at similar measurement rates [2,10,11,14]. This further motivates an experimental realization of the quantum circuit to study MPTs.

B. Critical behavior

In order to study the transition between the volume-law and the area-law phase in more detail and determine the critical measurement rate, we consider the scaling behavior of the von Neumann entanglement entropy. The constant entanglement entropy after T = 2N time cycles is expected to show the finite-size scaling form [2,10,11,13]

$$S_1(N, p, T) = \alpha(p_c) \ln(N) + F[(p - p_c)N^{1/\nu}], \qquad (11)$$



FIG. 4. The scaling behavior of the half-chain von Neumann entanglement entropy [Eq. (13)] for different system sizes $N \in \{6, 8, \ldots, 36\}$, for circuits with projective measurements P_i^{\pm} . The critical measurement density p_c and scaling exponent ν are adjusted to produce the best collapse, resulting in $p_c = 0.17 \pm 0.02$ and $\nu = 1.4 \pm 0.2$. All data points result from averaging the entanglement entropy after T = 2N time cycles of 500 circuit runs, and error bars are smaller than data points. Inset: The scaling behavior for additional applications of $\tilde{P}_i = |0\rangle\langle 0| + |0\rangle\langle 1|$ after each measurement, with $N \in \{6, 8, \ldots, 36\}$. An optimal collapse gives $p_c = 0.17 \pm 0.02$, $\nu = 1.4 \pm 0.2$ in this case; see Sec. V A.

with a factor $\alpha(p_c)$, scaling function *F*, and critical exponent ν . At criticality, we furthermore expect [2,10,11]

$$S_1(N, p_c, T) = \alpha(p_c)\ln(N) + b, \qquad (12)$$

with a constant offset *b*, and hence F(0) = const. We can determine the critical measurement rate p_c and the critical exponent v by eliminating the logarithmic term of the scaling form,

$$S_1(N, p, T) - S_1(N, p_c, T) = \tilde{F}[(p - p_c)N^{1/\nu}], \quad (13)$$

and comparing the expected behavior with numerical calculations.

We perform tensor network simulations for different measurement rates p and system sizes N and average the entanglement entropy after T = 2N time cycles over 500 circuit runs with randomly chosen single-qubit rotations and measurement placements. Due to the large entanglement entropies in the volume-law phase, we consider smaller system sizes for small measurement rates.

Figure 4 shows the resulting collapse of the entanglement entropies to the scaling form in Eq. (13), where we find $p_c =$ 0.17 ± 0.02 and $v = 1.4 \pm 0.2$ by optimizing the collapse. The clearly observable universal scaling behavior at criticality confirms the expectation of an entanglement phase transition. The determined critical exponent agrees with observations for different circuits in related works [4,5,10–12,20,23]. Apart from the scaling collapse, we analyze the scaling form at criticality in Eq. (12). We estimate the critical measurement rate p_c by fitting the scaling form to the numerical data at different measurement rates p and considering the fit quality, which is maximized at p_c . The resulting data are shown in the inset of Fig. 5, where we quantify the fit quality via the



FIG. 5. Scaling function of the entanglement entropy close to the critical measurement rate p_c . $S_1(N, 0.17, t)$ is plotted as function of ln(t) for fixed system sizes N = 34 (triangles) and N = 36(circles), and as a function of $\ln(x)$, $x = 2N/\pi$, for system sizes $N \in \{6, 8, \dots, 36\}$ after t = 80 time cycles, distinguishing even (circles) and odd (triangles) system sizes due to open boundary conditions [60,61]. All data points are averaged over 500 circuit runs, and error bars are smaller than the data points. Black lines show functions $g(N) = \alpha(0.17)\ln(x) + \tilde{b}$ fitted to the data points, giving fit parameters $\alpha(0.17) = 0.26 \pm 0.01$, $\tilde{b} = 0.33 \pm 0.03$ for even and $\alpha(0.17) = 0.28 \pm 0.01$, $\tilde{b} = 0.1 \pm 0.03$ for odd subsystem sizes. Shaded regions show fit uncertainties. Inset: Mean-squared error, Eq. (14), of fitted scaling forms according to Eq. (12) vs measurement rate p for odd and even subsystem sizes. A clear minimum around p = 0.15 to p = 0.17 indicates the critical measurement rate. Same data as in Fig. 4.

mean-squared error,

$$\mathcal{R}(p) = \langle [S_1(N, p, T) - f_p(N)]^2 \rangle_N, \tag{14}$$

with fitted function $f_p(N) = \alpha(p_c)\ln[N] + b$. The average is taken over the data for different system sizes N, and we distinguish even and odd subsystem sizes, since odd subsystem sizes show smaller entanglement entropies due to the open boundary conditions [60,61]. A clear minimum is observed around p = 0.15 to p = 0.17, which agrees well with the previously observed critical measurement rate and confirms the MPT in this regime.

The main plot in Fig. 5 shows that a similar logarithmic scaling form is observed in the dynamical behavior of the entanglement entropy in the critical regime for fixed system sizes. We plot $S_1(N, 0.17, t)$ as a function of $\ln(t)$ for fixed system sizes N = 34, N = 36 to represent odd and even subsystem sizes. Additionally, we plot the entanglement entropy as a function of $\ln(2N/\pi)$ for different subsystem sizes $N_A = N/2$ after t = 80 time cycles, where we include an offset in the scaling form which is expected from conformal field theories (CFTs) [13,43,61].

We observe that both scaling forms, in time and in system size, behave similarly when independently considering odd and even subsystem sizes. We further fit functions $g(N) = \alpha(p_c)\ln(2N/\pi) + \tilde{b}$ and $g(t) = \alpha_t(p_c)\ln(t) + \tilde{b}_t$ to the corresponding data, where we take the uncertainty of $p_c = 0.17 \pm 0.02$ into account and consider measurement rates in this regime. We observe similar coefficients $\alpha(p_c) = 0.3 \pm 0.15$ and $\alpha_t(p_c) = 0.3 \pm 0.15$ for both even and odd subsystem sizes. The ratio of the coefficients, $z = \alpha_t(p_c)/\alpha(p_c) = 1.04 \pm 0.04$ describes a dynamical critical exponent whose value indicates space-time conformal invariance [10,12,13,43,61]. Thus, our simulations demonstrate the emergence of a CFT at the studied MPT.

V. FURTHER EXPERIMENTAL CONSIDERATIONS

In the above, we have provided numerical evidence for an MPT in a model of a hybrid circuit using native trapped-ion gates. Despite the experimental feasibility of the MS and single-qubit gate set [30], as well as recent demonstrations of high-precision optical addressing that enables measurements to be performed throughout the circuit [33], several considerations remain before such an MPT might be observed in real trapped-ion hardware. In this section, we address what we believe are the major experimental challenges that remain, with a particular focus on a possible implementation in realistic trapped-ion experiments.

A. Measurement corrections from states outside the Hilbert space

Projective measurements of trapped-ion qubit states often populate states that are outside of the qubit Hilbert space. Such processes do not necessarily reduce the fidelity of state measurements, as the leakage onto unwanted states occurs only for one of the two qubit states. For example, during the spin-dependent fluorescent measurement in ¹⁷¹Yb⁺ ions, the state $|F = 1, m_F = 0\rangle$, representing the qubit state $|1\rangle$, leaks to $|F = 1, m_F = \pm 1\rangle$ states, while the state $|F = 0, m_F = 0\rangle$, representing the qubit state $|0\rangle$, is preserved [34]. Reinitializing the ion in the projected qubit state may be experimentally demanding, requiring real-time decision logic or transferring population via other metastable atomic states [62].

A potential experimental implementation can hence be simplified by applying an additional operator $\tilde{P}_i = |0\rangle\langle 0| + |0\rangle\langle 1|$ after each projective measurement operator P_i^{\pm} . Such an operation is routinely implemented with high fidelity in trapped-ion experiments, using optical pumping techniques. Any measured qubit then continues to evolve from the $|0\rangle$ state, which avoids the outcome-dependent operation of bringing the ion onto the $|1\rangle$ state. In the numerical simulations the wave function transforms according to

$$|\Psi\rangle \to \frac{\tilde{P}_i P_i^{\pm} |\Psi\rangle}{\|\tilde{P}_i P_i^{\pm} |\Psi\rangle\|},\tag{15}$$

which is implemented via the same probabilistic process as discussed earlier, followed by the application of \tilde{P}_i .

A comparison of numerical simulations using these additional operations after measurements with the previously discussed results (see Sec. IV B) demonstrates that the MPT is not strongly influenced by this replacement. The inset in Fig. 4 shows the resulting scaling collapse according to Eq. (13), whose optimization gives $p_c = 0.17 \pm 0.02$ and v = 1.4 ± 0.2 , which agrees with our observations for projective measurements. This result demonstrates a flexibility in adapting the measurement procedure to experimental restrictions and provides an important step towards efficient experimental realizations of MPTs.

B. Measurement crosstalk and SPAM errors

The realization of constant measurement rates in trappedion hybrid circuits can be possible using recent advances in single-ion optical addressing [32,33]. However, such an implementation comes with a source of error that directly affects the critical measurement rate. Due to the micron-scale spatial separation of the trapped ions, the addressing of a single ion can leak onto direct neighbors (crosstalk error), resulting in unwanted measurement of the neighboring qubit. This process increases the probability with which each ion is measured and thus effectively increases the measurement rate, shifting the critical measurement rate p_c to smaller values. Specifically, if a single ion is measured with probability p at a measurement layer in the circuit, and for each measurement the crosstalk error causes the additional measurement of an adjacent ion with probability p_d , the effective measurement rate can be considered as

$$(1 + p_{\rm d} - pp_{\rm d})p.$$
 (16)

Here the last term takes into account that an additional measurement caused by crosstalk errors has no effect if the ion is already measured. We further assume that left and right neighbors are addressed with equal probabilities and that the crosstalk probability is small, so that terms of higher order in p_d , resulting from crosstalk errors on both neighboring ions, are neglected.

In state-of-the-art trapped-ion addressing schemes, the relative intensity cross-talk error can be suppressed below a level of 10^{-4} at the neighboring ion [32,33]. The state measurement time (spin-dependent fluorescence) itself can be limited to 10 μ s using high resolution imaging systems [63]. Thus, the probability of an accidental measurement of the neighboring qubits (p_d) can be kept at a few percent level.

For example, if the state detection laser beam is at the saturation intensity of the ${}^{2}S_{1/2} - {}^{2}P_{1/2}$ transition on the target 171 Yb⁺ qubit, the probability of accidental measurement of the neighboring qubit via spontaneous emission from 10^{-4} level cross-talk can be kept at $p_{d} \approx 10^{-4} \times (2\pi \times 20 \text{ MHz/6}) \times 10 \ \mu \text{s} \approx 0.02$ [63,64]. Here $2\pi \times 20 \text{ MHz}$ is the natural linewidth of the ${}^{2}P_{1/2}$ state used in the measurement process. Accidental measurement of neighboring qubits can also be mitigated by using separate atomic species (or separate isotopes) [65] as processing qubits and read-out qubits, where the qubit state is transferred from a processing qubit to a read-out qubit before measurement.

However, a fundamental limit to the probability of accidental measurements is given by the absorption of photons emitted from an addressed ion by neighboring ions. The probability of this effect depends strongly on the spatial distance *d* between two ions and scales as $\sim 1/d^2$. In a typical experiment of ¹⁷¹Yb⁺ ions, a spatial distancing of $d \approx 5 \,\mu\text{m}$ can be achieved, so that this error is limited to $\lesssim 0.5\%$; see the Appendix for details. The error can be further reduced by orders of magnitude using e.g. quantum charge-coupled devices (QCCDs) [66].

With the estimate in Eq. (16) we expect a shift of p_c which is much smaller than the uncertainty in its value resulting from our computational methods. We confirmed via numerical studies of circuits including additional measurements with a relatively large leakage rate of $p_d = 0.1$ that the shift in p_c is not detectable within numerical uncertainties, emphasizing the robustness of the model to experimental errors.

Apart from the crosstalk errors in the single-qubit measurements, further imperfections appear when preparing the initial state, applying MS and single-qubit gates, or performing global measurements of the final state. Such SPAM errors are on the order of a few percent in existing trapped-ion hardware [30] and can also influence the critical measurement rate. However, since these errors are of the same order as measurement crosstalk errors, we expect the effects on the MPT to be similarly small.

Even though a noticeable shift of the critical measurement rate can result from the combination of multiple small imperfections, we expect the resulting shift of p_c to be small enough so that the MPT remains at finite measurement rates, $p_c > 0$. Thus, the MPT is robust against experimental imperfections and motivates an experimental realization.

C. Estimators and postselection

Beyond the above experimental considerations, an important difference between our tensor network simulations and real-world experiments is the procedure by which the longtime quantum wave function is characterized. The evaluation of observables in trapped-ion hardware is based on a statistical average of measurement outcomes, typically performed on the final quantum state. Most relevant for characterizing an MPT, the second Rényi entropy can be extracted from statistical correlations between local randomized measurement outcomes [59,67]. It can hence be evaluated from final-state measurement averages, much like any conventional observable. However, the probabilistic behavior of random circuit (monitored) measurements causes further limitations. This comes from the usual requirement of repeated state preparation and final-state measurements (possibly in multiple bases) in order to reconstruct any observable. In the case of hybrid circuits created by monitored measurements spread throughout the circuit, this necessity of repeated state preparation and measurement adds an additional requirement on postselection, which naively results in an exponential scaling overhead. In order to prepare a specific quantum state multiple times, the outcomes of all monitored single-qubit measurements need to be identical in the considered circuit runs. The probabilistic projections in the circuit for p > 0 thus require the postselection of the final quantum states of multiple circuit runs based on the binary measurement outcomes. For a hybrid quantum circuit of depth T acting on N ions, on average $\sim 2^{2pNT}$

circuit runs are needed in order to prepare a specific quantum state, causing the exponential overhead in circuit runs [7,12,21].

While microsecond-scale gate and measurement times of the trapped-ion hardware will allow experiments with $N \approx$ 10 qubits with measurement rates up to $p \approx 0.05$ within reasonable run times, the exponential postselection overhead will pose a challenge for larger systems and higher measurement rates. The achievable measurement rates are hence much smaller than the numerically observed critical measurement rate where the MPT happens in the studied circuit. Reducing the critical measurement rate p_c based on alternative circuit setups built of native ion gates might be a promising approach to reveal MPTs in a larger system.

Accidental measurements appearing in experimental circuit runs due to crosstalk errors as discussed in Sec. VB cause further issues in the postselection procedure. Since the outcomes of these stochastically appearing measurements are not monitored in the experiment, two similarly prepared final states can differ even when postselecting the monitored measurement outcomes. This strengthens the necessity of minimizing crosstalk effects in trapped-ion hardware, which can be achieved, e.g., by incorporating coupling of ions as in QCCD architectures [66]. We numerically estimate the average loss of fidelity due to stochastic crosstalk measurements compared to the ideally prepared state without leakage errors to be $\leq 10\%$ for relevant system sizes and the cross-talk rates discussed in Sec. VB. Understanding the exact effect of this fidelity loss on the sample-based evaluation of entanglement entropies is an interesting topic for future studies.

While this postselection procedure presents a clear bottleneck for experimental realizations of MPTs, recent works suggest ways to bypass this measurement overhead. Such proposals use additional ancilla qubits which are initially maximally entangled with the circuit qubits as a reference to locally probe the MPT [5,12,35]. A local order parameter can be associated with the average entropy of the reference qubits, which quantifies the amount of information shared with the initial state and is hence large in the volume-law phase and small in the area-law phase, undergoing the same transition as the entanglement entropy. The introduction of such local order parameters allows for the detection of the MPT at the critical measurement rate while reducing the necessary amount of global measurements and avoiding the exponential postselection overhead. Furthermore, the number of additional qubits can be kept as small as a single reference qubit, making the approach feasible for experimental realizations [12,35]. Apart from improvements of this reference qubit approach based on cross-entropy estimators to reduce the amount of experimental measurements [7], other proposals to circumvent the postselection process are based on mappings onto spacetime duals, which describe unitary dynamics and allow for an efficient measurement of Rényi entanglement entropies [36]. Such theoretically developed techniques propose experimentally feasible ways to overcome the postselection bottleneck by only slightly increasing the number of necessary qubits and may enable realizations of MPTs in trapped-ion hardware.

VI. DISCUSSION AND OUTLOOK

In this paper we have introduced a hybrid circuit model consisting of native trapped-ion gates that is a strong candidate for realizing a measurement-induced phase transition (MPT). In order to induce entanglement growth, the circuit uses a combination of two-qubit Mølmer-Sørensen gates and random single-qubit rotations. To compete with this entanglement growth, measurements are induced throughout the circuit at a constant rate p. Such measurements should be realizable via recent experimental advances in high-precision, single-ion optical addressing schemes [32,33].

The native gates studied in this work go beyond the Clifford group and can hence not be simulated efficiently with a scalable classical algorithm. In order to study this hybrid circuit, we therefore employ state-of-the-art tensor network simulations, evolving a matrix product state wave function using a time-evolving block decimation algorithm. These simulations demonstrate two clear entanglement regimes in the long-time behavior of the circuit, corresponding to volume-law entanglement for small *p* and area-law entanglement for large *p*. Further, our numerical results provide clear evidence for an MPT at a critical measurement density $p_c \approx 0.17$.

In the area-law regime, our tensor network simulations are able to converge typical hybrid circuits with up to 44 qubits and more than 10 000 unitary gates. Such sizes are sufficiently large to resolve the critical properties of the MPT in great detail. Collapsing our data near p_c to a critical scaling function reveals the universal exponent $v = 1.4 \pm 0.2$, which is within numerical uncertainty of the value of v = 4/3 given by firstpassage percolation in two dimensions [2,4,5,10,11,13,14,20]. In addition, the entanglement entropy shows clear logarithmic scaling at p_c , which agrees with previous studies of MPTs that show broad features of conventional conformal field theories [2,4,10–14,61]. The coefficients of the logarithmic fits in the space and time direction directly at p_c reveal coefficients that are the same to within numerical errors, indicating the apparent emergence of conformal symmetry.

Our numerical results suggest that modern trapped-ion experiments with a reasonable number of qubits are wellequipped to realize such MPTs. We have argued that even in the presence of realistic experimental errors, such as measurement crosstalk errors, the MPT is sufficiently robust to be induced with native entangling gates. We have further shown that the qubits can be brought into outcome-independent states after the projective measurements without affecting the MPT, which suggests a decision-logic free experimental implementation. However, we also find that without significant modification of the gate set architecture, experimental studies may be limited to measurement rates far in the volume-law phase. This is due to a postselection process required for evaluating physical observables, including the second Rényi entropy. A significant body of theoretical work is currently motivated by finding solutions to overcome this limitation [5,7,12,21,35,36].

Despite these challenges, our results indicate that trappedion quantum computers are compelling platforms to realize MPTs and could therefore conceivably be among the first engineered quantum systems to realize this exciting new

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APPENDIX: FUNDAMENTAL LIMITATIONS OF CROSSTALK ERRORS

The probability of crosstalk errors induced by absorption of photons emitted from an addressed ion can be kept at a small level. This probability depends strongly on the spatial distance between two ions and can be estimated via the photon absorption cross section σ of the neighboring ion. For the considered detection line of the ${}^{2}S_{1/2} - {}^{2}P_{1/2}$ transition, the cross section scales as $\sigma \approx \lambda^2/(2\pi)$ with the detection wavelength $\lambda = 369.5$ nm for ¹⁷¹Yb⁺. This corresponds to a cross section radius of $a = \sqrt{\sigma/\pi} \approx 80$ nm. Thus, a fraction proportional to $(1 - \cos(\theta))/(8\pi) \approx 6.4 \times 10^{-5}$ of the photons emitted by the targeted ion is absorbed, with the angle $\theta = \arctan(a/d)$ subtended by the photon absorption cross section area. For an optimal scattering rate of 7×10^6 Hz for the target Yb⁺ ion and detection times of 10 μ s, this gives a crosstalk error probability of $p_d \approx 5 \times 10^{-3}$. This probability can be increased when applying the corrective measurement scheme discussed in Sec. VA, since the additional pumping of an ion emits further photons.

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