

Kaleidoscope of quantum phases in a long-range interacting spin-1 chain

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(Received 13 October 2015; revised manuscript received 3 February 2016; published 11 May 2016)

Motivated directly by recent trapped-ion quantum simulation experiments, we carry out a comprehensive study of the phase diagram of a spin-1 chain with XXZ-type interactions that decay as $1/r^\alpha$, using a combination of finite and infinite-size DMRG calculations, spin-wave analysis, and field theory. In the absence of long-range interactions, varying the spin-coupling anisotropy leads to four distinct and well-studied phases: a ferromagnetic Ising phase, a disordered XY phase, a topological Haldane phase, and an antiferromagnetic Ising phase. If long-range interactions are antiferromagnetic and thus frustrated, we find primarily a quantitative change of the phase boundaries. On the other hand, ferromagnetic (nonfrustrated) long-range interactions qualitatively impact the entire phase diagram. Importantly, for $\alpha \lesssim 3$, long-range interactions destroy the Haldane phase, break the conformal symmetry of the XY phase, give rise to a new phase that spontaneously breaks a $U(1)$ continuous symmetry, and introduce a possibly exotic tricritical point with no direct parallel in short-range interacting spin chains. Importantly, we show that the main signatures of all five phases found could be observed experimentally in the near future.

DOI: [10.1103/PhysRevB.93.205115](https://doi.org/10.1103/PhysRevB.93.205115)

The study of quantum phase transitions in low-dimensional spin systems has been a major theme in condensed matter physics for many years [1]. A well-known implication of Mermin and Wagner's famous results [2] on finite temperature quantum systems is that, for a large class of one-dimensional quantum spin systems, long-range order is forbidden even at zero temperature. This absence of classical order promotes quantum fluctuations to a central role, and they often determine the qualitative properties of the quantum ground state. An important example, first conjectured by Haldane [3,4], is that a spin-1 antiferromagnetic Heisenberg chain possesses a disordered phase with an energy gap to bulk excitations, later identified as a symmetry protected topological phase [5,6]. Its spin-1/2 counterpart, despite possessing the same classical limit, has a disordered ground state with gapless excitations, and is described by a conformal field theory (CFT) [7].

Experimentally, such quantum phase transitions have been explored in quasi-1D materials, and more recently using artificial materials designed through the careful control of atomic, molecular, and optical (AMO) systems [8–11]. These AMO systems are usually well isolated from the environment, offer considerable tunability of system parameters, and make possible both measurement and control at the individual lattice-site level. A distinctive feature of AMO systems is that interactions between particles are often long ranged, decaying as a power law with distance ($1/r^\alpha$). The exponent α varies widely amongst different AMO systems, ranging from $\alpha = 6$ for van de Waals interactions in Rydberg atoms, to $\alpha = 3$ for polar molecules and magnetic atoms, to $\alpha = 0$ for atoms coupled to cavities [11–19]. The effect of long-range interactions can be tuned by either changing the dimensionality of the system, e.g., for neutral atoms or molecules in optical lattices, or by directly (and often continuously) altering the

value of α , e.g., in trapped ions or cold atoms coupled to photonic crystals [14]. The availability of tunable long-range interactions creates an entirely new degree of freedom—absent in typical condensed-matter systems—for inducing quantum phase transitions, and can potentially lead to novel quantum phases [20–23].

While long-range interacting classical models have been studied in considerable detail for some time [24–28], there is a relative lack of in-depth studies of quantum phase transitions in long-range interacting systems, despite the emerging experimental prospects for studying both their equilibrium and nonequilibrium properties [15–18,29–35]. One reason is that many analytically solvable lattice models become intractable when interactions are no longer short ranged, a well-known example being the spin-1/2 XXZ model. Thus exact analytical studies are either restricted to noninteracting bosonic and fermionic systems with long-range hopping and pairing [33,35–37], or to certain contrived long-range interacting spin models which are difficult to realize in real systems [38–41]. In addition, to properly incorporate long-range interactions in low-energy effective theories, existing field theoretic treatments need to be modified and usually become more complicated [42,43]. While spin-wave theories can be useful in treating long-range interactions [44,45], they are unable to distinguish major differences in quantum phases between integer and half-integer spin chains. Exact numerical studies for long-range interacting spin models are restricted to small system sizes and usually inconclusive [46–49], since the correlation length is generally divergent [32,50]. Approximate numerical techniques such as the density matrix renormalization group (DMRG) method have been adapted to treat long-range interactions [51], but determining complete diagrams with large-system-size calculations remains challenging, and those that exist are primarily for spin-1/2 chains [20,29,52,53].

In this paper, we carry out a comprehensive study of a spin-1 chain with tunable XXZ interactions that decay

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monotonically as $1/r^\alpha$, for all $\alpha > 0$. Our study is largely motivated by imminent trapped-ion-based experiments that can simulate this model with widely tunable index α [54–56]. In the absence of long-range interactions, the choice of spin-1 over spin-1/2 allows us to have four distinct quantum phases by varying the anisotropy of the interactions: a ferromagnetic (FM) phase and an antiferromagnetic (AFM) Ising phase that are both gapped and long-range ordered, a disordered gapless phase (the XY phase), and a gapped and topologically ordered phase (the Haldane phase). By using a combination of DMRG calculations, spin wave analysis, and field theory, we obtain the phase diagram for arbitrary anisotropy and all $\alpha > 0$, with both ferromagnetic and antiferromagnetic interactions. Our key observation is that when interactions in all spatial directions are antiferromagnetic, long-range interactions are frustrated, leading to primarily quantitative changes to the phase boundaries compared to the short-range interacting chain. Interestingly, we find that the topological Haldane phase is robust under long-range interactions with any $\alpha > 0$ [48,49,57]. However, when the interactions in the x - y plane become ferromagnetic, we find a number of significant modifications to the phase diagram: (1) The Haldane phase is destroyed at a finite α due to a closing of the bulk excitation gap; (2) the gapless XY phase, described by a CFT with central charge $c = 1$, disappears when $\alpha \lesssim 3$ due to a breakdown of conformal symmetry [33,35]; (3) the disappearance of the XY phase heralds the emergence of a new phase at $\alpha \lesssim 3$ (continuous-symmetry breaking, or CSB) in which the spins order in the xy plane, spontaneously breaking a $U(1)$ symmetry and possessing gapless excitations (Nambu-Goldstone modes); (4) Novel tricritical points, with no direct analog in short-range interacting 1D models, appear at the intersection of the Haldane, CSB, and XY/AFM phases.

The paper is organized as follows. In Sec. I, we introduce the model Hamiltonian and present complete phase diagrams for the ferromagnetic and antiferromagnetic cases. In Sec. II, we study the boundary of the FM phase, where a spin-wave approximation is found to be asymptotically exact in the large-system limit. In Sec. III, we determine both the XY-to-Haldane and Haldane-to-AFM transition lines accurately using DMRG calculations and use field theory arguments to explain the effect of long-range interactions on the boundary of the Haldane phase. In Sec. IV, we introduce the CSB phase and explain its emergence using spin-wave theory. The boundary between the CSB and XY phases is determined by a numerical calculation of central charge. In Sec. V, we show that all five phases possess distinct signatures that could be observed in near-future trapped ion quantum simulations with chains of 16 spins. Finally, we conclude the paper in Sec. VI and comment on a number of open questions.

I. MODEL HAMILTONIAN AND PHASE DIAGRAMS

We consider the following spin-1 Hamiltonian with long-range XXZ interactions in a 1D open-boundary chain:

$$H = \sum_{i>j} \frac{1}{(i-j)^\alpha} [J_{xy}(S_i^x S_j^x + S_i^y S_j^y) + J_z S_i^z S_j^z]. \quad (1)$$

Here $J_z \in (-\infty, \infty)$ and $\alpha \in (0, \infty)$ are allowed to vary continuously, and we consider both the $J_{xy} = 1$ (antiferromagnetic) and $J_{xy} = -1$ (ferromagnetic) cases. We note that, for $0 < \alpha < 1$, Eq. (1) does not have a well-defined thermodynamic limit when J_{xy} and/or J_z is ferromagnetic, since the ground-state energy density diverges. To make the ground-state energy extensive, we may impose an energy renormalization factor $N^{\alpha-1}$, first introduced by Kac [58], when taking the thermodynamic or continuum limit (here N is the chain length). For finite-size numerical calculations, we do not need to implement the Kac renormalization for $0 < \alpha < 1$ since ground-state properties are unaffected by energy renormalization [59].

Figure 1 shows our full phase diagram for both $J_{xy} = 1$ and $J_{xy} = -1$, with actual phase boundaries plotted using

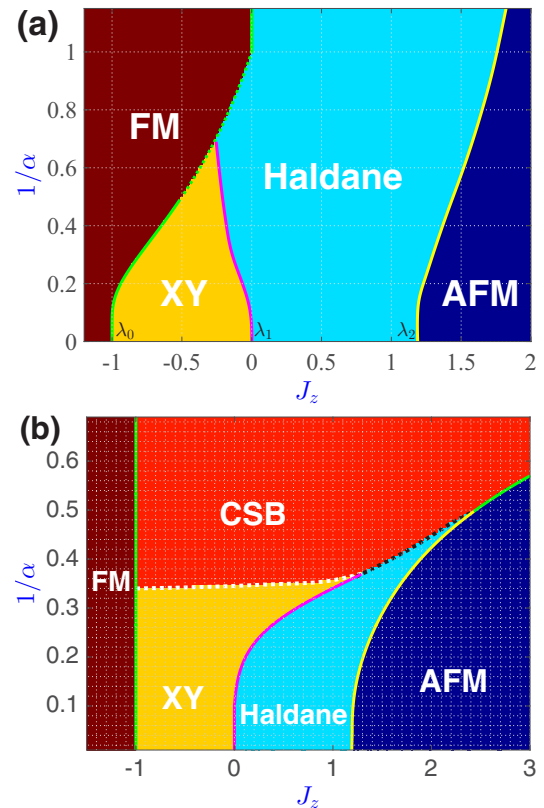


FIG. 1. Proposed phase diagram for (a) $J_{xy} = 1$ and (b) $J_{xy} = -1$. Five different phases are identified: a ferromagnetic (FM) Ising phase, an antiferromagnetic (AFM) Ising phase, a disordered XY phase, a topological Haldane phase, and a continuous symmetry breaking (CSB) phase. At $\alpha = \infty$, the transition points are denoted by $J_z = \lambda_{0,1,2}$ in (a). The FM-to-XY, FM-to-CSB, and CSB-to-AFM transitions are first order (green line); the XY-to-Haldane transition is BKT type with central charge $c = 1$ (purple line); the Haldane-to-AFM transition is second order with $c = 0.5$ (yellow line); the CSB-to-XY transition (white dotted line) has $c = 1$, but is a BKT-like transition corresponding to a universality class different from the XY-to-Haldane transition [83]; the CSB-to-Haldane transition (black dotted lines) appears to be an exotic continuous phase transition not described by a 1+1D CFT. The location of solid transition lines are expected to be accurate in the thermodynamic limit, while the location of dotted transition lines may have a small uncertainty.

the results of calculations discussed in the following sections. The nearest-neighbor interaction limit is achieved at $\alpha \rightarrow \infty$ ($1/\alpha = 0$). In this limit, the Hamiltonian in Eq. (1) with $J_{xy} = -1$ is equivalent to the one with $J_{xy} = 1$, as can be seen by performing a local unitary transformation that flips every other spin in the x - y plane while preserving the spin commutation relations: $S_i^{x,y} \rightarrow (-1)^i S_i^{x,y}$. The different ground-state phases of this short-range Hamiltonian have been well studied [60–62]. Notably, Haldane first conjectured [3,4] that for $\lambda_1 < J_z < \lambda_2$, a disordered gapped phase (the Haldane phase) will emerge. At $J_z = \lambda_2$, the ground state undergoes a second-order phase transition from the Haldane phase to an AFM phase, which belongs to the same universality class as the 2D Ising model. The value $\lambda_2 \approx 1.186$ has been found by various numerical techniques including Monte-Carlo [63], exact diagonalization [64], and DMRG [65–67]. At $J_z = \lambda_1$, a Berezinskii-Kosterlitz-Thouless (BKT) transition intervenes between the Haldane phase and a gapless disordered XY phase at $J_z < \lambda_1$. The value of λ_1 is theoretically predicted to be exactly zero after mapping Eq. (1) (for $\alpha = \infty$) to a field theory model using bosonization [68]. This prediction is supported by conformal field theory arguments [69] and a level spectroscopy method based on a renormalization group analysis and the $SU(2)/Z_2$ symmetry of the BKT transition [61,70–72]. Numerically, $\lambda_1 \approx 0$ has been verified via finite-size scaling [64,73,74] and DMRG [65]. Finally, at $J_z = \lambda_0 = -1$, a first-order phase transition from the XY phase to a ferromagnetic Ising phase takes place [61,66,75].

We now introduce our results for the long-range interacting case ($1/\alpha > 0$). For $J_{xy} = 1$ and $J_z > 0$, long-range interactions are frustrated and the Haldane-to-AFM phase transition point $\lambda_2(\alpha)$ increases moderately as α decreases, without changing the universality class of the transition. For sufficiently small $J_z < 0$, the ferromagnetic long-range interactions along the z direction eventually favor a ferromagnetic ground state, inducing a first-order transition at $\lambda_0(\alpha)$. The magnitude of the critical coupling, $|\lambda_0(\alpha)|$, decreases monotonically from 1 (at $\alpha = \infty$) to 0 (for all $\alpha \leq 1$) in the thermodynamic limit. The XY-to-Haldane phase boundary $\lambda_1(\alpha)$ becomes negative for finite α , eventually terminating in a tricritical point at the intersection of FM, Haldane, and XY phases. The entire XY phase (including the XY-to-Haldane phase boundary) has conformal symmetry with $c = 1$, and the XY-to-Haldane phase boundary remains a BKT transition until it terminates at the tricritical point. We note that the phase diagram [Fig. 1(a)] is similar to Fig. 1 in Ref. [76], which studies the XXZ spin-1 chain with next-nearest-neighbor interactions of tunable strength. This is partially because the frustrated long-range interactions in the x - y plane effectively cancel each other at different ranges, so their influences on the phase boundary are somewhat similar to those from next-nearest-neighbor interactions. However, we point out that the full $1/r^\alpha$ long-range interactions, frustrated or not, will result in power-law decaying correlation functions in the gapped phases (see Ref. [57] for details); such correlations are absent in models with next-nearest-neighbor interactions [76–79].

For $J_{xy} = -1$, where long-range interactions in the x - y plane are not frustrated, the phase diagram [Fig. 1(b)] shows a number of important qualitative differences from the nearest-neighbor phase diagram as α is decreased. First, the XY-to-

Haldane phase boundary bends significantly toward positive J_z , and we find the Haldane phase to terminate at $\alpha \approx 3$ for $J_z = 1$. Second, we expect the XY phase to disappear for $\alpha \lesssim 3$ due to the breakdown of conformal symmetry [33,35]. Third, for $\alpha \lesssim 3$ a new CSB phase emerges—this is not in violation of the Mermin-Wagner theorem, as it no longer applies for this range of interactions [2,52,80–83]. The CSB-to-AFM phase transition is expected to be first-order, since at large J_z and small α , quantum fluctuations play negligible roles for both the Néel-ordered state and the ordered CSB state. This behavior is similar to the transition between the AFM phase and the large- D phase (where a large positive anisotropy term $D \sum_i (S_i^z)^2$ causes all spins to stay in the $|S_i^z = 0\rangle$ state) reported in Refs. [65,66,75]. The Haldane phase has a $c = 1$ critical phase boundary with the XY phase, a $c = 0.5$ phase boundary with the AFM phase [67], and a possibly exotic phase boundary with the CSB phase, a boundary that is not described by a 1+1D CFT.

II. FM PHASE AND ITS BOUNDARY

Because the ferromagnetic state with all spins polarized along $\pm z$ (or an arbitrary superposition of these two states) is an exact eigenstate of the Hamiltonian for any value of α and J_z , we expect a first-order quantum phase transition at the boundary of the FM phase. The FM state has an energy $E_{\text{FM}} = J_z \sum_{i>j} (i-j)^{-\alpha}$, and the phase transition out of this state, defining the critical line $J_z = \lambda_0(\alpha)$, occurs when some other eigenstate with no ferromagnetic order appears with a lower energy. The dependence of λ_0 on α can be estimated using the following intuitive argument. For a given $J_z < 0$, the energy density of the ferromagnetic state in the thermodynamic limit is given by $\epsilon_{\text{FM}} = J_z \zeta(\alpha)$ [$\zeta(\alpha) \equiv \sum_{r=1}^{\infty} r^{-\alpha}$ is the Riemann zeta function], which diverges as $\alpha \rightarrow 1$. For $J_{xy} = 1$, the magnitude of the energy density arising from the term $\sum_{i>j} (S_i^x S_j^x + S_i^y S_j^y)/(i-j)^\alpha$ can be at most $\eta(\alpha) \equiv \sum_{r=1}^{\infty} (-1)^{r-1}/r^\alpha$ (the Dirichlet eta function), with this value obtained for any state that is Néel-ordered along some direction in the x - y plane. The competition between the energy of these two classical states gives a critical point $J_z \approx -\eta(\alpha)/\zeta(\alpha)$, which smoothly varies from $J_z = -1$ at $\alpha = \infty$ to $J_z = 0$ at $\alpha = 1$. For $J_{xy} = -1$, the situation is quite different, because the polarized state along any direction in the x - y plane has an energy density equal to $-\zeta(\alpha)$, and thus we naively expect the phase boundary to be at $J_z = -1$ for all $\alpha > 0$.

More formally, the boundary can be calculated via a spin-wave analysis. We treat the spin state that is polarized along the $+z$ direction as the vacuum state with no excitations and apply the Holstein-Primakoff transformation (for spin 1) to map spin excitations (spin-waves) into bosons: $S_i^z = 1 - a_i^\dagger a_i$, $S_i^+ \equiv S_i^x + i S_i^y = \sqrt{2} a_i^\dagger (1 - a_i^\dagger a_i/2)^{1/2}$. In the weak excitation limit, $\langle a_i^\dagger a_i \rangle \ll 1$, we can approximate $S_i^+ \approx \sqrt{2} a_i^\dagger$, and our Hamiltonian becomes

$$H_{\text{sw}} \approx \sum_{i>j} \frac{-J_z(a_i^\dagger a_i + a_j^\dagger a_j) + J_{xy}(a_i^\dagger a_j + a_j^\dagger a_i)}{(i-j)^\alpha}, \quad (2)$$

where we have ignored the interaction terms $a_i^\dagger a_j a_j^\dagger a_i$ since $\langle a_i^\dagger a_i \rangle, \langle a_j^\dagger a_j \rangle \ll 1$ is assumed. Assuming for the moment periodic boundary conditions, this quadratic Hamiltonian can be diagonalized by Fourier transformation, $H_{\text{sw}} = 2 \sum_k \omega_k c_k^\dagger c_k$, with the following dispersion relation ($q \equiv 2\pi k/N$) for an infinite system

$$\omega(q) = -J_z \sum_{r=1}^{\infty} r^{-\alpha} + J_{xy} \sum_{r=1}^{\infty} \cos(qr)/r^\alpha. \quad (3)$$

If $\omega_{\text{min}} \equiv \min \omega(q) > 0$, then the ground state of H_{sw} is the vacuum state of all modes k , and $\langle a_i^\dagger a_i \rangle = 0$ for all i , consistent with the approximation $\langle a_i^\dagger a_i \rangle \ll 1$. If $\omega_{\text{min}} < 0$, then the ground state has an extensive number of spin excitations and the spin-wave approximation should break down, and we do not expect the polarized state in the z direction to be the quantum ground state. The $\omega_{\text{min}} = 0$ condition thus sets the phase boundary for H_{sw} . For $J_{xy} = 1$, $\omega_{\text{min}} = \omega(q = \pi) = -J_z \zeta(\alpha) - \eta(\alpha)$, leading to a critical line of $J_z = -\eta(\alpha)/\zeta(\alpha)$. For $J_{xy} = -1$, $\omega_{\text{min}} = \omega(q = 0) = (1 - J_z)\zeta(\alpha)$, leading to a critical line at $J_z = -1$, independent of α . These results exactly match with the previous intuitive arguments.

We now compare the above spin-wave theory prediction with infinite-size DMRG calculations [84,85] for $J_{xy} = 1$. As seen in Fig. 2, the numerical results agree well with the spin-wave theory at large α , and the spin-wave prediction of $\lambda_0(\alpha)$ is asymptotically exact as $\alpha \rightarrow \infty$. However, a small but increasing difference in $\lambda_0(\alpha)$ is seen as α decreases. For $\alpha \gtrsim 1.5$, our infinite-size DMRG calculations converge well (see appendix A for our numerical treatment of long-range interactions), and we conjecture that it is the spin-wave approximation that starts to break down when α decreases. This is possibly due to stronger effects of interactions between spin-wave excitations as α becomes smaller, so that the spin-wave approximation (which ignores interactions) becomes less

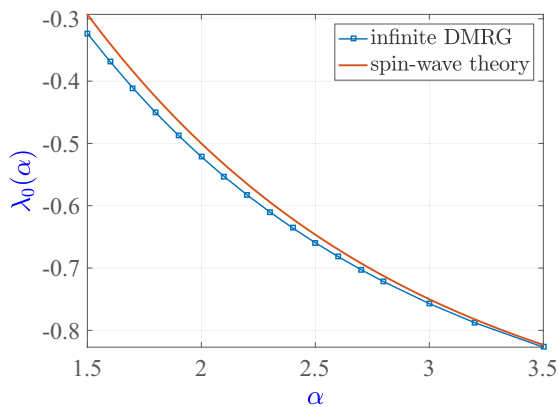


FIG. 2. Comparison of the (first-order) transition point $\lambda_0(\alpha)$ out of the FM phase calculated using infinite-size DMRG and spin-wave theory for $J_{xy} = 1$. The spin-wave theory predicts $\lambda_0(\alpha) = -\eta(\alpha)/\zeta(\alpha)$. The infinite-size DMRG calculations use a bond dimension $\chi = 100$, and increasing χ to 200 does not yield results distinguishable within the resolution of the plot. The transition point is numerically determined by finding the value of J_z at which the ground state energy density obtained from infinite-size DMRG calculations is equal to that of the FM state.

and less accurate. While our infinite-size DMRG calculations do not converge well for $\alpha \lesssim 1.5$, the spin-wave prediction should be asymptotically exact as $\alpha \rightarrow 1$, since the FM state's energy is super extensive for $\alpha \leq 1$ and $\lambda_0(\alpha \leq 1) = 0$. As a result, in Fig. 1(a) we have adopted the spin-wave prediction for the FM phase's boundary, but made the boundary line dotted for $1 < \alpha < 2$ to represent a small uncertainty in the transition point [for $\alpha > 2$ the uncertainty of the transition point is well below the resolution of Fig. 1(a)]. For $J_{xy} = -1$, our infinite-size DMRG calculations provide exactly the same transition point $\lambda_0(\alpha) = -1$ as the spin-wave theory, independent of α [Fig. 1(b)].

III. HALDANE PHASE AND ITS BOUNDARY

The existence of the Haldane phase in a spin-1 XXZ chain makes the phase diagram much richer than that of a spin-1/2 XXZ chain. We focus first on the XY-to-Haldane phase boundary $\lambda_1(\alpha)$. The transition out of the Haldane phase is signaled by a vanishing of the string-order correlation function $\mathcal{S}_{ij}^\xi \equiv \langle S_i^\xi S_j^\xi \prod_{i < k < j} (-1)^{S_k^\xi} \rangle$ ($\xi = x, y, z$) when $|i - j| \rightarrow \infty$. However, because the phase transition is of the BKT type, \mathcal{S}_{ij}^ξ changes rather smoothly with J_z and α for a finite $|i - j|$, and it is very challenging to find the exact transition point numerically. Finite-size scaling using exact diagonalization on small chains must be performed very carefully due to logarithmic corrections in system size [61,86–88], and infinite-size DMRG yields a phase transition point that depends strongly on the bond dimension χ (the dimension of the matrix product states used [89]), since the ground state bipartite entanglement entropy S grows logarithmically with system size N according to CFT: $S = c \log N + \text{const}$ [90]. As seen in Fig. 3, for $\chi = 100$ and at $\alpha = \infty$, the string-order correlation function \mathcal{S}_{ij}^z appears to start vanishing at $J_z \approx 0.3$, instead of at $J_z = 0$ as predicted by field theory [68]. However, this is consistent with previous infinite-size DMRG calculation results [65,66]. To extract a more accurate phase boundary, we perform a scaling of χ ranging from 50 to 200 near the XY-to-Haldane phase boundary, following a procedure similar to that in Ref. [65]. We then extract the XY-to-Haldane phase boundary (white line in Fig. 3) by determining the location where $\mathcal{S}_{ij}^z(\chi \rightarrow \infty)$ vanishes, which now correctly yields $J_z \approx 0$ at $\alpha = \infty$. However, we expect a few percent uncertainty in the transition point due to the use of \mathcal{S}_{ij}^z at a finite separation $|i - j|$, and due to the error in extrapolating $\mathcal{S}_{ij}^z(\chi \rightarrow \infty)$.

To explain why long-range interactions bend the XY-to-Haldane phase boundary in opposite directions for ferromagnetic and antiferromagnetic J_{xy} , we use an effective field theory first proposed by Haldane [3] and developed by Affleck [91]. The proper inclusion of long-range interactions within this field theoretic approach was discussed in detail in Ref. [57]. Here, we give a brief review of this field-theory treatment. Consider first the case of $J_{xy} = J_z = 1$. In this case, each spin-1 is mapped to a staggered field $\mathbf{n}(2i + \frac{1}{2}) = (\mathbf{S}_{2i} - \mathbf{S}_{2i+1})/2$ and a uniform field $\mathbf{l}(2i + \frac{1}{2}) = (\mathbf{S}_{2i} + \mathbf{S}_{2i+1})/2$. Importantly, we observe that the classical ground state of H is always Néel-ordered for any $\alpha > 0$, with $\mathbf{n}^2(x) = 1$ and $\mathbf{l}(x) = 0$ for any position x . The intuition behind this

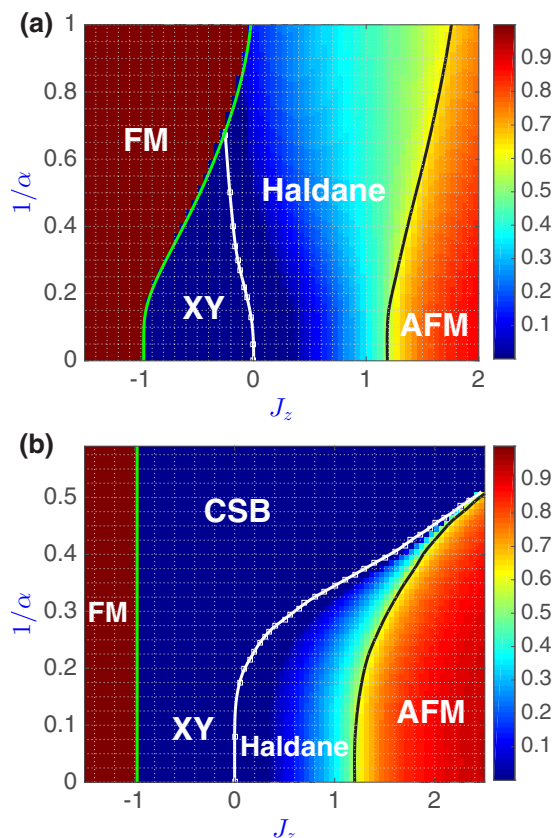


FIG. 3. Infinite-size DMRG calculation of $S_{ij}^z \equiv \langle S_i^z S_j^z \prod_{i < k < j} (-1)^{S_k^z} \rangle$ for a separation of $|i - j| = 500$. $S_{ij}^z = 1$ in the FM phase and $S_{ij}^z \approx 1$ deep in the AFM phase for any i and j . As $|i - j| \rightarrow \infty$, S_{ij}^z is finite for the Haldane phase and zero for the XY phase, thus we can use it to locate the XY-to-Haldane phase boundary. (a) $J_{xy} = 1$. The FM phase boundary (green line) is given by the spin-wave prediction $J_z = -\eta(\alpha)/\zeta(\alpha)$. (b) $J_{xy} = -1$. The FM phase boundary (green line) is exactly at $J_z = -1$. For both (a) and (b), we vary the bound dimension χ to accurately determine the XY-to-Haldane phase boundary, determining the value of J_z at which S_{ij}^z vanishes (for a large but finite $|i - j|$) and then extrapolating to the $\chi \rightarrow \infty$ limit (white squares fitted by the white line). The black line is the Haldane-to-AFM phase boundary, which is determined from $\langle S_i^z S_j^z \rangle$ (see Fig. 4).

decomposition is that, in the quantum ground state, $\mathbf{n}(x)$ should have only long-wave-length variations with $\mathbf{n}^2(x) \approx 1$, while $\mathbf{l}(x) \approx 0$ should represent long wave-length perturbations to the direction of $\mathbf{n}(x)$ due to quantum fluctuations. Therefore, when working with the Fourier-transformed fields $\mathbf{n}(q)$ and $\mathbf{l}(q)$, we can expand the Hamiltonian in powers of the momentum q and keep only the leading order terms.

The effective Hamiltonian in the continuum limit and momentum space reads (the lattice spacing is set to 1 for simplicity)

$$H_{\text{eff}} \approx \int dq [\omega(q)|\mathbf{n}(q)|^2 + \Omega(q)|\mathbf{l}(q)|^2], \quad (4)$$

where the cross terms between \mathbf{n} and \mathbf{l} are ignored because they involve $\mathbf{n}(q)$ near $q = \pi$. The dispersion relations $\Omega(q)$

and $\omega(q)$ can be expanded at small q as [92]:

$$\begin{aligned} \omega(q) &\equiv 2 \sum_{r=1}^{\infty} (-1)^r \frac{\cos(qr)}{r^\alpha} \\ &\approx -2\eta(\alpha) + \eta(\alpha - 2)q^2 + O(q^4), \\ \Omega(q) &\equiv 2 \sum_{r=1}^{\infty} \frac{\cos(qr)}{r^\alpha} \approx 2\zeta(\alpha) + \zeta(\alpha - 2)q^2 + O(q^4) \\ &\quad + 2\Gamma(1 - \alpha) \cos\left[\frac{\pi}{2}(\alpha - 1)\right] |q|^{\alpha-1}. \end{aligned} \quad (5)$$

For the \mathbf{n} field, we need to keep the q^2 term since the zeroth-order term gives a constant due to the approximation $\mathbf{n}^2(x) \approx 1$. The zeroth-order term in q for the \mathbf{l} field is the dominant source of quantum fluctuations, and we can ignore higher-order terms in determining whether H_{eff} is gapped or not (they do contribute to the long-distance behavior of correlation functions though [57]). Thus the Hamiltonian is approximately given by $H_{\text{eff}} \approx \int dq [\eta(\alpha - 2)q^2 |\mathbf{n}(q)|^2 + 2\zeta(\alpha) |\mathbf{l}(q)|^2]$. In a coherent-spin-state path-integral representation, the action is quadratic in the field \mathbf{l} and it can be integrated out [1,93]. The remaining path integral over the staggered field \mathbf{n} defines a 1+1D $O(3)$ nonlinear sigma model, with Lagrangian density (nonlinear constraint $\mathbf{n}^2(x) = 1$ implied)

$$\mathcal{L}(x) \approx \frac{1}{g} (|\partial \mathbf{n} / \partial t|^2 - v^2 |\partial \mathbf{n} / \partial x|^2). \quad (6)$$

Here the effective coupling g and spin-wave velocity v depend both on α and the lattice spacing (their exact values are not important to us). The coupling strength g flows towards infinity under renormalization group for the above Lagrangian [1,93], suggesting a disordered ground state with an excitation gap. This is corroborated by the $SU(n)$ variant of the Hamiltonian in the $n \rightarrow \infty$ limit, which can be analytically solved and contains a mass gap [91,94]. Now we adopt a phenomenological treatment [95,96] of the above Lagrangian (Eq. 6): The nonlinear constraint $\mathbf{n}^2 = 1$ can be approximately removed by introducing a mass gap Δ_α and a renormalized spin-wave velocity v_α . We thereby arrive at a free field theory with the Lagrangian density (written in momentum space)

$$\mathcal{L}(q) \propto |\partial \mathbf{n} / \partial t|^2 - (\Delta_\alpha^2 + v_\alpha^2 q^2) |\mathbf{n}(q)|^2. \quad (7)$$

Since $\Delta_\alpha \rightarrow \infty \approx 0.41$ [97,98] and $\Delta_\alpha \rightarrow 0 = 1$ (where the Hamiltonian becomes integrable), we infer that Δ_α should increase as α decreases. This speculation is confirmed by accurate finite-size DMRG calculations of Δ_α in Ref. [57].

Next, we consider the case of $J_{xy} = 1$ but $J_z < 1$. We can then write

$$H = \sum_{i>j} \frac{1}{(i-j)^\alpha} \mathbf{S}_i \cdot \mathbf{S}_j - (1 - J_z) \sum_{i>j} \frac{1}{(i-j)^\alpha} S_i^z S_j^z. \quad (8)$$

Following Refs. [76] and [91], the anisotropy term above can be treated as a negative mass term $(1 - J_z) f_\alpha n_z^2(q)$ to the Lagrangian density $\mathcal{L}(q)$ in Eq. (7). The precise value of the renormalization factor f_α is not important to us, but we expect it to continuously decrease as α becomes smaller, since the staggered field dominates in the Haldane phase and long-range interactions $[\sum_{i>j} \frac{1}{(i-j)^\alpha} S_i^z S_j^z$ in Eq. (8)]

are increasingly frustrated as α decreases. The mass gap for the field n_z is now smaller than for n_x and n_y , and reads $\Delta_\alpha(J_z) = \sqrt{\Delta_\alpha^2 - (1 - J_z)f_\alpha}$. Combined with the above discussion that Δ_α should increase with decreasing α , we require progressively more negative J_z to close the gap and transition into the XY phase as α decreases, thus explaining the shape of the XY-to-Haldane phase boundary in Fig. 3(a).

For $J_{xy} = -1$ and $J_z < 1$, the classical ground state is no longer Néel ordered and the field theory employed above is not valid. However, by rotating every other spin by π about the z axis, we generate a transformed Hamiltonian

$$H' = \sum_{i>j} \frac{(-1)^{i-j-1}}{(i-j)^\alpha} \mathbf{S}_i \cdot \mathbf{S}_j + \sum_{i>j} \frac{J_z - (-1)^{i-j-1}}{(i-j)^\alpha} S_i^z S_j^z. \quad (9)$$

Now the classical ground state is Néel ordered (along any direction for $J_z = 1$). The first term above is isotropic, and gets mapped to

$$\sum_{i>j} \frac{(-1)^{i-j-1}}{(i-j)^\alpha} \mathbf{S}_i \cdot \mathbf{S}_j \approx \int dq [\Omega(q)|\mathbf{n}(q)|^2 + \omega(q)|\mathbf{l}(q)|^2], \quad (10)$$

where the roles of $\omega(q)$ and $\Omega(q)$ are swapped as compared to Eq. (4). For $\alpha < 3$, $\Omega(q)$ in Eq. (5) is now dominated by the nonanalytic term $|q|^{\alpha-1}$ at small q , and we can no longer obtain the simple free Lagrangian in Eq. (7). In Ref. [57], it is shown that the $|q|^{\alpha-1}$ term in the dispersion of $\mathbf{n}(q)$ in Eq. (10) leads to a renormalization group flow towards a gapless ordered phase spontaneously breaking an $SU(2)$ symmetry for $\alpha < \alpha_c \lesssim 3$. For our complete Hamiltonian H' in Eq. (9), the anisotropy leads instead to a $U(1)$ continuous symmetry breaking phase for $\alpha < \alpha'_c$ (see the next section for further discussions, where α'_c is estimated to be 2.9 at $J_z = 1$). Our infinite-size DMRG calculations in Fig. 3(b) suggest that the Haldane phase terminates at a critical α around 3.1 for $J_z = 1$, and the XY phase is expected to exist in between the CSB phase and the Haldane phase at $J_z = 1$ for $\alpha'_c < \alpha < \alpha_c$.

For $\alpha > 3$, $\Omega(q)$ is dominated by q^2 and we can once again reduce H' to the free field Lagrangian Eq. (7), but with a different mass gap Δ'_α and spin-wave velocity v'_α . The anisotropy term in Eq. (9) changes the gap to $\Delta'_\alpha(J_z) = \sqrt{\Delta_\alpha^2 - (g_\alpha - J_z h_\alpha)}$. Here g_α is a renormalization factor due to nonfrustrating long-range interactions $\frac{(-1)^{i-j-1}}{(i-j)^\alpha} S_i^z S_j^z$ in Eq. (9), and should thus increase as α decreases, while h_α is a renormalization factor due to frustrating long-range interaction $\frac{1}{(i-j)^\alpha} S_i^z S_j^z$ in Eq. (9), and should decrease as α decreases. Together with the expectation that the gap Δ'_α should decrease with α [47,57] due to the appearance of gapless continuous symmetry breaking phase at $\alpha \lesssim 3$, we conclude that the gap closes at a point with J_z strictly larger than zero in the presence of long-range interactions, again consistent with our numerical results.

We point out that a different field theoretic approach based on non-Abelian bosonization [57,68] can also be employed to predict the qualitative changes to the XY-to-Haldane phase boundary. This technique has been used to predict the XY-to-Haldane phase boundary of a spin-1 XXZ chain with next-nearest-neighbor interactions [76], which is a reasonable approximation to our model when α is large enough that

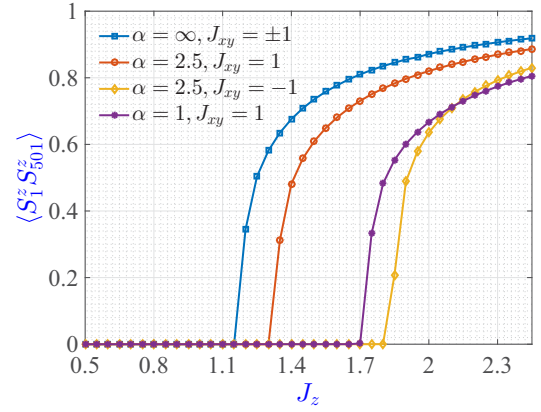


FIG. 4. $\langle S_1^z S_{501}^z \rangle$ as a function of J_z calculated using infinite-size DMRG for a few different sets of α and J_{xy} . The Haldane-to-AFM phase transition is clearly observed and we locate the transition point by finding the critical J_z (restricted to $J_z > 0$) above which $\langle S_1^z S_{501}^z \rangle > 0.1$. The curves shown look nearly identical when we increase the bond dimension used from 100 to 200.

next-nearest-neighbor interactions dominate over the next-longer-range interactions.

We end this section with a brief discussion of the boundary between the Haldane and AFM phases. Both the Haldane and AFM phases are gapped and have finite entanglement entropy in the infinite-system-size limit [99]. Thus we see well-converged results for bond dimensions of $\chi \geq 100$ in our infinite-size DMRG calculations. We extract the Haldane-to-AFM phase boundaries using the spin-spin correlation functions $C_{ij}^z \equiv \langle S_i^z S_j^z \rangle$ (see Fig. 4), and plot them as black lines in Figs. 3(a) and 3(b). Good agreement with existing literature [63–66] is found for the Haldane-to-AFM transition point at $\alpha = \infty$ ($1.15 < J_z < 1.2$). The bending of the Haldane-to-AFM phase boundary toward larger J_z for both $J_{xy} = 1$ and $J_{xy} = -1$ in the presence of long-range interactions can be understood via simple energetic considerations. In the AFM phase, the spins are (nearly) antialigned in the z direction; long-range interactions are strongly frustrated, and the energy $E = \sum_{i>j} \langle S_i^z S_j^z \rangle / (i-j)^\alpha$ at $\alpha \rightarrow 0$ is only half of the $\alpha = \infty$ case for a perfectly Néel ordered state. In the Haldane phase, the AFM order of spin correlations $\langle \mathbf{S}_i \cdot \mathbf{S}_j \rangle$ decays exponentially (followed by a small power-law tail [57]), and thus the ground state energy $E = \sum_{i>j} \langle \mathbf{S}_i \cdot \mathbf{S}_j \rangle / (i-j)^\alpha$ is much less frustrated by the long-range interactions. As a result, we expect the disordered ground state in the Haldane phase to have progressively lower energy than an AFM ordered state as α decreases at a given J_z , and hence a larger (but always finite even for $\alpha \rightarrow 0$) J_z is needed to make the transition from the Haldane phase into the AFM phase.

IV. CSB PHASE AND ITS BOUNDARY

The celebrated Mermin-Wagner theorem rigorously rules out continuous symmetry breaking in 1D and 2D quantum and classical spin systems at finite temperature, as long as the interactions satisfy the convergence condition $\sum_{i>j} J_{ij} r_{ij}^2 < \infty$ in the thermodynamic limit (r_{ij} and J_{ij} are respectively the distance and coupling strength between sites i and j) [2].

The long-distance properties of 1D spin systems at zero temperature can often be related to those of a 2D classical model at finite temperature; however, in the process of this mapping, the long-range interactions are only inherited by one of the two spatial directions in the classical model, and the Mermin-Wagner convergence condition will be satisfied for interactions decaying faster than $1/r^3$. Thus we expect no continuous symmetry breaking in the ground state of our Hamiltonian Eq. (1) for $\alpha > 3$. Indeed, we have found exclusively disordered or discrete (Z_2) symmetry breaking phases for $\alpha > 3$ in our phase diagrams (Fig. 1). Continuous symmetry breaking can (and does) appear when $\alpha < 3$. To gain a better understanding of the robustness of symmetry breaking states to quantum fluctuations, below we carry out a spin-wave analysis [100]. Similar analysis can be also found in Ref. [44] and [45] for Heisenberg chains with long-range interactions.

We start by considering the $J_{xy} = -1$ case, and take the state with all spins polarized along the $+x$ direction as the vacuum state. With this choice of vacuum, and assuming that the density of spin waves is small ($\langle a_i^\dagger a_i \rangle \ll 1$ in the following expressions), the Holstein-Primakoff mapping is now $S_i^x = 1 - a_i^\dagger a_i$, $S_i^y \approx (a_i^\dagger + a_i)/\sqrt{2}$, $S_i^z \approx (a_i^\dagger - a_i)/i\sqrt{2}$. Under this mapping, and dropping terms that are quartic in bosonic operators (again based on the assumption that $\langle a_i^\dagger a_i \rangle \ll 1$), H becomes

$$H_{\text{swx}} = \sum_{k=-N/2}^{N/2} (a_k^\dagger \quad a_{-k}) \begin{pmatrix} \omega_k & \mu_k \\ \mu_k & \omega_k \end{pmatrix} \begin{pmatrix} a_k \\ a_{-k} \end{pmatrix}; \quad (11)$$

$$\omega_k = \sum_{r=1}^{N/2} \frac{1}{r^\alpha} + \frac{J_z - 1}{2} \sum_{r=1}^{N/2} \frac{1}{r^\alpha} \cos\left(\frac{2\pi k}{N} r\right), \quad (12)$$

$$\mu_k = -\frac{J_z + 1}{2} \sum_{r=1}^{N/2} \frac{1}{r^\alpha} \cos\left(\frac{2\pi k}{N} r\right), \quad (13)$$

where $a_k = \frac{1}{\sqrt{N}} \sum_j e^{i2\pi jk/N} a_j$. H_{swx} can be diagonalized with a Bogoliubov transformation, yielding noninteracting Bogoliubov quasiparticles with a spectrum ν_k . Importantly, when $|\omega_k| > |\mu_k|$, $\nu_k > 0$ and the vacuum is dynamically stable. When $|\omega_k| < |\mu_k|$, ν_k is imaginary and the system is dynamically unstable indicating that we have made the wrong choice of a classical ground state. Using the expressions for ω_k and μ_k in Eqs. (12) and (13), we find that $|\omega_k| > |\mu_k|$ is satisfied for all $k \neq 0$ modes if and only if $-1 \leq J_z < \zeta(\alpha)/\eta(\alpha)$. This is because when $J_z < -1$, the classical ground state is ferromagnetic in z direction, and when $J_z > \zeta(\alpha)/\eta(\alpha)$ the classical ground state is Néel ordered along the z direction.

Because the Bogoliubov quasiparticles consist of both particles and holes, the ground state of H_{swx} can have a finite or divergent density of spin excitations, measured by

$$\begin{aligned} \langle a_i^\dagger a_i \rangle &= \frac{1}{N} \sum_{k \neq 0} \frac{1}{2} \left([1 - \mu_k^2/\omega_k^2]^{-1/2} - 1 \right) \\ \xrightarrow{N \rightarrow \infty} &= \frac{1}{4\pi} \int_{-\pi}^{\pi} dq [1 - \mu^2(q)/\omega^2(q)]^{-1/2} - 1. \end{aligned} \quad (14)$$

Expanding the integrand $[1 - \mu^2(q)/\omega^2(q)]^{-1/2}$ above around $q = 0$ to the lowest order in $1/q$, we find that

$$\begin{aligned} &[1 - \mu^2(q)/\omega^2(q)]^{-1/2} \\ &\approx \sqrt{\frac{(1 + J_z)\zeta(\alpha)}{2\zeta(\alpha - 2)q^2 - 4\Gamma(1 - \alpha)\cos[\frac{\pi}{2}(\alpha - 1)]|q|^{\alpha-1}}}, \end{aligned} \quad (15)$$

where $\Gamma(x)$ is the gamma function. For $\alpha > 3$, since $[1 - \mu^2(q)/\omega^2(q)]^{-1/2} \propto 1/|q|$ to leading order in $1/q$, the spin-wave density $\langle a_i^\dagger a_i \rangle \sim \ln(N)$ diverges as $N \rightarrow \infty$ according to Eq. (14). This means the long-range ferromagnetic order along the x direction is destroyed by quantum fluctuations in the thermodynamic limit; we expect that $\lim_{|i-j| \rightarrow \infty} \langle S_i^+ S_j^- \rangle = 0$, and the system will be disordered (either Haldane or XY). For $\alpha < 3$, instead we have $[1 - \mu^2(q)/\omega^2(q)]^{-1/2} \propto 1/|q|^{(\alpha-1)/2}$ to leading order in q , and the excitation density $\langle a_i^\dagger a_i \rangle$ converges to a finite constant. As a self-consistency condition, we also require $\langle a_i^\dagger a_i \rangle < 1$ to prevent the breakdown of the spin-wave approximation [44,52]. We expect a CSB phase in the parameter region of $\langle a_i^\dagger a_i \rangle < 1$, with nonvanishing spin order in the x - y plane (i.e., $\lim_{|i-j| \rightarrow \infty} \langle S_i^+ S_j^- \rangle \neq 0$), and a disordered phase when $\langle a_i^\dagger a_i \rangle > 1$. By numerically evaluating Eq. (14), which gives $\langle a_i^\dagger a_i \rangle$ in the infinite-size limit [101], we have obtained a phase diagram for $J_{xy} = -1$ under spin-wave approximation (Fig. 5).

For $J_{xy} = 1$ and $|J_z| < 1$, classically the spins prefer to antialign in the x - y plane. Expanding around this classical state, the spin-wave approximation leads to the same Hamiltonian

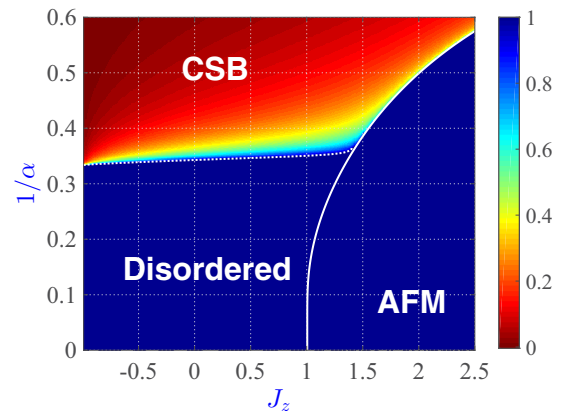


FIG. 5. Spin-wave excitation density $\langle a_i^\dagger a_i \rangle$ calculated using Eq. (14) for an infinite-size chain. For $J_z > \zeta(\alpha)/\eta(\alpha)$ (region to the right of the white line), imaginary frequencies appear in the Bogoliubov spectrum, indicating a classical instability toward the AFM phase. The region above the dotted and solid white lines has $\langle a_i^\dagger a_i \rangle \leq 1$, and is associated with the CSB phase. The remaining region in the plot has $\langle a_i^\dagger a_i \rangle > 1$, and is expected to be disordered. The disordered phase can be either the XY or the Haldane phase, but the spin-wave theory cannot distinguish one from the other. For better visibility, we have set $\langle a_i^\dagger a_i \rangle = 1$ for regions without CSB in the plot.

in Eq. (11) except that we have

$$\omega_k = \sum_{r=1}^{N/2} \frac{(-1)^{r-1}}{r^\alpha} + \frac{J_z - 1}{2} \sum_{r=1}^{N/2} \frac{(-1)^{r-1}}{r^\alpha} \cos\left(\frac{2\pi k}{N} r\right), \quad (16)$$

$$\mu_k = -\frac{J_z + 1}{2} \sum_{r=1}^{N/2} \frac{(-1)^{r-1}}{r^\alpha} \cos\left(\frac{2\pi k}{N} r\right). \quad (17)$$

As a result, now both $\mu(q)$ and $\omega(q)$ become fully analytic (in the $N \rightarrow \infty$ limit) due to the alternating sign $(-1)^r$ in Eqs. (16) and (17). Expanding around $q = 0$, we have

$$[1 - \mu^2(q)/\omega^2(q)]^{-1/2} \approx \sqrt{\frac{(1 + J_z)\eta(\alpha)}{2\eta(\alpha - 2)q^2}}. \quad (18)$$

As a result, $\langle a_i^\dagger a_i \rangle$ will be divergent for any $\alpha > 0$ due to the $1/|q|$ singularity in the integrand of Eq. (14). Thus continuous symmetry breaking is forbidden for all $\alpha > 0$ for $J_{xy} = 1$.

Now we confirm the spin-wave prediction of the CSB phase's boundary using DMRG calculations. Naively, one should calculate the CSB phase's order parameter $\lim_{|i-j| \rightarrow \infty} \langle S_i^+ S_j^- \rangle$. However, we find that in the XY phase $\langle S_i^+ S_j^- \rangle \sim 1/|i-j|^\eta$ decays with a rather slow power law (e.g., $\eta = 0.25$ at $J_z = 0$ and $\alpha = \infty$). At the maximum separation that we can calculate accurately using either finite or infinite-size DMRG, $\langle S_i^+ S_j^- \rangle$ only shows a crossover from the XY phase to the CSB phase. To faithfully determine the boundary of the CSB phase, we instead calculate the effective central charge c_{eff} as a function of α and J_z . We obtain c_{eff} by calculating the half-chain entanglement entropy S for two chains with different total lengths N_1 and N_2 using a finite-size DMRG algorithm [102]. Explicitly, for large N_1 and N_2 , we have

$$c_{\text{eff}} \approx 6 \frac{S(N_1) - S(N_2)}{\ln(N_1) - \ln(N_2)}. \quad (19)$$

In the XY phase (including its boundaries) and at the boundary between the Haldane and AFM phases, we expect 1+1D conformal symmetry in the underlying field theory model [67,69], with c_{eff} being the actual central charge representing the conformal anomaly [90]. In the Haldane, FM, and AFM phases, no 1+1D conformal symmetry exists due to the presence of a gap. Although the CSB phase is gapless, we expect a breakdown of 1+1D conformal symmetry due to the $1/r^\alpha$ long-range interactions that become relevant in the RG sense for $\alpha \lesssim 3$ [33,35,83]. We emphasize that in phases with no conformal symmetry, c_{eff} does not have the meaning of the central charge and is used only as a diagnostic here to numerically find phase boundaries.

We identify the XY-to-CSB phase boundary in Fig. 6 as the place where c_{eff} starts to become appreciably (5–10%) larger than 1. Here we find good agreement with the XY-to-CSB phase boundary predicted by spin-wave theory in Fig. (5) for $-1 < J_z \lesssim 1$. Together with perturbative field theory calculations presented in Ref. [83], we expect the phase boundary in Fig. 6 to be accurate within a few percent. The accuracy of the calculated c_{eff} can be further improved by finite-size scaling, which is however beyond the scope of the

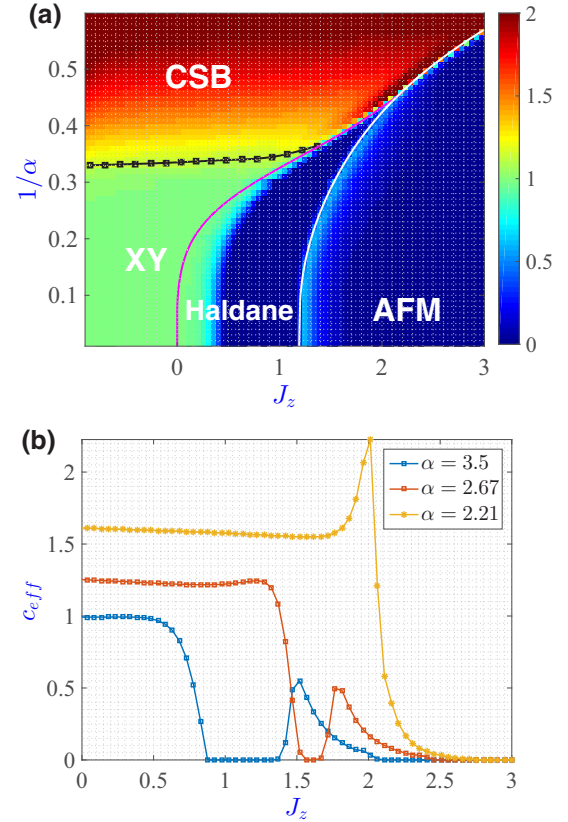


FIG. 6. Calculation of the effective central charge c_{eff} as a function of J_z and α for $J_{xy} = -1$, extracted from finite-size DMRG calculations with $N_1 = 100$, $N_2 = 110$, and a maximum bond dimension of 500. (a) The black squares (fitted by the black line) show where c_{eff} starts to deviate from 1 when going from the XY to the CSB phase. The purple line and white line are from Fig. 3, and show the boundaries of the Haldane phase. (The calculation of c_{eff} is inaccurate in predicting the location of the XY-to-Haldane transition due to strong finite-size effects [61,86–88].) For better contrast, locations with $c > 2$ are shown with the color corresponding to $c = 2$. (b) For our finite-size chains, the XY-to-Haldane BKT phase transition is signaled by a continuous drop of c_{eff} from 1 to 0 ($\alpha = 3.5$). The Haldane-to-AFM phase transition is identified by a peak with value around 0.5 in c_{eff} ($\alpha = 3.5$ and $\alpha = 2.67$). The CSB-to-Haldane transition is expected to be continuous and not associated with a central charge ($\alpha = 2.67$). The CSB-to-AFM transition has a sharp peak in c_{eff} ($\alpha = 2.21$), an indication of a first-order transition [67].

current study. The location of the CSB-XY-Haldane tricritical point is estimated to be at $\alpha \approx 2.75$ and $J_z \approx 1.35$.

From Ref. [83], it follows that the XY-to-CSB transition is a BKT-like transition that belongs to a universality class different from the XY-to-Haldane BKT transition. The Haldane-to-CSB transition is somewhat exotic, because the Haldane phase maps to a high-temperature disordered phase in a 2D classical model [93], and *in the absence of long-range interactions*, the CSB phase exists in 2D only at zero temperature [2] and is unlikely to undergo a phase transition directly to a high-temperature disordered phase. We also argue that the CSB-to-Haldane transition is not described by a 1+1D CFT, as supported by our numerical calculations shown in Fig. 6(b),

where c_{eff} changes smoothly (at least for finite chains) from a value larger than 1 to 0 during the CSB-to-Haldane transition.

The CSB-to-AFM phase transition is very likely to be first order, similar to the transition between the large- D and AFM phases studied in Refs. [66] and [75], despite the existence of quantum fluctuations in both phases. As shown in Fig. 6, we observe a sharp peak in c_{eff} at small α s when J_z is varied, indicating a first order transition [67], with further evidence that includes jumps in sublattice magnetization and spin-spin correlation across the CSB-to-AFM transition (not shown).

V. EXPERIMENTAL DETECTION

It was theoretically proposed in Refs. [54] and [55] that the Hamiltonian we consider can be simulated (for widely tunable J_z and $0 < \alpha < 3$) by using microwave field gradients or optical dipole forces to induce spin-spin interactions in a chain of trapped ions. The simulation of Eq. (1) with $J_{xy} = 1$ and $J_z = 0$ was experimentally demonstrated for a few ions with α tuned around 1 [56], where the ground state was adiabatically prepared by slowly ramping down an extra single-ion anisotropy term $D(t) \sum_i (S_i^z)^2$, with $D(t) > 0$. As the system size increases, the energy gap separating the ground state from the rest of the spectrum will become progressively smaller near the point where a phase transition between the “large- D ” phase and the XY/Haldane/FM/AFM phase occurs in the thermodynamic limit [75]. To avoid a slow ground state preparation process, we can adiabatically ramp down a staggered magnetic field in the z direction, $h(t) \sum_{i=1}^N (-1)^i S_i^z$, with $h(t) > 0$ [54,55]. By preparing an initial state that is the highest excited state of the staggered field Hamiltonian, the same adiabatic ramping process will lead us to the ground state of the Hamiltonian Eq. (1) with the opposite sign of both J_{xy} and J_z . As discussed in Ref. [55], the spin correlation functions $\langle S_i^z S_j^z \rangle$ and the string-order correlation $S_{ij}^z \equiv \langle S_i^z S_j^z \prod_{i < k < j} (-1)^{S_k^z} \rangle$ can be measured for any i and j , since one can obtain the complete statistics of all spins’ magnetization using spatially resolved measurements. Together with arbitrary single-spin rotations performed with microwave or optical Raman transitions, we can measure these correlations along any direction. Near-future experiments will most likely be limited to a few tens of spins. Although this limitation makes it difficult to probe continuous phase transitions, one can nevertheless observe important signatures of all five phases discussed in the manuscript by tuning J_z/J_{xy} and α deep into each phase. These signatures are summarized below and shown in Fig. 7.

FM phase [Fig. 7(a)]: Within the FM phase, $\langle S_i^z S_j^z \rangle = 1$ and $\langle S_i^x S_j^x \rangle = 0$ for any i and j , thus confirming perfect alignment of spins along the z direction.

AFM phase [Fig. 7(b)]: For sufficiently large $J_z > 0$, we have $\langle S_i^z S_j^z \rangle \approx (-1)^{i-j}$, showing a near perfect antialignment of spins along the z direction. In contrast, $\langle S_i^x S_j^x \rangle$ vanishes over a separation of just a few sites.

Haldane phase [Fig. 7(c)]: S_{ij}^z converges quickly to a nonzero constant as $|i - j|$ increases. In contrast, $\langle S_i^z S_j^z \rangle$ and $\langle S_i^x S_j^x \rangle$ vanish over a separation of just a few sites.

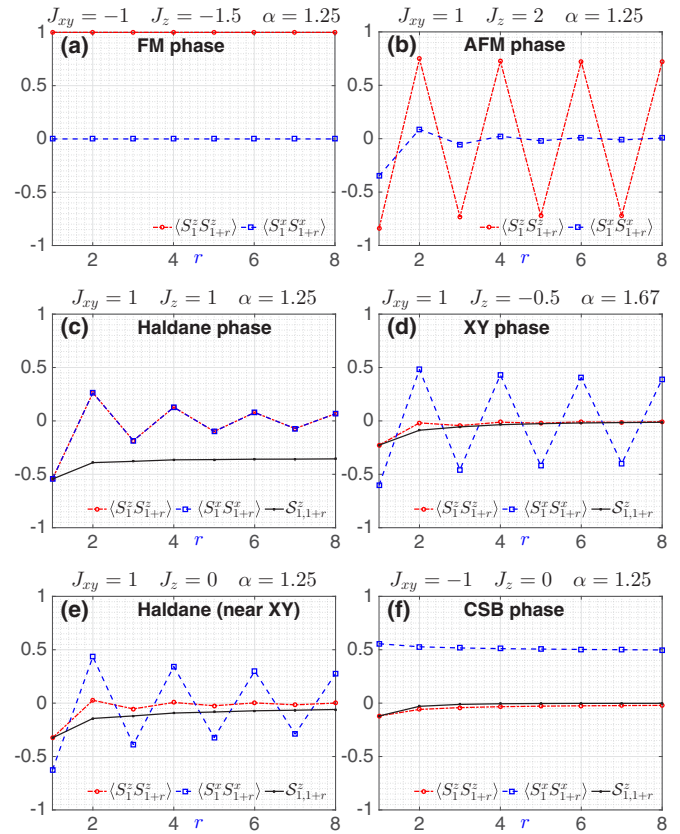


FIG. 7. Signatures of all five phases for a $N = 16$ spin chain. Except for (e), we tune J_{xy} , J_z , and α to set the ground state deep into each phase. Each phase is distinguished from the other phases by different behaviors in various spin-spin correlation functions.

XY phase [Fig. 7(d)]: We consider the XY phase for $J_{xy} = 1$ since the XY phase hardly exist for $\alpha < 3$ and $J_{xy} = -1$. S_{ij}^z and $\langle S_i^z S_j^z \rangle$ both decay quickly to zero as $|i - j|$ increases. $\langle S_i^x S_j^x \rangle$ oscillates and its amplitude decays very slowly (the slow decay reflects a relatively small value of the critical exponent associated with the correlation function decay).

CSB phase [Fig. 7(f)]: As in the XY phase, both S_{ij}^z and $\langle S_i^z S_j^z \rangle$ decay quickly to zero. However, $\langle S_i^x S_j^x \rangle$ converges quickly to approximately 0.5 at large $|i - j|$, showing a near perfect ordering of spins in the $x - y$ plane. Note that we are not explicitly breaking $U(1)$ symmetry here, so $\langle S_i^x S_j^y \rangle = \langle S_i^y S_j^x \rangle = \frac{1}{2} \langle S_i^+ S_j^- \rangle$. This is done because it is desirable for the experiment to operate within the $\sum_{i=1}^N S_i^z = 0$ subspace, where magnetic field noise and unwanted phonon couplings are suppressed [55,56].

Finally, we point out that, even in the experimental setup already demonstrated in Ref. [56], for which $J_z = 0$, one can still explore the two most interesting phases studied in this paper: the Haldane phase and the CSB phase. Note that, for $J_{xy} = 1$, $J_z = 0$ lies close to the Haldane-to-XY phase boundary, and thus one observes signatures of both phases, as in Fig. 7(e). Here the Haldane phase is identified via bulk correlations, but one can alternatively confirm the existence the Haldane phase by preparing edge excited states and measure edge excitation amplitudes [57].

VI. CONCLUSION AND OUTLOOK

By tuning the anisotropy $J_z/|J_{xy}|$ and the power-law exponent α , we have explored a rich variety of quantum phases—and the transitions between them—in a long-range interacting spin-1 XXZ chain. For $J_{xy} = -1$, long-range interactions give rise to a rather unusual phase diagram due to the emergence of a continuous symmetry breaking phase in one spatial dimension. Because the CSB phase cannot happen in a short-range interacting 1D spin-system, the nature of the phase transitions into and out of it is rather interesting; an in-depth study of the universality class of the CSB-to-XY transition was carried out in a separate work [83], where a similar transition in a long-range interacting spin-1/2 XXZ chain is analyzed. On the other hand, the CSB-to-Haldane transition, absent in spin-1/2 chains, requires further study to be understood thoroughly. The CSB-Haldane-AFM tricritical point is reminiscent of the tricritical point at the intersection of the large- D , Haldane and AFM phases, which has been related to the integrable Takhtajan-Babujian model described by an $SU(2)_2$ Wess-Zumino-Witten (WZW) model with central charge $c = 3/2$ [67,103–106]. Additional numerical calculations are needed to accurately determine the central charge at the CSB-Haldane-AFM tricritical point. Generalizations of our model to include single-ion anisotropy and a magnetic field are readily achievable in current trapped-ion experiments [55,56]. Understanding these exotic quantum phase transitions—induced by long-range interactions that are highly tunable in current experiments—requires the confrontation of numerous theoretical and numerical challenges, and motivates experimental quantum simulation of the model using AMO systems.

ACKNOWLEDGMENTS

We thank G. Pupillo, D. Vodola, L. Lepori, A. Turner, J. Pixley, M. Wall, P. Hess, A. Lee, J. Smith, A. Retzker and I. Cohen for helpful discussions. This work was supported by the ARO, the AFOSR, NSF PIF, NSF PFC at the JQI, and the ARL. M.F.-F. thanks the NRC for support.

APPENDIX: NUMERICAL TREATMENT OF LONG-RANGE INTERACTIONS

In our infinite-size and finite-size DMRG code [85], the $1/r^\alpha$ long-range interactions are represented as a

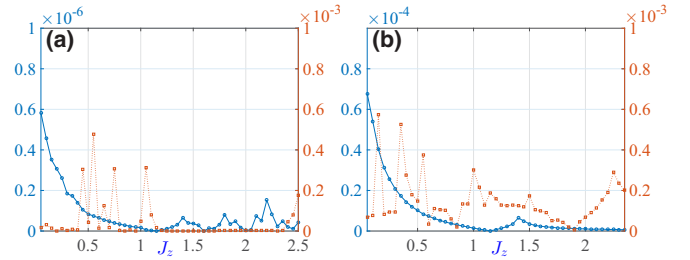


FIG. 8. Relative differences of $S_{1,501}^z$ (blue) calculated for $J_{xy} = 1$, $\alpha = 2$ and $\langle S_{1,501}^{\pm} \rangle$ (red) calculated for $J_{xy} = -1$, $\alpha = 2$ caused by (a) increasing L from 5000 to 10^4 (with K increasing from 9 to 10) and (b) decreasing ϵ_f from 10^{-10} to 10^{-11} (with K increasing from 9 to 10).

matrix product operator by fitting the power law to a sum of exponentials [51]. Specifically, we fit $f_r = 1/r^\alpha$ to $f'_r = \sum_{k=1}^K c_k e^{-r/\xi_k}$ for $r = 1, 2, \dots, L$. For a given L , we numerically find the minimum number of exponentials K that satisfy $\sum_{r=1}^L (f_r - f'_r)^2 \leq \epsilon_f$, with ϵ_f denoting the residual tolerance. The maximal range L is set to the chain length in our finite-size DMRG calculations, and to 5000 in our infinite-size DMRG calculations (much larger than the 500 site separation of correlations calculated in Fig. 3). ϵ_f is set to 10^{-12} in our finite-size DMRG calculations, and 10^{-10} in our infinite-size DMRG calculations. For all the calculations shown in the main text, we find no distinguishable differences within the resolution of our plots if we further increase L or decrease ϵ_f .

As an example, we show in Fig. 8 relative differences of $S_{1,501}^z$ (for $J_{xy} = 1$, $\alpha = 2$) and $\langle S_{1,501}^{\pm} \rangle$ (for $J_{xy} = -1$, $\alpha = 2$) caused by increasing L from 5000 to 10^4 and by decreasing ϵ_f from 10^{-10} to 10^{-11} . In all cases, the relative differences in the calculated observables are below 10^{-3} .

We have avoided the use of DMRG results if $\alpha < 1.5$ and interactions are unfrustrated in one or more directions ($J_{xy} = -1$ or $J_z < 0$ or both), because of the slow convergence of $\sum_{r=1}^L 1/r^\alpha$ with L as $\alpha \rightarrow 1$. Nevertheless, we do not expect new phase transitions in these situations based on Fig. 1, and we can instead infer the phases of the system there from the presented calculations.

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