

Hydrogenic states of monopoles in diluted quantum spin ice

Olga Petrova,¹ Roderich Moessner,¹ and S. L. Sondhi^{1,2}

¹Max-Planck-Institut für Physik komplexer Systeme, Nöthnitzer Strasse 38, 01187 Dresden, Germany

²Department of Physics, Princeton University, Princeton, New Jersey 08544, USA

(Received 14 January 2015; revised manuscript received 17 August 2015; published 1 September 2015)

We consider the effect of adding quantum dynamics to a classical topological spin liquid, with a particular view of how to best detect its presence in experiment. For the Coulomb phase of spin ice, we find quantum effects to be most visible in the gauge-charged monopole excitations. In the presence of weak dilution with nonmagnetic ions we find a particularly crisp phenomenon, namely, the emergence of hydrogenic excited states in which a magnetic monopole is bound to a vacancy at various distances. Via a mapping to an analytically tractable single particle problem on the Bethe lattice, we obtain an approximate expression for the dynamic neutron scattering structure factor.

DOI: [10.1103/PhysRevB.92.100401](https://doi.org/10.1103/PhysRevB.92.100401)

PACS number(s): 75.40.Gb, 75.10.Kt

The quest for spin liquids is an important enterprise in strongly correlated many-body physics in an era when a huge amount of theoretical interest has focused on forms of order outside the canonical broken symmetry paradigm [1–4]. The search involves identifying relatively simple Hamiltonians that host spin liquids and finding experimental systems and signatures—the latter being more elusive than in Landau ordered systems. Indeed, at this point the list of experimental systems where there is strong evidence of spin liquid behavior is small. Among them is the celebrated spin ice system, arising in some rare earth pyrochlore magnets, which exhibits a U(1) spin liquid and excitations which are condensed matter analogs of Coulombically interacting magnetic monopoles [5]. Spin ice is truly special at this point in hosting a three-dimensional spin liquid, but, owing to large magnetic moments, is limited to the classical regime, in which coherent quantum dynamics appears to play little role.

Logically, much recent interest has focused on looking for quantum generalizations of spin ice. There are several candidate materials for *quantum* spin ice behavior, such as Tb₂Ti₂O₇ [6–9], Yb₂Ti₂O₇ [10–12], and Pr₂Zr₂O₇ [13], but an unambiguous experimental signature of quantum spin ice has been lacking. Logically, much recent theoretical work has focused on looking for quantum generalizations of spin ice in which quantum fluctuations can lead to a fully quantum U(1) spin liquid [14–20].

Here we investigate the addition of quantum fluctuations to spin ice but in a different limit which is, plausibly, of relevance to existing materials. Fundamentally, we wish to understand the leading order effects of adding quantum dynamics about the classical spin ice limit. As we will detail below, this has a parametrically larger effect on monopole motion than on monopole-free ground states so the leading manifestations of quantum fluctuations appear when monopoles are present.

We begin this program by studying the simplest manifestation of the quantum mechanics of monopoles—a striking effect that appears in the response of quantum spin ice to the introduction of a vacancy or missing spin. We find that the lowest lying excited states in the vicinity of the vacancy resemble those of hydrogen modulo lattice induced mixing—they involve a magnetic monopole bound to the impurity site into an infinite set of levels. In the presence of a dilute set of such impurities, these states give rise to a characteristic signature in neutron

scattering at low temperatures which we discuss. Readers may note the family resemblance of these hydrogenic monopole states to hydrogenic states in doped semiconductors [21, and references therein], although we caution that the details have crucial differences. We also note that the response of spin liquids to impurities is of broad interest as a diagnostic of their internal dynamics [22,23]: what happens when you dope a spin liquid is the fundamental question of the resonating valence bond theory of high temperature superconductivity [24].

In the balance of this Rapid Communication we begin by briefly reviewing how the dynamics of quantum spin ice can be formulated as the quantum mechanics of monopoles. We then concentrate our attention on the problem of a vacancy spin and describe how it can be mapped to a good approximation to a monopole on a Bethe lattice interacting with a fixed Coulombic charge. This model leads to a family of hydrogenic bound states of the monopole along with a continuum band. In the technical heart of this Rapid Communication we solve this problem and obtain an exact closed form solution for the on-site Green's functions. We use these results to obtain the signature of the hydrogenic states in the structure factor of spin ice containing a dilute set of vacancy spins. We conclude with some comments and pointers to future work.

Quantum dipolar spin ice. Our model Hamiltonian

$$H_{\text{QDSI}} = H_{\text{DSI}} + \sum_i \mathbf{t} \cdot \mathbf{S}_i \quad (1)$$

consists, first, of the classical dipolar spin ice Hamiltonian, defined for Ising spins \mathbf{S}_i living on the sites of pyrochlore lattice and pointing along the local easy axis joining centers of neighboring tetrahedra [25,26, and references therein]:

$$H_{\text{DSI}} = \frac{\mu_0 \mu^2}{4\pi} \sum_{i < j} \left[\frac{\mathbf{S}_i \cdot \mathbf{S}_j}{r_{ij}^3} - \frac{3(\mathbf{S}_i \cdot \mathbf{r}_{ij})(\mathbf{S}_j \cdot \mathbf{r}_{ij})}{r_{ij}^5} \right]. \quad (2)$$

The second term in Eq. (1) is the transverse field, oriented perpendicular to the local easy axis, which adds the simplest quantum dynamics in the form of single spin flips. This simple form is convenient for a first theoretical analysis; for a more complete symmetry-based analysis of quantum terms in the Hamiltonian, see [17,27].

Ghost spins and the Bethe lattice. At this point we switch from a spin description to that referred to as the *dumbbell model* [5], which we quickly review. Each spin is replaced by a pair of magnetic charges $\pm q_m = \mu/a_d$ of opposite sign, where \bar{a}_d is a vector pointing between the centers of neighboring tetrahedra. By summing up the net charge at the center of each tetrahedron ($Q_\alpha \equiv \sum_{i \in \alpha} q_i = 0, \pm 2q_m, \pm 4q_m$), we can replace the dipolar piece of the spin Hamiltonian (2) by

$$H = \frac{\mu_0}{4\pi} \sum_{\alpha < \beta} \frac{Q_\alpha Q_\beta}{r_{\alpha\beta}} + \frac{v_0}{2} \sum_{\alpha} Q_\alpha^2, \quad (3)$$

Coulomb interactions between charges on the diamond lattice, with $v_0 = (2 + \sqrt{6})\mu_0/(4\pi a_d)$ the cost of creating a monopole, which can be shifted by a nearest-neighbor exchange term. In spin ice ground states, $Q_\alpha \equiv 0$. Flipping a spin in a ground state yields a pair of magnetic monopoles of charges $\pm 2q_m$ on adjoining tetrahedra which can then move apart via further spin flips at finite cost in energy. Each charge has three *majority* spins that are all pointing in or out of the tetrahedron, and a single *minority* spin pointing in the opposite direction. We note that the $\pm 4q_m$ excitations are explicitly excluded from the present analysis.

With this in hand let us discuss the energetics of substituting a magnetic ion on the pyrochlore lattice by a nonmagnetic impurity (see Fig. 1). Removing a spin from a classical spin ice state [Fig. 1(a)] leaves behind two monopoles of charges $\pm q_m$ [28] [Fig. 1(b)]. A bulk monopole with charge $\pm 2q_m$ can be “emitted” by the vacancy via flipping one of the two majority spins at each of the tetrahedra adjacent to the ghost spin [Fig. 1(c)]. The bulk monopole can then move around in the system. We define a quantity that can be thought of as the ionization energy of the vacancy: $I = \frac{\mu_0 \mu^2}{2\pi a_d^3} + 2v_0 \frac{\mu^2}{a_d^2}$. This is the total energy cost of an emitted monopole moved out to infinity. Emitting another monopole into the bulk would cost

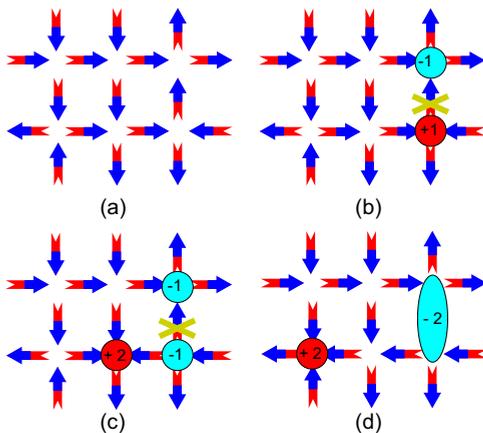


FIG. 1. (Color online) Spin ice projected onto a plane, with each vertex of the resulting square lattice in (a) corresponding to a pyrochlore tetrahedron. (b) A missing spin gives rise to a +1 and a -1 charge; (c) flipping one of the majority spins adjacent to the vacancy creates a bulk charge +2, inverting the sign of one of the vacancy charges; (d) the bulk charge propagates in the system through further spin flips, while the net charge of vacancy is approximated as a single -2 monopole.

additional energy of the order of $2v_0 \frac{\mu^2}{a_d^2}$, so the lowest energy charged excitation in the presence of an isolated vacancy is a single monopole with charge $\pm 2q_m$. Once the monopole is emitted into the bulk, it is free to hop through fluctuation-induced spin flips, while a net charge of opposite sign $\mp 2q_m$ remains at the vacancy [Fig. 1(d)].

Adding quantum dynamics via the transverse field in Eq. (1) has only a weak effect on the ground states, as connecting two of them requires flipping spins in closed loops, *minimally* six of them on a hexagon of the pyrochlore lattice. Near the classical limit, $v_0 \gg t$, such processes come with a prohibitively small energy scale, $\sim t^6/v_0^5$. By contrast, for a state containing a monopole, the lowest order effect—a monopole hopping onto a neighboring tetrahedron by flipping a majority spin—is parametrically stronger: *linear* in t !

Thus, in experiment, the most promising place to see quantum effects in spin ice is in the gauge-charged monopole excitations, rather than its gauge-neutral gapless emergent photons. Analogous considerations apply in the proximity of a vacancy, where we focus on the case of a monopole emitted into the bulk [Fig. 1(d)], also with low-order signatures. For this reason, here we perform a quantum calculation for the monopole states, and do a thermal sum over the nearly degenerate spin ice configurations.

We treat the problem as that of two Coulombic charges, one of which is stationary. As the charge propagates through the bulk, it changes the spin ice background. This process is difficult to capture exactly, but fortunately it is possible to make considerable progress via an effective model that we describe next. From this, we are able to extract the bound states in considerable detail, followed by a continuum band, much as we would expect for the hydrogen atom.

In order to investigate the problem of an isolated vacancy that has emitted a free monopole into the bulk, we switch to the *state lattice* description [29]. First, consider a new basis of the following (classical) states: a spin ice state with a vacancy, which we label $|0\rangle$, and states with an emitted monopole in the bulk, connected to $|0\rangle$ through single spin flips. Next, each site of the state lattice represents one of the basis states $|n\rangle$; while bonds connect those sites whose corresponding states are connected by single spin flips. Apart from site 0 (representing $|0\rangle$), the state lattice is trivalent. It can be shown that the smallest closed cycle in the state lattice of disordered pyrochlore spin ice has length 20 [30]. We therefore approximate the state lattice by a cycle-free infinite Cayley tree (the Bethe lattice) rooted at site 0 (Fig. 2). The monopole propagating in real space corresponds to a single particle hopping on this lattice

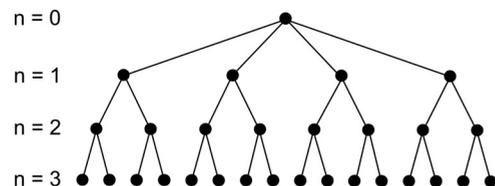


FIG. 2. The Bethe lattice describing state space. Its root, $n = 0$, corresponds to the unionized vacancy state $|0\rangle$.

in the presence of the Coulomb potential:

$$H|0\rangle = -t \sum_{m=1}^4 |0_m\rangle; \quad H|k\rangle = \left(I + \frac{C}{d_n}\right)|k\rangle - t \sum_{m=1}^3 |k_m\rangle, \quad (4)$$

where k labels a site at the n th generation of the Bethe lattice and the sums run over states reached by flipping majority spins of the monopole. d_n denotes the distance between the vacancy and the monopole in the bulk in units of a_d , such that C/d_n is the attractive Coulomb potential ($C = -\frac{\mu_0 \mu_g^2}{\pi a_d^2}$) between the two charges. In conventional spin ice, the cost of having a monopole is larger than the magnitude of the Coulomb interaction between two charges: $I > |C|$, where a_d has been set to unity. For concreteness, we use $C = -I/3$; $t = I/10$ in the following. Since we restrict ourselves to a particular starting spin ice configuration and omit other degenerate ice states from the discussion, the mapping from Eq. (1) to (4) is accurate to $O(t^5/v_0^4)$. Our final approximation concerns the distance between monopoles. Since the four sites at the first generation of the Bethe lattice correspond to the bulk monopole being one spin flip away from the vacancy, it is natural to approximate d_n in Eq. (4) by the generation of the Bethe lattice n . This definition fails to be exact already beyond $O(t^2)$, but should work sufficiently well in the $I, |C| \gg t$ regime, when the bulk monopole prefers not to move too far. In return for these approximations, we are able to solve exactly our idealized model, that of a single particle hopping on the Bethe lattice in the presence of a Coulomb potential $I + C/n$ for $n > 0$.

The Bethe lattice problem. We calculate the diagonal elements $\mathcal{G}_{ii}(\omega)$ of the lattice Green's function to infinite order in t [31,32]. We find [30] that each $\mathcal{G}_{ii}(\omega)$ can be written down in terms of a finite number of $\mathcal{G}_k^F(\omega)$, infinite sums involving particle hopping from a site at generation k to sites at generations $g > k$. The latter have a closed form expression in terms of the Gauss hypergeometric functions $F_1^2(a, b, c, z)$ [33]:

$$\mathcal{G}_k^F(\omega) = \frac{2k/\omega}{\sqrt{1+x^2} + 1} \frac{1}{k - \frac{C/\omega}{\sqrt{1+x^2}}} \times \frac{F_1^2\left(1 - \frac{C/\omega}{\sqrt{1+x^2}}, k+1, k+1 - \frac{C/\omega}{\sqrt{1+x^2}}, \frac{1-\sqrt{1+x^2}}{1+\sqrt{1+x^2}}\right)}{F_1^2\left(1 - \frac{C/\omega}{\sqrt{1+x^2}}, k, k - \frac{C/\omega}{\sqrt{1+x^2}}, \frac{1-\sqrt{1+x^2}}{1+\sqrt{1+x^2}}\right)}, \quad (5)$$

where $x^2 = -\frac{8t^2}{\omega^2}$. This yields the exact expression for any of the diagonal elements of the Green's function; for instance, at the root site

$$\mathcal{G}_{00}(\omega) = [\omega - 4t^2 \mathcal{G}_1^F(\omega - I)]^{-1}.$$

The full Green's function yields the energy levels via its poles and the local densities of states for each Bethe lattice generation, proportional to its imaginary part. The local density of states at site 0 in Fig. 3 indicates that indeed there are bound states followed by the continuum energy band. While the classical ground state (a spin ice state with a vacancy) would have zero energy, the ground state energy of the quantum problem ω_0 is lowered due to the hopping t . Low-lying excited

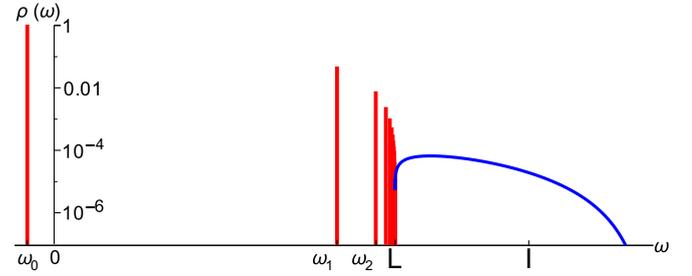


FIG. 3. (Color online) Local density of states at site 0 of the Coulomb problem on the Bethe lattice, for $C = -I/3$; $t = I/10$. Bound states (red) appear as sharp peaks; the lower edge of the continuum (blue) is labeled L , the classical ground state energy by 0, and the ionization energy by I .

states are separated from the ground state by a gap, which is also decreased from the classical value I through hopping and Coulomb attraction. They accumulate below the edge of the continuum band, located at $L = I - \sqrt{8t^2}$. In the Bethe lattice problem, the band of the extended states, of width linear in t , is confined to the region $I - \sqrt{8t^2} < \omega < I + \sqrt{8t^2}$. (Introducing closed cycles into the lattice has the effect of adding band tails, extending beyond these edges).

Signatures of monopoles in neutron scattering. One of our central results is the dynamic structure factor, defined as

$$S(\vec{q}, \Delta\omega) = \sum_f \delta(E_f - E_i - \Delta\omega) \left| \sum_{\vec{R}} \langle f | S_{\vec{R}}^+ | i \rangle e^{i\vec{q} \cdot \vec{R}} \right|^2. \quad (6)$$

In order to extract the information that is most relevant to spin ice experiments from the Bethe lattice model, we calculate a one-dimensional version of Eq. (6), averaged over all directions of \vec{q} . Such a quantity, $S(q, \Delta\omega)$, can be measured directly in a powder averaged neutron scattering experiment. The details of our calculation, carried out in the limit of dilute nonmagnetic impurities, are given in the Supplemental Material [30]. The dynamic structure factor $S(q, \Delta\omega)$, plotted in Fig. 4(a), has sharp features signaling the presence of bound states. The structure of the lines gives direct information about the character of the ground and excited states. The most visible signatures show up in elastic scattering and at the energy transfer equal to the difference between the first excited state and the ground state. For a well-localized ground state, the matrix elements between $S_{\vec{R}}^+ | i \rangle$ and excited states at higher energies (bound to the vacancy at distant radii) give rise to peaks whose structure is essentially identical up to a scale factor, as shown in Fig. 4(b). Note that for $t = 0$, the signals corresponding to $n \neq 1$ would be absent, vanishing as powers of t . Their presence thus yields direct evidence of the existence of quantum dynamics. Observing such hydrogenic bound states may well serve as the diagnostic of quantum spin ice which has thus far been lacking. An ideal compound would have a large value of total angular momentum J , and hence appreciable Coulomb splitting of the bound states, in addition to substantial transverse terms for multiple bound states to appear in neutron scattering. Moving away from the classical spin ice limit, much progress has been made very recently [34,35]. One needs substantial subleading

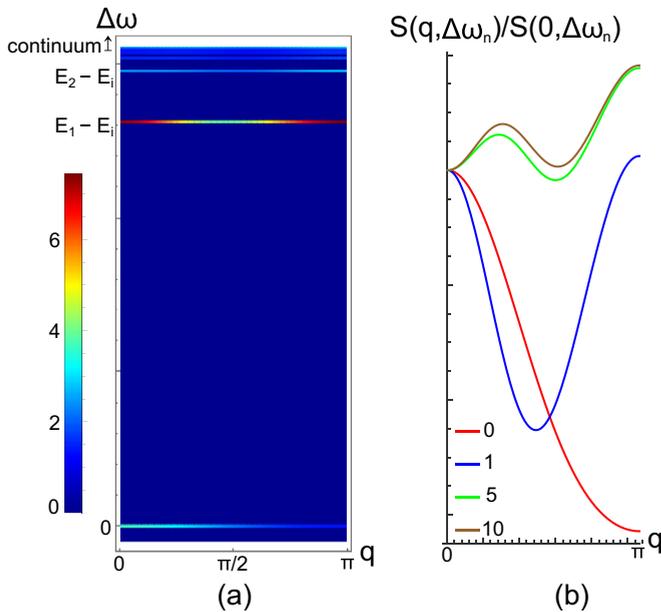


FIG. 4. (Color online) Left: Dynamic structure factor $S(q, \Delta\omega)$ for powder averaged neutron scattering. Each line is multiplied by $\exp[0.7|n - 1|]$, where n labels the all-even energy levels with $n = 0$ the ground state. Right: Line shapes $S(q, \Delta\omega_n)/S(0, \Delta\omega_n)$ for $n = 0, 1, 5, 10$. $C = -I/3$; $t = I/10$ throughout.

admixtures to the spin's crystal field ground state doublet, as suggested [34] to be the case in CdEr_2Se_4 [36]. Another origin of transverse terms is through interactions of the ion with surrounding magnetic moments [35]. In $\text{Dy}_2\text{Ti}_2\text{O}_7$, a transverse magnetic field on the order of 0.5 T [37,38] exists in the vicinity of a monopole. However, once projected onto the crystal-field ground state doublet, the Ising nature of the effective spin-1/2 still suppresses quantum fluctuations. Regarding compounds whose magnetic ions are not strictly Ising, one of the best-known quantum spin ice candidates, $\text{Yb}_2\text{Ti}_2\text{O}_7$, has an easy-plane anisotropy. While the exchange forces the magnetic moments in the (111) direction, there are

transverse terms of the same order as the (111) exchange [11]. Another easy-plane compound where a similar scenario may apply is CdDy_2Se_4 [39], whose large J may put this spinel in the right parameter range.

Conclusion and outlook. We have studied in detail the properties of magnetic monopoles in dipolar quantum spin ice. We have demonstrated that these are the prime indicators of the presence of quantum dynamics. In the presence of nonmagnetic impurities we have found both sharp hydrogenic bound states as well as a broad continuum energy band. While we believe these results to be robust, there is clearly much scope for further, presumably numerical, modeling taking into account the detailed lattice structure, as well as any material specific single-ion physics and terms in the quantum Hamiltonian.

The quantum dynamics of a pair of monopoles presents a more difficult problem due to the pair's center-of-mass motion. We are planning to address this issue, as well as clarify the detailed character of the continuum band of states in the vacancy problem, in future work. Additionally, despite neutron scattering being the method of choice for investigating magnetic materials, local disorder is an attractive subject for other types of experimental probes, such as nuclear magnetic resonance. While such techniques are beyond the scope of this work, our theoretical model can also be employed for calculating real space quantities accessible by the local probes.

Acknowledgments. The authors thank Subhro Bhattacharjee, Claudio Castelnovo, Radu Coldea, Michel Gingras, Siddharth Morampudi, Satoru Nakatsuji, Jeff Rau, Oleg Tchernyshyov, and Bruno Tomasello for useful discussions, and Yen Ting Lin and Nikos Bagis for pointing out the relevance of Ramanujan's work to the Green's function calculation. The authors acknowledge the support of the Helmholtz Virtual Institute *New States of Matter and their Excitations*, the US National Science Foundation under Grant No. DMR-1311781 (S.L.S.), the Alexander von Humboldt Foundation (O.P. and S.L.S.), and the German Science Foundation (DFG) via the Gottfried Wilhelm Leibniz Prize Programme at MPI-PKS as well as the DFG via SFB II43.

- [1] F. J. Wegner, *J. Math. Phys.* **12**, 2259 (1971).
- [2] P. Fazekas and P. W. Anderson, *Philos. Mag.* **30**, 423 (1974).
- [3] X. G. Wen and Q. Niu, *Phys. Rev. B* **41**, 9377 (1990).
- [4] X.-G. Wen, *Adv. Phys.* **44**, 405 (1995).
- [5] C. Castelnovo, R. Moessner, and S. L. Sondhi, *Nature(London)* **451**, 42 (2008).
- [6] H. R. Molavian, M. J. P. Gingras, and B. Canals, *Phys. Rev. Lett.* **98**, 157204 (2007).
- [7] J. P. C. Ruff, B. D. Gaulin, J. P. Castellan, K. C. Rule, J. P. Clancy, J. Rodriguez, and H. A. Dabkowska, *Phys. Rev. Lett.* **99**, 237202 (2007).
- [8] J. P. C. Ruff, Z. Islam, J. P. Clancy, K. A. Ross, H. Nojiri, Y. H. Matsuda, H. A. Dabkowska, A. D. Dabkowski, and B. D. Gaulin, *Phys. Rev. Lett.* **105**, 077203 (2010).
- [9] T. Fennell, M. Kenzelmann, B. Roessli, M. K. Haas, and R. J. Cava, *Phys. Rev. Lett.* **109**, 017201 (2012).
- [10] J. D. Thompson, P. A. McClarty, H. M. Rønnow, L. P. Regnault, A. Sorge, and M. J. P. Gingras, *Phys. Rev. Lett.* **106**, 187202 (2011).
- [11] K. A. Ross, L. Savary, B. D. Gaulin, and L. Balents, *Phys. Rev. X* **1**, 021002 (2011).
- [12] L. Pan, S. K. Kim, A. Ghosh, C. M. Morris, K. A. Ross, E. Kermarrec, B. D. Gaulin, S. M. Koohpayeh, O. Tchernyshyov, and N. P. Armitage, *Nat. Commun.* **5**, 4970 (2014).
- [13] K. Kimura, S. Nakatsuji, J.-J. Wen, C. Broholm, M. B. Stone, E. Nishibori, and H. Sawa, *Nat. Commun.* **4**, 1934 (2013).
- [14] R. Moessner and S. L. Sondhi, *Phys. Rev. B* **68**, 184512 (2003).
- [15] M. Hermele, M. P. A. Fisher, and L. Balents, *Phys. Rev. B* **69**, 064404 (2004).

- [16] C. L. Henley, *Ann. Rev. Condens. Matter Phys.* **1**, 179 (2010).
- [17] S. B. Lee, S. Onoda, and L. Balents, *Phys. Rev. B* **86**, 104412 (2012).
- [18] Y. Wan and O. Tchernyshyov, *Phys. Rev. Lett.* **108**, 247210 (2012).
- [19] O. Benton, O. Sikora, and N. Shannon, *Phys. Rev. B* **86**, 075154 (2012).
- [20] P. A. McClarty, O. Sikora, R. Moessner, K. Penc, F. Pollmann, and N. Shannon, [arXiv:1410.0451](https://arxiv.org/abs/1410.0451).
- [21] A. K. Ramdas and S. Rodriguez, *Rep. Prog. Phys.* **44**, 1297 (1981).
- [22] S. Sachdev, C. Buragohain, and M. Vojta, *Science* **286**, 2479 (1999).
- [23] S. Sachdev and M. Vojta, in *Proceedings of the XIII International Congress on Mathematical Physics*, edited by T. K. A. Fokas, A. Grigoryan, and B. Zegarlinski (International Press, Boston, 2000).
- [24] P. W. Anderson, *Science* **235**, 1196 (1987).
- [25] S. T. Bramwell and M. J. P. Gingras, *Science* **294**, 1495 (2001).
- [26] C. Castelnovo, R. Moessner, and S. Sondhi, *Ann. Rev. Condens. Matter Phys.* **3**, 35 (2012).
- [27] S. H. Curnoe, *Phys. Rev. B* **78**, 094418 (2008).
- [28] A. Sen and R. Moessner, *Phys. Rev. Lett.* **114**, 247207 (2015).
- [29] M. Chen, L. Onsager, J. Bonner, and J. Nagle, *J. Chem. Phys.* **60**, 405 (1974).
- [30] See Supplemental Material at <http://link.aps.org/supplemental/10.1103/PhysRevB.92.100401> for details.
- [31] W. F. Brinkman and T. M. Rice, *Phys. Rev. B* **2**, 1324 (1970).
- [32] J.-P. Gallinar, *Phys. Lett. A* **103**, 72 (1984).
- [33] B. C. Berndt, *Ramanujan's Notebooks, Part II* (Springer-Verlag, New York, 1989), p. 136.
- [34] J. G. Rau and M. J. P. Gingras, [arXiv:1503.04808](https://arxiv.org/abs/1503.04808).
- [35] B. Tomasello, C. Castelnovo, R. Moessner, and J. Quintanilla, [arXiv:1506.02672](https://arxiv.org/abs/1506.02672).
- [36] J. Lago, I. Živković, B. Z. Malkin, J. Rodriguez Fernandez, P. Ghigna, P. Dalmas de Réotier, A. Yaouanc, and T. Rojo, *Phys. Rev. Lett.* **104**, 247203 (2010).
- [37] B. Tomasello, Ph.D. thesis, University of Kent, 2014.
- [38] B. Tomasello and C. Castelnovo (private communication).
- [39] A. W. C. Wong, Z. Hao, and M. J. P. Gingras, *Phys. Rev. B* **88**, 144402 (2013).