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Repulsive Tomonaga-Luttinger liquid in the quasi-one-dimensional alternating spin- $\frac{1}{2}$ antiferromagnet NaVOPO₄

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We probe the magnetic field-induced Tomonaga-Luttinger liquid (TLL) state in the bond-alternating spin-1/2 antiferromagnetic (AFM) chain compound NaVOPO₄ using thermodynamic as well as local μ SR and ³¹P NMR probes down to mK temperatures in magnetic fields up to 14 T. The μ SR and NMR relaxation rates in the gapless TLL regime decay slowly following characteristic power-law behavior, enabling us to directly determine the interaction parameter *K* as a function of the magnetic field. These estimates are crosschecked using magnetization and specific heat data. The field-dependent *K* lies in the range of 0.4 < *K* < 1 and indicates the repulsive nature of interactions between the spinless fermions, in line with the theoretical predictions. This renders NaVOPO₄ the first experimental realization of TLL with repulsive fermionic interactions in hitherto studied *S* = 1/2 bond-alternating AFM-AFM chain compounds.

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Quantum phase transitions and quantum critical states in low-dimensional spin systems are an active research field in current condensed-matter physics [1]. Weakly coupled quasi-one-dimensional (1D) gapped spin-1/2 antiferromagnets (AFM) offer an excellent playground to study such quantum effects [2-6]. These gapped spin systems are highly vulnerable to the external magnetic field (H). Above the critical magnetic field (H_{C1}) of gap closing, a three-dimensional (3D) magnetic long-range-order (LRO) sets in, known as Bose-Einstein-condensation (BEC) of spin-1 triplons [7–9]. Further, above the 3D LRO, the gapless excitations can be universally described by a non-Fermi liquid type Tomonaga-Luttinger-liquid (TLL) theory [10] of interacting spinless fermions in 1D [2,6]. The low energy excitation of the TLL state is described by two parameters: the renormalized velocity of excitations u and the dimensionless interaction parameter K that determines the nature and strength of the interactions between the spinless fermions [2]. The TLL model has been successfully applied to a multitude of systems such as quantum nanowires [11], edge state of quantum Hall systems [12], carbon nanotubes [13], and most importantly magnetic insulators [14].

Although the physics of these diverse 1D systems has been explored in the framework of TLL model, the quantitative description of TLL parameters was missing until a long time. In this regard, recent notable achievements are the experimental observation of the TLL state with attractive (i.e., K > 1) and repulsive (i.e., K < 1) interactions in strongleg (C₇H₁₀N₂)₂CuBr₄ (DIMPY) [15–17] and strong-rung (C₅H₁₂N₂)₂CuBr₄ (BPCB) [8,17–19] two-leg spin ladders, respectively. Moreover, a remarkable control of the TLL properties by means of field tuning of the *u* and *K* parameters has also been demonstrated in these weakly coupled spin-ladder systems [16–18]. Here, *H* acts as a chemical potential that controls the filling of the fermion band and, therefore, tunes the dynamics of the TLL state. TLL with repulsive interaction is also demonstrated in spin-1/2 uniform Heisenberg/Ising spin chains [14,20–23].

Apart from spin-1/2 uniform Heisenberg/Ising spin chains and spin ladders, the Heisenberg bond-alternating spin-1/2 chains have been predicted to evince field-induced TLL state of interacting spinless fermions [24,25]. The nature of interactions between these fermions, attractive or repulsive, should be determined by the sign and hierarchy of the nearest-neighbor exchange couplings. One expects attractive interaction for ferromagnetic (FM)-AFM spin chains and repulsive interaction for AFM-AFM spin chains in the TLL state [25]. On the experimental side, only a very few bondalternating spin-1/2 chain systems are reported where TLL behavior is anticipated from the thermodynamic measurements. They include copper nitrate $Cu(NO_3)_2 \cdot 2.5D_2O$ [26] and pentafluorophenyl nitronyl nitroxide F₅PNN [27] with

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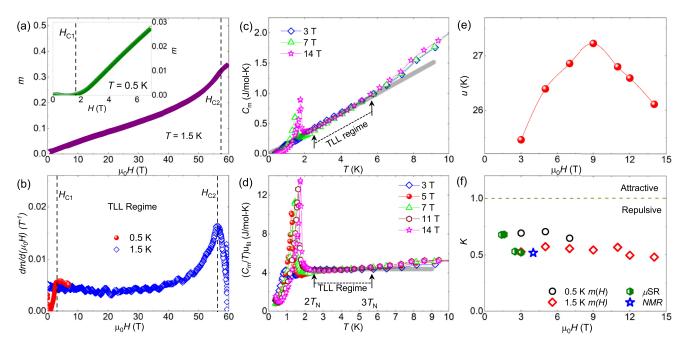


FIG. 1. (a) *m* vs *H* measured at T = 1.5 K. (Inset) *m* (after impurity subtraction) vs *H* at T = 0.5 K. (b) dm/dH vs *H* together for T = 1.5 and 0.5 K. (c) Magnetic specific heat (C_m) vs *T* for three selected fields. Dashed line guided by two upward arrows marks the *T*-linear TLL regime. (d) Normalized C_m vs *T* in different fields above H_{C1} . The linear TLL regime is marked by two arrows. (e) Normalized velocity *u* vs *H*, estimated from the C_m analysis. (f) TLL exponent *K* vs *H* deduced from magnetization, specific heat, μ SR, and NMR $1/T_1$ experiments.

AFM-AFM spin-chains, zinc-verdazyl complex [28] with FM-AFM spin-chains, etc. However, none of these reports present a detailed quantitative analysis and verification of the nature of interactions (attractive/repulsive) in the TLL state.

The recent discovery of an alternating spin-1/2 chain compound NaVOPO₄ [29] appears to be promising in this context. It features quasi-1D bond-alternating spin-1/2 AFM-AFM chains formed by VO₆ octahedra with leading exchange coupling $J/k_{\rm B} \simeq 39$ K along the chain direction with a weak alternation ratio $\alpha(=J'/J) \simeq 0.98$. The ground state is a singlet with a tiny zero-field spin-gap $\Delta_0/k_{\rm B} \simeq 2$ K that corresponds to a small accessible critical field of $\mu_0 H_{C1} \simeq 1.6$ T. The critical field of saturation is $\mu_0 H_{C2} \simeq 57$ T. An external field $H \ge H_{C1}$ closes the gap and gives rise to a 3D LRO (T_N) which is described well by 3D BEC of triplons. Since bondalternating spin chains are predicted to show evidence of both TLL and BEC phases, NaVOPO₄ gives a unique opportunity to study the crossover effect from 1D TLL at high-T s to 3D BEC at low T s. Moreover, the low value of H_{C1} provides an added advantage to explore the critical TLL behavior under laboratory fields.

In this Letter, we present a detailed study of the TLL state in NaVOPO₄. Using magnetization and specific heat and also powerful local tools such as nuclear magnetic resonance (NMR) and muon-spin relaxation (μ SR), we demonstrate that the interaction between spinless fermions is repulsive in nature within the TLL state, which is in accordance with the theoretical predictions [25]. Moreover, from the temperature-dependent NMR spin-lattice relaxation rate [1/*T*₁], we estimate the exact value of the TLL parameter (*K*)

employing the recent analytical results reported by Dupont *et al.* [23,30]. Our study not only links the 1D TLL theory to bond-alternating AFM-AFM Heisenberg spin-1/2 chain systems but also provides insight about the nature and strength of interactions between the spinless fermions in the gapless TLL state.

All the experiments were performed on the polycrystalline NaVOPO₄ sample. Details of the magnetization, specific heat, and ³¹P NMR measurements are described in Ref. [29]. μ SR measurements were carried out in zero-field (ZF) as well as in longitudinal-fields (LFs) down to 50 mK (see Ref. [31] for details).

We begin the discussion of prospective TLL behavior in NaVOPO₄ from the magnetization and specific heat. Figure 1(a) presents the reduced magnetization (m) vs H measured at T = 1.5 K, whereas the inset shows the corrected m vs H curve obtained after subtraction of the extrinsic paramagnetic contribution (see Ref. [29]). Here, the dimensionless quantity m is nothing but the expectation value of spin $\langle S \rangle$ [i.e., $m = \frac{M}{g\mu_{\rm B}} = \langle S \rangle$], calculated normalizing the measured magnetization by the Landé g factor $g \simeq 1.951$ (obtained from ESR experiments [29]). As expected for a gapped quantum magnet, at T = 0.5 K m remains nearly zero up to H_{C1} and dm/dH exhibits a rounded peak at $\mu_0 H_{C1} \simeq 1.6$ T as shown in Fig. 1(b). Similarly, at T = 1.5 K, m saturates in an applied field of $\mu_0 H_{C2} \simeq 57$ T where dm/dH shows another rounded maximum. These two rounded peaks in dm/dH [Fig. 1(b)] indicate the crossover in and out of the 1D TLL regime, respectively [17,26,32].

The magnetic specific heat (C_m) as a function of temperature is depicted in Fig. 1(c) at three representative fields

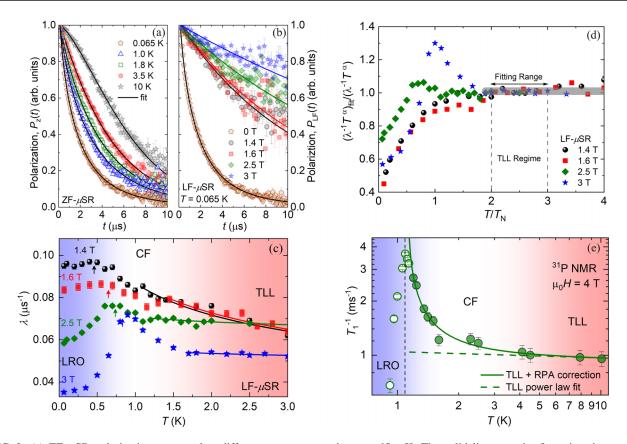


FIG. 2. (a) ZF- μ SR polarization measured at different temperatures down to 65 mK. The solid lines are the fits using the stretched exponential function, described in SM [31]. (b) μ SR polarization at 65 mK and in different LFs. The solid lines are the fits using single exponential function, described in the text. (c) Temperature dependent μ SR depolarization rate (λ) in different LFs. The solid lines are fits using Eq. (4). (d) Normalized $\lambda^{-1}T^{\alpha}$ vs T/T_N to show the collapse of data sets in different fields onto one line. (e) The RPA+TLL fit (solid line) using Eq. (5) to the $1/T_1(T)$ data at 4 T. The filled circles represent the data set considered for the fitting. The horizontal dashed line indicates the pure TLL contribution to the fit. The vertical dashed line denotes the T_N determined by the fit.

(for $H > H_{C1}$).¹ A striking feature observed in $C_m(T)$ is the appearance of a *T*-linear regime (2.5 to 7 K) above T_N which remains almost independent of field. Such a behavior is a robust signature of the TLL state [33]. In the TLL critical regime, the molar magnetic specific heat (C_m) is related to *u* as [2,17,34]

$$C_m = N_{\rm A} \frac{\pi k_{\rm B} T}{6u},\tag{1}$$

where N_A is the Avogadro's constant. To estimate u, we fitted $C_m(T)$ in the *T*-linear region by Eq. (1) and the corresponding u values obtained for different fields are plotted as a function of *H* in Fig. 1(e). In order to check the reliability of the estimation of u, C_m at different fields is normalized with respect to u [i.e., $(C_m/T)u_{fit}$] and plotted as a function of *T* in Fig. 1(d). Indeed, the $(C_m/T)u_{fit}$ curves above 2.5 K collapse on a flat line irrespective of the field value, which is

a clear demonstration of the TLL regime [17]. The variation of u in Fig. 1(e) is found to be reminiscent of the behavior reported for the spin-ladder compound BPCB with repulsive TLL interactions [17].

Another crucial parameter that characterizes TLL correlations is K. The relation between K and u can be expressed as [17]

$$\frac{\partial m}{\partial(\mu_0 H)} = \frac{g\mu_{\rm B}}{2\pi k_{\rm B}} \left(\frac{K}{u}\right). \tag{2}$$

The *u* values obtained from the specific heat analysis and differential magnetization $\partial m/\partial(\mu_0 H)$ data at 0.5 and at 1.5 K from Fig. 1(b) are used in Eq. (2) to determine *K*. The obtained *K* values plotted in Fig. 1(f) are in the range of 0.5 < K < 1. The value of *K* less than 1 indicates repulsive TLL interactions. Moreover, the value of *K* is also found to be more than 0.5, which suggests Heisenberg-type interactions with dominant transverse AFM fluctuations [23]. However, we note that the above method of estimating the TLL parameters may be biased by the quality of the experimental data [17]. Therefore, we back our conclusions using local probes like μ SR and NMR, both of which give direct access to the low-energy spin dynamics.

Figure 2(a) shows representative muon spin polarization curves measured at different temperatures from 10 K to 65 mK

 $^{{}^{1}}C_{\rm m}$ was obtained by subtracting the phonon contribution from the total specific heat. In order to estimate the phonon part, we fitted the high-temperature part of the total specific heat in zero field by a sum of multiple Debye terms as in Ref. [29] and the fit was extrapolated down to low temperatures. The same phonon contribution was also used to estimate $C_{\rm m}$ in different fields.

in zero field. At 10 K, it shows a Gaussian-like behavior. As the temperature is lowered, the shape of $P_z(T)$ eventually changes to exponential-like behavior. It exhibits neither coherent oscillations nor a loss of initial polarization down to 65 mK, ruling out the transition to a magnetic LRO state. The spectra also do not decay to 1/3 of the initial polarization at long times, characteristic of a quasistatic local magnetic field at the muon stopping site, thus excluding a spin-glass type transition [35,36]. These observations are consistent with a singlet/disordered ground state reported previously in zero field [29]. However, a weak increase in the muon depolarization rate (λ) below the temperature scale of $\Delta_0/k_B \simeq 2$ K can be ascribed to the quasi-free spins that arise from singlets broken by implanted muon or from a small amount of impurity/orphan spins (see Ref. [31]).

To look for the temperature-dependent electronic spin dynamics in NaVOPO₄ at higher fields $(H \ge H_{C1})$, we have measured μ SR response in four different LFs (i.e., 1.4, 1.6, 2.5, and 3 T). Magnetic field at the muon stop sites includes contributions from nuclear as well as electronic field fluctuations. In an applied longitudinal field, the nuclear contribution to the muon spin relaxation is expected to be fully quenched and thus one can reliably probe the electronic spin dynamics. Figure 2(b) presents the muon spin polarization at the lowest measured temperature of 65 mK in different LFs. The increasing slope of the curves with the applied field suggests a slowing down of electronic fluctuations with increasing field, in the field-induced ordered state [29]. Remarkably, the LF- μ SR polarization curve at the highest measured LF of 3 T also exhibits substantial relaxation, which is not completely quenched, implying electronic nature of the fluctuations. As the compound shows the onset of field-induced magnetic LRO above H_{C1} , one would expect distinct oscillations in the muon spin polarization below $T_{\rm N}$ or at least at the lowest measured temperature 65 mK. However, in our case we did not observe any such oscillations which could be due to the broad distribution of the local magnetic field at the muon site [37]. Similar muon response without visible oscillations has been observed in the ordered state of other gapped quantum AFMs [38,39].

The muon spin polarization curves over the entire measured temperature range and for all LFs follow a simple exponential relaxation [Fig. 2(b)]

$$P_{\rm LF}(t) = P_0 e^{-\lambda t} + P_b, \qquad (3)$$

where P_0 is the initial muon spin polarization and P_b is the nonrelaxing background component arising from the sample holder or the sample background. For each field, both P_0 and P_b are nearly temperature-independent. The estimated λ is plotted against temperature in Fig. 2(c) where we can identify three distinct temperature regimes. In the low-temperature regime, as the temperature rises, λ increases and shows an anomaly at $T_{\rm N}$, suggesting the onset of magnetic LRO. With increasing LF, the anomaly at T_N becomes more pronounced and shifts towards higher temperatures. The peak position in different fields traces the phase boundary (magnon BEC) shown in Ref. [29]. The region immediately above T_N is the critical fluctuation (CF) regime where thermal fluctuations dominate the relaxation. In the CF region, λ shows a decreasing behavior until the power-law behavior sets in above $\sim 2T_{\rm N}$ indicating TLL fluctuations. We note in passing that a peak in λ is observed even at $\mu_0 H = 1.4$ T, reflecting that the actual H_{C1} could be lower than 1.4 T.

In quantum spin chains, the spin-spin correlation decays following the power law, with the exponent being a simple function of K [40]. Hence, one can estimate the TLL parameter (K) by modeling the μ SR depolarization rate in the gapless TLL regime (above T_N) with a power law of the form [39]

$$\lambda = T_{\text{TLL}}^{-1}(T) \propto T^{1/2K-1}.$$
(4)

We fitted the data in Fig. 2(c) by Eq. (4) in the temperature window $2T_N \leq T \leq 3T_N$. To validate the choice of this temperature window, we normalized λ as λ/T^{α} (where $\alpha =$ 1/2K - 1) and plotted it against the reduced temperature T/T_N [see Fig. 2(d)]. Clearly, the data for all the four fields collapse on a straight line for $2T_N \leq T \leq 3T_N$ and also beyond $3T_N$, highlighting the universal power-law behavior in the TLL regime and confirming the reliability of the estimated α or K. Further, to ensure that the extracted K value is not affected by a particular choice of the fitting range, we fitted the data four times [39]: first over the full fitting window $2T_N \leq T \leq 3T_N$, then excluding either of the end points, and finally excluding both the end points. Consequently, the final value of K is taken as the inverse-variance weighted average and is plotted against H in Fig. 1(f). Thus, the obtained TLL parameter K < 1 once again proves the repulsive interactions between the spinless fermions in the TLL state. Moreover, the K value shows an increasing trend as the field is reduced. Conventionally, the value of K is expected to increase towards 1 with decreasing field since the system approaches the noninteracting and/or 1D quantum critical regime when the field is reduced to H_{C1} [16].

We note that a simple power-law dependence of the relaxation rate, which is predicted for a purely 1D spin system, underestimates the value of K since it does not take into account the 3D critical fluctuations that are predominant at low temperatures, especially in the region close to $T_{\rm N}$. In the past, attempts have been made to estimate K for the spinladder compound DIMPY from the temperature-dependent NMR $1/T_1$ using the generic power law [16,20]. However, the obtained values of K were not in good agreement with the theory. This mismatch was ascribed to the enhancement of the relaxation due to the 3D critical fluctuations in the vicinity of T_N [30]. Recently, the effect of this fluctuation near $T_{\rm N}$ has been theoretically modeled by Dupont *et al.* [30] using a random phase approximation (RPA). Thereafter, this approximation was used by Horvatić et al. [23] as a correction to the TLL power-law behavior that takes the form (RPA+TLL)

$$1/T_1 \propto T^{1/2K-1}\phi(K, T_N/T).$$
 (5)

Here, $\phi(K, T_N/T)$ is the RPA correction factor, which is a multiplicative correction function of *K* and T_N [23,30]. The validity of this method was independently tested on Ising spin-1/2 chain compound BaCo₂V₂O₈ and spin-ladder compound DIMPY [23]. The fitting window in this method can be extended to 1.2 T_N because the effect of critical fluctuations is included.

In Fig. 2(e), we show the temperature dependence of $1/T_1$ in a magnetic field of 4 T, which is well above H_{C1} . The sharp λ -type anomaly at $T_N \simeq 1.09$ K reflects the onset of a 3D magnetic LRO. The nonlinear fit using the analytical expression [Eq. (5)] returns $K \simeq 0.52$ and $T_N \simeq 1.11$ K. This K value clearly reflects that the interactions are repulsive in nature and are in line with our previous analysis. Moreover, the obtained T_N value is also in excellent agreement with our experimental data.

Quantum magnets demonstrating both BEC and TLL physics are rare. The majority of known materials feature very high critical fields or show the onset of field-induced LRO at relatively high temperatures. In this regard, NaVOPO₄ is a very promising candidate system in which both BEC and TLL physics can be realized due to the low critical field of $\mu_0 H_{C1} \simeq 1.4$ T. Indeed, we are able to tune the TLL parameter (K) in the gapless state by applying the magnetic field. For a comparison, we have plotted K vs H in Fig. 1(f) extracted from the specific heat, NMR, and μ SR data. The K values from the different experiments are consistent and remain below unity, indicating the repulsive nature of the interactions. Moreover, as we move towards the noninteracting regime, i.e., close to H_{C1} , the K value increases slowly. The overall behavior of u with H is symmetric and its shape is similar to the previously reported spin-ladder compound BPCB with repulsive TLL interactions [3,15,17]. Our results for NaVOPO₄ offer the first experimental observation of the TLL physics with repulsive interactions in bond-alternating spin chains. This finding is in an excellent agreement with theory [25].

In summary, we have obtained direct experimental confirmation for repulsive interactions between spinless fermions in the TLL state of NaVOPO₄. Several experimental methods consistently show repulsive interactions in the gapless TLL regime above H_{C1} and at temperatures above $\sim 2T_N$. The μ SR and NMR relaxation rates, which are measures of the correlation time, follow power-law behavior and provide direct estimates of the interaction parameter *K*. A more exact value of *K* is obtained from the analysis of $1/T_1$ using the recently proposed RPA+TLL method. Our results rigorously establish NaVOPO₄ as the first candidate bond-alternating spin-1/2 AFM-AFM chain system manifesting gapless TLL phase with repulsive interactions, validating the theoretical predictions.

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