Simulating large-size quantum spin chains on cloud-based superconducting quantum computers

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Quantum computers have the potential to efficiently simulate large-scale quantum systems for which classical approaches are bound to fail. Even though several existing quantum devices now feature total qubit numbers of more than 100, their applicability remains plagued by the presence of noise and errors. Thus, the degree to which large quantum systems can successfully be simulated on these devices remains unclear. Here, we report on numerical results of physics-motivated variational ansatzes, as well as cloud simulations performed on several of IBM's superconducting quantum computers to simulate ground states of spin chains having a wide range of system sizes up to 102 qubits. Our numerical analysis shows that the accuracy of the ground-state energy and fidelity improves substantially by increasing the number of layers used in the ansatzes. From the cloud experiments, we find that the ground-state energies extracted from realizations across different quantum computers and system sizes reach the expected values to within errors that are small (i.e., on the percent level), including the inference of the energy density in the thermodynamic limit from these values. We achieve this accuracy through a combination of physics-motivated variational ansatzes, and efficient, scalable energy-measurement and error-mitigation protocols, including the use of a reference state in the zero-noise extrapolation. By using a 102-qubit system, we have been able to successfully apply up to 3186 CNOT gates in a single circuit when performing gate-error mitigation. Our accurate, error-mitigated results for random parameters in the ansatz states suggest that a standalone hybrid quantum-classical variational approach for large-scale XXZ models considered in this work is feasible.

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I. INTRODUCTION

The notion of quantum computers traces back to the works of Benioff [1], Mannin [2], and Feynman [3]. In particular, Feynman suggested using quantum computers to simulate other quantum systems instead of using classical computers, giving rise to the notion of a universal quantum simulator [4]. A critical breakthrough was made by Shor, whose quantum factoring algorithm outperforms classical algorithms almost exponentially faster [5]. Experimental progress has come a long way, leading to the burgeoning of quantum devices, with the total qubit number now exceeding one hundred in the best cases. However, these devices are still regarded as noisy intermediate-scale quantum (NISQ) processors [6], not yet suitable for full-scale quantum error correction and fault-tolerant quantum computation. However, there are concurrent efforts to develop error mitigation techniques [7-15]and algorithms [16] for NISQ devices to realize their potential quantum advantage [17]. Recent notable experimental achievements include random quantum circuits of around 50 qubits [18,19], boson sampling with large numbers of photons [20,21], Hartree-Fock method implementation for quantum chemistry with 12 qubits [22], realization of the toric-code state with 31 qubits [23], and quantum walks on a 62-qubit processor [24].

Despite all this effort and accomplishment, the central question remains of whether NISQ computers can be of practical use for the simulation of large quantum systems and to extract accurate observable values, such as the energy of the simulated quantum states. So far, most experiments with quantitatively accurate results have been limited to small numbers of qubits, around ten or below [9-12,22], with a few others reaching beyond twenty [15,23]. None of them has demonstrated accurate results over a wide range of system sizes with the same model and across different devices. There are also challenges to overcome for largescale experiments (around or over one hundred qubits) with useful outcomes, including the need for high-fidelity gates and readout as well as scalable and efficacious approaches to mitigating the effects of noise and errors on the measured observables.

In this work, we have designed a family of physicsmotivated variational ansatzes and analyzed their accuracy for the Heisenberg and XXZ spin chains. For small numbers of qubits, our ansatzes can reach exact ground-state wave functions and for larger number of qubits, the accuracy of the ground-state energy and fidelity improves substantially by increasing the number of layers used in the ansatzes. For

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TABLE I. Results related to the open-chain Heisenberg model. The numerical calculation was done with the MPS method using a bond dimension $\chi = 64$. The "error" in the last column represents the relative error between the experimentally estimated value E_{exp} and the exact ground-state energy E_{gs} . *Note that these values were obtained by averaging results over different backends and/or different groups of physical qubits; see Tables S.2 and S.3 in the Supplemental Material [45] for the complete list of results.

N	θ^*_{even}	$ heta^*_{\mathrm{odd}}$	$E^*_{ m ansatz}$	$E_{\rm gs}$	ϵ	f	$E_{ m exp}$	error
4	0.151748	0.215765	-6.464102	-6.464102	0	1.0000	-6.5(1.6)	0.56%
6	0.141671	0.216088	-9.880996	-9.974309	0.94%	0.9923	$-9.9(1.9)^{*}$	0.19%
8	0.138569	0.216093	-13.299823	-13.499730	1.48%	0.9796	-13.2(2.2)	2.22%
10	0.13710	0.216102	-16.719307	-17.032141	1.84%	0.9639	$-16.7(1.3)^{*}$	1.95%
12	0.136248	0.216110	-20.139037	-20.568363	2.09%	0.9462	-20.3(2.1)	1.30%
14	0.135688	0.216115	-23.558885	-24.106899	2.27%	0.9271	-23.6(1.8)	2.10%
16	0.135293	0.216120	-26.978800	-27.646949	2.42%	0.9072	$-25.8(1.6)^{*}$	6.68%
18	0.134999	0.216123	-30.398756	-31.188044	2.53%	0.8867	$-30.7(0.7)^*$	1.56%
20	0.134773	0.216126	-33.818738	-34.729893	2.62%	0.8659	$-33.0(0.5)^{*}$	4.98%
30	0.134132	0.216134	-50.918850	-52.445423	2.91%	0.7614	$-50.2(2.0)^{*}$	4.28%
40	0.133832	0.216139	-68.019098	-70.165893	3.06%	0.6629	$-68.5(2.0)^{*}$	2.34%
50	0.133658	0.216141	-85.119397	-87.888441	3.15%	0.5737	-85.0(2.8)*	3.29%
60	0.133544	0.216143	-102.219721	-105.612060	3.21%	0.4946	-99(4)	6.26%
70	0.133464	0.216144	-119.320058	-123.336305	3.26%	0.4253	-125(7)	1.35%
80	0.133405	0.216145	-136.420403	-141.060947	3.29%	0.3649	-138.5(2.5)	1.82%
90	0.133359	0.216146	-153.520754	-158.785857	3.32%	0.3126	-153(5)	3.64%
98	0.133329	0.216146	-167.201038	-172.965924	3.33%	0.2760	-168.1(2.6)	2.81%
100	0.133323	0.216146	-170.621109	-176.510957	3.34%	0.2675	-173(9)	1.99%
102	0.133316	0.216146	-174.041180	-180.055995	3.34%	0.2592	-177.5(2.7)	1.42%

example, with six layers and N = 50 qubits in the Heisenberg model, the fidelity of our ansatz with the ground-state wave function is above 0.9 and the accuracy in the ground-state energy is above 99.75%. For potential realization, we have also developed efficient approaches for measuring energies, including the use of Bell measurement. Moreover, we have also improved scalable mitigation methods to extract accurate GS energy values for large systems, despite the presence of noise and errors in the gates and the readout. We have also introduced a reference state in the zero-noise extrapolation (rZNE) to further improve the accuracy of the results.

To demonstrate that noisy quantum processors can still be used to provide accurate results, we have utilized nine distinct cloud quantum computers, and carried over 40 different sets of cloud experiments for the Heisenberg model (see Tables I-III) and over 48 sets of them for the XXZ model (see the figures presenting the results below). We present realizations of approximate ground states (GS) of these spin chains having nineteen different system sizes, ranging from 4 to 102 qubits. To distinguish our work from experiments performed on in-house devices or customized physical apparatuses, we shall refer to our use of third-party hardware as "cloud experiments," as well as to make a distinction from numerical simulations. We report the extracted GS energies, accurate to within a few percent level of error, including the inference of the energy density in the thermodynamic limit from these values. We emphasize that these cloud experiments are not equivalent to numerical simulations, as the actual devices have substantial noise and errors and devices' condition can drift over time, and sometimes the same submitted jobs can fail. Nevertheless, cloudbased experiments offer a new paradigm for research and development. In addition, we have used our procedure to measure the energies of several ansatz states that have randomly chosen parameters, and obtained accurate mitigated energy values. Our work thus establishes a simple–yet substantially improved–quantum variational protocol with mitigation, and paves the way for massive use of large NISQ computers for fundamental physics studies of many-body systems, as well as for practical applications, including optimization problems.

The remaining structure of the paper is as follows. In Sec. II, we discuss the motivation that leads to our family of ansatzes for the Heisenberg and *XXZ* spin chains and give the decomposition of the ansatz gates in terms of elementary ones. In Sec. II, we analyze our ansatzes numerically and present results for various numbers of qubits and layers in the ansatz, including cases of both open and periodic conditions. In Sec. IV, we discuss scalable methods to measure energy and mitigate measured energy values and present cloud experimental results from real devices. In Sec. V, we present results of numerical analysis of our ansatz on a two-leg ladder. We conclude in Sec. VI. Additional supporting materials are presented in the Appendix.

II. HEISENBERG AND XXZ MODELS AND THE ANSATZ FOR GROUND STATES

Quantum spin systems, such as the Heisenberg [25] and XXZ models [26,27], have sparked analytical development and understanding of quantum phases and also served as a testbed for numerical techniques. The Hamiltonian of the spin-1/2 XXZ spin chain with the open boundary condition

TABLE II.	Various Heisenberg sp	in-chain cloud exp	periments perf	formed on the	127-qubit ibm_	washington backend	/device of IBM Q
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N	$E_{\rm exp}$	ϵ_{ans}	$\epsilon_{ m gs}$	QOD	Qubits used	Shots	Rep.
10	-16.9(3.4)	1.08%	0.78%	44(7)	[30,31,32,36,51,50,49,48,47,35]	40K	50
10	-16.5(1.4)	1.31%	3.12%	31(4)	[32,36,51,50,49,48,47,35,28,29]	40K	25
20	-33.8(1.4)	0.055%	2.67%	20.3(1.0)	[50,51,36,32,31,30,29,28,35,47,	40K	25
					48,49,55,68,69,70,74,89,88,87]		
30	-51.1(2.3)	0.36%	2.57%	19.6(1.1)	[115,116,117,118,110,100,101,102,92,83,84,85,73,	40K	50
					66,67,68,55,49,50,51,36,32,31,30,29,28,27,26,16,8]		
40	-69.2(1.6)	1.74%	1.38	12.1(0.8)	[2,1,0,14,18,19,20,33,39,38,37,52,56,57,58,59,60,61,62,	40K	50
					72,81,82,83,92,102,103,104,105,106,93,87,		
					86,85,73,66,67,68,55,49,48]		
50	-86.9(1.8)	2.09%	1.12%	13.19(0.28)	list of 40 qubits +[47,35,28,27,26,25,24,23,22,15]	40K	25
60	-99(4)	3.15%	6.26%	11.28(0.34)	[3,2,1,0,14,18,19,20,33,39,38,37,52,56,57,58,59,	40K	50
					60,53,41,42,43,34,24,25,26,27,28,29,30,31,32,		
					36,51,50,49,48,47,46,45,54,64,63,62,72,81,82,83,92,		
					102,103,104,105,106,107,108,112,126,125,124]		
70	-125(7)	4.76%	1.35%	13.1(0.8)	[3,2,1,0,14,18,19,20,21,22,23,24,25,26,27,28,29,30,	40K	50
					31,32,36,51,50,49,48,47,46,45,54,64,65,66,67,68,69		
					70,74,89,88,87,93,106,105,104,111,122,121,120,119,		
					118,110,100,101,102,92,83,82,81,80,79,		
					91,98,97,96,95,94,90,75,76,77]		
80	-138.5(2.5)	1.52%	1.82%	12.79(0.23)	[3,2,1,0,14,18,19,20,33,39,38,37,52,56,57,58,	40K	50
					59,60,53,41,42,43,34,24,25,26,27,28,29,30,31,32,36,		
					51,50,49,55,68,69,70,74,89,88,87,93,106,105,104,		
					111,122,121,120,119,118,110,100,101,102,92,83,84,85,73,		
					66,65,64,63,62,72,81,80,79,91,98,97,96,95,94,90,75]		
90	-153(5)	0.34%	3.64%	12.5(0.4)	[3,2,1,0,14,18,19,20,21,22,23,24,25,26,27,28,29,30,31,32,	40K	25
					36,51,50,49,48,47,46,45,44,43,42,41,40,39,38,37,52,56,		
					57,58,59,60,61,62,63,64,65,66,67,68,69,70,74,89,88,87,86,		
					85,84,83,82,81,80,79,78,77,76,75,90,94,95,96,97,98,99,		
					100,101,102,103,104,105,106,107,108,		
					112,126,125,124,123,122]		
98	-168.1(2.6)	0.54%	2.81%	12.32(0.27)	[3,2,1,0,14,18,19,20,33,39,40,41,42,43,34,24,23,22,15,4,5,	40K	75
					6,7,8,16,26,27,28,29,30,31,32,36,51,50,49,48,47,46,45,		
					54,64,63,62,61,60,59,58,71,77,76,75,90,94,95,96,97,98,91,		
					79,80,81,82,83,84,85,73,66,67,68,69,70,74,89,88,87,93,		
					106,107,108,112,126,125,124,123,122,111,104,103,102,		
					101,100,110,118,117,116,115,114]		
100	-173(9)	1.39%	1.99%	11.9(0.14)	[3,2,1,0,14,18,19,20,21,22,15,4,5,6,7,8,16,26,27,28,29,30,	40K	50
					31,32,36,51,50,49,48,47,46,45,44,43,42,41,40,39,38,37,52,		
					56,57,58,59,60,61,62,63,64,65,66,67,68,69,70,74,89,88,87,		
					86,85,84,83,82,81,80, 79,78,77,76,75,90,94,95, 96,97,98,		
					99,100,101,102,103,104,105,106,107,108,112,126,125,124,		
					123,122,121,120,119,118,117,116]		
102	-177.5(2.7)	1.99%	1.42%	12.23(0.17)	list of 100 qubits +[115,114]	40K	75

reads

$$\hat{H}_{XXZ}(\Delta) = \sum_{j=1}^{N-1} \hat{h}_{XXZ}^{[j,j+1]}(\Delta)$$

= $\sum_{j=1}^{N-1} \left(\sigma_x^{[j]} \sigma_x^{[j+1]} + \sigma_y^{[j]} \sigma_y^{[j+1]} + \Delta \sigma_z^{[j]} \sigma_z^{[j+1]} \right),$
(1)

where Δ represents the anisotropy in the coupling. For $\Delta = 1$, the model reduces to the isotropic antiferromagnetic Heisenberg chain. The model is known to possess three distinct

quantum phases: (i) a ferromagnetic phase for $\Delta < -1$, where classical states, such as $|\uparrow\uparrow\uparrow\ldots\rangle$ and $|\downarrow\downarrow\downarrow\ldots\rangle$, are ground states; (ii) a gapless, critical phase, for $-1 < \Delta < 1$; and (iii) an antiferromagnetic phase for $\Delta > 1$. We will mainly focus on the range of $\Delta > -1$ with nontrivial ground states.

In the following, we explain how we use adiabatic connection [28] to arrive at a physics-motivated ansatz, schematically shown in Fig. 1(a), and justify its validity by considering the gap structure through the adiabatic connection, as illustrated in Fig. 2. We also analyze how well the ansatz performs and how its accuracy is improved by increasing the number of layers. TABLE III. Heisenberg spin-chain cloud experiments performed on all available 27-qubit backends/devices and the 65-qubit ibmq_brooklyn of IBM Q.

	Backend						
N	E _{exp}	$\epsilon_{ m ans}$	$\epsilon_{ m gs}$	QOD	qubits used	shots	rep.
ibm_	_auckland						
20	-34.3(1.3)	1.42%	1.24%	13.4(0.5)	[13, 12, 10, 7, 4, 1, 2, 3, 5, 8, 11, 14, 16, 19, 22, 25, 24, 23, 21, 18]	100K	50
ibm	cairo						
16	-27.5(2.5)	1.93%	0.53%	13.5(1.1)	[1,2,3,5,8,11,14,16,19,22,25,24,23,18,15,12]	100K	50
18	-28.9(1.3)	4.93%	7.34%	13.4(0.4)	[1,2,3,5,8,11,14,16,19,22,25,24,23,18,15,12]	100K	50
18	-32.3(1.1)	6.25%	3.57%	13.51(1.1)	[6,7,4,1,2,3,5,8,11,14,16,19,22,25,24,23,21,18]	100K	50
20	-32.8(1.1)	3.01%	5.56%	12.47(0.35)	[6,7,4,1,2,3,5,8,11,14,16,19,22,25,24,23,21,18,15,12]	100K	50
ibm	hanoi						
20	-32.3(1.3)	4.49%	7.00%	17.0(0.7)	[5,3,2,1,4,7,10,12,15,18,21,23,24,25,22,19,16,14,11,8]	100K	100
ibma	q_kolkata						
20	-33.5(1.2)	0.94%	3.54%	18.6(0.8)	[26,25,22,19,16,14,11,8,5,3,2,1,4,7,10,12,15,18,21,23]	40K	25
ibmo	q_montreal						
6	-9.9(1.6)	0.75%	0.19%	13.1(0.6)	[20,19,22,25,24,23]	32K	100
20	-31.7(1.7)	6.26%	8.72%	13.0(0.5)	[12,10,7,4,1,2,3,5,8,11,14,16,19,22,25,24,23,21,18,17]	32K	50
ibmo	mumbai						
4	-6.5(1.6)	0.56%	0.56%	37(9)	[12,15,18,17]	8192	146
6	-9.9(3.5)	0.19%	0.75%	64(11)	[7,10,12,15,18,17]	8192	146
8	-13.2(2.2)	0.75%	2.22%	36(5)	[1,4,7,10,12,15,18,21]	8192	146
10	-16.6(3.3)	0.71%	2.54%	38(7)	[0,1,4,7,10,12,15,18,21,23]	8192	100
12	-20.3(2.1)	0.80%	1.30%	14.0(1.4)	[8,5,3,2,1,4,7,10,12,15,18,21]	8192	100
14	-23.6(1.8)	0.17%	2.10%	14.1(1.0)	[8,5,3,2,1,4,7,10,12,15,18,21,23,24]	8192	100
16	-24.0(2.1)	10.1%	7.42%	12.1(0.8)	[8,5,3,2,1,4,7,10,12,15,18,21,23,24,25,22]	8192	100
18	-29.7(1.3)	2.30%	4.77%	16.7(0.7)	[8,5,3,2,1,4,7,10,12,15,18,21,23,24,25,22,19,16]	8192	100
20	-31.6(2.3)	6.56%	9.01%	14.2(1.0)	[8,5,3,2,1,4,7,10,12,15,18,21,23,24,25,22,19,16,14,11]	8192	100
ibm	toronto						
18	-32.1(1.7)	5.60%	2.92%	13.1(0.6)	[0,1,4,7,10,12,15,18,21,23,24,25,22,19,16,14,11,8]	16K	50
	-32.3(1.4)	4.49%	7.00%	13.0(0.5)	[0,1,4,7,10,12,15,18,21,23,24,25,22,19,16,14,11,8,5,3]	32K	50
ibmo	brooklyn						
10	-16.8(1.6)	0.48%	1.36%	24.0(2.1)	[53,47,48,49,40,35,34,33,25,19]	100K	50
20	-33.4(0.9)	1.24%	3.83%	22.7(0.7)	[9.8.7.6.5.4.3.2.1.0.10.13.14.15.16.17.18.19.25.33]	100K	50
20	-34.0(3.1)	0.54%	2.10%	27.4(3.4)	[0,1,2,3,4,5,6,7,8,12,21,20,19,18,17,16,15,24,29,28]	20K	80
30	-49.3(3.3)	3.18%	6.00%	18.4(1.1)	[43,52,56,57,58,59,60,53,47,48,49,40,35,34,33,32,31,	100K	50
	.,				30.2.24,15,16,17,11,4,5,6,7,8,9]		
40	-68(4)	0.028%	3.09%	20.6(1.3)	[63.62,61,60.53,47,46,45,39,31,32,33,25,19,18,17,16,15,	100K	50
	. ,				14,13,10,0,1,2,3,4,5,6,7,8,12,21,22,23,26,37,36,35,40,49]		
50	-83(5)	2.49%	5.56%	18.7(1.5)	[43,52,56,57,58,59,60,61,62,63,64,54,51,50,49,40,35,36,	100K	50
				× /	37,26,23,22,21,12,8,7,6,5,4,3,2,1,0,10,13,14,15,16,		
					17,18,19,25,33,32,31,39,45,46,47,48]		

A. Gap structure of the interpolated Hamiltonian and the ansatz structure from adiabatic evolution

For the *XXZ* interaction $\hat{h}_{XXZ}^{[j,j+1]}(\Delta)$ on a bond involving two nearest-neighbor qubits, the singlet pair $|\Psi^-\rangle = (|01\rangle - |10\rangle)/\sqrt{2}$ has an energy value $-2 - \Delta$, the triplet $|\Psi^+\rangle = (|01\rangle + |10\rangle)/\sqrt{2}$ has energy $2 - \Delta$, and both $|00\rangle$ and $|11\rangle$ (or equivalently the two entangled triplets $|\Phi^{\pm}\rangle = (|00\rangle \pm |11\rangle)/\sqrt{2}$) have energy Δ . Note that we have used the notation $|0/1\rangle$ to replace $|\uparrow/\downarrow\rangle$, the eigenstates of the Pauli Z operator σ_z . Thus, the singlet is the ground state of the simple two-qubit *XXZ* interaction for $\Delta > -1$. This means that the following Hamiltonian with interaction only on odd numbers of bonds only even N,

$$\hat{H}_{\text{odd}} = \sum_{j=1}^{N/2-1} \left(\sigma_x^{[2j-1]} \sigma_x^{[2j]} + \sigma_y^{[2j-1]} \sigma_y^{[2j]} + \Delta \sigma_z^{[2j-1]} \sigma_z^{[2j]} \right),$$
(2)

has its unique ground state being the product of singlets over these odd bonds, i.e., a linear chain of valence-bond state,

$$|\psi_{\text{singlets}}\rangle = \frac{1}{\sqrt{2^{N/2}}} \prod_{j=1}^{N/2} (|01\rangle - |10\rangle)_{2j-1,2j}.$$
 (3)



FIG. 1. Variational ansatz and the layout of a 102-qubit quantum computer. (a) The variational ansatz structure, (b) the layout of the 127-qubit ibm_washington backend, where a chain of 102 qubits is illustrated by the thick, shaded line.

We then expect that $|\psi_{\text{singlet}}\rangle$ is adiabatically connected to the ground state of the *XXZ* model, by connecting \hat{H}_{odd} to the full *XXZ* Hamiltonian \hat{H}_{XXZ} via the following linear interpolation,

$$\hat{H}(s) = (1 - s)\hat{H}_{\text{odd}} + s\hat{H}_{XXZ} = \hat{H}_o(s) + \hat{H}_e(s).$$
(4)

We regroup it into interaction terms on even and odd bonds, denoted collectively by $\hat{H}_o(s)$ and $\hat{H}_e(s)$, respectively, and it is straightforward to see that $\hat{H}_o(s) = \hat{H}_{odd}$, but

$$\hat{H}_{e}(s) = s \sum_{j=1}^{N/2-1} \left(\sigma_{x}^{[2j]} \sigma_{x}^{[2j+1]} + \sigma_{y}^{[2j]} \sigma_{y}^{[2j+1]} + \sigma_{z}^{[2j]} \sigma_{z}^{[2j+1]} \right),$$
(5)

is a rescaled version of the XXZ model on even bonds.

We check the spectral properties of this Hamiltonian for small N and find that $\hat{H}(s)$ is gapped in the range $s \in$ [0, 1] for $\Delta > -1$; see Fig. 2 for two different Δ values using 8 qubits. This means that the product of singlets $|\psi_{\text{singlet}}\rangle$ is adiabatically connected to the ground state of the XXZ model via the evolution $|\psi(1)\rangle = \mathbf{U}_{\text{evo}}|\psi_{\text{singlets}}\rangle =$ $e^{-i\int_{s=0}^{1} ds \hat{H}(s)}|\psi_{\text{singlets}}\rangle$. Discretizing the evolution operator \mathbf{U}_{evo} , we have the following Trotterized approximation

$$\mathbf{U}_{\text{evo}} \approx \prod_{l=1}^{N_L} e^{-i\hat{H}(s_l)\delta s} \approx \prod_{l=1}^{N_L} \left(e^{-i\hat{H}_{\text{e}}(s_l)\delta s} e^{-i\hat{H}_{\text{o}}(s_l)\delta s} \right), \quad (6)$$

where N_L is the number of discretized time steps or layers and $\delta s = 1/N_L$ is the dimensionless step size. To allow for flexibility, we turn the discretized evolution into a variational form and arrive at the structure of the ansatz shown in Fig. 1(a)



FIG. 2. The energy gap of 8-qubit XXZ model with open-boundary condition in the Hamiltonian interpolation $\hat{H}(s)$ from one with interaction on odd bonds only to one with interaction on all bonds, for (a) $\Delta = 1$, i.e., the Heisenberg model, and (b) the XXZ model at $\Delta = -0.8$.

with gates in the *l*th layer being

$$U_{\text{even/odd}}^{(l)}(\{\theta\}) = \bigotimes_{j \in \text{even/odd}} \left[e^{-i\theta_{e/o,x}^{(l)} \sigma_x^{[j]} \sigma_x^{[j+1]} - i\theta_{e/o,y}^{(l)} \sigma_y^{[j]} \sigma_y^{[j+1]} - i\theta_{e/o,z}^{(l)} \sigma_z^{[j]} \sigma_z^{[j+1]}} \right],$$
(7)

where $\{\theta\}'s$ are a set of variational parameters, the subscripts e/o denote the association with even and odd bonds, respectively. Therefore, we arrive at the following N_L -layer variational ansatz state

$$|\psi_{\text{ansatz}}(\{\theta\})\rangle = \bigotimes_{l=1}^{N_L} \left[U_{\text{even}}^{(l)}(\{\theta_e\}) U_{\text{odd}}^{(l)}(\{\theta_o\}) \right] |\psi_{\text{singlets}}\rangle.$$
(8)

In Fig. 1(b), we illustrate a particular choice of a chain with 102 qubits, which we have used to implement the Heisenberg model. We remark that our construction from the adiabatic connection is similar to how the quantum approximate optimization algorithm (QAOA) ansatz originates from discretizing the Ising interaction and the transverse field [28]. But they differ in the goal: the QAOA aims to minimize the energy of a classical Ising Hamiltonian using the transverse-field part as a driver, whereas our goal is to optimize the energy of a quantum Hamiltonian as a whole. We note that similar physics-motivated ansatzes were used previously [29–31], for the Heisenberg mode on the kagome lattice was also studied numerically in Ref. [32].

B. Creation of singlets

Each singlet pair in $|\psi_{\text{singlets}}\rangle$ can be created from $|00\rangle$ by simple single-qubit gates (the Hadamard *H* and the Pauli *X*

gates) followed by a CNOT gate,

$$|\Psi^{-}\rangle = \frac{1}{\sqrt{2}}(|01\rangle - |10\rangle) = U_{\text{singlet}}|00\rangle$$
$$= \frac{|0\rangle}{|0\rangle} \underbrace{X}_{\text{H}} \underbrace{H}_{\text{H}}$$
(9)

and, thus, the product of such singlet pairs can be created in parallel with these circuits,

$$|\psi_{\text{singlets}}\rangle = U_{\text{init}}|0\dots0\rangle = \bigotimes_{i=1}^{N/2} U_{\text{singlet}}^{[i]}|0\dots0\rangle, \qquad (10)$$

with the superscript *i* denoting the pair of qubits for the singlet creation. We note that the reverse of the latter part corresponds to Bell measurement,

which can be used to measure the energy contribution of a pair of qubits; see below.

C. Gate decomposition

Let us define the essential two-qubit Rxyz gate that we need,

$$R_{xyz}(\theta_x, \theta_y, \theta_z) \equiv e^{-i(\theta_x/2)\sigma_x \otimes \sigma_x - i(\theta_y/2)\sigma_y \otimes \sigma_y - i(\theta_z/2)\sigma_z \otimes \sigma_z}, \quad (12)$$

where a factor of 1/2 is inserted in the definition of the Rxyz gate to match the convention of single-qubit rotation and we have used the tensor product notation " \otimes " to emphasize the two-qubit structure in the gate. We present a decomposition of the Rxyz gate that has a minimum number of CNOTs [33] (which is three) in the decomposition,

$$R_{xyz}(\theta_x, \theta_y, \theta_z) = \underbrace{R_z(\theta_z)}_{H} \underbrace{R_z(\theta_x + \frac{\pi}{2})}_{H} \underbrace$$

where *H* is the Hadamard gate, $R_{\alpha}(\theta) = e^{-i\theta\sigma_{\alpha}/2}$ is the singlequbit rotation around α -axis ($\alpha = x, y, z$) by an angle θ , and *S* is the one-qubit phase gate $S = e^{i\pi/4}R_z(\pi/2)$. The gate $U_{\text{even/odd}}(\theta) = R_{xyz}(2\theta, 2\theta, 2\theta)$ will be used for the Heisenberg model, and for the *XXZ* model, due to the ZZ anisotropy, we will allow $\theta_z = 2\theta_2$ parameter to be independent from $\theta_x = \theta_y = 2\theta_1$, and thus $R_{xyz}(2\theta_1, 2\theta_1, 2\theta_2)$ is needed.

Note that as the circuit action is symmetric with respect to swapping the two qubits, one can flip the circuit in the last line to fit the desired or natural direction of the CNOT gate. One can also replace the Hadamard gate *H* by a combination of the square root of *X* gate (or equivalently $R_x(\pi/2)$, which is

among the native gates in IBM Quantum Computers), and the phase gate *S* via the identity $H = SR_x(\pi/2)S$. Note that *S* is equivalent to $R_z(\pi/2)$ up to an irrelevant global phase factor, and, therefore, the circuit can be expressed entirely in terms of IBM Q's native gate set: { R_z , S_X , CNOT, *X*}, where S_X is the square root of *X*.

D. Boundary condition and even-odd choice of the total qubit number

We note that Fig. 1 mainly depicts the ansatz for the open boundary condition (OBC). For the periodic boundary



FIG. 3. Simulation results of the ansatz. (a) the fidelity of the optimal ansatz state with the ground state of the open-chain Heisenberg model vs the total number of qubits N for one to six layers in the ansatz; (b) the corresponding relative error in the GS energy. (c) The fidelity and (d) relative energy error of the optimal ansatz state and the exact ground state of the periodic-boundary-condition Heisenberg model with the total number of qubits N for one to six layers in the ansatz.

condition (PBC), our ansatz (in the sublayer labeled as even gates) should include the two-qubit gates that act on the last qubit and the first qubit. From the numerical results presented in Fig. 3 (and to be discussed below), we do see that our ansatz gives comparable ground-state energy accuracy in both cases. For the ground-state fidelity for the PBC case seems to have values less than those in the case of the OBC. For practical purposes, it is easier on the real devices to find a open-ended path with good CNOT links. To find one periodic path with good CNOT links along the way will limit the size of the loop. In our cloud experiments, we focus on the case of the OPC.

Throughout this paper, we focus on the total qubit number N being even, as we argued earlier that the product of singlets (requiring N being even) can be adiabatically connected to the ground state of the XXZ model for finite even N. For odd N, the XXZ model has doubly degenerate ground states. To see this we consider the antiferromagnetic Heisenberg chain, for example. It was known that the ground state is a singlet [34], but this is valid for even N, because if one has an odd number of qubits (spin-1/2), the state with smallest spin magnitude one can construct is 1/2, which indicates double degeneracy.

However, we also checked that for odd N, the interpolated Hamiltonian has a finite gap throughout the bulk range of the interpolating parameter s, except at the end s = 1, where the the final Hamiltonian is degenerate. If the goal is to target one ground state, then our ansatz can be modified by changing the initial state to be (N - 1)/2 singlet pairs plus one remaining qubit arbitrarily initialized. The subsequent variational gate structure can be made the same, but we may need to allow them with site dependence. However, we have not tested this extensively.

III. ANALYSIS OF THE ANSATZ

We have explained how we arrive at a physics-motivated ansatz,

$$|\psi_{\text{ansatz}}(\{\theta\})\rangle = \mathbf{U}(\{\theta\})|0\dots0\rangle$$

= $\bigotimes_{l=1}^{N_L} \left[U_{\text{even}}^{(l)}(\{\theta_e\})U_{\text{odd}}^{(l)}(\{\theta_o\}) \right] U_{\text{init}}|0\dots0\rangle,$ (14)

schematically shown in Fig. 1(a), where we have used U_{init} to denote the initialization step that takes $|0 \dots 0\rangle \equiv |\uparrow \dots \uparrow\rangle$ to a product of singlets or Bell pairs $|\psi_{\text{singlets}}\rangle$ in Eq. (10).

Variational ansatzes and trial wave functions are commonly used in physics. Well-known examples include the BCS wave function for superconductivity [35] and the Laughlin wave function for the fractional quantum Hall effect [36]. Despite not exactly representing the GS, they capture the essential physical properties. To analyze how well our ansatzes simulate the GS wave functions and their energy, we minimize $E_{\text{ansatz}}(\{\theta\})$ to obtain the optimal parameters $\{\theta^*\}$ and then compute the GS fidelity [37] $f \equiv |\langle \psi_{gs} | \psi_{ansatz}(\{\theta^*\}) \rangle|$ and the error in the GS energy $\epsilon \equiv |E_{\text{ansatz}}(\{\theta^*\}) - E_{\text{gs}}|/|E_{\text{gs}}|$ [38], where $|\psi_{\rm gs}\rangle$ denotes the GS and $E_{\rm gs}$ its exact energy. In particular, we find that in the case of the open boundary condition, the one-layer ansatz can reach exact ground states of the N = 4 chain and the three-layer ansatz can reach exact ground states of the N = 6 chain. (See Appendix A for the exact ground state wave function of 4-qubit XXZ model.) In the case of the periodic boundary condition, the one-layer ansatz can reach exact ground states of the N = 4 chain and the two-layer ansatz can reach exact ground states of the N = 6chain. For larger chains, exact preparation of ground states is possible but requires more layers of the ansatz [39]. We note that even for N = 3 electrons, Laughlin's droplet state is not an exact wave function [36,40], but his wave functions are the key to understand the fractional quantum effect. That variational ansatzes contain the exact ground states is a desired feature, as it can ascertain optimality of variational parameters.

In Table I, we show the optimized parameters, energy, and overlap with MPS diagonalized ground-state wave function using one layer of our ansatz for the Heisenberg model. We check that the results agree with the exact computation for the qubit number $N \leq 12$. As expected, the fidelity decreases with the number of qubits, but much slower than exponentially. In contrast, the approximate ground-state energy seems to reach about 3% of error even for large chains using just one layer in the ansatz (e.g., 3.33% even for N = 100).

Using one to six layers in the ansatz, we compare the fidelity and energy error versus N in Figs. 3(a) and 3(b) for the open chain. For large systems, we use matrix product states (MPS) [41–44] for these calculations. The results improve substantially with increasing layers: with six layers and N = 50, the fidelity is above 0.9 and the accuracy in the GS energy is above 99.75%. For the periodic chain [see results in Figs. 3(c) and 3(d)], given our ansatz breaks the translation invariance (down to two sites), it may take a few layers to approximately restore the invariance. Overall, we do see general improvement in both quantities as the number of layers in the ansatz increases.

For the numerical optimization, we used both the quasi-Newton algorithm (with Matlab's fminunc function) and the sequential quadratic programming (SQP) algorithm (with Matlab's fminunc). The optimized parameters for small chains are used for the initial guess of larger chains. Our six-layer ansatz for the Heisenberg model contains only 12 variational parameters. For more layers, we expect that techniques used in machine learning, such as stochastic gradients, should be helpful.

IV. CLOUD EXPERIMENTS, rZNE, AND RESULTS

We have performed cloud experiments by creating the (one-layer) ansatz states and measuring their energies on nine different backends of IBM Q, which contain 27, 65, and 127 qubits on three types of layouts (see Fig. 4, Table IV, and Appendix B). The cloud experimental results of the 19 different sizes (ranging from 4 to 102 qubits) of Heisenberg chains are also summarized in Table I; their relative errors with the ground-state energy values are within a few percentages. We will explain below how the experimental results and their mitigated values were obtained. For the purpose of demonstration, we mostly use the numerically optimized parameters to run the state creation circuits. But still, we also test the feasibility of the hybrid quantum-classical approach by performing cloud experiments with random parameters below.

We first discuss different approaches to measure the total energy and then the readout-error and gate-error mitigation methods to extract estimated values from experiments. In particular, we will introduce a reference state in doing the gate-error mitigation.

A. Measuring energy: Several approaches

Here we describe three approaches that we have used to measure the energy expectation value. Ideally the three approaches should give the same results and we have indeed tested all three experimentally and verified that they give the same results within a few percentages of errors for small systems; see Appendix D.

(1) Tomography-based approach. For the models we consider, the Hamiltonian terms are of the form $\sigma_{\alpha}^{[j]}\sigma_{\alpha}^{[j+1]}$, where j is the site number and α is the spin direction (x, y or z). Naively, if we can obtain the reduced density $\rho_{i,i+1}$ matrix for the pair (j, j + 1) then we can calculate the energy contribution from $\text{Tr}(\rho_{j,j+1}\sigma_{\alpha}^{[j]}\sigma_{\alpha}^{[j+1]})$. But this requires state tomography and seems to need to run $9N_{\text{bond}}$ different circuits for the total energy, where N_{bond} is the number of nearest pairs or bonds, e.g., N - 1 for an open chain and N for a periodic one. However, we can improve the efficiency by performing the state tomography in parallel. Doing so, we just need two sets of state tomography circuits (for even and odd bonds respectively) to obtain the reduced density matrices of neighboring pairs of qubits. We will later discuss the measurement mitigation on pairs of qubits associated with bonds to extract reliable energy contribution. Doing the tomography in parallel reduces the number of circuits to measure to 9×2 , which is independent of the model size. The benefit of this is that any one- and two-qubit observables are readily available, such as the local spin observables and the concurrence which quantifies nearest-neighbor entanglement, and it applies to all nearest-neighbor interacting spins. (Extension to finiteranged interaction is straightforward but requires more sets of measurements and multiqubit state tomography.) For our cloud experimental results on measuring the concurrence in a 8-qubit XXZ chain; see Appendix D.

(2) *XYZ* measurement. The second, slightly reduced measurement method is to measure separately the two neighboring qubits on each bond in basis α and then average over the classical results $\overline{\sigma_{\alpha}^{[k]}\sigma_{\alpha}^{[k+1]}}$, treating $\sigma_{\alpha} = \pm 1$ from the



FIG. 4. Illustration of the layout of some backends used in this work: (a) the layout of 27-qubit machines, such as ibm_auckland, ibm_cairo, ibm_hanoi, ibm_kolkata, ibmq_montreal, ibmq_mumbai, and ibm_toronto; (b) the 65-qubit layout of ibmq_brooklyn; and (c) the 127-qubit layout of ibm_washington. An edge between two qubits indicates that a direct CNOT gate can be executed between them. See Table IV for certain properties of these backends.

measurement outcome assignment. This naively requires $3N_{\text{bond}}$ different circuits for the total energy. But a much simplified implementation is to measure all qubits in α basis and calculate the average $\overline{\sigma_{\alpha}^{[k]}\sigma_{\alpha}^{[k+1]}}$ for all bonds. This only requires three different measurement settings to obtain the total energy. Such simplification is applicable for the same reason mentioned in the previous approach (1).

(3) Bell measurement. There is another method that uses Bell-state measurement on all bonds. It exploits the specific form of $\hat{h}_{XXZ}^{[j,j+1]}(\Delta)$, for which $|\Psi^-\rangle \equiv (|01\rangle - |10\rangle)/\sqrt{2}$ has energy $-2 - \Delta$, the triplet $|\Psi^+\rangle \equiv (|01\rangle + |10\rangle)/\sqrt{2}$ has energy $2 - \Delta$, and both $|00\rangle$ and $|11\rangle$ (or equivalently $|\Phi^{\pm}\rangle \equiv (|00\rangle \pm |11\rangle)/\sqrt{2}$) have energy Δ . We can identify the Bell state on a particular bond with Bell measurement. The energy contribution of that bond is the energy corresponding to the Bell state obtained from the measurement outcome. The total energy can be calculated by adding up every bond's energy contribution. In practice, the Bell measurement requires a short circuit including a CNOT gate, and the effect of its error can be mitigated; see below.

TABLE IV. Properties of various IBM Q backends used in this work, data taken in late June, 2022. Our cloud experiments were performed during the span from March 2022 to June 2022. (Detailed information of these on specific dates when a job was run could be obtained in the Supplemental Material [45]). Q. volume is the quantum volume and N_{tot} is the total number of qubits in the backend. The basis gate set of these backends includes CX, ID, RZ, SX, and X, where CX denotes the CNOT gate, ID is the identity gate, RZ is the *z*-rotation gate, and SX is the square root of the Pauli X gate. The volume of the detailed device information, such as properties of individual qubits and individual gate and readout errors, is too large to list in the table here. Moreover, the device properties may have shifted over time. We have included such detailed device properties in the Supplemental Material [45].

Backend (Q. volume)	N _{tot}	Processor type	Average frequency	Average CNOT error	Average readout error	Average T ₁ time	Average T ₂ time
ibm_auckland (64)	27	Falcon r5.11	4.97 GHz	1.042×10^{-2}	$1.439 imes 10^{-2}$	178.38 μs	152.09 μs
ibmq_brooklyn(32)	65	Hummingbird r2	5.13 GHz	2.842×10^{-2}	2.928×10^{-2}	74.35 μs	77.66 μs
ibm_cairo(64)	27	Falcon r5.11	5.13 GHz	7.969×10^{-2}	1.352×10^{-2}	101.71 μs	132.51 μs
ibm_hanoi (64)	27	Falcon r5.11	5.00 GHz	4.444×10^{-2}	1.357×10^{-2}	151.26 μs	116.79 μs
ibmq_kolkata(128)	27	Falcon r5.11	5.10 GHz	4.801×10^{-2}	1.556×10^{-2}	118.67 μs	96.82 μs
ibmq_montreal (128)	27	Falcon r4	5.00 GHz	1.943×10^{-2}	3.426×10^{-2}	119.39 μs	102.78 μs
ibmq_mumbai(128)	27	Falcon r5.1	4.89 GHz	7.984×10^{-2}	2.665×10^{-2}	135.78 μs	117.56 μs
ibmq_toronto (32)	27	Falcon r4	5.09 GHz	8.680×10^{-2}	6.050×10^{-2}	115.84 μs	104.92 μs
<pre>ibm_washington (64)</pre>	127	Eagle r1	5.07 GHz	4.734×10^{-2}	2.789×10^{-2}	94.38 μs	90.82 µs

Naively, this approach of measuring energy requires N_{bond} different measurement patterns appended at the end of state creation as readout for the total energy if each bond is measured separately. However, we can divide the bonds into even and odd groups, as above, and can perform the Bell measurement in parallel within each group. Then we only need to perform *two* different sets of measurements. This turns out to be the approach we used to perform large-system (up to N = 102 qubits) cloud experiments on real devices to obtain the total energy.

We mainly use the Bell-state approach to measure the system's energy, as it requires the least resource compared to two other approaches.

B. Error mitigation

Here we describe the approaches we have used to implement error mitigation on of readout and gate errors. For additional information, see Appendix C, and for example codes, see the Supplemental Material [45].

1. Measurement/readout error mitigation

For superconducting qubits, the readout error can be as large as 10% or more and it is therefore crucial to mitigate the measurement error to calculate the correct energy of the created state on the real device. Due to the expanding deployment of cloud quantum computers, the interest in the issue of state preparation and readout error has recently been rekindled [46–51]. The key idea is to first characterize the measurement pattern dependent on the state input, such as from the detector tomography or simply measuring the probability matrix \mathcal{M} that relates the input states to the measured outcomes, i.e., $\vec{P}_{\text{measured}} = \mathcal{M}\vec{P}_{\text{ideal}}$, where $\vec{P}_{\text{measured}}$ and \vec{P}_{ideal} represent respectively the measured and ideal probability distribution. By properly inverting the relation with the constraint that the outcome distribution P_{ideal} be nonnegative, one can obtain the mitigated distribution to evaluate the observables.

For N qubits, the complete matrix \mathcal{M} is of size $2^N \times 2^N$ and requires preparation of 2^N computational states, thus is not efficient and is only doable for a small number of qubits. As the models we consider here contain only nearest-neighbor interactions, we are mainly concerned with measurement mitigation for pairs of qubits in a bond, i.e., involved in the interacting Hamiltonian, and such simplification allows us to deal with large systems in a practical way. We can perform readout mitigation pairwise for the nearest-neighbor two qubits on all bonds. Similar to the energy measurement, this can be reduced to two sets of mitigation, i.e., on pairs of even and odd bonds, respectively. Each mitigation requires four different inputs from all two-qubit computational basis and measurement in the same basis gives rise to a 4×4 matrix \mathcal{M} , which we can then use to infer the ideal two-qubit measurement distribution so as to obtain the mitigated energy contribution.

2. Bell-measurement mitigation

In our cloud experiments with large numbers of qubits, the local energy for a pair of qubits is obtained by measuring in the Bell-state basis, which uses an inverse circuit for Bell-state preparation and involves CNOT gates. To mitigate potential errors caused by imperfect CNOT gates, we adopt the above readout mitigation for the Bell measurement. Specifically, for each pair in the bonds, we prepare the four Bell states and then immediately measure qubits pairwise in the Bell-state basis, such as the circuit shown previously, to obtain a 4×4 Bell-state assignment matrix \mathcal{M}_{Bell} for each pair. With this we can mitigate the outcome distribution and hence the energy value obtained from the Bell-state measurement.

3. Gate-error mitigation

By doing readout mitigation we are probing the properties of the state actually created on the quantum devices. However, the observable expectation is affected by gate errors as well that prevent us from obtaining the idealized value. To estimate the latter, prior works have considered pulse and gate-error mitigation by extrapolating to the zero-error limit [7-9], and this is an extrapolation of the physical observables, not the actual observable values associated with the quantum states created. Nevertheless, it is still important to see how well quantum computers can estimate these values despite the noise and errors, especially in the regime where direct classical calculations might not be feasible.

However, to perform accurate gate mitigation, one needs to have substantial access to the hardware performing pulse-level optimization and operations [9], which is still not practical for dealing with a large number of qubits. (Note that recent experiments have been carried on 26 qubits using pulse-level zero-noise extrapolation [15].) Instead, we will use the gate-level zero-noise extrapolation (ZNE) approach discussed in Refs. [10–13]. In particular, our approach builds on the idea in Ref. [13] and we prepare the circuits to create $|\psi_n\rangle = \mathbf{U}(\mathbf{U}^{-1}\mathbf{U})^n |0...0\rangle$, where *n* is a nonnegative integer and $\mathbf{U} = \mathbf{U}_{\text{var}}(\{\theta\})U_{\text{init}}$, as in Eq. (14) of the main text, denotes the circuit to prepare the ansatz state from the fiducial state $|0...0\rangle$, i.e., $|\psi_{\text{ansatz}}\rangle = \mathbf{U}(\{\theta\})|0...0\rangle$, and then use several forward-backward repetitions in U to evaluate the observable $\mathcal{O}_n = \langle \psi_n | \hat{O} | \psi_n \rangle$, as a function of *n*. Ideally, different n should give the same state and hence the same value for the observable \hat{O} . However, noise and errors spoil this and the state with larger n should be noisier. The extrapolation to the gate-level zero-noise limit is done by a fitting to \mathcal{O}_n with $m = 2n + 1 \rightarrow 0$ limit.

4. Reference-state gate-error mitigation

Building on this, we propose to use a reference state (or possibly multiple ones), which is contained in the ansatz family, for example, the product of Bell pairs that we use below (via setting all θ 's to zero, i.e., $|\psi_{\text{singlets}}\rangle = \mathbf{U}(\{\theta\} = 0)|0...0\rangle$), with a known exact energy value, to improve the extrapolation of the energy value or other observables. Running the energy cloud experiment for this reference state with the above gate-level mitigation, we obtain the naively-extrapolated experimental value and hence the possible mismatch with the exact value. Using such knowledge for the reference state as a calibration, we can estimate the expected value of the ansatz state from the naive experimental value. Combining both gate and readout error mitigation, we are able to reach the accuracy of the extrapolated energy with

a few percentages of the exact value for all ranges of the qubit number that we have tested on real devices. We expect that this reference-state ZNE (rZNE) may be applied to the general VQE platform. It does not require additional circuits from randomized compiling, as done, e.g., in Ref. [12,14], but averaging the results from these randomized circuits can be used to further improve the accuracy.

C. Cloud experimental results: Reference-state zero-noise extrapolation applied

From the 102-qubit Heisenberg-chain experimental data in Figs. 5(a) and 5(b), we fit the total energy of the optimal ansatz state and Bell pairs to a form $f_E(m) = a \exp(-bm) + c$ and obtain their respective ZNE values -199.2 and -169.8. The energy of Bell pairs with *N* qubits is known exactly, $-(2 + \Delta)N/2$, which is -153 when N = 102 and $\Delta = 1$. The two values for the Bell pairs enable us to naively correct the ansatz state energy from -199.2 to $-199.2/(169.8/153) \approx -179$, close to the numerical MPS value -174.04.

The noise in real devices is very complex, but a simplified model on neighboring two qubits is a depolarizing channel: $\rho_2 \rightarrow (1 - p_m)\rho_2 + p_m I \otimes I/4$. One naively expects that $p_m = 1 - e^{-bm}$ and this implies that $f_E(m) = Ee^{-bm}$. (In Appendix E, we provide further results of the decay constant b.) But in our fitting above, we observe a nonzero residual value c. This means that we should rescale only the drop (i.e., a) from m = 0 to $m \to \infty$. To be more precise, the rescale factor r is obtained via $a_B \cdot r + c_B = E_{\text{bell}}$, where E_{bell} is the exact Bell pairs energy and the subscript B in a and c denotes the parameters obtained from fitting the cloud experiments for Bell pairs. Assume the cloud experiments for the optimal ansatz experience similar noise and errors, as their circuit structure and depth are identical (except the rotation parameters), we obtain the extrapolated experimental ansatz energy to be $E_{exp} = a_E \times r + c_E$, where the subscript E denotes the parameters obtained from fitting the cloud experiments for the ansatz. Applying this to the 102-qubit cloud experiment, we obtain almost the same result (up to rounding): $E_{exp} =$ -179.1 ± 3.1 . This is how all the other reported data were obtained. From our experience, the results obtained this way do not differ much from the naive rescaling in most of our cloud experiments.

We note that in this set of cloud experiments, the maximal CNOT depth is 63 and the maxmial total number of CNOTs used is 3186. Similar experimental results on 50 qubits in ibmq_brooklyn are shown in Figs. 5(c) and 5(d). We also note that, in Ref. [12], $|00...0\rangle$ was used as a reference state in a circuit that is used to extract the depolarizing rate. Additional randomized instances (e.g., 448 copies) of each main circuit were needed for averaging [12], in addition to tripling and quintupling all CNOT gates for ZNE [11].

1. Cloud experimental results for the Heisenberg model on various sizes

Table I has a summary of cloud experimental results of Heisenberg chains with 19 different sizes (ranging from 4 to 102 qubits) and some were averaged over several different sets of qubits or different devices. We refer to Tables II and III for a comprehensive list of 39 mitigated results on nine different backends. These backends possess different qubit numbers, quantum volumes, and noise and error rates (see Table IV), but the success across all these backends (with varying numbers of qubits used) demonstrates the utility of such a simple and scalable rZNE approach. With these results, we can, for example, extract the energy per site in the thermodynamic limit (see also Fig. 6), which yields a value of -1.713 ± 0.046 that agrees with the exact Bethe ansatz calcuation, $4(\ln 2 - 1) \approx -1.773$ [27], within 3.4% of deviation.

2. Results using random parameters

To illustrate a proof-of-principle demonstration of the potential hybrid quantum-classical approach, we have also performed additional cloud experiments on 40 qubits with random parameters in the ansatz, and our rZNE method gives energy values agreeing well with the numerically calculated values; see Figs. 5(e) and 5(f). This demonstrates that it is feasible to use quantum devices to extract mitigated expectation values accurately, and, based on these, estimate the next iteration of the variational parameters by classical computers. Hence, there is no need to know the optimal variational parameters in advance, and the rZNE-mitigated variational algorithm can potentially become practical for large-scale NISQ devices.

3. Results for the XXZ model

As constructed, our approach works equally well for the *XXZ* model, and in Fig. 7, we present the two sets of cloud experimental results for a wide range of $\Delta \in [-0.8, 1.4]$ with N = 40 and N = 80 spins, respectively, carried out on two separate backends, ibmq_brooklyn and ibm_washington. The mitigated values agree well with the anticipated ansatz values. (The results for additional experiments with N = 8 chain are shown in Appendix D and fall within several percentages of the numerical ground-state energy values.)

V. BEYOND ONE DIMENSION

We expect that our protocol can be generalized to twodimensional structures. As a concrete example beyond the strict 1D, we consider a two-leg ladder of the XXZ model (see Fig. 8). We first prepare all vertical pairs of qubits in singlets, and apply the XXZ anastz gates to all horizontal odd bonds, even bonds, and then all vertical bonds. This constitutes a onelayer ansatz and can be repeated for multiple layers. We have performed numerically simulations and found that for N = 6(total number of spins in the ladder), the two-layer ansatz can achieve the exact ground state for $\Delta > -1$ up to machine precision. It is likely that for larger N, exact ground states can be achieved by using more layers. Our ansatz achieves very high GS fidelity, exceeding 0.95 even for N = 50 with just three layers, as well as high acurracy in the GS energy; see Fig. 9. There is an interesting phase diagram from this two-leg model [52,53], including a Haldane phase, which could potentially be implemented on a digital quantum processor.



FIG. 5. Zero-noise gate-error extrapolation for cloud experimental realizations of optimal ansatz states of Heisenberg spin chains, along with the use of Bell pairs. The data points presented were already processed by measurement-error migitation. (a), (b) 102-qubit Heisenberg chain on ibm_washington for the optimal ansatz state and Bell pairs, respectively; (c), (d), similarly for a 50-qubit Heisenberg chain on ibmq_brooklyn; (e), (f) results for two 40-qubit random ansatz states on ibmq_brooklyn. In panel (e), the parameters [θ_{even} , θ_{odd}] = [3.5, 0.7] were used and the exact ansatz energy is -16.0669. In panel (f), parameters [0.3,1.7] were used and the exact ansatz energy is -48.0625. Separate cloud experiments (results not shown in plots) with 40-qubit Bell pairs gives a naive extrapolation of the Bell pairs energy to be -67.0(4.0), whose ideal value is -60. The migtigated values with the reference state for panels (e) and (f) are -15.4 \pm 0.7 and -46.1 ± 2.4, respectively.



FIG. 6. Approximated ground-state energy per site vs the total number N in the spin chain obtained from cloud experiments performed on various IBM Q backends using the one-layer variational ansatz. The dashed line is a fit from the data: -1.713 + 0.393/N and the approximated ground-state energy at the $N \rightarrow \infty$ limit is -1.713 ± 0.046 , which is compared to the exact result from the Bethe ansatz solution $4(\ln 2 - 1) \approx -1.773$.

VI. SUMMARY

In this work, we have analyzed physics-motivated ansatzes for the Heisenberg and the XXZ spin chains and showed their potential use to simulate the ground-state wave function and extract the ground-state energy. The accuracy improves substantially by increasing the ansatz layers. We have also developed efficient and scalable methods for measuring the energy and mitigated the experimentally measured values. By carrying out the cloud realization and measurement of the ansatz states, we have demonstrated that short-depth variational circuits could be applied to large systems of qubits, with up to 102 qubits performed on real devices. Despite the presence of substantial noise and errors in current devices, we have been able to improve and implement efficient error mitigation schemes to deduce accurate ground-state energy from experiments on large systems. For example, from the data obtained from our cloud experiments, we were able to exact the estimated ground-state energy density value that agrees with the exact Bethe ansatz solution to within 4% of error. We note that for our Heisenberg and XXZ models, the reference state is chosen to be a product of singlet pairs. For other models, the choice of the reference states may be important.

Our work thus opens up the potential practical use of error mitigated VQE on large quantum computing backends for improved accuracy. One first applies our rZNE (combining readout mitigation and possibly further mitigation) to obtain the extracted observable value(s) and/or its gradients from quantum devices, then uses classical computers to search for variational parameters to be used in the subsequent iteration of experiments with mitigation. The procedure is then iterated until the mitigated observable value(s) converge to within certain accuracy. Such an error-mitigated, rZNE VQE approach, though not yet practical for large systems in the current cloud-based setting, due to limited allocated time and long job queues, seems plausible in dedicated experiments. Our cloud experiments using randomly chosen parameters already demonstrated agreement with the expected ansatz energy. To enter a regime where quantum advantage may be realized, we will likely need to go beyond one dimension, e.g., two dimensions, where classical simulations of quantum manybody systems become intractable as the system size increases. Toward this goal, we have also analyzed a two-leg ladder and showed the applicability of our ansatz.

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APPENDIX A: EXACT 4-QUBIT GROUND STATE FOR THE OPEN CHAIN

For four-qubit XXZ model with the open-boundary condition, we could assume the ground state to be of the form

$$\begin{aligned} |\psi_0\rangle &= a \frac{1}{\sqrt{2}} (|0101\rangle + |1010\rangle) + b \frac{1}{\sqrt{2}} (|0011\rangle + |1100\rangle) \\ &+ c \frac{1}{\sqrt{2}} (|1001\rangle + |0110\rangle), \end{aligned}$$
(A1)

and evaluate the energy, giving

$$\langle \psi_0 | H_{XXZ} | \psi_0 \rangle = 4a(b+2c) + (b^2 - c^2 - 3a^2)\Delta.$$
 (A2)

To minimize the energy with the constraint that $a^2 + b^2 + c^2 = 1$, we can introduce a Lagrange multiplier to enforce the constraint to the optimization of a quadratic function of *a*, *b* and *c*. It can be formulated as solving an eigenvalue problem for a 3×3 matrix, and we find that the exact ground-state energy $E_0(\Delta)$ can be obtained from the lowest real root of a third-order polynomial,

$$(12\Delta - 3\Delta^3) - (20 + \Delta^2)x + 3\Delta x^2 + x^3 = 0.$$
 (A3)

The coefficients (a, b, c) in the ground-state wave function can be obtained as follows:

$$\tilde{a}(\Delta) = [-\Delta + E_0(\Delta)]/2, \tag{A4}$$

$$\tilde{b}(\Delta) = 1, \tag{A5}$$

$$\tilde{c}(\Delta) = [-4 - 3\Delta^2 + 2\Delta E_0(\Delta) + E_0(\Delta)^2]/8,$$
 (A6)

with $(a, b, c) = (\tilde{a}, \tilde{b}, \tilde{c})/\sqrt{\tilde{a}^2 + \tilde{b}^2 + \tilde{c}^2}$.



FIG. 7. The mitigated energy results for XXZ spin chains. (a) A 40-qubit XXZ chain on ibmq_brooklyn and (b) an 80-qubit XXZ chain on ibm_washington. The information about which physical qubits were used is listed in the Appendix.

One can also write down the equation for the parameters in the one-layer ansatz to arrive at the above four-qubit state (A1), up to a global phase, and in principle solve for the optimal parameters for the ground state. However, we could not find a closed form expression for the parameters.

For the periodic boundary condition, one can use the anastz,

$$\begin{split} |\psi_0\rangle = a \frac{1}{\sqrt{2}} (|0101\rangle + |1010\rangle) + b \frac{1}{2} (|0011\rangle + |0110\rangle \\ + |1100\rangle + |1001\rangle), \end{split} \tag{A7}$$



FIG. 8. The anastz for the two-leg ladder. (a) We initialize the state in a product of singlets which are formed between the upper spins and the lower spins. One-layer ansatz includes (b)–(d), where the gates are indicated by shaded rectangles.

and the problem reduces to solving a 2×2 matrix, which was previously presented in the Supplemental Material of Ref. [54], and the analytic expression for the four-qubit ground state is available there. Given that we know from numerics that the one-layer ansatz contains the exact groundstate wave function, we can similarly solve for the optimal parameters by identifying the ansatz and exact wave functions.

APPENDIX B: PROPERTIES OF QUANTUM BACKENDS AND THE CHOICE OF QUBITS

The average properties of the nine quantum backends of IBM are listed in Table IV and there are three different layouts, as illustrated in Fig. 4. Seven of the backends have 27 qubits, the backend ibmq_brooklyn has 65 qubits, and ibm_washington has 127 qubits, with the last also shown in Fig. 1(b). As it is not possible to list all the detailed information in this Appendix, we have downloaded complete device properties for individual qubits, as well gate errors and readout errors in the Supplemental Material [45]. Before cloud experiments were performed, we examined the detailed error rates reported on the service website and chose a path with a desired total number of sites along those connected qubits so as to avoid CNOT links with high error rates. For large system sizes, it is inevitable that we encounter a few CNOT links that may have somewhat higher error rates than others. We note that the detailed noise and error rates may drift over time as the devices are regularly calibrated and this impact large paths more than small ones. For example, to perform the 80-qubit XXZ model cloud experiments in Fig. 7(b), we had to use a different path from the one used previously for the Heisenberg model (reported in Table II) to avoid certain CNOT links with large error rates.

APPENDIX C: DATA ANALYSIS

1. Measurement mitigation

In the main text, we have described details of measurement/readout mitigation, including the standard one and our Bell-state measurement mitigation. In the Supplemental Material [45], we have provided a Python



FIG. 9. The fidelity (a) and relative energy error (b) of the optimal ansatz state and the exact ground state of the two-leg Heisenberg model with the total number of qubits N using one to three layers in the ansatz. The bond dimension we used in the MPS is $\chi = 64$.

notebook that implements the Bell-measurement mitigation, as well as an exemplary data output from an actual cloud experiment. With sufficient repetitions of energy measurements (see Tables II and III for the numbers, ranging from 25 to 146 on different selections of qubits and backends), we can then estimate the average and uncertainty of the energy.

2. Zero-noise extrapolation

In fitting the energy data (e.g., after measurement mitigation is applied), as described in the main text, we use an exponential function

$$f_E(m) = a \exp(-bm) + c, \qquad (C1)$$

where m = 2n + 1 is the total number of **U** or **U**⁻¹ in the circuit to prepare the ansatz state. For most experiments we used *n* from 0 to 4; for some we had *n* from 0 to 5. In all the experiments we did, the uncertainty from extrapolations is significantly larger than the statistical uncertainty from the energy measurements. The error bars from the extrapolated values are estimated from the extreme fit parameters values within the 68.27% confidence interval, corresponding to one sigma standard deviation. In the Supplemental Material [45], we have provided a Mathematica notebook that illustrates the analysis.

3. Reference-state gate-error mitigation

After we get the fitted energy curves for both experiment energy $f_e(m)$ and reference-state energy $f_r(m)$

$$f_e(m) = a_e \exp(-b_e m) + c_e, \tag{C2}$$

$$f_r(m) = a_r \exp(-b_r m) + c_r, \tag{C3}$$

we use the theoretically known value of the reference-state energy E_{ref} to calibrate the "experimental" energy E_{exp} ,

$$E_{\rm exp} = a_e \frac{E_{\rm ref} - c_r}{a_r} + c_e, \qquad (C4)$$

where the uncertainty of the E_{exp} comes from combining the uncertainties of all the fit parameters in the equation. In the same Mathematica notebook, we also implement this analysis for two different sets of data.

APPENDIX D: ADDITIONAL CLOUD EXPERIMENTAL RESULTS

1. Results on Heisenberg chains

The nine IBM Q backends we use have three different layouts, as illustrated in Fig. 4. The complete list of the results from the cloud experiments for the Heisenberg model on various backends and with various number of qubits is shown in Tables II and III. These were carried out using the Bell-measurement approach.

2. Results on XXZ chains

We have also performed cloud experiments for 8-qubit *XXZ* model on ibmq_montreal, with Δ ranging from -0.8 to 1.4, and we use two different measurement methods to calculate the energy, as shown in the Fig. 10. The two methods of the *XYZ* measurement and of the Bell measurement agree with each other. In addition, we have also used quantum state tomography to measure the total energy at two different values of Δ (0 and 1); the energy values obtained from tomography also agree with the other two approaches. In particular, the energy results from the state tomography give -13.46 ± 0.31 at $\Delta = 1$ and -9.3 ± 0.8 at $\Delta = 0$.

3. Concurrence results

With the tomography approach, we have obtained additionally the concurrence for all the bonds, and the cloud experimental results are compared to those of the ansatzes and the exact solution in Fig. 11. Due to the open boundary condition, the concurrence alternates from large to small between odd and even bonds. The entanglement on all even bonds is identically zero for the one-layer ansatz. This is due to the initial state being product of singlet pairs on odd bonds



FIG. 10. The energy results for an 8-qubit XXZ chain on ibmq_montreal. (a) The energy is obtained using the Bell-measurement approach on physical qubits [15,12,13,14,16,19,22,25]; (b) The energy is obtained using the XYZ measurement approach on physical qubits [11,14,16,19,22,25,24,23]. We have also performed energy measurement using quantum state tomography for $\Delta = 0$ and 1. The three methods for measuring energy agree very well in our cloud experiments.

and the one-layer entangling operation on even bonds is not strong enough to make the pairs on even bonds entangled. For odd bonds, the concurrence values inferred from the clound experiments are {0.890372, 0.767076, 0.683096, 0.768255} at $\Delta = 1$ and {0.850059, 0.663988, 0.648982, 0.812279} at $\Delta = 0$. As the quantum phase transition at the Heisenberg point $\Delta = 1$ is infinite-order, the concurrence does not exhibit singularity across the transition, so we did not perform cloud experiments for the concurrence over a wide of Δ , but only for $\Delta = 0$ &1 as an illustration. These concurrence values were obtained by use our rZNE approach with the naive extrapolation using Bell pairs as the reference. In doing ZNE, we had to repeat (UU⁻¹) several times, but the resulting reduced density



4. Information for Fig. 7 in the main text

(a) A 40-qubit *XXZ* chain on ibmq_brooklyn with physical qubits being [38,41,42,43,52,56,57,58,59,60,53,47, 46,45,39,31, 30,29,24,15,16,17,11,4,5,6,7,8,12,21,20,19,25, 33,34,35,40, 49,50,51]. (b) An 80-qubit *XXZ* chain on ibm_washington, with physical qubits being [97,96,95, 94,90,75,76,77,71,58,57,56,52,37,38,39,33,20,21,22,23,24,25, 26,27,28,29,30,31,32,36,51,50,49,48,47,46,45,44,43,42,41,53, 60,61,62,63,64,65,66,67,68,69,70,74,89,88,87,93,106,107,108,





FIG. 11. Ground-state entanglement property—concurrence—for a chain of 8-spin XXZ model with the open-boundary condition: (a) $\Delta = 1$ and (b) $\Delta = 0$. The concurrence is calculated for two neighboring qubits (j, j + 1) on *j*th bond $(j \in [1, 7])$ using quantum states obtained from optimizing one-layer, two-layer, and three-layer ansatzes, as well as from exact diagonalization of the XXZ Hamiltonian and from the cloud experiment done on *ibmq_montreal* with the one-layer ansatz. Note that with just one layer, the concurrence on the even bonds is zero. From these we observe that the entanglement is decreasing from the one-layer optimal ansatz to two- and to the three-layer one, toward the exact solution. The reason is that the initial state of the ansatz is a product of singlet Bell states on odd bonds, which possesses a very high global entanglement. The gates on even bonds act to decrease the entanglement of Bell states (on odd bonds) to increase the entanglement on even bonds.



FIG. 12. The decay coefficient *b* from extracting the ansatz energy (b_E) and from the Bell pairs energy (b_B) vs the anisotropy parameter Δ in the 80-qubit *XXZ* model, performed in ibm_washington. The QOD is related to the decay rate *b* as QOD = 7/b.

112,126,125,124,123,122,111,104,103,102,101,100,110,118, 117,116,115,114].

APPENDIX E: QUANTUM OBSERVABLE DEPTH

In fitting the energy data, we use an exponential function $f_E(m) = a \exp(-bm) + c$, where m = 2n + 1 is the total number of **U** or \mathbf{U}^{-1} in the circuit to construct the state. We note that each **U** contains seven layers of CNOT gates. When such an exponential-decay fitting works, the quantity 7/b, roughly speaking, represents the decay depth in the quantum circuit for the total energy, which we will refer to as the quan-

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