

Confinement of Fractional Excitations in a Triangular Lattice Antiferromagnet

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High-resolution neutron and THz spectroscopies are used to study the magnetic excitation spectrum of Cs_2CoBr_4 , a distorted-triangular-lattice antiferromagnet with nearly XY -type anisotropy. What was previously thought of as a broad excitation continuum [L. Facheris *et al.*, *Phys. Rev. Lett.* **129**, 087201 (2022)] is shown to be a series of dispersive bound states reminiscent of “Zeeman ladders” in quasi-one-dimensional Ising systems. At wave vectors where interchain interactions cancel at the mean field level, they can indeed be interpreted as bound finite-width kinks in individual chains. Elsewhere in the Brillouin zone their true two-dimensional structure and propagation are revealed.

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In conventional magnetic insulators the dynamic response is typically dominated by coherent single-particle $S = 1$ excitations, also known as magnons or spin waves. In many low-dimensional and highly frustrated quantum spin systems elementary excitations carry fractional quantum numbers, be they spinons in Heisenberg spin chains [1–4], Majorana fermions in the now-famous Kitaev model [5–7], or pseudocharge defects in spin ice like Coulomb phases [8]. The physical excitation spectrum, such as that measured by neutron spectroscopy, is then dominated by broad multiparticle continua [9–13]. In addition to the continuum, fractional excitations may also form bound states due to attractive interactions between them. A spectacular new phenomenon emerges when interactions are *confining*, i.e., do not fall off with distance, much like strong forces that bind quarks in hadrons [14]. This produces a hierarchical series of bound states inside the resulting potential well. An example is the sequence of domain wall (kink) bound states in quasi-one-dimensional (quasi-1D) Ising spin chains [15–17]. The confining potential for this model is linear and results from 3D couplings, which generate an effective field acting on individual chains [16]. The binding energies are, in supreme mathematical elegance, spaced according to the negative zeros of the Airy function [15,17]. The best-known experimental examples of such “Zeeman ladder” spectra are the quasi-1D Ising ferromagnet CoNb_2O_6 [18] and antiferromagnet (AF) $\text{BaCo}_2\text{V}_2\text{O}_8$ [19], as well as the isostructural compound $\text{SrCo}_2\text{V}_2\text{O}_8$ [20], where as many as eight consecutive bound states are observed. Shorter sequences have been found in another prototypical Ising spin chain material, RbCoCl_3 [21].

Bound states are always easier to form in 1D, where even an arbitrary-shallow potential well provides at least one robust localized state for a propagating particle. Fractional

excitations can also form bound states in higher dimensional spin systems [22,23]. Their structure and propagation are much more complex than that of simple paired domain walls in spin chains. Unfortunately, no Zeeman ladder analogues, i.e., no hierarchical binding sequences of fractional excitations, have been found in higher dimensions to date. In the present work we report the observation of such a phenomenon in a quasi-2D distorted-triangular-lattice AF. That the quintessentially 1D physics of bound kinks survives in 2D is remarkable. We argue that it is “rescued” at certain special wave vectors by the intrinsic frustration in triangular lattice geometry. Elsewhere in the Brillouin zone the bound states are no longer restricted to single chains and are to be viewed as 2D objects propagating in the entire triangular plane.

The material in question, Cs_2CoBr_4 (space group $Pnma$, $a = 10.19$, $b = 7.73$, $c = 13.51$ Å), is a very interesting $J - J'$ model distorted-triangular-lattice AF [24,25]. Despite a prominent triangular motif in its structure, it demonstrates certain 1D features such as a field-induced incommensurate spin density wave with Tomonaga-Luttinger spin liquid type dynamics and a propagation vector controlled by a one-dimensional nesting in the spinon Fermi sea. Its true 2D nature is manifest in the presence of a robust $m = 1/3$ magnetization plateau, typical of a triangular AF. The model magnetic Hamiltonian is described in detail in Refs. [24,25]. The key structural features are chains of Co^{2+} ions that run along the crystallographic b axis of the orthorhombic lattice (see Fig. 1 in Ref. [24]). The chains are coupled in the (bc) plane in a zigzag fashion to form a distorted triangular network [inset of Fig. 1(d)]. Easy-plane single-ion anisotropy ensures that the low-energy physics of the spin-3/2 Co^{2+} ions can be described in terms of effective $S = 1/2$ pseudospins. The components of the effective exchange coupling constants are

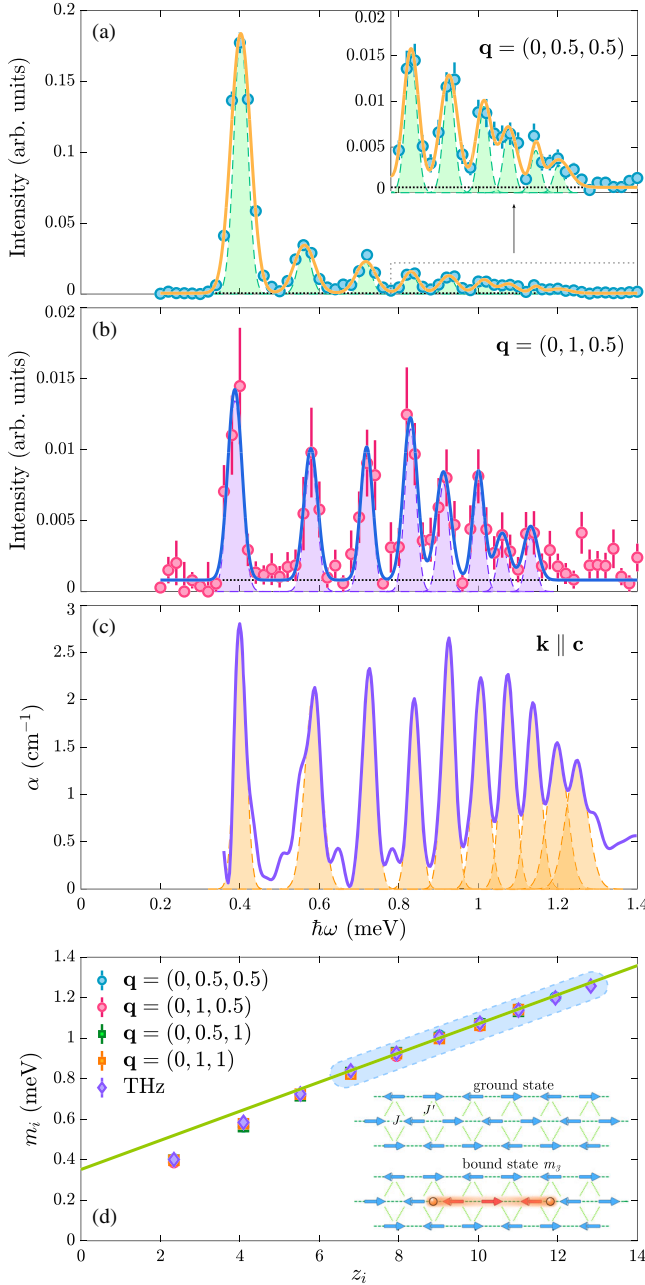


FIG. 1. (a),(b) Neutron scattering intensity (solid symbols) measured at $T = 40$ mK versus energy transfer at the 1D AF zone centers $\mathbf{q} = (0, 0.5, 0.5)$ and $\mathbf{q} = (0, 1, 0.5)$, respectively. The data are integrated fully along h direction and in ± 0.025 r.l.u. and ± 0.25 r.l.u. along k and l , respectively. Solid lines are fits to a series of Gaussian peaks. Dashed Gaussians represent the calculated experimental energy resolution. Black dotted lines indicate the fitted flat background. (c) Measured THz absorption (solid line) versus absorbed photon energy for light propagating along the \mathbf{c} axis at 0.2 K. Dashed areas highlight the Gaussian components of the fit to extract peak positions. (d) Measured excitation energy plotted versus the value of negative roots of the Airy function. The solid line is a linear fit as described in the text. The blue area highlights the points used for the fit. Inset: cartoons of the magnetic ground state and a representative $m = 3$ two-kink bound state.

subject to restrictions imposed by the pseudospin projection. A simplistic spin-wave analysis of previous inelastic neutron data provided a rough estimate for the nearest-neighbor in-chain AF exchange tensor components: $J^{XX} \sim J$, $J^{YY} \sim 1.1J$, $J^{ZZ} \sim 0.25J$, $J = 0.8$ meV [25]. Here Y is chosen along the \mathbf{b} crystallographic direction, and X and Z alternate between adjacent chains, where anisotropy planes are almost orthogonal. Note that this is practically a planar exchange anisotropy, with only a tiny in-plane Ising component to account for the $\Delta \sim 0.4$ meV spectral gap found in this system. The frustrated interchain coupling J' is significant, of the order of $0.45J$, and is of predominantly Ising (YY) character. Interplane interactions J'' are not frustrated. The material orders magnetically in a collinear stripe-type structure, with an ordering wave vector $(0, 1/2, 1/2)$ [see inset in Fig. 1(d)]. The Néel temperature $T_N = 1.3$ K allows us to estimate J'' . If this were the only coupling between chains with no additional frustration due to J' , we could expect $k_B T_N \sim 2\Delta / \ln(\Delta/J'')$ [26]. The actual value of J'' must be larger than thus obtained, as the in-plane frustration interferes with the emerging magnetic structure. A certain upper estimate is given by the mean field picture where $k_B T_N \sim 2J''S(S+1)$. This leads us to conclude that 3×10^{-4} meV $\lesssim J'' \lesssim 0.075$ meV $\ll J$, confirming the quasi-2D character of the material.

Our previous inelastic neutron scattering experiments indicated that the excitation spectrum in zero applied field is a gapped continuum of states, with intensity concentrated on its lower bound, and a strong dispersion along the chain axis [25]. The central finding of the present work is that this “continuum” is actually a sequence of at least nine sharp bound states that previously could not be observed due to poor experimental energy resolution. New neutron data were collected at the LET time-of-flight spectrometer at ISIS (UK), using 2.35 meV incident energy neutrons in repetition-rate-multiplication mode [27]. We used the same 1.16 g single crystal as in [25] mounted on a ^3He - ^4He dilution refrigerator. All measurements were performed at a base temperature of 40 mK. In the experiment the sample was rotated 180° around the \mathbf{a} axis in steps of 1° . The spectra were measured for ~ 10 -min counting time at each sample position.

We first focus on the 1D AF zone centers ($\mathbf{q}\mathbf{b} = 0, \pi$), where interchain interactions within the triangular planes cancel out at the mean field-RPA level, and where spin wave theory predicts no transverse dispersion or intensity modulation of excitations. Figures 1(a) and 1(b) show constant- \mathbf{q} cuts through the data at wave vectors $\mathbf{q} = (0, 0.5, 0.5)$ and $\mathbf{q} = (0, 1, 0.5)$, respectively. A sequence of sharp peaks is clearly apparent in both cases. A fit to the data using empirical Gaussian profiles yields an accurate measure of the peak positions and shows that their widths are essentially resolution limited. In Figs. 1(a) and 1(b) this is emphasized by the shaded Gaussians representing the computed experimental resolution [28].

Corroborative evidence is also obtained by THz spectroscopy. The experiment was performed with a

Martin-Puplett-type interferometer and a ^3He - ^4He dilution refrigerator with base temperature of 150 mK using a ^3He -cooled Si bolometer at 0.3 K. The sample was a circular plate approximately 1 mm thick in c direction and 4 mm in diameter. THz radiation propagating along the crystal c axis was unpolarized and the apodized instrumental resolution was 0.025 meV. The THz absorption spectrum is shown in Fig. 1(c). It is calculated as a difference of spectra measured at 0.2 K and 2 K, i.e., in the magnetically ordered phase and above T_N . The THz spectrum appears to have some features absent in the neutron spectrum [33], but all peaks found in the latter are also present here. The positions of these peaks were determined in Gaussian fits (shaded peaks) in a narrow range ± 0.025 meV near each peak value. Comparing data measured at low temperature to that collected at 2 K [28] reveals that the series of peaks is endemic to the ordered phase and disappears above T_N .

The spacing between the excitation peaks present in both measurements corresponds to confinement in an approximately linear one-dimensional potential. To demonstrate this, we plot the excitation energies deduced from neutron spectra at several wave vectors, as well as the positions of corresponding THz peaks, versus the negative roots z_i of the Airy function in Fig. 1(d). For a precise linear attractive potential $\lambda|x|$ between the dispersive particles [35], near the minimum $\epsilon(k) = m_0 + \hbar^2 k^2 / 2\mu$ we expect the excitation energies to be [17,18]

$$m_i = 2m_0 + (\hbar\lambda)^{2/3} \mu^{-1/3} z_i \quad \text{with } i = 1, 2, \dots \quad (1)$$

In the actual data, the linear dependence is apparent for all but the first few points. As will be addressed in more detail below, this slight deviation indicates that the confining force increases somewhat at short distances. From a linear fit to the higher-energy peaks we can immediately extract the slope 0.072(3) meV and the energy of a single particle $m_0 = 0.18(1)$ meV (half-intercept). Using the single-particle kinetic mass $\hbar^2/\mu = 0.39 \text{ meV} \times b^2$ [28], we estimate the confining force constant $\lambda = 0.031(2) \text{ meV}/b$ [36].

The next point that we make is that the observed bound states at the 1D AF zone center are essentially 1D objects. This is concluded by analyzing the neutron spectra shown in Figs. 2(a) and 2(b) and Supplemental Fig. 5(a) in [28]. The bound states do not propagate in either transverse direction and thus have an essentially flat dispersion. Moreover, their intensity shows a monotonic falloff and no obvious transverse modulation, as plotted for the first two modes in Figs. 2(e) and 2(f), and Supplemental Fig. 5(b) [28]. This implies that these excitations do not involve cross-chain correlations and are confined to a single chain.

For the first mode, the measured transverse wave vector dependence is precisely accounted for [orange solid lines in Figs. 2(e) and 2(f), and Supplemental Fig. 5(b)] by the combined effects of (i) the magnetic form factor of Co^{2+} and (ii) a neutron polarization factor for spin components

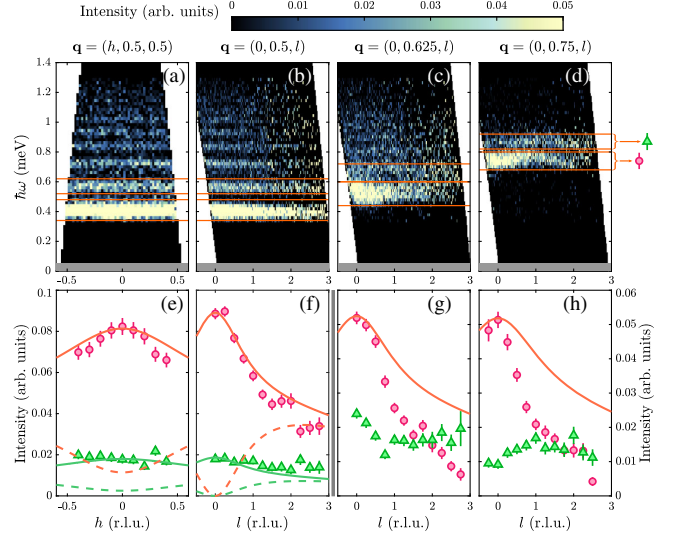


FIG. 2. (a)–(d) False color plot of neutron scattering intensity measured at $T = 40$ mK plotted versus energy transfer and momentum transfer transverse to the crystallographic b axis. Gray areas mask regions of elastic-incoherent scattering. Background subtraction has been performed as described in [28]. The orange regions represent energy-integration windows used to extract the cuts in panels below. (e)–(h) Intensity-momentum cuts (solid symbols) for the first two modes. The orange and green lines show the product of the magnetic form-factor-squared for Co^{2+} and calculated neutron polarization factor for excitations polarized perpendicular (solid) and parallel (dashed) to the ordered moment direction, respectively. The overall scale factor of the lines is arbitrary, but always consistent between the two polarizations and consistent between panels (e) and (f).

perpendicular to the chain axis (to the direction of ordered moments in the ground state). Longitudinal excitations would become stronger for scattering vectors deviating from the chain axis (orange dashed line). Already for the second mode the identification of polarization is ambiguous (green lines). Unlike in Ising chains, in Cs_2CoBr_4 the conservation of the chain-axis projection of angular momentum is not expected, meaning that excitations cannot be unambiguously classified as longitudinal nor transverse.

The 1D nature of zone-center excitations prompts a simple interpretation. Similarly to the situation in CoNb_2O_6 and $\text{BaCo}_2\text{V}_2\text{O}_8$, the observed modes are bound states of two kinks (domain walls) in individual chains. Naively, ignoring the kinetic energy of the kinks as in the purely Ising case, such an excitation can be visualized as in the cartoon in the inset of Fig. 1(d). The energy m_0 is to be associated with that of a single domain wall. As a consistency check, we can compare that to the computed energy of a domain wall in a classical spin chain. Using $J^{YY}/J^{XX} \sim 1.1$ as estimated for Cs_2CoBr_4 , with a trivial numerical classical-energy minimization procedure we get $m_0 \sim 0.9JS^2 = 0.18 \text{ meV}$, in excellent agreement with the measured value.

Geometric frustration ensures that at the magnetic zone center these strings of flipped spins within a single chain incur no energy cost due to interactions with adjacent chains within the triangular lattice. Moreover, any transverse dispersion is suppressed. At the same time, the interaction energy due to unfrustrated interlayer coupling is proportional to the string length, resulting in confinement. In this simplistic picture, the confining force is $\lambda = 2J''S^2/b$. This yields an interlayer coupling constant $J'' = 0.062(4)$ meV, inside the possible range deduced from T_N .

The apparent difference between our data and those for the Ising-chain AF $\text{BaCo}_2\text{V}_2\text{O}_8$ [19] is the presence of *two* Zeeman ladder series in the latter material. The two series correspond to transverse and longitudinal excitations, the first transverse mode being a spin flip, i.e., a conventional magnon. They are offset in energy by λb in our notation [28]. For Cs_2CoBr_4 this amounts to as little as 0.03 meV. The two series cannot be resolved within the resolution and counting statistics of our neutron experiment but may account for the extra shoulder peaks seen in the THz data in Fig. 1(c). A similar estimate for $\text{BaCo}_2\text{V}_2\text{O}_8$ yields a larger offset of about 0.34 meV that is easily resolved experimentally [19].

The deviation from linear-potential behavior at low energies is also readily explained. Since the material is almost planar, the domain walls are not confined to a single bond as in the ideal Ising case, but have a characteristic size l [37]. We can estimate that quantity in a classical spin chain using the above-mentioned anisotropy parameters: $l \sim 2b$. The energy of the first few bound states is thus modified due to a physical overlap of the two bounding domain walls. Once the kinks are separated by a distance of more than $\sim l$, this interaction becomes negligible and the confinement potential becomes linear.

In summary, at the 1D AF zone center the simplistic kink model from 1D Ising physics consistently explains the observed excitation energies. The first mode is transversely polarized and shows an intensity pattern very consistent with that of a magnon in spin wave theory [25], being strongly suppressed at the 1D structural zone centers (integer k). The intensities of higher modes are not that easily interpreted: they may have a mixed polarization and are less suppressed at integer k values [Figs. 1(a) and 1(b)].

Away from the 1D AF zone centers, the excitations are considerably more complex. This is very clear in the longitudinal dispersion of the bound states shown in Figs. 3(a) and 3(b). Other than at $\mathbf{q}\mathbf{b} = 0, \pi$ ($k = 0, 1/2$) the m_1 mode splits into two branches, each with an asymmetric dispersion relation. In fact, the m_1 state at $\mathbf{q}\mathbf{b} = \pi$ seems to be continuously connected to the m_2 excitations at $\mathbf{q}\mathbf{b} = 2\pi$ ($k = 1$) and vice versa. Fitting the dispersion of the strongest low-energy mode in the vicinity of $\mathbf{q}\mathbf{b} = \pi$ to a Lorentz-invariant relativistic form

$$(\hbar\omega_{\mathbf{q}})^2 = \hbar^2\Delta(\mathbf{q}\mathbf{b})^2/\mu + \Delta^2, \quad (2)$$

yields the value of kinetic mass quoted above.

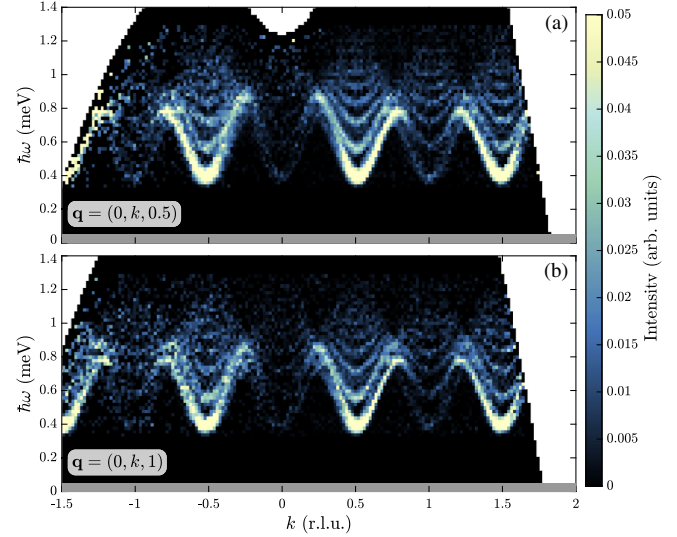


FIG. 3. (a),(b) False color plot of neutron scattering intensity measured at $T = 40$ mK plotted versus energy transfer and momentum transfer along $\mathbf{q} = (0, k, 0.5)$ and $\mathbf{q} = (0, k, 1)$, respectively. The data were fully integrated along h , and in the range ± 0.25 r.l.u. along l around the central value. The gray areas mask regions where the incoherent scattering dominates the signal. Background subtraction has been performed as described in [28].

A look at the intensities reveals that other than at the special wave vectors, the bound states can no longer be seen as strings in a single chain, but are “dressed” with correlations extending to several neighboring chains in the triangular plane. This conclusion is reached from Figs. 2(c) and 2(d), that show a transverse cut of the spectrum at $\mathbf{q}\mathbf{b} = 5\pi/4$ and $\mathbf{q}\mathbf{b} = 3\pi/2$, respectively. As plotted in Figs. 2(g) and 2(h), the measured intensity of the first two modes now shows a much steeper transverse wave vector dependence than computed from just the polarization and form factors (solid lines). The second mode even seems to show signs of intensity oscillations, with a possible mixing or hybridization and resulting intensity transfer with the first mode.

Our data reveal that away from the special wave vectors the bound states also *propagate* in two dimensions, albeit with a small bandwidth. Indeed, in Fig. 2(d) one can see that at $\mathbf{q}\mathbf{b} = 3\pi/2$ the bound states develop a nonzero dispersion along the c^* direction, in contrast to what is seen at $\mathbf{q}\mathbf{b} = 0, \pi$. Although the bandwidth of transverse dispersion, 0.08 meV, is at the limit of our experimental resolution, qualitatively one can say that $\mathbf{q}\mathbf{c} = 0, 4\pi$ are dispersion minima for the m_1 mode, while the maximum is at $\mathbf{q}\mathbf{c} = 2\pi$. That periodicity is consistent with having two chains per unit cell along the c -axis direction in the crystal structure.

Overall, the differences between our results and spectra of Ising spin chains [18,19] are striking. In the latter, all bound states, including the first one, are much less

dispersive than the lower edge of the entire spectrum, which approximately corresponds to the lower edge of the two-kink continuum in the absence of long-range order. As a result, each bound state persists only in a restricted area in the Brillouin zone. In contrast, in Cs_2CoBr_4 a few of the lower-energy bound states are highly dispersive and span across the entire zone.

In summary, we demonstrate that Zeeman ladders of confined fractional excitations can exist in a *bona fide* quasi-2D system. These states are inherently related to those in 1D models, as revealed at special wave vectors where 2D interactions are canceled by geometric frustration. However, elsewhere in reciprocal space their true 2D character is manifest. Once again, the distorted triangular lattice model provides a link between 1D and 2D quantum magnetism.

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- [36] This force of ~ 6 fN corresponds to the gravity pull between two average humans at a separation of 8 km.
- [37] We define the domain wall width in a spin- S chain as the distance over which the z -axis spin component changes from $S/2$ to $-S/2$ near its center.
- [38] L. Facheris *et al.*, Spin-density wave dynamics in a 2D distorted triangular lattice antiferromagnet, *STFC ISIS Neutron and Muon Source* (2022), [10.5286/ISIS.E.RB2210048](https://doi.org/10.5286/ISIS.E.RB2210048).