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Measuring the Rényi entropy of a two-site Fermi-Hubbard model on a trapped ion quantum computer

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The efficient simulation of correlated quantum systems is a promising near-term application of quantum computers. Here, we present a measurement of the second Rényi entropy of the ground state of the two-site Fermi-Hubbard model on a 5-qubit programmable quantum computer based on trapped ions. Our work illustrates the extraction of a non-linear characteristic of a quantum state using a controlled-swap gate acting on two copies of the state. This scalable measurement of entanglement on a universal quantum computer will, with more qubits, provide insights into many-body quantum systems that are impossible to simulate on classical computers.

One of the striking differences between classical and quantum systems is the phenomenon of entanglement. Analyzing large entangled states is of considerable interest for quantum computing applications. This is particularly relevant to quantum chemistry, quantum electrodynamics, and materials science simulations involving interacting fermions [1, 2], small versions of which have been simulated on few-qubit quantum computers [3–6]. Recently, a quantum algorithm was developed to construct the entanglement spectrum of an arbitrary wave function prepared on a quantum computer via measurement of the Rényi entropies [7]. In this Letter we measure the second Rényi entropy in a 5-qubit circuit by implementing a controlled-swap (C-Swap) gate, and mitigate experimental errors by exploiting the symmetry properties of this gate. We note that previous measurements of the Rényi entropy such as [8] were not implemented on universal machines and may not be easily generalizable to arbitrary Hamiltonians or scalable to larger systems.

For a many-body quantum system ideally described by the state $|\Psi\rangle$ and composed of two subsystems A and B , the n th Rényi entropy is given by $S_n = \frac{1}{1-n} \log(R_n)$, where

$$R_n = \text{Tr}(\hat{\rho}_A^n) \quad (1)$$

is the trace of the n th power of the reduced density matrix $\hat{\rho}_A = \text{Tr}_B(|\Psi\rangle\langle\Psi|)$. For non-zero entanglement we have $R_2 < 1$, which has the same universality properties as the von Neumann entropy $S = -\text{Tr}(\hat{\rho}_A \log(\hat{\rho}_A))$. Both are measures of the entanglement between A and B , and provide valuable information about the underlying physics of the system. For example, the Rényi entropy can be used to study out-of-equilibrium scaling of entanglement [9–14], which includes distinguishing many-body localized states from thermalized states [15–

19] through their time dependence and dimensional scaling law [20], and to study topological order [21, 22] and quantum critical systems [23–25].

The system under investigation for this work is the two-site Fermi-Hubbard model, which describes interacting electrons on a lattice [26, 27]. Despite its simplicity, it has been postulated as a model for complex phenomena such as high-temperature superconductivity. Since its behavior in the thermodynamic limit remains inaccessible to classical numerical techniques, it has become a prime candidate for simulation by quantum computers [28, 29].

Our work consists of several co-designed theoretical and experimental steps. First, we find an efficient mapping from the electronic problem to the qubit space. Second, we develop a circuit for digitized adiabatic evolution to prepare the ground state of the model, parametrized by the Trotter step size and the total evolution time. Based on available experimental resources and estimated Trotter errors, we choose a set of parameters that best corresponds to the result from exact diagonalization. Third, we realize the C-Swap gate, which is the key to efficiently extracting the second Rényi entropy. Finally, we integrate all of these elements into one circuit (Fig. 1) that prepares and evolves two copies of the Fermi-Hubbard system, and measures the second Rényi entropy with the help of an ancilla qubit. Importantly, in this step, we use the symmetry properties of the C-Swap gate to post-process the information contained in the four data qubits and reduce experimental errors. We note that other methods to measure entanglement entropy have been proposed using either multiple copies of the system [30–33], or random quenches [34].

We execute the relevant quantum circuits on a trapped-ion system, that constitutes a programmable five-qubit quantum computer with full connectivity [35, 36]. Qubits are realized in the hyperfine-split $^2\text{S}_{1/2}$ ground level of $^{171}\text{Yb}^+$ ions confined in a Paul trap. Single- and two-qubit gate fidelities are typically

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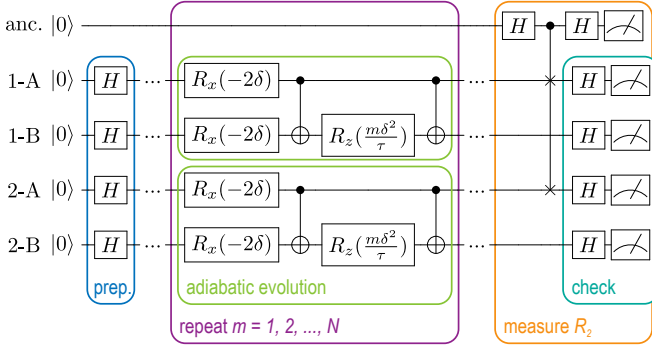


FIG. 1. The quantum circuit for the adiabatic evolution of two copies (labeled 1 and 2) of the two-site Fermi Hubbard model (each realized in two qubits A and B), and the measurement of the second Rényi entropy. The non-interacting ground state is prepared via the application of Hadamard gates, followed by the digitized adiabatic evolution to a finite value of U by repeated application of the central code block. m is an integer referring to the m -th step of the adiabatic evolution; δ and τ are given in the text. The angles in the rotation gates are in radians. The Rényi entropy R_2 is measured by applying a C-Swap gate controlled by an ancilla qubit (anc.) on subsystems A and subsequent detection of the state of the ancilla. By applying additional Hadamard gates and measuring the system qubits, experimental errors can be detected and discarded.

99.1(5)% and 98.5(5)%, respectively. Typical gate times are $10 \mu\text{s}$ for single- and $210 \mu\text{s}$ for two-qubit gates. The computational gates such as H, CNOT, and C-Swap are generated in a modular fashion by a compiler which breaks them down into constituent physical-level single- and two-qubit gates from the library (see Appendix A).

The Fermi-Hubbard Hamiltonian is

$$\hat{H} = -J \sum_{\langle i,j \rangle, \sigma} (\hat{c}_{i,\sigma}^\dagger \hat{c}_{j,\sigma} + \hat{c}_{j,\sigma}^\dagger \hat{c}_{i,\sigma}) + U \sum_{i=1}^N \hat{n}_{i\uparrow} \hat{n}_{i\downarrow} \quad (2)$$

where $\hat{c}_{i,\sigma}^\dagger$ and $\hat{c}_{i,\sigma}$ are the electron creation and annihilation operators, respectively, for spin state $\sigma \in \{\downarrow, \uparrow\}$ and site i , and $\hat{n}_{i,\sigma} = \hat{c}_{i,\sigma}^\dagger \hat{c}_{i,\sigma}$ is the electron number operator. Here J is the hopping strength and U is the on-site interaction. We consider the smallest non-trivial version of the model, which involves two sites and two electrons with total spin $S_z = 0$ along the z -axis.

Typically, when mapping electronic problems to qubits, the Jordan-Wigner [37] or Bravyi-Kitaev [2] mappings are used, both of which work in the second quantized basis. Here, the number of qubits N_q is equal to the number of single-electron states N_s , which would require $N_q = N_s = 4$ qubits.

However, in many problems, there are additional symmetries or conservation laws that can reduce the size of the Hilbert space. The Hamiltonian above conserves both the number of electrons N_e and the total spin along z , S_z . Therefore, a first quantization mapping, in which the Hilbert space sizes of the qubit system and the many-

electron problem are equal, uses qubits most efficiently. This is an important optimization for near-term quantum hardware.

In first quantization, the Hilbert space size is 4, which can be mapped to two qubits as $|00\rangle = \{1\uparrow 1\downarrow\}$, $|01\rangle = \{1\uparrow 2\downarrow\}$, $|10\rangle = \{2\uparrow 1\downarrow\}$, and $|11\rangle = \{2\uparrow 2\downarrow\}$. $\{i_{\sigma_i} j_{\sigma_j}\}$ represent the Slater determinants, which satisfy the number and spin conservation laws of the Hamiltonian. In this mapping, one qubit represents the up spin space and the other the down spin space. In this basis, the Hamiltonian is

$$\hat{H} = \begin{bmatrix} U & -J & -J & 0 \\ -J & 0 & 0 & -J \\ -J & 0 & 0 & -J \\ 0 & -J & -J & U \end{bmatrix}. \quad (3)$$

This can be written in terms of operators acting on the two qubits (up to a constant, and scaling energy by J) [38, 39]

$$\hat{H} = -(\hat{X}_1 + \hat{X}_2) + \frac{U}{2} \hat{Z}_1 \hat{Z}_2, \quad (4)$$

where \hat{X}_i and \hat{Z}_i are Pauli matrices.

To prepare the ground state of \hat{H} at finite U , we use digitized adiabatic evolution from the zero-interaction ground state. The time-dependent Fermi-Hubbard Hamiltonian $\hat{H}(t)$ is

$$\hat{H}(t) = -(\hat{X}_1 + \hat{X}_2) + \frac{t}{2\tau} \hat{Z}_1 \hat{Z}_2 \quad (5)$$

from $t = 0$ to $t = U\tau$ by linear interpolation. At $U = 0$, the ground state is $(|0\rangle + |1\rangle) \otimes (|0\rangle + |1\rangle)$, which can be prepared with Hadamard gates (see Fig. 1).

The first order Trotterization for a Hamiltonian with two non-commuting terms $\hat{H}_{a,b}$ is $\exp(-i(\hat{H}_a + \hat{H}_b)\delta) = \exp(-i\hat{H}_a\delta)\exp(-i\hat{H}_b\delta) + \hat{\mathcal{O}}(\delta^2)$. For larger evolution time τ and smaller step size δ , the approximation of the ground state is more accurate (see Appendix B). Here, at time $t = m\delta$, where m is an integer, we set $\hat{H}_a = -(\hat{X}_1 + \hat{X}_2)$, $\hat{H}_b = \frac{m\delta}{2\tau} \hat{Z}_1 \hat{Z}_2$. Putting together Trotterization and digitized adiabatic evolution, we obtain the following sequence of unitary operations to prepare the ground state:

$$\hat{V} = \prod_{m=1}^M \left[\exp(i\delta \hat{X}_1) \exp(i\delta \hat{X}_2) \exp\left(-im \frac{\delta^2}{2\tau} \hat{Z}_1 \hat{Z}_2\right) \right]. \quad (6)$$

Following [40], we use the convention that $\hat{R}_x(\theta) = \exp(-i\theta \hat{X}/2)$, and similarly for \hat{Z} . This leads us to the circuit for the digitized adiabatic evolution as shown in figure 1.

We simulate the evolution on a classical computer to investigate the scaling of the error in R_2 in the first order Trotter approximation (see Appendix B).

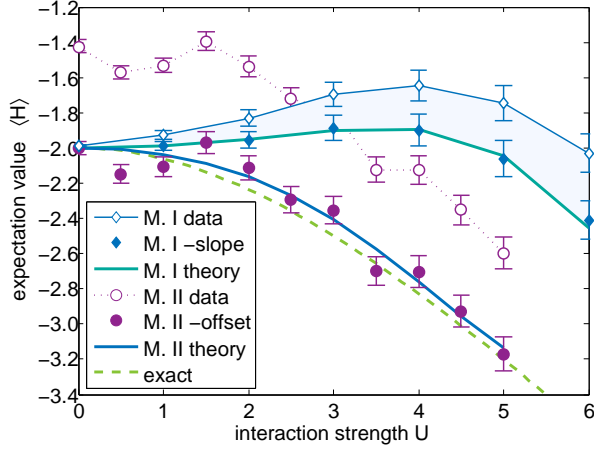


FIG. 2. Expectation value of the Hamiltonian $\langle \hat{H} \rangle$ of the Fermi-Hubbard model after adiabatic evolution, to the values of U given by the abscissa. Method I (M. I, open diamonds) sees an increase in experimental error with U (shaded area) since N_{steps} increases with U . Subtracting a straight line of slope 0.063 best matches the data points to the theory curve (filled diamonds). For method II (M. II, open circles), $N_{\text{steps}}=5$ is constant and the experimental error in $\langle \hat{H} \rangle$ is seen as an offset. Even for large systems, this can be measured for the integrable case $U = 0$ and then subtracted (filled circles). The corrected results (measured offset 0.58) follow the theory curve well. Method II is a better match to the exact solution.

For the experimental implementation, the memory depth of the arbitrary waveform generator used to control the laser pulses limits the length of a gate sequence to 8ms, bounding the number of Trotter steps to $N_{\text{steps}} \leq 6$. We implement two different methods for evolving to finite U .

Method I: For a fixed τ and δ , one can sample at intervals of τ/δ and $N_{\text{steps}} = U\tau/\delta$. In order to evolve from $U = 0$ to $U = 6$, we choose $\tau/\delta = 1$ and go from $N_{\text{steps}} = 0$ to $N_{\text{steps}} = 6$. As a result, only finitely-spaced values of U can be chosen. Here, smaller values of U involve fewer gates and will be less affected by experimental errors.

Method II: We fix $N_{\text{steps}} = 5$, and can hence sample any value of U up to N_{steps} . In this case, $\tau = N_{\text{steps}}\delta/U$. Here, the same number of gates are performed at every value of U and so the magnitude of the experimental error should be similar at every point.

Based on simulations (see Appendix C), we choose the parameters that seem to most closely follow the results from exact diagonalization. For method I, $\delta = 0.1$ and $\tau = 0.1$, while for method II, $\delta = 0.25$ and $\tau = 1.25/U$.

We first implement the digitized adiabatic evolution by itself for both methods in a two-qubit experiment on our system. We prepare the qubits along the x -axis of the Bloch sphere, in the ground state of the non-interacting Hamiltonian ($U = 0$) using Hadamard gates (see Fig. 1). Then we evolve the state in steps from $U = 0$ to $U = 6$ for method I and $U = 5$ for method II. Finally we measure along the z -axis and separately (with the help

of additional Hadamard gates) along the x -axis to calculate the expectation value $\langle \hat{H} \rangle$ from $\langle \hat{X}_1 \rangle$, $\langle \hat{X}_2 \rangle$, and $\langle \hat{Z}_1 \hat{Z}_2 \rangle$. The results are shown in figure 2. We see that method II shows an offset from the ideal value. If we subtract the value measured for $U = 0$ (no evolution) from all data points, they match the theoretical expectation closely. For method I, the number of gates and hence the error incurred grows with U . Subtracting a straight line of slope 0.063 models this increase well. This is consistent with the small-circuit error model developed in [41]. In a larger system, where the dynamics are unknown, this correction cannot be easily determined. The offset seen in method II, however, corresponds to the error in the eigenvalue of a non-interacting and hence easily integrable system, which can be more generally applied. Alternatively, zero-noise extrapolation as recently done in [42] can be employed, if the extra circuit depth needed is available. We note that while both methods provide good results when simulating the Rényi entropy (see below), only method II matches the exact evolution of $\langle \hat{H} \rangle$ well. A particular Trotterization doesn't in general reproduce all observables with similar accuracy.

To measure R_2 , we follow the technique outlined in [7], which requires two copies of the state $|\Psi\rangle = \sum_{i,j} c_{ij} |a_i\rangle |b_j\rangle$, where $|a_i\rangle$ and $|b_j\rangle$ are the basis states of subspace A and B , respectively. R_2 is given by the expectation value of the Swap operator on subspace A ,

$$R_2 = \langle \Psi | \langle \Psi' | \text{Swap}_A | \Psi \rangle | \Psi' \rangle, \quad (7)$$

where the apostrophe indicates states belonging to subsystem 2, and the operator Swap_A acts as follows:

$$\text{Swap}_A |\Psi\rangle | \Psi' \rangle = \sum_{i,j} \sum_{i',j'} c_{ij} c_{i'j'} |a_{i'}\rangle |b_j\rangle |a_i\rangle |b_{j'}\rangle. \quad (8)$$

To extract R_2 experimentally, we apply the Swap-gate to the subsystems A of two copies of the adiabatically evolved state, conditional on the state of an ancilla qubit. The ancilla qubit is prepared and measured in the x -basis by applying a Hadamard gate before and after the C-Swap gate (see Fig. 1). Repeating the simulation and measurement, and averaging the results allows us to determine the probability P_a to find the ancilla qubit in state $|0\rangle$ or $|1\rangle$ from which R_2 is calculated as $R_2 = P_a(0) - P_a(1)$.

The C-Swap or Fredkin gate [43] has been experimentally implemented in NMR [44] and photonic systems [45, 46]. Our work is the first implementation of a C-Swap gate with trapped ions. Its evolution operator is

$$\hat{U}_{\text{C-Swap}} = \begin{pmatrix} 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 1 \end{pmatrix}. \quad (9)$$

We realize this gate by programming our quantum compiler to break it down into gates from our native library [35, 36]. It requires seven entangling gates and fourteen

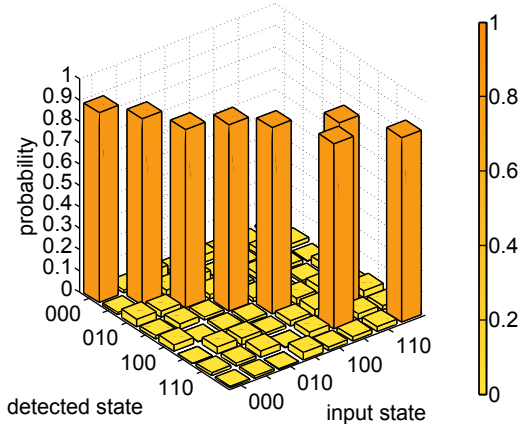


FIG. 3. Implementation of the C-Swap or Fredkin gate. The plot shows results for all eight input states, reproducing the evolution operator (eqn. 9) with an average success probability of 86.8(3)%, while the state of the control qubit is correct with 94.0(2)% probability. The results have been corrected for $\sim 1\%$ state-preparation and measurement error.

single-qubit rotations. A circuit diagram detailing its modular implementation is shown in Appendix D. We apply each logical input state and record the output state probabilities. The results are shown in figure 3. Compared to the ideal evolution operator shown in equation 9, the average success probability is 86.8(3)%. The control qubit, on which the measurement of R_2 hinges, is found to be in the correct state with 94.0(2)% probability. We also apply the gate on a balanced superposition of input states and measure 85.2(7)% correct outcomes. From these measurements, we derive an estimate for the process fidelity of 83(3)%, using the method described in [47].

With all elements of the circuit in place, we now implement the entire algorithm to adiabatically evolve two copies of the two-qubit Fermi-Hubbard system and measure the second Rényi entropy using the top qubit as the ancilla (see Fig. 1). The results are shown as open symbols for method I and II in figure 4(a) and (b), respectively. Both figures include a curve showing the theoretically expected values for the chosen Trotter step (δ) and evolution time (τ), as well as the exact solution for comparison. The exact solution shows R_2 monotonically decreasing with increasing interaction, implying increasing entanglement between the qubits. For method I the curve starts just below 0.9. At this point, no adiabatic evolution is applied, and this value is expected from the performance of the C-Swap gate. The deviation from the theoretically expected curve increases as more evolution steps are taken. For method II, 27 entangling gates have to be performed regardless of the value of U and we observe a systematically lowered value of R_2 . It is clear that the measurement outcomes of the ancilla qubit are determined by both the dynamics of the model and the errors in the physical gates.

To distinguish between these two phenomena, we de-

velop a method to detect erroneous runs using the additional information available in the four data qubits representing the two systems. The eigenvalues of the Swap operator are ± 1 with the corresponding eigenstates being even and odd functions of the qubits being swapped. For any operator \hat{V} , with eigenvalues λ_m and eigenvectors $|m\rangle$ such that $\hat{V}|m\rangle = \lambda_m|m\rangle$, and a state $|\Phi\rangle = \sum_m d_m|m\rangle$, the expectation value $\langle\Phi|\hat{V}|\Phi\rangle$ can be obtained by applying \hat{V} conditional on the state of an ancilla qubit. The ancilla is prepared along x before the controlled- \hat{V} operation and measured in the X-basis thereafter. Just before measurement, the qubits are in the following state, where the first ket in each term describes the ancilla qubit:

$$|0\rangle \sum_m (1 + \lambda_m) d_m |m\rangle + |1\rangle \sum_m (1 - \lambda_m) d_m |m\rangle. \quad (10)$$

This is essentially the circuit used to measure the expectation value of the Swap_A operator to determine the Rényi entropy. Therefore we can make the observation that the probability of $|1\rangle|a_i\rangle|b_j\rangle|a_{i'}\rangle|b_{j'}\rangle$ is 0 if $i' = i$ or if $j' = j$. This implies that twelve of the 32 possible output states of the 5-qubit register should have zero-weight when it comes to evaluating R_2 (see Appendix E). We re-analyze the data after discarding such outcomes and find the values given by the filled symbols in figure 4. The data points now follow the theoretical curves, showing that the method succeeds in substantially reducing experimental errors. This technique is general since it only depends on the symmetry of the C-Swap operation and is independent of the model Hamiltonian under investigation. The yield, or fraction of data runs that are kept under this method, is $\sim 84\%$ for method II, and drops with U from 94% to 83% for method I (see Appendix E). For the deeper circuits, the technique is not able to discard all errors, however, as can be seen in figure 4(a) at $U = 6$.

In summary, we have demonstrated a complete chain of steps (i.e. the ‘full quantum computing stack’) to simulate a model Hamiltonian on a quantum computer and measure bipartite entanglement using a circuit with up to 31 entangling and 78 single-qubit gates. Each step is scalable to a larger system of qubits on the trapped-ion hardware platform. The technique can also be implemented on different quantum computing architectures. The error-mitigation method we have used is expected to scale well to larger system sizes since the number of states it discards scales with the size of the Hilbert space of the system, but it will break down at larger circuit depths. Therefore it is a good technique for the NISQ (Noisy Intermediate-Scale Quantum) era [48], where we might expect short-depth circuits acting on several qubit systems without fault tolerance to be useful for certain applications.

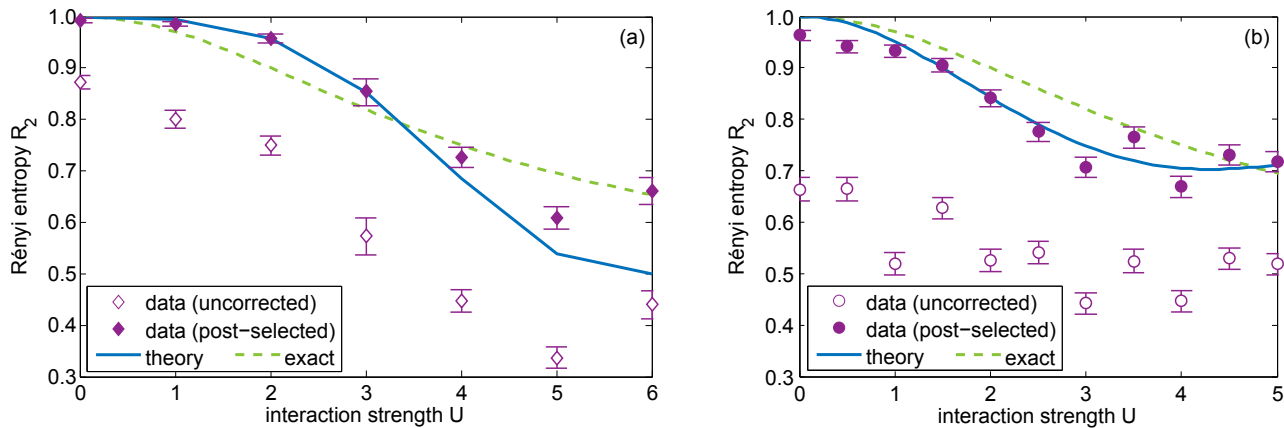


FIG. 4. Results of the measurement of R_2 after the digitized adiabatic evolution to different interaction strengths U according to method I (a) and II (b) compared to the expected curve (solid line) and the exact solution (dashed line). The open symbols show the original data while the filled symbols show the same data after post-selection based on the symmetry of the C-Swap gate.

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Appendix A: Experimental system

We perform the experiment on a quantum computer consisting of a chain of five $^{171}\text{Yb}^+$ ions confined in a Paul trap and laser cooled near the motional ground state. The hyperfine-split $^2S_{1/2}$ ground level with an energy difference of 12.642821 GHz provides a pair of qubit states, $|0\rangle = |0,0\rangle$ and $|1\rangle = |1,0\rangle$ with quantum numbers $|F, m_F\rangle$, which are magnetic field independent to first order. The $1/e$ -coherence time of this so-called “atomic clock” qubit is 1.5(5) s in our system, which can be extended by reducing magnetic field noise. Optical pumping is used to initialize the state of all ions, and the final states are measured collectively via state-dependent fluorescence detection [49]. Each ion is mapped to a distinct channel of a photomultiplier tube (PMT) array. The average state detection fidelity is 99.4(1)% for a single qubit, while a 5-qubit state is typically read out with 95.7(1)% average fidelity, limited by channel-to-channel crosstalk. These state detection and measurement (SPAM) errors are characterized in detail by measuring the state-to-state error matrix. For averaged data such as the one shown in this work, they are corrected by re-normalizing the averaged state vector by the inverse of this matrix. Quantum operations are achieved by applying two Raman beams from a single 355nm mode-locked laser, which form beat notes near

the qubit frequency. The first Raman beam is a global beam applied to the entire chain, while the second is split into individual addressing beams, each of which can be switched independently to target any single qubit [35]. Single qubit gates are generated by driving resonant Rabi rotations (R-gates) of defined phase, amplitude, and duration. Two-qubit gates (so-called XX-gates) are realized by illuminating two ions with beat-note frequencies near the motional sidebands and creating an effective spin-spin (Ising) interaction via transient entanglement between the state of two ions and all modes of motion [50–52]. To ensure that the motion is left disentangled from the qubit states at the end of the interaction, we employ a pulse shaping scheme by modulating the amplitude of the global beam [53, 54]. The signal to drive each ion is generated by an individual Arbitrary Waveform Generator (AWG) which allows us to efficiently apply single-qubit Z-rotations as classical phase advances.

Appendix B: Trotter error and resource analysis

Figure 5 shows the effect of varying the Trotter step size δ and the adiabatic evolution time τ on the accuracy of the measured R_2 as compared to the exact value between $U = 0$ and $U = 10$. We only consider the first order Trotter approximation because with only two non-commuting terms in the Hamiltonian, the first and second order approximations are asymptotically equivalent. We see that increasing τ increases the amplitude of the oscillations around the exact result, whereas increasing δ leads to an increase in a constant offset from the exact value. When δ becomes too large (compared to the largest eigenvalue of the Hamiltonian), the Trotter approximation breaks down.

Figure 6 shows how the error from the first-order Trotter approximation varies with different parameters. The error in R_2 is defined as the chi-squared goodness of fit of

the output of the quantum algorithm to the actual data between $U=0$ to $U=10$, that is

$$\epsilon_{R_2} = \overline{\left(\frac{(R_2^{\text{exact}} - R_2^{\text{sim}})^2}{R_2^{\text{exact}}} \right)}, \quad (\text{B1})$$

where the average is over $U = 0$ to $U = 10$ measured at intervals of $U = 0.1$. As expected from the adiabatic evolution theorem, $\epsilon_{R_2} \sim \tau^{-2}$, which is the behavior seen in figure 6(a).

The error in the wavefunction, defined as the projection of the wavefunction orthogonal to the ground state, averaged between $U=0$ to $U=10$, is given by

$$\epsilon_\Psi = 1 - \overline{\langle \Psi^{\text{exact}} | \Psi^{\text{sim}} \rangle^2}. \quad (\text{B2})$$

The data show that the error in the wavefunction and the error in the actual quantity of interest can scale differently.

Figure 7 shows the circuit depth as a function of δ and τ scales linearly with each, which is to be expected from the first-order Trotter approximation and adiabatic evolution.

Appendix C: Parameter choices for the experiment

Figure 8 shows theoretical calculations of the outcome of the R_2 measurement for various parameters that satisfy the experimental restriction $N_{\text{steps}} \leq 6$. For the experiment, the following sets were chosen: For method I, $\delta = \tau = 0.1$, and for method II, $\delta = 0.25$ and $\tau = 1.25/U$.

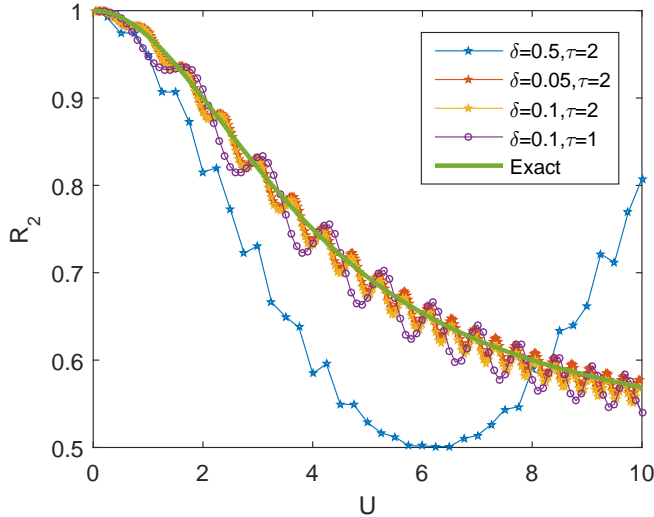


FIG. 5. The value of R_2 output from the quantum algorithm compared to the exact value at different step values of the adiabatic evolution time τ and Trotter step size δ for the first-order Trotter approximation.

Appendix D: The C-Swap or Fredkin gate

The compiler breaks down the Fredkin (C-Swap) gate into native R- and XX-gates as given by the circuit in figure 9.

Appendix E: Error reduction

Using the convention that the top-most qubit in the five-qubit output state after the circuit shown in figure 1 is the most significant bit in the binary representation of the computational basis, we can assign each state from $|00000\rangle$ to $|11111\rangle$ a decimal value from 0 to 31. The observation about the symmetry of the state after the C-Swap gate implies that the states numbered 16, 17, 18, 20, 21, 23, 24, 26, 27, 29, 30, and 31 should have zero-weight at the end of the circuit. This can be used to post-select the results by discarding runs that result in these outcomes. Figure 11 shows the results before and after this step. It also gives the yields associated with each experiment. We see that we discard less than 20% of the data and achieve an improvement in the results of over 40% (see Fig. 4).

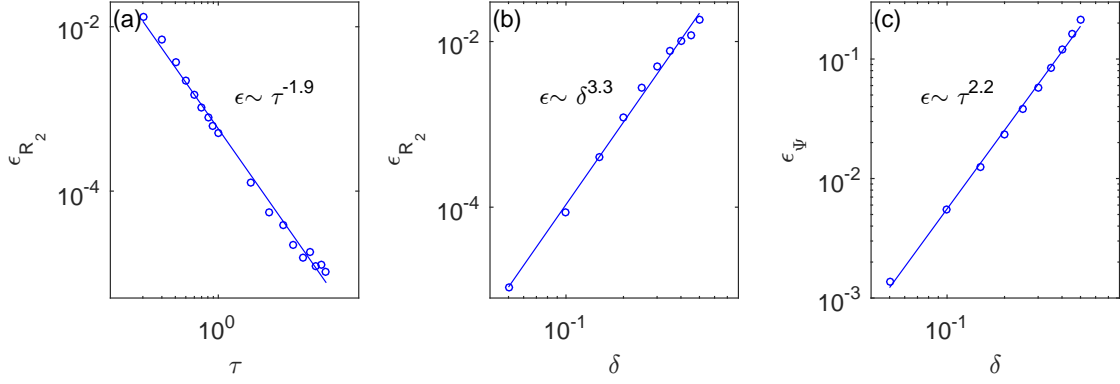


FIG. 6. (a) The error in R_2 , ϵ_{R_2} (defined in the text) as a function of evolution time τ at fixed $\delta = 0.05$ fits well to a line of slope -1.9. (b) ϵ_{R_2} as a function of Trotter step size δ at fixed $\tau = 10$ fits well to lines of slope 3.3. (c) The error in the wavefunction ϵ_Ψ (defined in the text) at fixed $\tau = 10$ as a function of Trotter step size δ fits well to a line of slope 2.2.

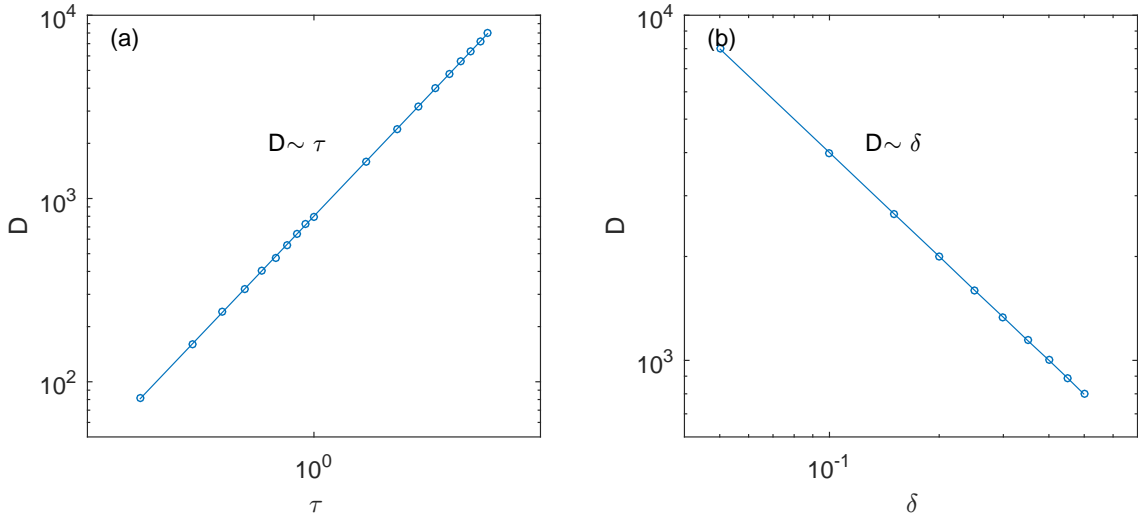


FIG. 7. In first Trotter approximation, (a) the parallel circuit depth D as a function of evolution time τ at fixed $\delta = 0.05$ fits well to lines of slope 1, and (b) D as a function of Trotter step size δ at fixed $\tau = 10$ fits well to lines of slope -1. Here D is defined as the circuit depth for adiabatic evolution to $U = 10$.

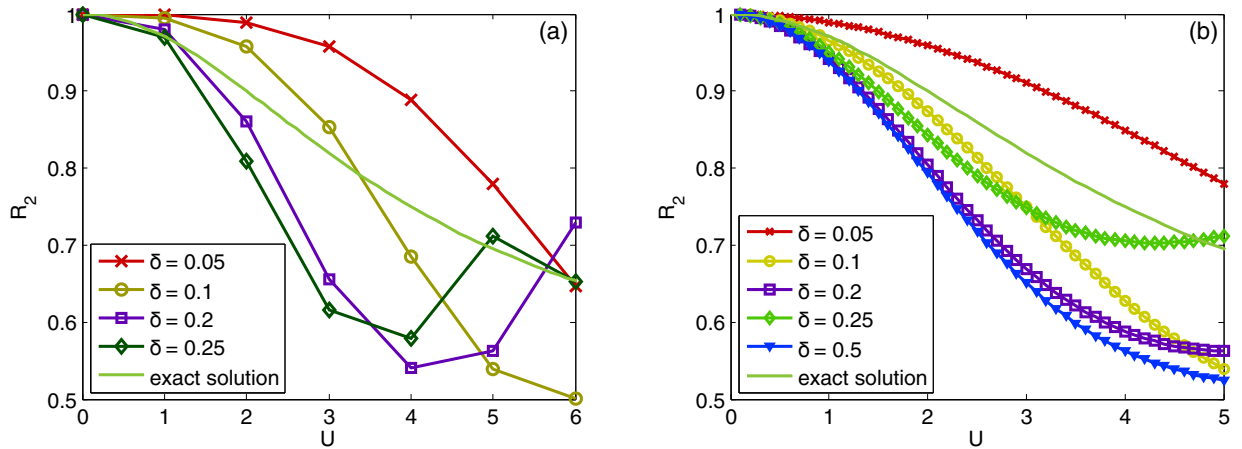


FIG. 8. The value of R_2 resulting from the quantum algorithm compared to the exact value for different values of the adiabatic evolution time τ and Trotter step size δ for (a) method I with fixed $\tau/\delta = 1$, and (b) method II with fixed $N_{\text{steps}} = 5$.

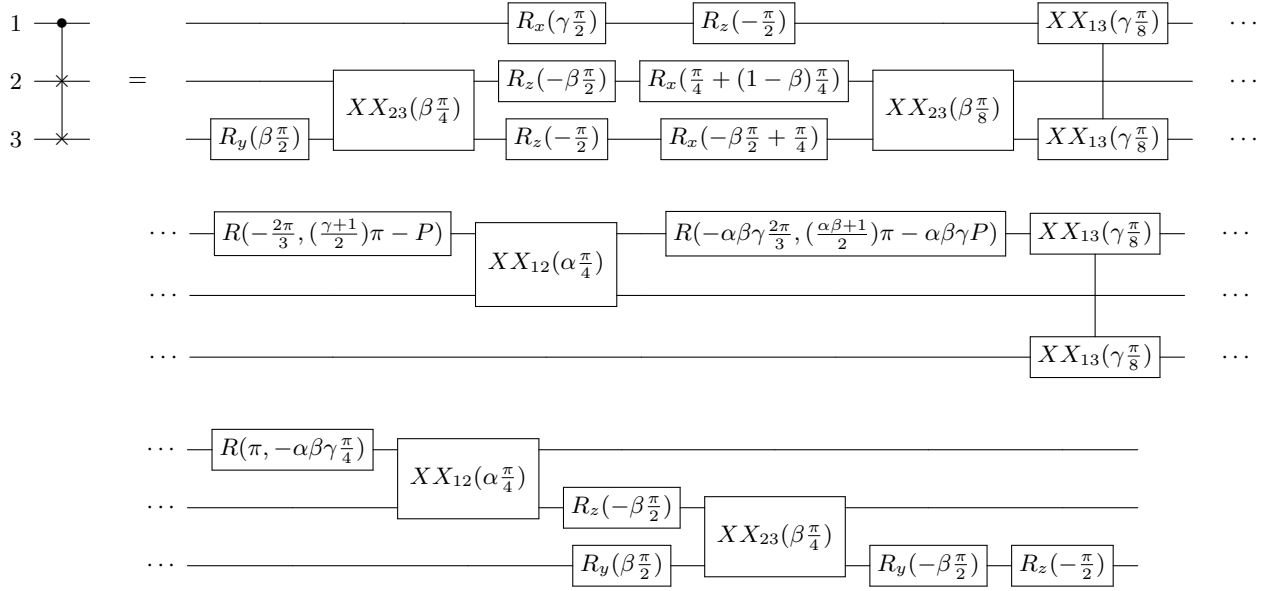


FIG. 9. Controlled-Swap (or Fredkin) gate implementation using entangling $XX(\chi)$ gates, and single qubit rotations $R_x(\theta)$, $R_y(\theta)$, and $R(\theta, \phi)$. The parameters are $\alpha = \text{sgn}(\chi_{12})$, $\beta = \text{sgn}(\chi_{23})$, $\gamma = \text{sgn}(\chi_{13})$, and $P = \arcsin \sqrt{\frac{2}{3}}$.

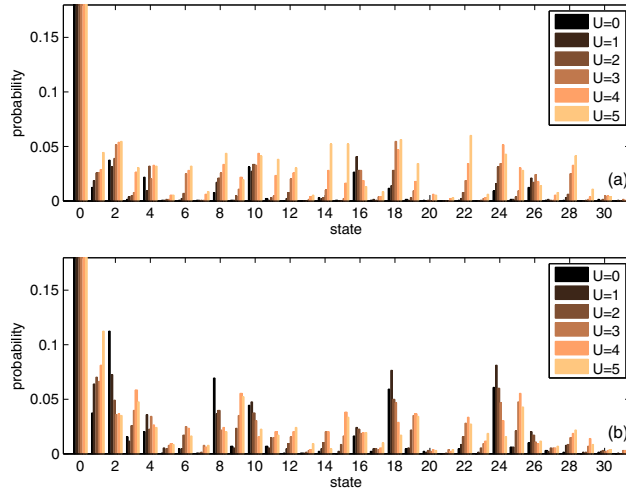


FIG. 10. The original output of the circuit shown in figure 1 averaged over 2000-2500 runs, corrected for SPAM errors. Method I is shown in (a) and method II in (b).

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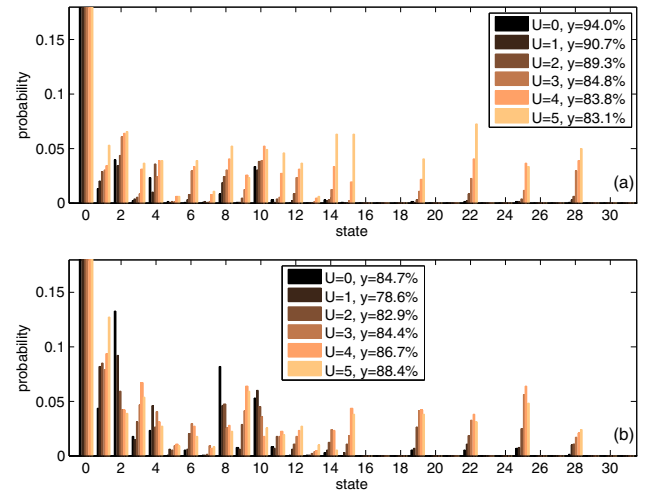


FIG. 11. The output of the circuit shown in figure 1 averaged over 2000-2500 runs, corrected for SPAM errors and after discarding in post-selection any runs that resulted in the states given in the text. Method I is shown in (a) and method II in (b). The legend gives the yields, i.e. the fraction of experimental runs that was kept.

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