I. INTRODUCTION

Low-dimensional frustrated spin systems have attracted great interest due to their importance for the understanding of emergent phenomena, such as reentrant phase transitions [1], flat-band physics [2], anomalous robustness of topological order [3–5], and spin liquid phases [6–10]. Frustrated spin ladders have received special attention for several reasons: (i) due to their low dimensionality, the interplay of frustration and quantum correlations leads to a variety of topological and nontopological phases such as the Haldane phase, dimer order and various spin liquids [11–26]; (ii) they are quasi-one-dimensional, and show characteristics of both one- and two-dimensional systems [27], and (iii) powerful numerical and analytical methods are available for studying their low-energy properties [28,29].

Most studies of spin ladders have considered a single type of spin, in particular, σ = 1/2. Ladders with two types of spin (τ > σ), so-called mixed-spin ladders, have received less attention, although this heterogeneity produces qualitatively new effects [30–35]. In this paper, we study the ground state phase diagram of a mixed-spin (1, 1/2) ladder with coupling constants \( J_l, J_r, J_d \) (illustrated in Fig. 1), focusing on the effects of diagonal interactions \( (J_d) \). We limit ourselves to \( J_l \geq 0, J_d \geq 0 \) but admit both ferromagnetic and antiferromagnetic rung couplings \( J_r \).

Some knowledge about the ground state can be gained thanks to the Lieb-Mattis theorem [36], which fixes the total spin \( S \) for cases where the lattice can be subdivided into sublattices \( \mathcal{A} \) and \( \mathcal{B} \) in such a way that \( J_{ij} = 0 \) if \( i, j \in \mathcal{A} \) or \( i, j \in \mathcal{B} \) and \( J_{ij} \geq 0 \) if \( i \in \mathcal{A} \) and \( j \in \mathcal{B} \). Then \( S = |S_\mathcal{A} - S_\mathcal{B}| \), where \( S_\mathcal{A} \) and \( S_\mathcal{B} \) are the largest possible values of the spin on sublattices \( \mathcal{A} \) and \( \mathcal{B} \), respectively. Figure 2 shows three cases where this separation is particularly simple. In the special case of vanishing diagonal coupling, \( J_d = 0 \) (but \( J_l > 0, J_r > 0 \)),\n
The detailed nature of the RS-H transition in the frustrated spin 1/2 ladder has been a matter of debate. A first-order transition is well established in a wide region of parameter space, but for weak interchain coupling an intermediate phase, a “columnar dimer phase,” has been conjectured [38]. Numerous studies have since tried to find evidence for this phase, with little success [23,25,26,39,40]. More recent calculations

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fully agree with a single transition, of first order for not too small couplings [41–43]. Similarly the RS-H transition found for the mixed-spin model has been interpreted as a single first-order transition due to a level crossing of the RS and H singlet ground states, at the same time “some peculiarities of the RS-H transition in the mixed-spin system, as compared to the uniform-spin case” were noticed [37].

Our aim is to shed light on the RS-H transition in the mixed-spin ladder, using tools of quantum information theory, which are better suited for investigating subtle details of ladders [44] than, e.g., the dependence of bond energies on coupling parameters, as used in Ref. [37]. The remarkable concept of entanglement combined with new algorithms has improved our understanding of many-body systems in general and quantum phase transitions in particular [45–52]. The best-known measure of entanglement is the von Neumann entanglement entropy (EE) which is widely used to detect the ground state is an exact dimer state, a product of singlet-singlet ground states, at the same time “some peculiarities of the RS-H transition in the mixed-spin system, as compared to the uniform-spin case” were noticed [37].

FIG. 1. Pictorial representation of a mixed-spin ladder with different intra- and interleg exchange couplings. The blue and red dots represent the σ and τ spins, respectively. The index i numbers the unit cells.

FIG. 2. Sublattices A (full symbols) and B (empty symbols) for three special cases: (a) \( J_d = 0, J_i > 0, J_r > 0 \), (b) \( J_1 = 0, J_r > 0, J_d > 0 \), and (c) \( J_i = 0, J_r > 0, J_d > 0 \).
where \( n \) and \( n' \) label the legs, and the summations \( \sum_i \) run over unit cells (see Fig. 1). Here, \( J_i \) is the intra-leg exchange interaction between spins \( \sigma \) and \( \tau \), and the other two couplings refer to the inter-leg interactions, \( J_r \) on the rungs and \( J_d \) across the diagonals.

The Hamiltonian (1) has several symmetries, including SU(2), time-reversal, “leg-swap”, inversion and discrete translations. Some of them may be explicitly broken by boundary conditions. Nevertheless, for long enough ladders these symmetries may be partially restored (well inside the ladder if a correlation length exists which is much smaller than the number of unit cells).

If some symmetry is spontaneously broken we can define appropriate order parameters. Important additional information about the ground state can be gained from spin correlation functions. Those across the rungs are defined as

\[
S_n(i) := (\sigma_i^{(1)} \cdot \tau_i^{(2)}) \quad S_r(i) := (\tau_i^{(1)} \cdot \tau_i^{(2)}),
\]

where \( \langle \ldots \rangle \) denotes the expectation value with respect to the ground state. The “columnar” correlation functions (those along legs) are conveniently labeled by rung numbers \( \ell \) (instead of cell numbers \( i \))

\[
S_m(\ell) := \begin{cases} 
(\sigma^{(n)}_i \cdot \tau^{(n)}_i), & \ell = 2i - 1 \\
(\tau^{(n)}_i \cdot \sigma^{(n)}_{i+1}), & \ell = 2i 
\end{cases},
\]

where \( n = 1, 2 \) numbers the legs and \( \ell \) runs from 1 to \( L = 2N \). Similarly, the “diagonal” correlation functions are

\[
S_d1(\ell) := \begin{cases} 
(\sigma^{(1)}_i \cdot \tau^{(2)}_i), & \ell = 2i - 1 \\
(\tau^{(1)}_i \cdot \sigma^{(2)}_{i+1}), & \ell = 2i
\end{cases},
\]

and

\[
S_d2(\ell) := \begin{cases} 
(\sigma^{(2)}_i \cdot \tau^{(1)}_i), & \ell = 2i - 1 \\
(\tau^{(2)}_i \cdot \sigma^{(1)}_{i+1}), & \ell = 2i
\end{cases}.
\]

These correlation functions satisfy the inequalities

\[
-\frac{3}{4} \leq S_n(i) \leq \frac{1}{4}, \quad -2 \leq S_r(i) \leq 1,
\]

\[
-1 \leq S_m(\ell) \leq \frac{1}{4}, \quad a = l, d, \quad n = 1, 2.
\]

If the spatial symmetries are preserved in the ground state, the rung correlations are independent of the cell number \( i \), and both columnar and diagonal correlations do not depend on the rung number \( \ell \) (nor on \( n \)).

III. FRUSTRATED MIXED-SPIN PLAQUETTE

It is instructive to consider first the building block of the mixed-spin ladder, consisting of four spins on the corners of a square, two with spin 1 and two spin 1/2. Eigenstates of both the total spin and the Hamiltonian can be deduced analytically. The results may serve as a starting point for the construction of effective low-energy Hamiltonians.

We consider the general case where the rung couplings can be different \( (J_r \rightarrow J_r, J_t) \). The Hamiltonian

\[
H = J_s(\sigma_1 \cdot \sigma_2 + J_t \tau_1 \cdot \tau_2 + J_r(\tau_1 \cdot \sigma_1 + \tau_2 \cdot \sigma_2))
\]

\[
+ J_d(\tau_1 \cdot \sigma_2 + \tau_2 \cdot \sigma_1),
\]

and is invariant with respect to an interchange of spin operators, \( \tau_1 \leftrightarrow \tau_2, \sigma_1 \leftrightarrow \sigma_2 \) (“leg-swap symmetry”).

We classify the states of the 36-dimensional Hilbert space according to the eigenstates of \( S^2, S_z \) and the leg-swap operation. This basis is constructed explicitly in Appendix A. For \( S_z = 0 \) we obtain five states which are even under the leg-swap operation (one with \( S = 3 \), one with \( S = 2 \), three with \( S = 1 \)) and five states which are odd (two with \( S = 2 \), one with \( S = 1 \), two with \( S = 0 \)). The Hamiltonian is block-diagonal in this basis, with a 3 \( \times \) 3 matrix as largest block.

The ground state symmetry depends on the coupling constants. If all couplings are equal, \( J_t = J_d = J_r = J \), the Hamiltonian is directly related to the total spin,

\[
H = \frac{J}{2} \left( S^2 - \frac{11}{2} \right),
\]

and the ground state has \( S = 3 \) for \( J < 0 \) and \( S = 0 \) for \( J > 0 \).

The diagonalization of the Hamiltonian (Appendix A) yields the phase diagram of Fig. 3. In this paper, we concentrate ourselves mostly on the parameter region \( J_t = J_r = J_r = 1 \). Fig. 4 shows the energy spectrum for this case. Remarkably, all eigenvalues are linear functions of \( J_d \). At the fully symmetric point the lines collapse to the values predicted by Eq. (9). The ground state is always a spin singlet, but of different character above and below \( J_d = 1 \), where the singlet levels cross. It is worthwhile to add that for generic parameter sets the two singlet levels repel each other and are separated by a gap.

The spin correlation functions defined by Eqs. (2) to (4) provide valuable insight into the character of the eigenstates. They are readily evaluated using the Hellman-Feynman theorem. As a simple example we first consider the eigenstate with \( S = 3 \), which has an energy \( J_1 + J_d + \frac{1}{2} J_r + J_r \) and therefore correlation functions \( S_1 = S_0 = 1/2, S_0 = 1/4 \), and \( S_r = 1 \). These are just the upper bounds of Eq. (6). We now turn to the ground state. The correlation functions are calculated using

\[
\]
IV. WEAK AND STRONG RUNG COUPLINGS

We return now to the mixed-spin ladder. In this section, we investigate the limits of weak and strong rung couplings, which can be understood analytically. We limit ourselves to the unfrustrated ladder, $J_l = 0$, and consider three special cases, weak antiferromagnetic rung coupling, ferromagnetic rung coupling and strong antiferromagnetic rung coupling.

Our mixed-spin ladder is invariant under the exchange of leg and diagonal couplings [40], i.e., $H(J_l, J_r, J_d) = H(J_d, J_r, J_l)$, therefore the entire discussion below can be applied to the case of $J_l = 0$ and $J_d > 0$, by replacing $J_l$ by $J_d$.

A. Weak antiferromagnetic rung coupling

In the limit of vanishing rung coupling, the ladder is decoupled into two equivalent mixed-spin (1,1/2) chains. According to the Lieb-Mattis theorem [36], each chain has a total spin $S_{tot} = N/2$, where $N$ is the number of unit cells, and thus exhibits ferrimagnetic long-range order. Since the elementary cell (of a chain) consists of two spins, linear spin wave theory yields two types of magnons, a gapless “acoustic” branch with dispersion $\omega_k^a / J_l = -\frac{1}{4} + \left(\frac{1}{4} + 2 \sin^2 k\right)^{1/2}$ ($\sim k^2$ for small $k$), and a gapped “optical” branch with dispersion $\omega_k^o / J_l = 1/4 + \left(\frac{1}{4} + 2 \sin^2 k\right)^{1/2}$ [67–69]. When an antiferromagnetic rung coupling is switched on, spin wave theory predicts a linear dispersion of the gapless mode, reflecting the antiferromagnetic character of the ladder system, whereas the optical mode moves upward [70]. However, DMRG calculations show that a spin gap opens, which first increases quadratically as a function of $J_r$ up to $J_r \sim 0.3$, and then grows linearly [70].

B. Ferromagnetic rung coupling

For $J_r < 0$ and $J_d = 0$, the $\sigma$ spins form rung triplets and the $\tau$ spins form rung quintets, therefore the ground state is a product of rung-triplet and rung-quintet states. Low-energy excited states are rung singlets for $\sigma$ spins and rung triplets for $\tau$ spins, separated from the ground state by energy gaps of $J_r$ and $2J_r$, respectively. As soon as leg couplings are switched on, when $J_l \ll |J_r|$, the ladder behaves like a ferrimagnetic spin (1,2) chain, with long-range order, a total spin $S_{tot} = N$, an acoustic mode and an optical mode (at $\omega_0^a = 2J_l$). In the opposite limit of weak ferromagnetic rung coupling ($|J_l| \ll J_r$), the magnetic moments of the two chains are aligned, giving again $S_{tot} = N$, while the low-energy excitation spectrum remains essentially that of two independent chains, with acoustic and optical modes as described above. Hence we expect the two limits of weak and strong ferromagnetic rung couplings to be smoothly connected.

C. Strong antiferromagnetic rung coupling

For $J_r > 0$ and $J_d = 0$, the ground state is a product of local rung singlets, with an energy (per unit cell) of $-(11/4)J_r$. The first excited states are triplets, separated from the ground state by a finite energy gap of size $J_r$. With increasing leg coupling $J_r$ the energy gap decreases monotonically. Appendix B shows that first-order perturbation theory in $J_r$ does not give any contribution from quintets ($S = 2$) on the
Fig. 5. MPSs representation for the mixed-spin (1,1/2) ladder. The rungs with spins $\sigma = 1/2$ ($\tau = 1$) are indexed by $A$ and $C$ ($B$ and $D$). A spin-1/2 (1) rung has a local Hilbert space with dimension $d^A = d^C = 4$ ($d^B = d^D = 9$).

V. NUMERICAL METHODS

A. Matrix product states

Matrix product states (MPSs) [71,72] provide an efficient representation of the ground state of one-dimensional systems obeying the area law, for which the entanglement entropy grows with the boundary of a specific area rather than its volume [73]. To use MPSs for our model, we map the ladder onto a chain and consider each rung as a supersite with a larger Hilbert space. In Vidal’s representation [46], a generic state of a one-dimensional system is described in terms of two sets of matrices. We choose four pairs of matrices $(\lambda_A, \lambda_B, \lambda_C, \lambda_D)$, which allow us to include phases showing a doubling of the unit cell, such as a dimer solid or the ferrimagnetic state [2, 1, 1, ..., 2, 1, 1, 1] [37]. In this representation, illustrated in Fig. 5, an arbitrary state of our ladder is

$$|\psi\rangle = \sum_{i_1, \ldots, i_N} \left[ \cdots \Gamma^{n_4}_{A} \lambda^{A}_{i_4} \Gamma^{n_{4}+1}_{B} \lambda^{b}_{i_4} \Gamma^{n_{4}+2}_{C} \lambda^{C}_{i_4} \Gamma^{n_{4}+3}_{D} \lambda^{D}_{i_4} \cdots \right] |i_1\rangle \otimes \cdots \otimes |i_N\rangle,$$  \hspace{1cm} (12)

where $i_n$ numbers the states of rung $n$ ($i_n = 1, \ldots, 4$ for $\sigma$-rungs and $i_n = 1, \ldots, 9$ for $\tau$-rungs). The $\lambda$’s are diagonal matrices with the non-negative “Schmidt coefficients” $\lambda_i$ on the diagonal, and $\lambda_i^2$ are the eigenvalues of the reduced density matrix (pres = $T_{FB(A)}|\psi\rangle\langle\psi|$), where $A$ and $B$ are two halves of the ladder). The EE is directly connected to these eigenvalues through

$$S = -\sum_i \lambda_i^2 \ln \lambda_i^2.$$  \hspace{1cm} (13)

The matrices $\Gamma$ in (12) correspond to transformations between different Schmidt bases. The dimension $\chi$ of the matrices is a key parameter in tensor network states, called bond dimension, and the accuracy of the state (12) is controlled by this quantity. For weakly entangled states, a fairly small bond dimension is sufficient to obtain sensible results. But there are situations where very large bond dimensions are required, such as gapless critical systems [74].

B. Time-evolving block decimation

One of the efficient MPS-based algorithms for simulating one-dimensional quantum many-body systems is the iTEBD technique [46,48,75–77]. In iTEBD, using the imaginary time evolution of a quantum state we can find the ground state of the Hamiltonian $H$ through the relation

$$|GS\rangle = \lim_{\beta \to \infty} \exp(-\beta H)|\psi_0\rangle,$$  \hspace{1cm} (14)

where $|\psi_0\rangle$ is an initial “guess state,” chosen in the form of Eq. (12). If $|\psi_0\rangle$ is orthogonal to $|GS\rangle$, for instance because it has a different symmetry, the ground state cannot be reached by this method. Practical difficulties may appear in the vicinity of a continuous phase transition, where a judicious choice of the guess state is of crucial importance. However, for a first order transition between two gapped phases, such as that between RS and H phases, the method is expected to work well.

To proceed, we first rewrite the Hamiltonian (1) as

$$H = \sum_i \left( h_i^{(\sigma)} + h_i^{(\tau)} \right),$$  \hspace{1cm} (15)

where

$$h_i^{(\sigma)} = J_1 \sigma_1^{(i)} \cdot \sigma_1^{(i+1)} + J_2 \sum_{n \neq n'} \sigma_1^{(i)} \cdot \sigma_{1+n}'^{(i)} + J_3 \sum_{n \neq n'} \sigma_1^{(i)} \cdot \sigma_{1+n}'^{(i)},$$

$$h_i^{(\tau)} = J_4 \tau_1^{(i)} \cdot \tau_1^{(i+1)} + J_5 \sum_{n \neq n'} \tau_1^{(i)} \cdot \tau_{1+n}'^{(i)} + J_6 \sum_{n \neq n'} \tau_1^{(i)} \cdot \tau_{1+n}'^{(i)}.$$  \hspace{1cm} (16)

If the “time” $\beta$ is divided into a large number of intervals of width $\delta$ one can use the (first-order) Suzuki-Trotter decomposition

$$e^{-\beta H} \approx \prod_i e^{-\delta h_i^{(\sigma)}} e^{-\delta h_i^{(\tau)}}.$$  \hspace{1cm} (17)

Starting with an initial guess state of the form (12), we apply the operator $e^{-\delta H}$ iteratively to update the matrix product representation, until the ground state energy or the entropy converges. In our ladder system, we use the second-order Suzuki-Trotter decomposition. The simulations are started with a time step $\delta = 0.5$, which gradually is decreased to $\delta = 10^{-5}$. A key feature of the iTEBD algorithm is that it directly treats the infinite system by exploiting translational invariance, therefore it is free of finite-size effects.

C. DMRG

We also use the DMRG technique, especially to study the ground state in the vicinity of transition points. Unlike iTEBD, DMRG is a variational method. In other respects, the two methods have many steps in common. We iteratively optimize the MPSs of two neighboring sites to minimize the ground state energy, and then project the Hamiltonian onto a variational space. We use an iterative algorithm such as Lanczos to lower the energy. The two-site update is repeated for each pair of neighboring sites until the wave function converges to the ground state.
we introduce the dimensionless parameter $\chi$. In order to obtain the complete ground state phase diagram, we consider the regime with large magnetic moments on the rungs. We find in the entire range of bond dimensions, $\chi = 100$ and 120. The results for the two cases are almost the same. The inset of Fig. 6 confirms that the energy has well converged for $\chi = 120$. The cusps at $R = 0$ points to a phase transition of first order. Some insight on the magnetic order parameters are zero). The degeneracy of the ES levels offers a versatile tool both for identifying different phases and for locating quantum phase transitions [57,58,78–80]. As seen in Fig. 8, the level degeneracy of the ES is odd (1 or 3) in the RS phase ($0 < R < 1$), while in the F$_1$ phase ($-1 < R \leq 0$) the ES levels are nondegenerate.

A. iTEBD

We first discuss results obtained by the iTEBD method. The ground state energy is shown in Fig. 6 for two different bond dimensions, $\chi = 100$ and 120. The results for the two cases are almost the same. The inset of Fig. 6 confirms that the energy has well converged for $\chi = 120$. The cusps at $R = 0$ points to a phase transition of first order. Some insight on the two sides of the transition can be gained by calculating the magnetic moments on the rungs. We find in the entire range of bond dimensions, $\chi = 100$ and 120. The results for the two cases are almost the same. The inset of Fig. 6 confirms that the energy has well converged for $\chi = 120$. The cusps at $R = 0$ points to a phase transition of first order. Some insight on the magnetic order parameters are zero).

VI. UNFRUSTRATED LADDER

In the absence of diagonal interactions, our mixed-spin ladder (1) is unfrustrated and, as discussed in Sec. IV, its ground state is ferrimagnetically long-range ordered in the limit of strong antiferromagnetic rung couplings and magnetically disordered in the limit of strong antiferromagnetic rung couplings. In order to obtain the complete ground state phase diagram, we introduce the dimensionless parameter $R = J_l/(J_l + J_r)$. The limits $R = \pm 1$ correspond to the strong rung coupling regimes, whereas $R = 0$ is the limiting case of two decoupled mixed-spin chains.

A. iTEBD

We first discuss results obtained by the iTEBD method. The ground state energy is shown in Fig. 6 for two different bond dimensions, $\chi = 100$ and 120. The results for the two cases are almost the same. The inset of Fig. 6 confirms that the energy has well converged for $\chi = 120$. The cusps at $R = 0$ points to a phase transition of first order. Some insight on the two sides of the transition can be gained by calculating the magnetic moments on the rungs. We find in the entire range of bond dimensions, $\chi = 100$ and 120. The results for the two cases are almost the same. The inset of Fig. 6 confirms that the energy has well converged for $\chi = 120$. The cusps at $R = 0$ points to a phase transition of first order. Some insight on the magnetic order parameters are zero).

B. DMRG

We now discuss results obtained with DMRG, which is less sensitive to initial guess states than iTEBD. Computations were performed for ladders of different sizes with open boundary conditions. For greater efficiency we increased the bond dimension to 1000. The ground state energies obtained with DMRG (after extrapolation) match perfectly those calculated with iTEBD, except in the tiny region of $R$ discussed above, where the DMRG values are in general smaller.

The EE is obtained by cutting the ladder into two halves, and tracing out the degrees of freedom of one of the halves. The results presented in Fig. 7 exhibit a discontinuity at $R = 0$ where a phase transition of first order occurs from the ferrimagnetically ordered F$_1$ phase to the RS phase (where all magnetic order parameters are zero).

The degeneracy of the ES levels offers a versatile tool both for identifying different phases and for locating quantum phase transitions [57,58,78–80]. As seen in Fig. 8, the level degeneracy of the ES is odd (1 or 3) in the RS phase ($0 < R < 1$), while in the F$_1$ phase ($-1 < R \leq 0$) the ES levels are nondegenerate.

VII. FRUSTRATED LADDER

We now consider the case where all three exchange couplings are finite. Some regions in parameter space can be understood without detailed calculations, especially those with no or small frustration. Thus for $J_d = J_l$ our ladder model can be mapped onto a generalized mixed-spin Heisenberg chain. Its ground state can be guessed in a simple way both for
For $J_d < J_j$ and for $0 < J_j \ll J_d$. Regions with strong frustration, such as those where the three coupling parameters are all positive and of a similar size, are of course difficult to handle using qualitative arguments. We therefore resort to numerical calculations, limiting ourselves to the region $J_j = J_d = 1$.

A. Generalized Heisenberg chain

For $J_d = J_j$, it is convenient to introduce rung spins

$$S_i := \sigma_i^{(1)} + \sigma_i^{(2)}, \quad T_i := \tau_i^{(1)} + \tau_i^{(2)},$$

(18)

because in this case the Hamiltonian can be written in terms of these operators,

$$H = \frac{J_j}{2} \sum_i (S_i^2 + T_i^2) + J_d \sum_i (S_i \cdot T_i + T_i \cdot S_{i+1}),$$

(19)

where we have neglected an additive constant. The rung spin operators do not have a fixed “length”, but can assume the values $S_i = 0, 1, T_i = 0, 1, 2$. For $J_j < 0$, the first term has the lowest eigenvalue if both $S_i$ and $T_i$ are as large as possible, i.e., $S_i = 1$ and $T_i = 2$. But this is also true for the second term (it is obvious for classical spins or for the Néel state). Therefore the ground state is expected to be ferrimagnetic ($F_1$) if $J_j < 0$ (and $J_d = J_j > 0$). For positive $J_j$, the first term in the Hamiltonian favors singlet rung spins, in contrast to the second term which is lowest for maximal rung spins. The system is frustrated. However, for very small positive values of $J_j$ the first term can be neglected and we obtain again a ferrimagnetic phase of type $F_1$.

Our numerical analysis of nearest-neighbor-spin correlations on the rungs imply that $\langle T_i^2 \rangle \approx \langle S_i^2 \rangle$ for $0 \leq J_d \leq 0.86$, which implies that the rung-quantum states on the spin-1/2 rungs have negligible weight and can be projected out. This lends support to the mapping of the frustrated mixed-spin $(1, 1/2)$ ladder onto a frustrated uniform spin-1/2 ladder, as proposed in Ref. [37].

B. iTEBD

We have studied numerically the ground state energy, both the EE and the Schmidt gap (the difference between the two largest eigenvalues of the reduced density matrix [81]), as well as the level degeneracy of the ES. We therefore present the overall picture obtained with iTEBD. More details will be given in Sec. VIII for DMRG results, in particular regarding the intermediate phase.

Figure 9 shows the ground state energy as a function of $J_d$. Cusps at certain values of $J_d$ (indicated by the dashed vertical lines), are indicative for phase transitions. Corresponding jumps in the EE are more pronounced, and found to occur at $J_d \approx 0.74, 0.86, 1.17$. The gross features of the ground state energy – an almost linear increase for $J_d \lesssim 0.7$ and a linear decrease for $J_d \gtrsim 0.7$ – match the behavior found for the plaquette (Fig. 4). The different values of the “critical points” can be attributed to the fact that in the ladder a site is connected to two neighbors by $J_d$ and only to one in the plaquette. The situation is reminiscent of that encountered for the antiferromagnetic XX chain in a transverse field $h$, where the transition to the ferromagnetic state occurs at $J = 2h$ for two sites, but already at $J = h$ for the chain. An additional similarity between the plaquette and the ladder is the vanishing of the excitation gap at criticality, observed as a level crossing in the case of the plaquette and by numerical evidence in the case of the ladder (for the same specific coupling parameters as used here) [37].

We have also studied both the ES and local order parameters in the different phases. No magnetic order was found both for $0 \leq J_d < 0.71$ and for $0.74 < J_d \leq 0.86$. The low-lying levels of the ES are odd-degenerate for $0 \leq J_d < 0.71$ (as in the RS phase [57,78]), and even-degenerate for $0.74 < J_d \leq 0.86$ (as in the H phase). For $J_d > 0.86$ the ES is nondegenerate and magnetic order does exist. A sharp first-order transition at $J_d \approx 1.17$ separates two different magnetic phases, the $F_1$ phase corresponding to the sector $[1, 2, \ldots, 1, 2]$ for $J_d \gtrsim 1.17$, and the $F_2$ phase corresponding to the sector $[1, 1, 1, 2, \ldots, 1, 1, 1, 2]$ for $0.86 < J_d \lesssim 1.17$. The $F_2$ phase breaks the translational symmetry and can be considered an intermediate phase between the $H$ and
F₁ phases. This picture agrees with the phase diagram of Ref. [37], although the transition point between F₁ and F₂ phases is somewhat different.

In the narrow interval 0.71 ≤ J_d < 0.74, the ES levels are not found to exhibit any characteristic degeneracy, neither of the RS-type nor of the H-type, and we often detect magnetic order. We attribute the erratic data to limitations of the method. In fact, the spin gap seems to be very small in this region [37], therefore one would need both a large parameter β and a large bond dimension χ to obtain consistent results, very much like in the unfrustrated case for J_d = 0, |J_r| ≪ J_l. Moreover, in our ansatz we included a possible period doubling, which allowed us to reproduce the F₂ phase, but excluded ground states with longer periods, for which we find good evidence on the basis of DMRG data (to be discussed below).

VIII. INTERMEDIATE SINGLET PHASE

As discussed above, for 0 < J_d ≤ 0.86 the ground state of our model appears to be well represented by that of a frustrated spin-1/2 ladder (with couplings J_l′, J_r′, J_d′) [37], which has been intensively investigated using both analytical and numerical methods [18,25,26]. An intermediate columnar-dimer phase has been reported only in a narrow neighborhood of J_r′ = 0.38 (for J_l′ = 1 and J_d′ = 0.2) [26], i.e., in a “weak coupling” region in parameter space which does not correspond to the “strong coupling” region of our intermediate phase. This apparent discrepancy is less serious if one keeps in mind that the low-energy excitations are quite different in the two models (the spin gap remains finite at the RS-H transition in the spin-1/2 ladder [40], in contrast to the softening observed in the mixed-spin ladder [37]). Therefore the (approximate) mapping between the ground states of the two models cannot be used to rule out a strong-coupling intermediate phase in the mixed-spin ladder.

We now discuss DMRG data for the frustrated mixed-spin ladder, focusing on the parameter region J_l = J_r = 1, 0.4 < J_d < 1. All magnetic order parameters ⟨τ_i^(n)⟩ and ⟨σ_i^(n)⟩ vanish for J_d ≲ 0.86, and the existence of an intermediate phase is confirmed on the basis of spatially modulated spin correlations.

A. EE, Schmidt gap and ES

We first use tools of quantum information theory for pinpointing the intermediate singlet phase. Results for the EE as well as the Schmidt gap are presented in Fig. 10. The sharp changes of the EE at J_d ≃ 0.707 and J_d ≃ 0.73 indeed are clear signatures for a distinct phase, squeezed in between RS and H phases. It will be shown later that dimerization plays a role, and therefore we refer to this new ground state as the D phase. At the first boundary, J_d = 0.707, the EE changes continuously, which points to a second-order transition from the RS to the D phase. At the second boundary, J_d = 0.73, a discontinuous jump indicates a first-order transition from the D to the H phase. The Schmidt gap is almost constant in the RS phase, decreases gradually in the D phase, and tends to 0 in the H phase.

The ES is illustrated in Fig. 11. The low-lying ES levels have odd and even degeneracies in the RS and H phases, respectively, and are nondegenerate in the ferrimagnetic phases. In the D phase, the ES levels have a mixed even-odd degeneracy.

FIG. 10. EE (top) and Schmidt gap (bottom) vs J_d, for different sizes, L = 120, 180, and 240, obtained by DMRG with χ gradually increased from 300 to 1000.

FIG. 11. DMRG results for the low-lying ES levels and their degeneracies (marked by circles).
FIG. 12. Spin correlations across τ-rungs (upper panel) and across σ-rungs (lower panel), as functions of the cell number $i$, obtained by DMRG.

B. Short-range spin correlations

Figure 12 displays the rung correlations as functions of the cell number $i$ for different values of $J_d$. Clearly, boundary effects are limited to a few cells. Both for $J_d \leq 0.707$ and for $J_d > 0.73$, the data do not depend on $i$; for small $J_d$ they are close to their lower (singlet) bounds while for $J_d \gtrsim 0.8$, they approach values corresponding to spin 1. In the D phase, $0.707 < J_d < 0.73$, the rung correlations are no longer constant but oscillate with a $J_d$-dependent wave vector.

Both columnar and diagonal correlation functions, depicted in Fig. 13, are $\ell$-independent outside the D phase and show incommensurate modulations inside. However, there is an additional rapid oscillation—a weak dimerization—which breaks the inversion symmetry. Both correlation functions are found to be independent of $n$, therefore the leg-swap symmetry is not broken. The translational symmetry is of course broken both by the slow incommensurate modulations and by the dimerization.

We have repeated our computations for ladders of larger sizes (for example, $L = 240$) and also for ladders with odd $L$. The main features, in particular the oscillations in the bulk, remain the same.

Similar incommensurate oscillations have been seen in DMRG results for an SU(3) spin ladder, where they appear in an intermediate phase between a valence-bond crystal for small rung couplings and a critical Luttinger liquid for large rung couplings [82].

C. Interpretation

To quantify the rapid oscillations of columnar and diagonal correlation functions, we introduce staggered order parameters

$$S_a(i) := \sum_n [S_{an}(2i - 1) - S_{an}(2i)], \quad a = l, d. \quad (20)$$

Both $S_l(i)$ and $S_d(i)$ vanish in the RS and H phases, but are finite in the D phase. Figure 14 shows that these functions oscillate, actually with the same wave vectors $q$ as the rung correlations (Fig. 12). The oscillations are sinusoidal and therefore we could also state that the original correlation functions $S_{an}(\ell)$ have an oscillatory component with wave vector $\pi - 2q$. This reminds us of the BOW in nearly half-filled Peierls systems, where a commensurate-incommensurate transition from bond alternation at half filling to an incommensurate harmonic oscillation away from half filling occurs. Very close to half filling the bond-order wave is not simply sinusoidal but has the form of a “soliton lattice,” consisting of relatively wide regions with constant order parameter and narrow domain walls in which the order parameter changes rapidly. We do not see any evidence for domain walls, maybe simply because when $q$ approaches 0 the amplitude also tends to 0.

We have deduced the wave vector $q$ for the staggered order parameters by fitting sinusoidal functions to the data points far from the edges of the ladder. The result shown in Fig. 15
FIG. 14. Staggered bond orders along legs (top) and diagonals (bottom).

is consistent with a slow increase from zero at the lower boundary of the intermediate phase and a rapid fall to zero at the upper boundary, similarly to the behavior reported in Ref. [82]. Figure 15 also shows the “BOW parameter” $B$, the oscillation amplitude at the center of the ladder. Clearly $B$ is only nonzero in the D phase.

An important question is whether finite size effects are responsible for the intermediate phase. This has been an issue in the context of the Hubbard model on the honeycomb lattice, for which Meng and coworkers reported an intermediate spin liquid phase for relatively small system sizes [83], while subsequent work by Sorella and others showed that this phase disappears in larger systems [84]. Therefore we have investigated the size dependence of the oscillation amplitudes. The results for the BOW parameter $B$, shown in Fig. 15, indicate that the intermediate phase survives in the thermodynamic limit. However, the limiting value of $B$ is rather small and therefore additional studies for larger system sizes would be very useful for strengthening the case for these incommensurate spin patterns.

IX. SUMMARY AND OUTLOOK

In this paper, we have described our study of a frustrated mixed-spin ladder, consisting alternatively of spin $1/2$ and spin 1 rungs. We have used tools from quantum information, in particular the entanglement entropy (EE), the Schmidt gap and the entanglement spectrum (ES), as well as correlation functions to characterize the various phases. Three types of interactions were taken into account, one along legs (coupling $J_l$), one on rungs ($J_r$) and one on diagonals ($J_d$). We limited ourselves mostly on the region $J_l = J_r = 1, J_d \geq 0$. Three distinct nonmagnetic phases were identified, a gapped RS phase with odd degeneracy of the ES levels (for $J_d \lesssim 0.7$), a gapped H phase with even degeneracy of the ES levels (for $J_d \gtrsim 0.7$), and an intermediate D phase with mixed even-odd degeneracies of the ES levels (for $J_d \approx 0.7$).

The overall behavior of short-range correlation functions is depicted in Fig. 16. Three bonds are quite prominent, the singlet bond on the $\tau$ rung and the “magnetic” bonds on the diagonals, both in the RS phase, as well as the triplet bond on the $\sigma$ rung in the H phase. Therefore, with one exception,

FIG. 16. Pictorial representation of spin correlations in the unit cell (defined in Fig. 1) for coupling constants $J_l = J_r = 1$ and $J_d = 0.5$ (left figure, RS phase), $J_d = 0.715$ (middle figure, D phase), $J_d = 0.85$ (right figure, H phase). The grayscale reaches from the white shade (lower bound of the correlations – singlet bonds) to the black shade (upper bound of the correlations – magnetic bonds).
the bonds are far from being singlets, the favorite state of two antiferromagnetically coupled spins. This is a clear signature for frustration.

The most striking features of our study are the incommensurate spatial oscillations of spin correlation functions, observed only in the D phase. Their wave vector depends on the coupling strength. Their amplitude can serve as an order parameter; its behavior close to the phase boundaries suggests that the transition is continuous at the RS-D boundary and of first order at the D-H boundary.

Our choice of coupling parameters is quite special, as became apparent in our calculations for the elementary plaquette (Sec. III). In fact, for this choice the energy spectrum of the plaquette is “integrable” (level crossing but no level repulsion), for other couplings it is generically nonintegrable (level repulsion but no level crossing). Therefore it would be desirable to study the ladder in a wider region of parameter space.

We have used relatively small system sizes. It is true that the extrapolation to the thermodynamic limit did preserve a finite order parameter in the D phase, at the same time its small value is worrisome. Clearly, additional calculations for larger systems would be very useful.

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APPENDIX A: DIAGNOSTICIZATION OF THE PLAQUETTE HAMILTONIAN

The ten eigenstates of the Hamiltonian (7) with $S_z = 0$ can be grouped into even and odd states (with respect to the leg-swap operation), namely,

$$|a\rangle_{\pm} = \frac{1}{\sqrt{2}}(|00 \downarrow\downarrow \rangle \pm |00 \uparrow\uparrow \rangle),$$

$$|b\rangle_{\pm} = \frac{1}{\sqrt{2}}(|+ - \uparrow\downarrow \rangle \pm |- + \downarrow\uparrow \rangle),$$

$$|c\rangle_{\pm} = \frac{1}{\sqrt{2}}(|- + \uparrow\downarrow \rangle \pm |+ - \downarrow\uparrow \rangle),$$

$$|d\rangle_{\pm} = \frac{1}{\sqrt{2}}(|0- \uparrow\uparrow \rangle \pm |0 + \downarrow\downarrow \rangle),$$

$$|e\rangle_{\pm} = \frac{1}{\sqrt{2}}(|00 \uparrow\downarrow \rangle \pm |00 \downarrow\uparrow \rangle).$$  \hspace{1cm} \text{(A1)}

Once the eigenstates in this subspace are determined, those for $S_z \neq 0$ are easily obtained by applying $S_z$ and $S_\pm$. These new states are uninteresting as long as we are just searching for the eigenvalues of $S^2$ and $H$, because if $|\Psi\rangle$ is an eigenstate of $S^2$ and $H$, the same holds for $S_z|\Psi\rangle$, with the same eigenvalues.

It is straightforward to calculate eigenstates and eigenvalues of $S^2$ in this basis. We find one state with $S = 3$, one with $S = 2$ and three with $S = 1$ from the even basis states,

$$|3, 0\rangle = \frac{1}{\sqrt{5}} \left[ |a\rangle + |d\rangle + \frac{1}{\sqrt{2}} (|b\rangle + |c\rangle) + \sqrt{2} |e\rangle \right],$$

$$|2, 0\rangle_+ = \frac{1}{\sqrt{2}} (|a\rangle - |d\rangle),$$

$$|1, 0\rangle_{(1)} = \frac{1}{\sqrt{2}} (|b\rangle - |c\rangle),$$

$$|1, 0\rangle_{(2)} = \frac{1}{2} (|a\rangle + |d\rangle - \sqrt{2} |e\rangle),$$

$$|1, 0\rangle_{(3)} = \frac{1}{\sqrt{3}} \left[ \frac{1}{2} (|a\rangle + |d\rangle) - \sqrt{2} (|b\rangle + |c\rangle) \right] + \frac{1}{\sqrt{2}} |e\rangle. \hspace{1cm} \text{(A2)}$$

The odd basis states yield two eigenstates of $S^2$ with $S = 2$, one with $S = 1$ and two singlet states,

$$|2, 0\rangle_{(1)} = \frac{1}{\sqrt{6}} \left[ |a\rangle + |d\rangle + \sqrt{2} (|b\rangle + |c\rangle) \right],$$

$$|2, 0\rangle_{(2)} = \frac{1}{\sqrt{6}} (|b\rangle - |c\rangle + 2 |e\rangle),$$

$$|1, 0\rangle_- = \frac{1}{\sqrt{2}} (|a\rangle - |d\rangle),$$

$$|0, 0\rangle_{(1)} = \frac{1}{\sqrt{3}} \left[ |a\rangle + |d\rangle - \frac{1}{\sqrt{2}} (|b\rangle + |c\rangle) \right],$$

$$|0, 0\rangle_{(2)} = \frac{1}{\sqrt{3}} (|b\rangle - |c\rangle - |e\rangle). \hspace{1cm} \text{(A3)}$$

We now determine the eigenstates and eigenvalues of the Hamiltonian using these basis states. We find for the “nongenerate” states

$$H|3, 0\rangle = \left( J_t + J_d + \frac{1}{2} J_\tau \right) |3, 0\rangle,$$

$$H|2, 0\rangle_+ = \left[ - \frac{1}{2} (J_t + J_d) + \frac{1}{2} J_\tau \right] |2, 0\rangle_+,$$

$$H|1, 0\rangle_- = \left[ - \frac{1}{2} (J_t + J_d) + \frac{1}{2} J_\tau - J_\tau \right] |1, 0\rangle_- \hspace{1cm} \text{(A4)}$$

For the “doubly degenerate” states, we obtain

$$H|2, 0\rangle_{(1)} = \left[ \frac{1}{2} (J_t + J_d) + 4 J_\tau \right] |2, 0\rangle_{(1)} + \frac{1}{\sqrt{2}} (J_t - J_d) |2, 0\rangle_{(2)}, \hspace{1cm} \text{(A5)}$$

$$H|2, 0\rangle_{(2)} = \frac{1}{\sqrt{2}} (J_t - J_d) |2, 0\rangle_{(1)} - \frac{3}{4} J_\tau \right] |2, 0\rangle_{(2)}.$$
and therefore we have just to diagonalize $2 \times 2$ matrices. For the singlet states, we find

$$E = \frac{1}{2} \left( h_{11} + h_{22} \pm \sqrt{(h_{11} - h_{22})^2 + 4h_{12}^2} \right),$$  \hspace{1cm} (A7)

where

$$h_{11} = -J_i - J_d + \frac{1}{2} J_\sigma - J_\tau,$$

$$h_{22} = \frac{1}{2} J_\sigma - 2 J_\tau,$$

$$h_{12} = \sqrt{2}(J_d - J_\tau).$$  \hspace{1cm} (A8)

The remaining “triply degenerate” states satisfy the eigenvalue equation

$$H(1, 0)^{(a)}_+ = 3 \sum_{\beta=1}^3 h_{a\beta} |1, 0)^{(b)}_+ , \alpha = 1, 2, 3$$  \hspace{1cm} (A9)

with matrix elements

$$h_{11} = -\left( \frac{3}{4} J_\sigma + J_\tau \right),$$

$$h_{12} = h_{21} = \frac{1}{2} (J_d - J_\tau),$$

$$h_{13} = h_{31} = \frac{\sqrt{2}}{2} (J_d - J_\tau),$$

$$h_{22} = \frac{1}{4} [-5(J_i + J_d) + J_\sigma + 2 J_\tau],$$

$$h_{23} = h_{32} = \frac{\sqrt{2}}{4} (-J_i - J_d + 2 J_\tau),$$

$$h_{33} = \frac{1}{4} (-J_i - J_d + J_\sigma - 6 J_\tau).$$  \hspace{1cm} (A10)

Mmiraculously, for the particular parameter set $J_i = J_d = J_\sigma = J_\tau = 1$ all eigenvalues are simple linear functions of $J_d$, as shown in Table II and illustrated in Fig. 4.

### APPENDIX B: EXPANSION IN POWERS OF $J_i$ AND $J_d$ FOR $J_i > 0$

We use perturbation theory to evaluate the ground state for antiferromagnetic rung coupling, $J_i > 0$, and small leg coupling, $J_d \ll J_i$. A rung with $\sigma$-spins has four states, a singlet ($S = 0$)

$$|s_\alpha\rangle = \frac{1}{\sqrt{2}} (|\uparrow\rangle|\downarrow\rangle - |\downarrow\rangle|\uparrow\rangle),$$  \hspace{1cm} (B1)

and three triplet states ($S = 1$)

$$|t^+\rangle = |\uparrow\rangle|\uparrow\rangle, \quad |t^-\rangle = |\downarrow\rangle|\downarrow\rangle,$$

$$|t^0\rangle = \frac{1}{\sqrt{2}} (|\uparrow\rangle|\downarrow\rangle + |\downarrow\rangle|\uparrow\rangle),$$  \hspace{1cm} (B2)

where we have used the convention that in the product states the first ket is on chain 1, the second on chain 2.

A rung with $\tau$-spins has nine states, a quintet ($S = 2$)

$$|q^{\pm}\rangle = |\pm\rangle|\pm\rangle,$$

$$|q^{\pm\pm}\rangle = \frac{1}{\sqrt{2}} (|\pm\rangle|0\rangle + |0\rangle|\pm\rangle),$$

$$|q^0\rangle = \frac{1}{\sqrt{6}} (|+\rangle|-\rangle + 2|0\rangle|0\rangle + |-\rangle|+\rangle),$$  \hspace{1cm} (B3)

a triplet ($S = 1$)

$$|t^\pm\rangle = \pm \frac{1}{\sqrt{2}} (|\pm\rangle|0\rangle - |0\rangle|\pm\rangle),$$

$$|t^0\rangle = \frac{1}{\sqrt{2}} (|+\rangle|-\rangle - |-\rangle|+\rangle),$$  \hspace{1cm} (B4)

and a singlet ($S = 0$)

$$|s_\tau\rangle = \frac{1}{\sqrt{3}} (|0\rangle|0\rangle - |+\rangle|-\rangle - |-\rangle|+\rangle),$$  \hspace{1cm} (B5)

where $|\pm\rangle$ and $|0\rangle$ are the eigenstates of $\tau_z$ with eigenvalues ±1 and 0, respectively.

For $J_i = J_d = 0$ and $J_\tau > 0$ is a product of rung singlets $\bigotimes_i |s_\sigma_i\rangle|s_\tau_i\rangle$. To first order in $J_i$ we have to apply the operator $\sigma^{(1)} \cdot \tau^{(1)} + \sigma^{(2)} \cdot \tau^{(2)}$ to each pair of adjacent rungs. We find

$$\langle \sigma^{(1)} \cdot \tau^{(1)} + \sigma^{(2)} \cdot \tau^{(2)} | s_\tau \rangle = \sqrt{2} \left( |t^+\rangle|t^0\rangle - |t^-\rangle|t^0\rangle - |t^0\rangle|t^+\rangle + |t^0\rangle|t^-\rangle \right).$$  \hspace{1cm} (B6)

Triples appear in first-order perturbation theory for the ground state, but there are no quintets.

To second order in $J_i$ three different terms are generated. If the operator $\sigma^{(1)} \cdot \tau^{(1)} + \sigma^{(2)} \cdot \tau^{(2)}$ acts on four distinct rungs, the expression (B6) simply appears twice and no quintets are produced. If the same operator is applied twice to a single pair of rungs, quintets neither appear. This is easily understood by applying $\sigma^{(1)} \cdot \tau^{(1)} + \sigma^{(2)} \cdot \tau^{(2)}$ to Eq. (B6). In the third case, where three neighboring rungs are involved, say, at sites $i - 1, i, i + 1$, quintet states do appear.
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