

The Effects of Three Magnons Interactions in the Magnon-Density Waves of Triangular Spin Lattices

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Abstract

We investigate the magnon-density waves proposed as the longitudinal excitations in triangular lattice antiferromagnets by including the cubic and quartic corrections in the large-s expansion. The longitudinal excitation spectra for the two-dimensional (2D) triangular antiferromagnetic model and quasi-one-dimensional (quasi-1D) antiferromagnetic materials have been obtained for a general quantum spin number s. For the 2D triangular lattice model, we find a significant reduction (about 40%) in the energy spectra at the zone boundaries due to both the cubic and quartic corrections. For the quasi-1D antiferromagnets, since the cubic term comes from the very weak couplings on the hexagonal planes, they make very little correction to the energy spectra, whereas the major correction contribution comes from the quartic terms in the couplings along the chains with the numerical values for the energy gaps in good agreement with the experimental results as reported earlier (Merdan and Xian in Phys Rev B 87:174434, 2013).

Keywords Quantum antiferromagnets · Longitudinal excitation · Spin-wave excitation

1 Introduction

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Since Haldane [1] predicted difference between the excitations of integer spin and half-odd-integer spin chains, the nature of excitations of quantum Heisenberg antiferromagnets has attracted both experimental and theoretical attentions. In particular, for the spin-1 chains, the singlet ground state is separated from the triplet excitation

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states by an energy gap. This theoretical prediction has been confirmed experimentally in the quasi-1D spin-1 antiferromagnetic compounds such as CsNiCl₃ and RbNiCl₃ [2]. Haldane's conjecture was also supported by some other experiments [2–6] and theoretical studies [7–11]. Furthermore, a longitudinal excitation has been proposed by Affleck for explanation of a gapped excitation mode observed in very low temperature in these quasi-1D hexagonal antiferromagnetic compounds CsNiCl₃ and RbNiCl₃ which possess Néel order at low temperature [12,13]. This longitudinal mode describes the fluctuations of the long-range order parameter and is beyond the spin-wave theory (SWT) which predicts only the transverse spin-wave excitations (magnons).

On the other hand, the triangular lattice Heisenberg antiferromagnet is the prototype system of geometrically frustrated magnets and has been under intensive investigation for fundamentally different types of ground and excited states [14–16]. It is now widely accepted that the ground state of the antiferromagnet on a triangle lattice has the long-range noncollinear Néel-like order with the 120° magnetic three-sublattice structure as predicted by various methods [17,18], including a SWT based on three sublattices [19-29]. The interaction between spin-wave excitations in antiferromagnetic materials of collinear spin configuration is depicted by higher-order anharmonicities beginning with the quartic term [30,31]. The higher-order anharmonicities of antiferromagnetic systems with noncollinear spin configuration begin with the cubic term which describes the coupling between transverse (one-magnon) and longitudinal (twomagnon) fluctuations [32,33], in addition to the quartic term. This cubic term is similar to those that describe the interaction between one- and two-particle states of phonons in crystals [34] and excitations in superfluid bosonic systems [35]. In noncollinear antiferromagnets, the cubic term comes from products of the spin operator components S^z and $S^{x,y}$, which are not present in collinear lattices. For the correction in spin-wave spectrum, the cubic term has been included in perturbation theory and represents the coupling of the transverse fluctuations in one sublattice to the longitudinal ones in the others [22,26,32,33,36-38].

For a generic quantum spin-s antiferromagnetic Hamiltonian system with a Néel order, a microscopic theory of the longitudinal modes has been proposed [39]. In this theory, the longitudinal excitations are identified as the collective modes of the magnon-density waves, and the corresponding wave functions are constructed by employing the magnon-density operator S^z in similar fashion to Feynman's theory on the low-lying excited states of the helium-4 superfluid where the particle density operator is used [40]. In our earlier calculations for the quasi-1D hexagonal structures of CsNiCl₃ and RbNiCl₃ and tetragonal structure of KCuF₃, we find that, after inclusion of the higher-order contributions from the quartic terms in the large-s expansion, the energy gap values at the magnetic wavevector are in good agreement with experimental results [41,42].

Although there is no report of direct experimental observations of longitudinal modes in 2D triangle antiferromagnetic lattices, a theoretical investigation of dynamic structure factors does find some broad peaks in the two-magnon continuum and a massive contribution from the longitudinal fluctuations to the high energy spectral weight, clearly indicating the strong magnon-magnon interactions in the system [32]. In this article, we extend our preliminary investigation of the longitudinal modes in the 2D triangle antiferromagnetic model [43], focusing now on the higher-order calculations



by including both the cubic and quartic terms. Our results show a significant reduction in the energy spectra due to the higher-order corrections. We also examine the cubic term contribution to the energy spectrum correction for the quasi-1D hexagonal systems of CsNiCl₃ and RbNiCl₃, not considered in our earlier study [41]. We find that in these systems, the cubic term contribution is negligible, mainly due to the very weak coupling on the triangular planes of the systems.

We organize this article as follows. Section 2 outlines the main results of the spin-wave theory for the triangular lattice model using the bosonization approach. In Sects. 3 and 4, we review our microscopic theory for the longitudinal excitations, including the higher-order corrections from the cubic and quartic terms and using the approximated ground state from SWT, and apply to the 2D triangular antiferromagnetic model. In Sect. 5, we re-examine our calculation of the higher-order corrections in the quasi-1D hexagonal systems where there are several experimental results for comparison. We notice that the energy gap changes very little after the inclusion of the cubic contribution, mainly due to the small coefficient for the plane Hamiltonian when compared with the coefficient of the perpendicular (chain) Hamiltonian. In Sect. 6, we conclude this article by a summary and a critical discussion of the longitudinal modes in 2D triangle lattices.

2 Spin-Wave Formalism for Triangular Lattice Model

The Heisenberg antiferromagnet on a triangular lattice is described by Hamiltonian with spin operator S,

$$H = J \sum_{\langle i,j \rangle} \mathbf{S}_i \cdot \mathbf{S}_j, \tag{1}$$

where J>0 is the coupling parameter and the sum on $\langle i,j\rangle$ runs over all the nearest-neighbor pairs of the triangular lattice once. The classical ground state of the antiferromagnetic Heisenberg model on a triangular lattice consists of three sublattices where the direction of each spin on one sublattice forms an angle of 120° from those on the other two sublattices. We choose the direction of classical orientation in the xz-plane at the one i-sublattice surrounded by six j-sublattices. The Hamiltonian of Eq. (1) can be transformed into a rotating local basis as

$$S_i^x \to S_i^x \cos \theta_i + S_i^z \sin \theta_i,$$

$$S_i^y \to S_i^y,$$

$$S_i^z \to S_i^z \cos \theta_i - S_i^x \sin \theta_i,$$
(2)

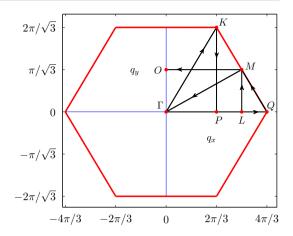
where $\theta_i = \mathbf{Q} \cdot \mathbf{r}_i$ and $\mathbf{Q} = (4\pi/3, 0)$ is the magnetic ordering wavevector of the hexagonal Brillouin zone of the triangular lattice as shown in Fig. 1.

The Hamiltonian operator of Eq. (1) after this transformation is given by

$$H = J \sum_{\langle i,j \rangle} \left[\cos(\theta_i - \theta_j) (S_i^x S_j^x + S_i^z S_j^z) + \xi S_i^y S_j^y + \sin(\theta_i - \theta_j) (S_i^z S_i^x - S_i^x S_j^z) \right],$$
(3)



Fig. 1 The hexagonal first Brillouin zone of a triangular lattice in reciprocal space. The coordinates of the labeled points are $\Gamma=(0,0)$, $P=(2\pi/3,0)$, $L=(\pi,0)$, $Q=(4\pi/3,0)$, $M=(\pi,\pi/\sqrt{3})$, $K=(2\pi/3,2\pi/\sqrt{3})$ and $O=(0,\pi/\sqrt{3})$ (Color figure online)



where we have also introduced an anisotropy parameter $\xi (\leq 1)$ along the y-axis. The Holstein–Primakoff transformation which transforms spin operators into bosons is used for the spin-wave calculations such that

$$S_i^z = s - a_i^{\dagger} a_i, \quad S_i^+ = \sqrt{2s} f_i a_i, \quad S_i^- = \sqrt{2s} a_i^{\dagger} f_i,$$
 (4)

where $f_i = \sqrt{1 - a_i^{\dagger} a_i/2s}$, s is the spin quantum number and $S_i^{\pm} = S_i^x \pm i S_i^y$. Substituting Eqs. (4) into (3) and approximating the expansion of the square root in f_i to the first order in $a_i^{\dagger} a_i/2s$, we obtain the following Hamiltonian

$$H = H_0 + H_1 + H_2 + H_3 + H_4, (5)$$

where $H_0 = -3/2JNs^2$ is the classical ground-state energy $\mathcal{O}(s^2)$, H_2 is the harmonic part of the linear SWT (LSWT) correction $\mathcal{O}(s)$, H_3 is the cubic anharmonic term $\mathcal{O}(s^{1/2})$ and H_4 is the quartic anharmonic term $\mathcal{O}(s^0)$. The LSWT depicts the harmonic approximation or noninteracting magnons. The quadratic terms in H_2 can be written as

$$H_{2} = \frac{1}{4} Js \sum_{\langle i,j \rangle} \left[2(n_{i} + n_{j}) - (1 + 2\xi)(a_{i}a_{j} + a_{i}^{\dagger}a_{j}^{\dagger}) - (1 - 2\xi)(a_{i}a_{j}^{\dagger} + a_{i}^{\dagger}a_{j}) \right], \tag{6}$$

where $n_i = a_i^{\dagger} a_i$ and $n_j = a_j^{\dagger} a_j$ are number operators. After Fourier transformation for the boson operators with the Fourier component operators $a_{\bf q}$ and $a_{\bf q}^{\dagger}$, and performing the diagonalization of H_2 by the canonical Bogoliubov transformation, $a_{\bf q} = u_{\bf q} \alpha_{\bf q} + v_{\bf q} \alpha_{\bf -q}^{\dagger}$, the linear spin-wave Hamiltonian now reads



$$H' = H_0 + H_2 = -\frac{3}{2}JNs(s+1) + \sum_{\mathbf{q}} \mathcal{E}_{\mathbf{q}} \left(\alpha_{\mathbf{q}}^{\dagger} \alpha_{\mathbf{q}} + \frac{1}{2} \right), \tag{7}$$

where $\mathcal{E}_{\mathbf{q}} = 3Js\,\omega_{\mathbf{q}}$ is the spin-wave excitation spectrum with the dimensionless spectrum $\omega_{\mathbf{q}}$ given by

$$\omega_{\mathbf{q}} = \sqrt{A_{\mathbf{q}}^2 - B_{\mathbf{q}}^2} = \sqrt{(1 - \gamma_{\mathbf{q}})(1 + 2\xi\gamma_{\mathbf{q}})},$$
 (8)

with $A_{\mathbf{q}}$ and $B_{\mathbf{q}}$ defined by

$$A_{\mathbf{q}} = 1 + \left(\xi - \frac{1}{2}\right)\gamma_{\mathbf{q}}, \quad B_{\mathbf{q}} = \left(\xi + \frac{1}{2}\right)\gamma_{\mathbf{q}},$$
 (9)

respectively, and γ_q defined by

$$\gamma_{\mathbf{q}} = \frac{1}{z} \sum_{\varrho} e^{i\mathbf{q} \cdot \mathbf{r}_{\varrho}} = \frac{1}{3} \left(\cos q_x + 2 \cos \frac{q_x}{2} \cos \frac{\sqrt{3}}{2} q_y \right), \tag{10}$$

with the summation over the nearest-neighbor index ϱ and the coordination number z=6 for the triangular lattice.

The cubic term exists in the triangular lattice because of the coupling of S^z and S^x spin components. In terms of boson operators, the cubic term reads

$$H_3 = J\sqrt{\frac{s}{2}} \sum_{\langle i,j \rangle} \sin(\theta_i - \theta_j) \left[(a_i + a_i^{\dagger}) n_j - n_i (a_j + a_j^{\dagger}) \right]. \tag{11}$$

We notice that for the collinear spin lattices, $\sin(\theta_i - \theta_j) = 0$ and the cubic terms vanish and that H_1 with one boson terms always cancel out. Furthermore, the LSWT ground-state expectation value of the three-boson operators is always zero. This cubic term has been included in the perturbation theory with the contribution of order $1/s^2$. In more details, after performing Fourier and Bogoliubov transformations, we obtain

$$H_{3} = Jzi\sqrt{\frac{3s}{8N}} \sum_{\mathbf{q},\mathbf{k}} \left[\frac{1}{2!} F_{1}(\mathbf{q},\mathbf{k})_{\mathbf{q}\mathbf{k}-\mathbf{q}\mathbf{k}}^{\dagger\dagger} + \frac{1}{3!} F_{2}(\mathbf{q},\mathbf{k}) \alpha_{\mathbf{q}}^{\dagger} \alpha_{\mathbf{k}+\mathbf{q}}^{\dagger} \alpha_{\mathbf{k}}^{\dagger} + \text{H.c.} \right], \tag{12}$$

with $F_1(\mathbf{q}, \mathbf{k})$ and $F_2(\mathbf{q}, \mathbf{k})$ given by

$$F_{1}(\mathbf{q}, \mathbf{k}) = \bar{\gamma}_{\mathbf{q}}(u_{\mathbf{q}} + v_{\mathbf{q}})(u_{\mathbf{k}}u_{\mathbf{q}-\mathbf{k}} + v_{\mathbf{k}}v_{\mathbf{q}-\mathbf{k}}) + \bar{\gamma}_{\mathbf{k}}(u_{\mathbf{k}} + v_{\mathbf{k}})(u_{\mathbf{q}}u_{\mathbf{q}-\mathbf{k}} + v_{\mathbf{q}}v_{\mathbf{q}-\mathbf{k}}) - \bar{\gamma}_{\mathbf{q}-\mathbf{k}}(u_{\mathbf{q}-\mathbf{k}} + v_{\mathbf{q}-\mathbf{k}})(u_{\mathbf{q}}u_{\mathbf{k}} + v_{\mathbf{q}}v_{\mathbf{k}}),$$
(13)



and

$$F_{2}(\mathbf{q}, \mathbf{k}) = \bar{\gamma}_{\mathbf{q}}(u_{\mathbf{q}} + v_{\mathbf{q}})(u_{\mathbf{k}}v_{\mathbf{q}+\mathbf{k}} + v_{\mathbf{k}}u_{\mathbf{q}+\mathbf{k}})$$

$$+ \bar{\gamma}_{\mathbf{k}}(u_{\mathbf{k}} + v_{\mathbf{k}})(u_{\mathbf{q}}v_{\mathbf{q}+\mathbf{k}} + v_{\mathbf{q}}u_{\mathbf{q}+\mathbf{k}})$$

$$- \bar{\gamma}_{\mathbf{q}+\mathbf{k}}(u_{\mathbf{q}+\mathbf{k}} + v_{\mathbf{q}+\mathbf{k}})(u_{\mathbf{q}}v_{\mathbf{k}} + v_{\mathbf{q}}u_{\mathbf{k}}), \tag{14}$$

where u_i and v_i are Bogoliubov parameters and the function $\bar{\gamma}_{\mathbf{q}}$ is given by

$$\bar{\gamma}_{\mathbf{q}} = \frac{1}{3} \left(\sin q_x - 2 \sin \frac{q_x}{2} \cos \frac{\sqrt{3}}{2} q_y \right). \tag{15}$$

The first term in Eq. (12) is called "decay" which describes the interaction between one- and two-magnon states, and it is symmetric under permutation of two outgoing momenta. The second term is called "source," and it is symmetric under permutation of three outgoing momenta [44]. The $1/s^2$ contribution from the second-order perturbation of H_3 is evaluated by Miyake [22,36] such that

$$\delta E_3 = -\frac{z^2 J^2 s}{16N} \sum_{\mathbf{q}, \mathbf{k}} \frac{F_2(\mathbf{q}, \mathbf{k})^2}{\mathcal{E}_{\mathbf{q}} + \mathcal{E}_{\mathbf{k}} + \mathcal{E}_{\mathbf{q} + \mathbf{k}}}.$$
 (16)

The quartic anharmonic term in Eq. (5) reads

$$H_4 = \frac{1}{4} J \sum_{\langle i,j \rangle} \left[-n_i n_j + \frac{1}{4} (1 + 2\xi)(n_i + n_j) a_i a_j + (1 - 2\xi) \{ a_j^{\dagger} (n_i + n_j) a_i + \text{H.c.} \right].$$
 (17)

For simplicity, we define the following Hartree–Fock averages (the LSWT ground-state expectation values) of the triangular lattice

$$\rho = \langle a_l^{\dagger} a_l \rangle = \frac{1}{N} \sum_{\mathbf{q}} \rho_{\mathbf{q}}, \quad \mu_{\varrho} = \langle a_l^{\dagger} a_{l+\varrho} \rangle = \frac{1}{N} \sum_{\mathbf{q}} e^{i\mathbf{q}\cdot\varrho} \rho_{\mathbf{q}},
\Delta_{\varrho} = \langle a_l a_{l+\varrho} \rangle = \frac{1}{N} \sum_{\mathbf{q}} e^{i\mathbf{q}\cdot\varrho} \Delta_{\mathbf{q}}, \quad \delta = \langle a_l a_l \rangle = \frac{1}{N} \sum_{\mathbf{q}} \Delta_{\mathbf{q}},$$
(18)

with $\Delta_{\mathbf{q}}$ and $\rho_{\mathbf{q}}$ defined as

$$\Delta_{\mathbf{q}} = \frac{1}{2} \frac{B_{\mathbf{q}}}{\sqrt{A_{\mathbf{q}}^2 - B_{\mathbf{q}}^2}}, \quad \rho_{\mathbf{q}} = \frac{1}{2} \left(\frac{A_{\mathbf{q}}}{\sqrt{A_{\mathbf{q}}^2 - B_{\mathbf{q}}^2}} - 1 \right).$$
(19)

The ground-state expectation value of the quartic H_4 of Eq. (17) can be calculated first by applying Fourier transformation and then Bogoliubov transformation using Wick's



theorem. The ground-state energy correction in terms of the Hartree–Fock averages is given by

$$\delta E_4 = -\frac{1}{4} J N z \left[\rho^2 + \mu_{\varrho}^2 + \Delta_{\varrho}^2 - (1 + 2\xi) \left(\rho \Delta_{\varrho} + \frac{1}{2} \mu_{\varrho} \delta \right) - (1 - 2\xi) \left(\rho \mu_{\varrho} + \frac{1}{2} \Delta_{\varrho} \delta \right) \right]. \tag{20}$$

Thus, the total ground-state energy can be calculated from all these contributions for the isotropic case $\xi = 1$ as

$$E = -\frac{1}{4}JNzs^{2} \left[1 + \frac{I_{2}}{s} + \frac{(I_{3} + I_{4})}{(2s)^{2}} \right], \tag{21}$$

where I_2 is related to harmonic part H_2 with numerical value given by

$$I_2 = 1 - \frac{1}{N} \sum_{\mathbf{q}} \omega_{\mathbf{q}} = 0.218412,$$
 (22)

and the other constants I_3 and I_4 are related to δE_3 and δE_4 , respectively, with numerical values calculated at $\xi = 1$

$$I_3 = \frac{2}{N^2} \sum_{\mathbf{q}, \mathbf{k}} \frac{F_2(\mathbf{q}, \mathbf{k})^2}{\omega_{\mathbf{q}} + \omega_{\mathbf{k}} + \omega_{\mathbf{q} + \mathbf{k}}} = 0.2756(2), \tag{23}$$

$$I_{4} = 4\left(\rho^{2} + \mu_{\varrho}^{2} + \Delta_{\varrho}^{2} - 3(\rho\Delta_{\varrho} + \frac{1}{2}\mu_{\varrho}\delta) + \left(\rho\mu_{\varrho} + \frac{1}{2}\Delta_{\varrho}\delta\right)\right) = -0.25429.$$
(24)

These numerical results have been obtained by Miyake [22]. The integration of I_3 is a four-dimensional integral and has been calculated by Monte Carlo integration using Mathematica software.

The sublattice magnetization M in general can be written in terms of the magnon-density ρ as

$$M = s - \rho. \tag{25}$$

Within the linear spin-wave approximation, ρ is given by $\rho_0 = 0.261303$. The higher-order correction to the sublattice magnetization can be expressed as $M = s - \rho_0 + \frac{\delta s_2}{2s}$. Miyake first calculated $\delta s_2 = 0.0110$ [22], later confirmed by Chernyshev and Zhitomirsky [44] using a different method. But Chubukov [26] obtained $\delta s_2 = 0.027$, perhaps because of an integration problem.



3 Longitudinal Excitations Formalism

In antiferromagnetic quantum systems with a Néel-like long-range order, the longitudinal excitations correspond to the fluctuations in the order parameter. We identify the longitudinal modes as the magnon-density waves (MDW), well defined only in the systems where the interactions between the transverse magnons are significant. In the low-dimensional systems, the magnon density is high enough to support the longitudinal waves, such as the cases of the quasi-1D systems mentioned in Sect. 1. It remains questionable whether or not the interaction between magnons in pure 2D systems such as the triangle antiferromagnet is strong enough to support the longitudinal modes, although there is some indication this may be so [32].

The magnon-density operator is given by spin operator S^z , and so S^z can be used to construct the wave function of longitudinal excitation state in a similar fashion as Feynman's theory of the phonon–roton excitation state of the helium superfluid, where the density operator is the usual particle density operator [40,45]. The longitudinal excitation state is thus constructed by applying the magnon-density fluctuation operator X_q on the ground-state $|\Psi_g\rangle$ as

$$|\Psi_e\rangle = X_{\mathbf{q}}|\Psi_g\rangle,\tag{26}$$

where $X_{\mathbf{q}}$ is given in terms of the Fourier transformation of S^z operator as,

$$X_{\mathbf{q}} = \frac{1}{\sqrt{N}} \sum_{l} e^{i\mathbf{q} \cdot \mathbf{r}_{l}} S_{l}^{z}, \quad q > 0, \tag{27}$$

with index l running over all lattice sites. The condition q > 0 in Eq. (27) guarantees the orthogonality to the ground state. The energy spectrum of longitudinal excitation is given by [46]

$$E(\mathbf{q}) = \frac{N(\mathbf{q})}{S(\mathbf{q})},\tag{28}$$

where $N(\mathbf{q})$ is the ground-state expectation value of a double commutator such that

$$N(\mathbf{q}) = \frac{1}{2} \langle [X_{-\mathbf{q}}, [H, X_{\mathbf{q}}]] \rangle_g, \tag{29}$$

and $S(\mathbf{q})$ is the normalization integral or the structure factor of the lattice model

$$S(\mathbf{q}) = \langle \Psi_e | \Psi_e \rangle = \frac{1}{N} \sum_{l,l'} e^{i\mathbf{q}.(\mathbf{r}_l - \mathbf{r}_{l'})} \langle S_l^z S_{l'}^z \rangle_g.$$
 (30)

We apply the SWT for the approximation of the ground-state $|\Psi_g\rangle$ in the following sections to calculate these expectation values.



4 Magnon-Density Waves in 2D Triangular Lattice

The one-sublattice Hamiltonian of Eq. (3) after the rotation for the triangular lattice is used to obtain the double commutator of Eq. (29) as

$$N(\mathbf{q}) = \frac{1}{4}sJ \sum_{\rho} \left[(1 + 2\xi)(1 + \gamma_q)\tilde{g}_{\varrho} + (1 - 2\xi)(1 - \gamma_q)\tilde{g}_{\varrho}' + \frac{I_3}{8s} \right], \quad (31)$$

where I_3 contains cubic terms and is defined in Eq. (23), and the transverse correlation functions \tilde{g}_{ϱ} and \tilde{g}'_{ϱ} are defined as

$$\tilde{g}_{\varrho} = \frac{1}{2s} \langle S_{l}^{+} S_{l+\varrho}^{+} \rangle_{g}, \quad \tilde{g}'_{\varrho} = \frac{1}{2s} \langle S_{l}^{+} S_{l+\varrho}^{-} \rangle_{g}.$$
 (32)

Due to the lattice translational symmetry, both correlation functions are independent of index l. These functions contain the contribution from quadratic and quartic terms and both given in terms of the Hartree–Fock averages of Eq. (18) as

$$\tilde{g}_{\varrho} = \Delta_{\varrho} - \frac{2\rho \, \Delta_{\varrho} + \mu_{\varrho} \, \delta}{2s},$$

$$\tilde{g}'_{\varrho} = \mu_{\varrho} - \frac{2\rho \, \mu_{\varrho} + \Delta_{\varrho} \delta}{2s}.$$
(33)

We obtain the numerical results at the isotropic point $\xi=1$ as $\tilde{g}_{\varrho}=0.12598$ and $\tilde{g}'_{\varrho}=0.03642$ for all the six nearest neighbors. As it can be seen, $N(\mathbf{q})$ is dominated by \tilde{g}_{ϱ} . The structure factor is independent of s, and is given by

$$S(\mathbf{q}) = \rho + \frac{1}{N} \sum_{\mathbf{q}'} \rho_{\mathbf{q}'} \rho_{\mathbf{q}+\mathbf{q}'} + \frac{1}{N} \sum_{\mathbf{q}'} \Delta_{\mathbf{q}'} \Delta_{\mathbf{q}+\mathbf{q}'}.$$
 (34)

We notice that the calculations of both Eqs. (33) and (34) involve up to four boson operators of the quartic terms, but not the cubic term. We then calculate the longitudinal excitation spectrum $E(\mathbf{q})$ given by Eq. (28). From the numerical calculation, we found that this spectrum of the longitudinal mode is gapless in the thermodynamic limit, as $E(\mathbf{q}) \to 0$ at both $\mathbf{q} \to 0$ and $\mathbf{q} \to \pm \mathbf{Q}$. Two longitudinal modes for the triangular lattice antiferromagnets due to the noncollinear nature of the Néel-like order can be obtained by folding of the wavevector. We denote one as L_{+} with the spectrum $E(\mathbf{q} + \mathbf{Q})$ and the other as L_{-} with the spectrum $E(\mathbf{q} - \mathbf{Q})$. We plot both spectra at isotropic case $\xi = 1$ in Figs. 2 and 3. We find that the energy values of the spectra reduce significantly by about 40% at the zone boundaries after the inclusion of the quartic and cubic corrections, and that the two longitudinal modes are nearly degenerate, only differ by a few percents on the zone boundaries. For example, the L_{+} energy value at P is 0.6545zs J in the first-order calculation, reduces to 0.3829zs J after including the higher-order corrections with the cubic contribution of 0.0933zsJ and the quartic contribution of -0.3648zsJ. The spectrum for both modes is still gapless at Γ point, where $\gamma_{\mathbf{q}} = 1$, and at the points K and Q where $\gamma_{\mathbf{q}} = 1/2$.



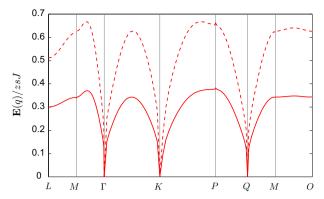


Fig. 2 The excitation spectrum of the longitudinal mode L_+ along $(LM\Gamma KPQMO)$ of the BZ with isotropic case $\xi=1$. It is gapless at Γ , K and Q points. The longitudinal spectra calculated from the first-order and higher-order approximations are indicated by the dash and solid lines, respectively (Color figure online)

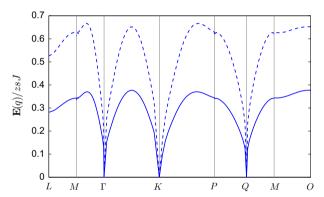


Fig. 3 The excitation spectrum of the longitudinal mode L_- along $(LM\Gamma KPQMO)$ of the BZ with isotropic case $\xi=1$. It is gapless at Γ , K and Q points. The longitudinal spectra calculated from the first-order and higher-order approximations are indicated by the dash and solid lines, respectively (Color figure online)

The numerical calculation demonstrates that the gapless spectra of L_- and L_+ modes are expected due to the slow logarithmic divergence in both the second and third terms in the structure factor $S(\mathbf{q})$ of Eq. (34). More precisely, near Γ , K and Q points, we find that $S(\mathbf{q}) \propto -\ln q$, and hence, the excitation spectrum $E(\mathbf{q}) \propto -1/\ln q$ with different coefficients.

The logarithmic behavior of the structure factor and the energy spectrum of the triangular lattice model is similar to that of the square lattice model investigated earlier [39,42,46,47]. We have identified these gapless modes of the 2D models as quasi-gapped modes because any finite size effect or anisotropy will induce a large energy gap when compared with the counterparts of the spin-wave spectrum. The effect of anisotropy can be investigated by considering the value of ξ parameter in Eq. (3) differing from unity. For example, for a tiny anisotropy such as $\xi = 1 - 1.5 \times 10^{-4}$,



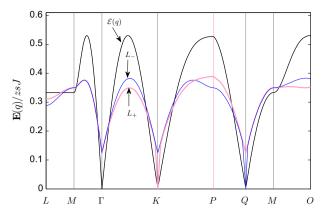


Fig. 4 The longitudinal modes L_- and L_+ together with spin-wave excitation spectrum $\mathcal{E}(q)$ of the 2D triangle lattice along $(LM\Gamma KPQMO)$ of the BZ with an anisotropy $\xi=1-1.5\times 10^{-4}$. The longitudinal gap values at Γ , K, and Q points for both modes, L_- and L_+ , respectively, are 0.2030zsJ in the first-order approximation and 0.1242szJ after including the higher-order correction. The transverse spin-wave gap is 0.0075zsJ (Color figure online)

at Γ point for both modes, we obtain the energy gap value of 0.2030zsJ in the first-order approximation and 0.1242szJ after including the higher-order corrections with the cubic contribution of 0.0407zsJ and quartic contribution of -0.1195zsJ. The gap value of the corresponding spin-wave spectrum at the same value of anisotropy is 0.0075zsJ, much smaller. In particular, we find that the longitudinal energy gap value is proportional to $1/[-\ln(1-\xi)]$, in comparison with the spin-wave gap which is proportional to $\sqrt{1-\xi}$, when $\xi \to 1$. In order to make further comparisons between the longitudinal mode and the transverse spin-waves mode, we plot both the spectra with $\xi = 1 - 1.5 \times 10^{-4}$ in Fig. 4 along the path $(LM\Gamma KPQMO)$ of the BZ. The different gap values for the longitudinal and transverse mode at Γ , K and Q points can be clearly seen. The spin-wave spectra at Γ point are still gapless where $\gamma_{\bf q} = 1$, whereas both longitudinal modes have the gap value of 0.1242szJ. The L_+ mode has the same gap at Q point, but it is gapless at K point, and vice versa for the L_- mode.

Before we turn to the quasi-1D systems in the next section, we like to mention that although we do find the significant energy reduction in the two longitudinal modes after the inclusion of the higher-order terms, we cannot at the moment directly relate our values to the peak structures of dynamic structure calculated in Ref. [32] based on SWT.

5 Magnon-Density Waves in Quasi-1D Triangular Lattices

We now turn to the longitudinal modes for the quasi-1D hexagonal antiferromagnetic systems, modeled by the following Heisenberg Hamiltonian with a strong interaction J along the chains and weak interaction J' on the hexagonal planes,



Table 1 The numerical results for the energy gap of the L_{\perp} mode with and without the cubic term contribution at the magnetic wavevector for the three quasi-1D materials

Quasi-1D materials	L_{-} mode before	L_{-} mode after
CsNiCl ₃	(0.490721)2J	(0.490837)2J
RbNiCl ₃	(0.718999)2J	(0.719291)2J
$CsMnI_3$	(1.19189)2J	(1.19191)2J

$$H = 2J \sum_{\langle i,j \rangle}^{\text{chain}} \mathbf{S}_i \cdot \mathbf{S}_j + 2J' \sum_{\langle i,j \rangle}^{\text{plane}} \mathbf{S}_i \cdot \mathbf{S}_j.$$
 (35)

An energy gap about 0.41(2J) has been observed by the neutron scattering experiments for CsNiCl₃ [2] with spin s = 1, J = 0.345 and J' = 0.0054 THz. This energy gap does not belong to the transverse spin-wave spectra, but belong to the longitudinal modes, as first proposed by Affleck [12,13]. Following Affleck [12,13], we earlier calculated the energy gap of the lower longitudinal mode L_{-} at the point $\mathbf{Q} = (4\pi/3, 0, \pi)$ using Eq. (28) but including only the quartic correction and obtained a value of (0.4907)2J [41], in reasonable agreement with the experimental result. Now after including the cubic contribution (as described by I_3 of Eq. 31), we obtain a value of (0.4908)2J for this gap, with a very small change. For RbNiCl₃ also with s=1but J'/J = 0.0295, the experimental result of the gap value is about 0.51 THz [48]; our result is 0.6974 THz with only quartic correction and 0.6977 after including cubic correction. The compound CsMnI₃ has spin quantum number s = 5/2 and the very small ratio of couplings $\epsilon = J'/J \approx 0.005$, for which the SWT approximation for its ground state is very poor; our result for the gap value of 0.47199 THz with only quartic correction and 0.47200 THz after including the cubic correction is in very poor comparison with the experimental results of about 0.1 THz by Harrison et al. [49]. Clearly in the case of CsMnI₃, we need better ground state in order to obtain better results for the energy gap of the longitudinal modes as mentioned before.

In general, as we can see from Table 1, the contribution from the cubic term is tiny for the energy spectra of the longitudinal modes. This is mainly due to the small value of $\epsilon = J'/J$ in these systems, namely the coupling J' on the hexagonal plans with non-vanishing cubic contribution is much smaller than the coupling J along the chains for which the cubic term vanishes. The energy spectra of the longitudinal modes for such quasi-1D systems of Eq. (35) can be expressed as sum of the chain and plane parts as,

$$E_{\pm} = L_{\pm}^{\text{chain}} + L_{\pm}^{\text{plane}}.$$
 (36)

In Table 2, we present the numerical results for the energy gap due to the planar term of Eq. (36) before and after including cubic corrections for the three quasi-1D materials, and where we define the cubic contribution as $\Delta L_{-}^{\rm plane}$. We can see that the cubic correction relative to the quartic contribution is similar in ratio to that of the 2D triangular lattice model discussed in Sect. 4.



Table 2 The numerical results for the L-energy gap of the planar term of the Hamiltonian Eq. (35) with and without the cubic term contribution at the magnetic wavevector for the three quasi-1D materials

Quasi-1D materials	L_{-}^{plane} before	L_{-}^{plane} after	$\Delta L_{\perp}^{ m plane}$
CsNiCl3	0.0309086 zs J'	$0.0333744 \ zsJ'$	$0.00246585 \ zsJ'$
RbNiCl3	0.0544659 zsJ'	$0.0577724 \ zsJ'$	$0.00330649 \ zsJ'$
CsMnI3	0.0231435 zsJ'	$0.0237096 \ zsJ'$	$0.00056611 \ zsJ'$

6 Conclusion

In this paper, we have investigated the longitudinal excitations of the 2D triangular antiferromagnetic lattice and the quasi-1D hexagonal systems after including the higher-order corrections. For the 2D triangular model, we find significant reduction of about 40% in the energy spectra from the higher-order contributions of the cubic and quartic terms. For the quasi-1D hexagonal materials, we find that the cubic corrections are negligible when compared with the quartic corrections which was calculated earlier [41]. This is mainly because of the weak coupling on the triangular planes.

Our numerical values for the energy gap of the longitudinal modes after including the higher-order corrections are in reasonable agreement with the experimental results for the spin-1 compounds CsNiCl₃ and RbNiCl₃, but are poor for the spin-5/2 compound CsMnI₃ because the approximate ground state by SWT is poor for this compound which is very close to a quantum critical point. Clearly, a better ground state for this compound will be needed in our calculation of the longitudinal modes in order to make reasonable comparison with the experiment.

Another point that needs addressing is the question of how well defined are the longitudinal modes in 2D triangular antiferromagnets since the magnon-density ρ in the order parameter of Eq. (25) may not be high enough to support the longitudinal modes. This is similar to the case of the 2D antiferromagnet on a square lattice. As we mentioned earlier, although there is no direct experimental evidence of these longitudinal modes in 2D triangular antiferromagnet, theoretical investigation of dynamic structure factors does find some broad peaks in the two-magnon continuum [32], indicating the strong longitudinal collective fluctuations. It will also be desirable to investigate the spontaneous decay of the longitudinal modes due to the coupling to the magnons as represented by the cubic terms in the Hamiltonian [33], and we wish to report our investigation in the future.

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