The quantum J_1 – J'_1 – J_2 spin-1/2 Heisenberg model: influence of the interchain coupling on the ground-state magnetic ordering in two dimensions

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Abstract

We study the phase diagram of the two-dimensional (2D) J_1 – J'_1 – J_2 spin-1/2 Heisenberg model by means of the coupled cluster method. The effect of the coupling J'_1 on the Néel and stripe states is investigated. We find that the quantum critical points for the Néel and stripe phases increase as the coupling strength J'_1 is increased, and an intermediate phase emerges above the region at $J'_1 \approx 0.6$ when $J_1 = 1$. We find indications for a quantum triple point at $J'_1 \approx 0.60 \pm 0.03$, $J_2 \approx 0.33 \pm 0.02$ for $J_1 = 1$.

1. Introduction

The spin-1/2 Heisenberg antiferromagnet has been much studied in recent years due to the discovery or successful syntheses of such new magnetic materials as the layered-oxide high-temperature superconductors. Also of much interest has been the interplay between frustration and quantum fluctuations in two-dimensional (2D) quantum spin systems that can lead to quantum phase transitions between magnetically ordered semiclassical and novel quantum paramagnetic ground-state phases, see, e.g. [1, 2].

For example, the frustrated 2D antiferromagnetic J_1 – J_2 model with nearest-neighbour (J_1) and next-nearest-neighbour (J_2) bonds has attracted much attention both theoretically (see, e.g. [3–10] and references cited therein) and experimentally [11, 12]. It is now well accepted that the model exhibits two phases displaying semiclassical magnetic long-range order (LRO) at small and at large J_2 , separated by an intermediate quantum paramagnetic phase without magnetic LRO in the parameter region $J_{c1} < J_2 < J_{c2}$, where $J_{c1} \approx 0.4J_1$ and $J_{c2} \approx 0.6J_1$. The ground state (gs) for $J_2 < J_{c1}$ exhibits Néel magnetic LRO, whereas for $J_2 > J_{c2}$ it exhibits collinear stripe LRO.

In real systems deviations from the ideal 2D J_1 – J_2 model such as spin anisotropies [8, 13] or interlayer coupling may be

relevant [10]. In addition, a three-dimensional (3D) version of the J_1 – J_2 model [14] has also been considered.

An interesting generalization of the pure J_1 – J_2 model has been introduced recently by Nersesyan and Tsvelik [15]. They consider a spatially anisotropic spin-1/2 2D J_1 – J'_1 – J_2 model, where the nearest-neighbour bonds have different strengths J_1 and J'_1 in the x and y directions. This model has been further studied by other groups using the exact diagonalization (ED) of small lattice samples of $N \leq 36$ sites [16], and the continuum limit of the model [17]. Both groups support the prediction of the resonating valence-bond state by Nersesyan and Tsvelik [15] for $J_2 = 0.5J'_1 \ll J_1$, and the limit of small interchain coupling extends along a curve nearly coincident with the line where the energy is maximum. The model has also been studied by Moukouri [18] using a two-step density-matrix renormalization group approach.

Our aim here is to further the study of this model by using the coupled cluster method (CCM). The CCM (see, e.g. [19–21] and references cited therein) is one of the most powerful and universally applicable techniques of quantum many-body theory. It has been applied successfully to calculate with high accuracy the ground- and excited-state properties of many lattice quantum spin systems (see, e.g. [10, 21–26] and references cited therein). A particularly important result from

Figure 1. (a) J_1 – J_1' – J_2 model; — J_1 ; --- J_1' ; ··· J_2 ; (b) Néel state, (c) stripe state—columnar and (d) stripe state—row. Arrows in (b)–(d) represent spins situated on the sites of the square lattice (indicated by • in (a)).

our calculations is the indicated existence of a quantum triple point (QTP) at nonzero (positive) values of J_1 , J_1' and J_2 .

2. The model

The J_1 – J_1' – J_2 model is a spin-1/2 Heisenberg model on a square lattice with three kinds of exchange bonds, with strength J_1 along the row direction, J_1' along the column direction, and J_2 along the diagonals, as shown in figure 1. All exchanges are assumed positive here, and we set $J_1 = 1$. The Hamiltonian of the model is described by

$$H = J_1 \sum_{i,l} \mathbf{s}_{i,l} \cdot \mathbf{s}_{i+1,l} + J_1' \sum_{i,l} \mathbf{s}_{i,l} \cdot \mathbf{s}_{i,l+1} + J_2 \sum_{i,l} (\mathbf{s}_{i,l} \cdot \mathbf{s}_{i+1,l+1} + \mathbf{s}_{i+1,l} \cdot \mathbf{s}_{i,l+1}), \quad (1)$$

where the index (i, l) labels the x (row) and y (column) components of the lattice sites.

This model has two types of classical gs, namely, the Néel (π, π) state and stripe states (columnar stripe $(\pi, 0)$ and row stripe $(0, \pi)$), the spin orientations of which are shown in figures 1(b)–(d) respectively. There is clearly a symmetry under the interchange of rows and columns, $J_1 \rightleftharpoons J_1'$, which implies that we need only consider the range of parameters with $J_1' < J_1$. The ground-state (gs) energies of the various classical states are given by

$$\begin{aligned} \frac{E_{\text{Néel}}^{\text{cl}}}{N} &= \frac{1}{4}(-J_1 - J_1' + 2J_2), \\ \frac{E_{\text{columnar}}^{\text{cl}}}{N} &= \frac{1}{4}(-J_1 + J_1' - 2J_2), \\ \frac{E_{\text{row}}^{\text{cl}}}{N} &= \frac{1}{4}(J_1 - J_1' - 2J_2). \end{aligned} \quad (2)$$

We take $J_1 = 1$ and $J_1' < 1$. Clearly, from (2), the classical gs is then either the Néel state (if $J_1' > 2J_2$) or the columnar stripe state (if $J_1' < 2J_2$). Hence, the (first-order) classical phase transition between the Néel and columnar stripe states occurs at $J_2^c = \frac{1}{2}J_1'$, $\forall J_1 > J_1'$.

3. The coupled cluster method

The CCM formalism is now briefly described (and see [19–26] for further details). The starting point for the CCM calculation is to select a normalized model state $|\Phi\rangle$. It is often convenient to take the classical ground state as the model state for spin

Table 1. Number of fundamental LSUB n configurations ($\sharp$ f.c.) for the Néel and stripe states of the spin-1/2 J_1 – J_1' – J_2 model.

Method	$\sharp$ f.c.	
	Néel	stripe
LSUB2	2	1
LSUB4	13	9
LSUB6	146	106
LSUB8	2 555	1 922
LSUB10	59 124	45 825

systems. Hence our model states are the Néel state and the columnar stripe state. In order to treat each site identically, we perform a mathematical rotation of the local axes of the spins such that all spins in the reference state align along the negative z -axis. The Schrödinger ground-state ket and bra CCM equations are $H|\Psi\rangle = E|\Psi\rangle$ and $\langle\tilde{\Psi}|H = E\langle\tilde{\Psi}|$ respectively. The CCM employs the exponential ansatz, $|\Psi\rangle = e^S|\Phi\rangle$ and $\langle\tilde{\Psi}| = \langle\Phi|\tilde{S}e^{-S}$. The correlation operator S is expressed as $S = \sum_{I \neq 0} S_I C_I^+$ and its counterpart is $\tilde{S} = 1 + \sum_{I \neq 0} \tilde{S}_I C_I^-$. The multi-spin creation operators $C_I^+ \equiv (C_I^-)^\dagger$, with $C_0^+ \equiv 1$, are written as $C_I^+ \equiv s_{j_1}^+ s_{j_2}^+ \cdots s_{j_n}^+$. The gs energy is $E = \langle\Phi|e^{-S} H e^S|\Phi\rangle$; and the staggered magnetization M in the rotated spin coordinates is $M \equiv -\frac{1}{N} \langle\tilde{\Psi}|\sum_{j=1}^N s_j^z|\Psi\rangle$.

The ket- and bra-state correlation coefficients $(S_I, \tilde{S}_I)$ are calculated by requiring the expectation value $\bar{H} \equiv \langle\tilde{\Psi}|H|\Psi\rangle$ to be a minimum with respect to all parameters $(S_I, \tilde{S}_I)$ such that $\langle\Phi|C_I^- e^{-S} H e^S|\Phi\rangle = 0$ and $\langle\Phi|\tilde{S}(e^{-S} H e^S - E_0)C_I^+|\Phi\rangle = 0$; $\forall I \neq 0$.

4. Approximation scheme

The CCM formalism is exact if all spin configurations are included in the S and $\tilde{S}$ operators. In practice, however, truncations are needed. As in much of our previous work, we employ here the localized LSUB n scheme [6, 10, 21–26], in which all possible multi-spin-flip correlations over different locales on the lattice defined by n or fewer contiguous lattice sites are retained. The numbers of such fundamental configurations (namely, those that are distinct under the symmetries of the Hamiltonian and of the model state $|\Phi\rangle$) that are retained for the Néel and stripe states of the current model in various LSUB n approximations are shown in table 1. In order to solve the corresponding coupled sets of CCM bra- and ket-equations we use parallel computing [27].

Figure 2. Extrapolated CCM LSUB n results for the gs energy per spin, E/N , for $J_1' = 0.2, 0.4, 0.6, 0.8, 1.0$, using the Néel and stripe states of the $s = 1/2$ J_1 - J_1' - J_2 model. The LSUB n results are extrapolated in the limit $n \rightarrow \infty$ using the set $n = \{4, 6, 8, 10\}$. The NN exchange coupling $J_1 = 1$.

(This figure is in colour only in the electronic version)

5. Extrapolation scheme

In practice one needs to extrapolate the raw LSUB n data to the $n \rightarrow \infty$ limit. Based on previous experience [10, 23, 26] we use the following empirical scaling laws for the extrapolations of the gs energy per spin,

$$E/N = a_0 + a_1 n^{-2} + a_2 n^{-4}, \quad (3)$$

and the staggered magnetization of frustrated models,

$$M = b_0 + b_1 n^{-\nu}, \quad (4)$$

where the exponent ν is a fitting parameter. In order to fit to any fitting formula that contains n unknown parameters, one desirable rule is to have at least $(n + 1)$ data points ($n + 1$ rule) to obtain a robust and stable fit. In our results below the LSUB n results for $n = \{4, 6, 8, 10\}$ are extrapolated.

6. Results

In figure 2 we show the gs energy per spin as a function of J_2 . For each value of J_1' two curves are shown, one (for smaller values of J_2) using the Néel state, and the other (for larger values of J_2) using the stripe state as the CCM model state. Both sets of curves have the natural termination points [6, 21, 22] shown. For $J_1' \lesssim 0.6$ the two curves for a given value of J_1' cross (or, in the limit, meet) very smoothly near their maxima, at a value of J_2 slightly larger than the classical transition point of $0.5J_1'$. This behaviour is indicative of a second-order quantum phase transition, by contrast with the first-order classical transition from (2). For $J_1' \gtrsim 0.6$ the curves no longer cross at a physical value (namely, where the calculated staggered magnetization is positive), indicating the opening up of an intermediate quantum phase between the semiclassical Néel and stripe phases.

In figure 3(a) we show the equivalent staggered magnetization, M , for the Néel state, extrapolated using (4).

Figure 3. Extrapolated CCM LSUB n results for the gs staggered magnetization, M , for $J_1' = 0.6, 0.65, 0.7$ for the Néel state of the $s = 1/2$ J_1 - J_1' - J_2 model. (a) Results using (4), $M = b_0 + b_1 n^{-\nu}$. (b) Results using (5), $M = c_0 + n^{-0.5}(c_1 + c_2 n^{-1})$. The LSUB n results are extrapolated in the limit $n \rightarrow \infty$ using the set $n = \{4, 6, 8, 10\}$. The NN exchange coupling $J_1 = 1$.

We observe that the extrapolation scheme produces smooth and physically reasonable results, except for a very narrow anomalous ‘shoulder’ region near the points where M vanishes for $0.6 \lesssim J_1' \lesssim 0.75$. This critical regime is undoubtedly difficult to fit with the simple two-term scheme of (4). Our method for curing this problem and stabilizing the curves is to make efficient use of the information we obtain in (4) to extract the exponent ν , and then to use that value to infer the next term in the series. We find, very gratifyingly, that the value for ν fitted to (4) turns out to be very close to 0.5 for all values of J_1' and J_2 except very close to the critical point. Therefore, we use the form of (5),

$$M = c_0 + n^{-0.5}(c_1 + c_2 n^{-1}). \quad (5)$$

The use of (5) now removes the anomalous shoulder, as shown in figure 3(b). Henceforth, in all of the results we discuss, we use (5) for the staggered magnetization.

The raw LSUB n data for $n = \{2, 4, 6, 8, 10\}$ and $n = \{6, 8, 10\}$ are also extrapolated. The results for $n = \{2, 4, 6, 8, 10\}$ and $n = \{4, 6, 8, 10\}$ are similar as they both

Figure 4. The extrapolated CCM LSUB n results for the staggered magnetization, M , for $J_1' = 0.2, 0.4, 0.6, 0.8, 1.0$ of the $s = 1/2 J_1 - J_1' - J_2$ model. The LSUB n results are extrapolated in the limit $n \rightarrow \infty$ using the set $n = \{4, 6, 8, 10\}$. The NN exchange coupling $J_1 = 1$.

obey the $n + 1$ rule as mentioned above. This adds credence to the validity and stability of our results. The extrapolated curves for $n = \{6, 8, 10\}$ have a very minor ‘shoulder’ which is undoubtedly due to using only three data points to fit the three unknown terms, thus violating the $n + 1$ rule.

In figure 4 we show our equivalent results for the staggered magnetization to those in figure 2 for the gs energy. We note the surprising result that M vanishes for both the quantum Néel and stripe phases at almost exactly the same critical value of J_2 , for a given J_1' , so long as $J_1' \lesssim 0.6$. Conversely, for $J_1' \gtrsim 0.6$ there exists an intermediate region between the critical points at which $M \rightarrow 0$ for these two phases. The order parameters of both the Néel and the stripe phases vanish continuously both below and above the correspondingly indicated quantum triple point (QTP), as is again typical of second-order transitions. We note, however, that there exists some evidence from such other sources as ED calculations for the $J_1 - J_2$ model (i.e. the present model with $J_1' = J_1$) in two dimensions [5], that the transition between the stripe and intermediate phases is first-order, while others [28] have argued it may be close to second-order. We discuss these points further in section 7 below.

We show in figure 5 the zero-temperature phase diagram of the 2D spin-1/2 $J_1 - J_1' - J_2$ model, as obtained from our extrapolated results for both the gs energy and the order parameter. The phase diagram using the extrapolated LSUB n results based on $n = \{2, 4, 6, 8, 10\}$ is very similar to that of figure 5, which again adds credence to the validity and stability of our results.

Within the high-order CCM that we have used, our results certainly seem to provide clear and consistent evidence for a QTP at $J_1' \approx 0.6$ for $J_1 = 1$. For $J_1' \lesssim 0.6$ there exist only the Néel and stripe phases, with a second-order transition between them. For $J_1' \gtrsim 0.6$ there exists an intermediate (magnetically disordered, i.e. paramagnetic) quantum phase, which requires further investigation. Although the nature of the intermediate phase is still under discussion, a valence-bond solid phase seems to be the most favoured [4, 9]. On the other hand,

Figure 5. The extrapolated CCM LSUB n results for the gs phase diagram of the $s = 1/2 J_1 - J_1' - J_2$ model. The LSUB n results are extrapolated in the limit $n \rightarrow \infty$ using the set $n = \{4, 6, 8, 10\}$. The NN exchange coupling $J_1 = 1$. QTP $\equiv$ quantum triple point.

another possibility for the paramagnetic phase is the resonating valence-bond (RVB) phase [16].

7. Discussion and conclusions

In conclusion, our most important result is the evidence presented for a QTP for which our best estimate is $J_1' \approx 0.60 \pm 0.03$, $J_2 \approx 0.33 \pm 0.02$ for $J_1 = 1$. Below this point we predict a second-order phase transition between the quantum Néel and stripe phases, whereas above it these two phases are separated by an intermediate phase. Although the order parameters for the Néel and stripe phases vanish on the Néel-intermediate and stripe-intermediate phase boundary lines, respectively, above the QTP, we are unable to conclude more about the nature of the transitions at those boundaries, since the present calculations have not addressed at all the nature of the intermediate phase.

Other calculations on this model [16, 17] differ predominantly by giving a QTP at $J_1' = 0 = J_2$ for $J_1 = 1$. We believe that the difference arises essentially from the nature of the alternative methods used. For example, due to the small size of the lattices used, the ED calculations of [16] might miss the longer-range correlations that become increasingly important the nearer one approaches the QTP. By contrast, at *any* level of truncation the CCM always incorporates some long-range correlations through the important exponentiated parametrization of the wavefunction that lies at the heart of the method, as described in section 3. Also, the results of Starykh and Balents [17] are based on an analysis of a continuum version of the $J_1 - J_1' - J_2$ model rather than on the discrete lattice model itself. This could easily account for our differing predictions for the position of the QTP.

As we have discussed above (and see figures 4 and 5), our results indicate that there is a continuous (second-order) transition below the QTP between the two ordered semiclassical phases with different broken symmetries. This is undoubtedly a rather novel and surprising result that has few precedents, and one that seemingly violates the Landau

criterion of symmetry change. We note, however, that a similar scenario has been discussed recently in the related, but different, context of deconfined quantum criticality, as we explain more fully below.

Thus, it has been argued recently by other authors [29] that there also exists a continuous (i.e. second-order) phase transition for the spin-1/2 pure J_1 - J_2 model on the 2D square lattice (i.e. our case with $J'_1 = J_1$) between the Néel state and what we have called here the intermediate state, and which those authors identify as a paramagnetic valence-bond solid (VBS) state. Such direct second-order quantum phase transitions between two states with different broken symmetries, and hence characterized by two seemingly independent order parameters, are difficult to understand within the standard Ginzburg–Landau critical theory, as we now discuss.

Thus, the competition between two such distinct kinds of quantum order associated with different broken symmetries would lead generically in the Ginzburg–Landau scenario to only one of three possibilities: (i) a first-order transition between the two states, (ii) an intermediate region of co-existence between both phases with both kinds of order present or (iii) a region of intermediate phase with neither of the orders of these two phases present. A direct second-order transition between states of different broken symmetries is only allowable within the standard Ginzburg–Landau critical theory if it arises by an accidental fine-tuning of the disparate order parameters to a multicritical point. Thus, for the spin-1/2 pure J_1 - J_2 model and its quantum phase transition suggested by Senthil *et al* [29], it would require the completely accidental coincidence (or near coincidence) of the point where the magnetic order parameter (i.e. the staggered magnetization) vanishes for the Néel phase with the point where the dimer order parameter vanishes for the VBS phase. Since each of these phases has a different broken symmetry (namely, spin-rotation symmetry for the Néel phase and lattice symmetry for the VBS phase), one would naively expect both that each transition is described by its own *independent* order parameter (i.e. the staggered magnetization for the Néel phase and the dimer order parameter for the VBS phase) and that the two transitions should therefore be *independent* of each other.

By contrast, the ‘deconfined’ type of quantum phase transition postulated by Senthil *et al* [29] permits direct second-order quantum phase transitions between such states with different forms of broken symmetry. In their scenario the quantum critical points still separate phases characterized by order parameters of the conventional (i.e. in their language, ‘confining’) kind, but their proposed new critical theory involves fractional degrees of freedom (namely, spinons for the spin-1/2 J_1 - J_2 model on the 2D square lattice) that interact via an emergent gauge field. For the specific example of the spin-1/2 J_1 - J_2 model, the order parameters of both the Néel and VBS phases discussed above are represented in terms of the spinons, which themselves become ‘deconfined’ exactly at the critical point. That the spinons are the fundamental constituents of *both* order parameters then offers a natural explanation for the direct second-order phase transition between two states of the system that otherwise seem very different on the basis of their broken symmetries.

Despite the compelling nature of the arguments posited by Senthil *et al* [29], we should mention, however, that other authors believe the phase transition in the J_1 - J_2 model not to be due to a deconfinement of spinons. For example, Sarker *et al* [9] have given arguments, based both on numerical results from series expansion analyses and on spin-wave theory, that the spin-1/2 J_1 - J_2 transition is not of the above second-order deconfined type but, in their view, is more likely to be a (weakly) first-order transition between the Néel phase and a VBS phase with columnar dimerization.

One should also note that other, less radical, mechanisms have also been proposed to explain such second-order phase transitions and their seeming disagreement (except by accidental fine-tuning) with Ginzburg–Landau theory. What seems clearly to be minimally required is that the order parameters of the two phases with different broken symmetry should be related in some way. Thus, a Ginzburg–Landau-type theory can only be preserved if it contains additional terms in the effective theory that represent interactions between the two order parameters. For example, just such an effective theory has been proposed for the 2D spin-1/2 J_1 - J_2 model on the square lattice by Sushkov *et al* [30], and further discussed in [9].

Whether or not the deconfined phase transition theory of Senthil *et al* [29] survives the controversy that still surrounds it, we note that one of the motivations that led to it was the existence of various other numerical calculations in recent years that also point to a direct second-order quantum phase transition between phases of different broken symmetry that are characterized by seemingly independent order parameters. Examples include the quantum phase transitions between: (i) an antiferromagnetic Mott insulator and a $d_{x^2-y^2}$ superconductor in a 2D Hubbard model on a square lattice, based on quantum Monte Carlo (QMC) simulations on lattices up to size 16×16 [31]; and (ii) superfluid and stripe-order phases in a 2D square lattice spin-1/2 XY model with both a nearest-neighbour coupling term (J) and a four-spin ring exchange term (K), based on a QMC method (namely, a stochastic series expansion technique) with lattices up to size 64×64 [32]. While most previous such numerical evidence for continuous (second-order) quantum phase transitions between states with different broken symmetry has come from QMC simulations, the well-known ‘sign problem’ inherent to QMC techniques has meant that it has not been easy to apply that method to frustrated spin-lattice systems of the type considered here. We believe that the use of the CCM for such systems, as reported here, opens up a new arena and sheds a new spotlight on this fascinating and still unresolved larger field.

Returning to our own results for the spin-1/2 J_1 - J'_1 - J_2 model on the 2D square lattice, of course, one may also argue that what we have observed as a continuous (second-order) transition below the QTP may in reality be a very weak first-order transition, which is thereby not in violation of the Landau symmetry change criterion. Our completely independent calculations for the two semiclassical phases can, obviously, never entirely preclude this latter possibility. Nevertheless, it is clear from our results (and see figures 2 and 4) that the data

below the QTP are fully consistent only with a transition which if it is not continuous is *very* weakly first-order for *all* values $0 \leq J'_1 \lesssim 0.60$ below the QTP.

Finally, one could also argue that in this same range there might exist a very narrow strip of intermediate quantum disordered phase, which would then reconcile our results (at least qualitatively) with those from the exact diagonalization of small clusters by Sindzingre [16]. While this possibility, again, can never be entirely ruled out by any numerical calculation such as ours, we have clearly demonstrated that our own extrapolation schemes are both robust and internally consistent enough to rule out any but an *extremely* narrow strip of intermediate disordered phase for $0 \leq J'_1 \lesssim 0.60$.

We end by noting that two of the unique strengths of the CCM are its ability to deal with highly frustrated systems as easily as unfrustrated ones, and its use from the outset of infinite lattices, which leads in turn to its ability to yield accurate phase boundaries even near a possible QTP. Our own results for the ground-state energy and staggered magnetization provide a set of independent checks that lead us to believe that we have a self-consistent and coherent description of this extremely challenging system. We also believe that the present high-order CCM results are among the numerically most accurate for this and related spin–lattice models containing frustration. Nevertheless, our suggested novel phase scenario certainly needs further confirmation by the application of alternative high-order methods. It would also be of considerable interest to repeat the investigation for the computationally more challenging case of the same system for spin-1 particles, and we intend to report on this system in the future.

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