Diffusive high-temperature transport in the one-dimensional Hubbard model

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We consider charge and spin transport in the one-dimensional Hubbard model at infinite temperature, halffilling, and zero magnetization. Implementing matrix-product-operator simulations of the nonequilibrium steady states of boundary-driven open Hubbard chains for up to 100 sites we find clear numerical evidence of diffusive transport for weak driving and any (nonzero and finite) value of the interaction U. In addition, nonequilibrium steady states are characterized by nonzero spin-up spin-down density-density correlations. For maximal driving, on the other hand, we find an insulating behavior and cosine-shaped density profiles.

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I. INTRODUCTION

The one-dimensional (1D) Hubbard model is the simplest model of strongly interacting dynamics of spinful fermions on a lattice within a single-band approximation. Its Hamiltonian reads

$$H = -t \sum_{i=1}^{L-1} \sum_{s \in \{\uparrow,\downarrow\}} (c_{i,s}^{\dagger} c_{i+1,s} + \text{H.c.}) + U \sum_{i=1}^{L} n_{i\uparrow} n_{i\downarrow}, \quad (1)$$

where the operators $c_{i,s}$, $c_{i,s}^{\dagger}$, for spin $s \in \{\uparrow,\downarrow\}$ and site $i \in \{1, \ldots, L\}$ are standard fermionic annihilation and creation operators and $n_{i,s} := c_{i,s}^{\dagger} c_{i,s}$ are the density operators. The model is solvable by a coordinate Bethe ansat z^1 and possesses an infinite number of conservation laws.² While stationary properties of the 1D Hubbard model are well understood, see Ref. 3, much less is known about its dynamics, for instance, about the transport behavior. In thermodynamically the most interesting regime, at half-filled band and zero magnetization $\sum_{i=1}^{L} \langle n_{i,s} \rangle = L/2$, studied in the present paper, the model is gapped for charge excitations and gapless for spin excitations, for any $U \neq 0$. At zero temperature, it is, therefore, an example of a Mott (charge) insulator and ballistic (ideal) spin conductor. Transport can be qualitatively characterized by a Drude weight—a linear response indicator of ballistic transport, defined as the weight of zero-frequency singular term $\delta(\omega)$ in the real part of conductivity $\sigma(\omega)$. Spin and charge Drude weights at zero temperature have been calculated in Ref. 4, with finite-size corrections given in Ref. 5, while a regular part of $\sigma(\omega)$ is studied in Ref. 6. At nonzero temperature, on the other hand, no rigorous result is known and there is no consensus between numerical and Bethe ansatz based results. Thermodynamic Bethe ansatz suggests⁷ that, even at half-filling, the charge Drude weight is finite, so the model was predicted to exhibit ideal charge transport; for a similar conclusion, see also quantum Monte Carlo calculation in Ref. 8. Analytical calculations at large U on the other hand support vanishing charge Drude weight;⁹ for a study of low-energy excitations, see also Ref. 10. Exact numerical simulations of small systems, again at half-filling and at high/infinite temperature, suggest $^{11} \sim 1/L$ scaling of the charge Drude weight. For temperatures much smaller than the gap, semiclassical arguments together with field-theoretical scattering rates predict diffusive transport.¹² Vanishing finitetemperature Drude weight in thermodynamic limit (TL) $L \rightarrow \infty$ offers the possibility of an insulating or diffusive behavior.

However, at high or infinite temperatures, nonequilibrium transport properties of 1D Hubbard model in either charge or spin sector are not known as there has been up to date no analytical or numerical method capable of reliably treating this regime. In this paper, we employ nonequilibrium steady state simulations¹³ using an efficient matrix product ansatz¹⁴ for the time-dependent density matrix and provide a clear evidence of diffusive transport for both attractive U < 0 and repulsive U > 0 cases at infinite temperature and weak driving (corresponding to the regime of linear response). Namely, we show 1/L scaling of charge as well as of spin current and clear linear density profiles. We also briefly discuss density-density correlations, exotic transport behavior in the regime of strong driving, and ballistic transport away from half-filling.

II. BOUNDARY DRIVEN HUBBARD CHAIN

Using the Jordan-Wigner transformation, we can map the 1D Hubbard model (1) to a spin-1/2 ladder system. Namely, writing $c_{i\uparrow} = P_{i-1}^{(\sigma)}\sigma_i^-$, where $P_i^{(\sigma)} = \sigma_1^z \cdots \sigma_i^z$ for spin-up fermions, and $c_{i\downarrow} = P_L^{(\sigma)}P_{i-1}^{(\tau)}\tau_i^-$, where $P_i^{(\tau)} = \tau_1^z \cdots \tau_i^z$ for spin-down fermions, one can verify that fermionic operators $c_{i,s}, c_{i,s}^{\dagger}$ satisfy canonical anticommutation relations provided σ_i^{α} and τ_i^{α} are two sets of Pauli matrices [and $\sigma_j^{\pm} := (\sigma_j^x \pm i\sigma_j^y)/2$, $\tau_j^{\pm} := (\tau_j^x \pm i\tau_j^y)/2$]. Writing the Hubbard Hamiltonian (1) in spin-ladder form, one obtains

$$H = -\frac{t}{2} \sum_{i=1}^{L-1} \left(\sigma_i^{x} \sigma_{i+1}^{x} + \sigma_i^{y} \sigma_{i+1}^{y} + \tau_i^{x} \tau_{i+1}^{x} + \tau_i^{y} \tau_{i+1}^{y} \right) + \frac{U}{4} \sum_{i=1}^{L} \left(\sigma_i^{z} + 1 \right) \left(\tau_i^{z} + 1 \right).$$
(2)

The spin-1/2 ladder system consists of two XX chains in two legs and a Z-Z type interchain coupling along the rungs. For numerical simulations of the Hubbard model, we shall use this ladder formulation (2).

To induce a nonequilibrium situation, two legs are coupled to independent reservoirs. Their action is decribed in an effective way via the Lindblad equation¹⁵ for the density matrix ρ of the ladder system:

$$\frac{d}{dt}\rho = \mathbf{i}[\rho, H] + \mathcal{L}^{\text{dis}}(\rho), \tag{3}$$

where the dissipative term is expressed in terms of Lindblad operators L_k , as

$$\mathcal{L}^{\text{dis}}(\rho) = \sum_{k} ([L_k \rho, L_k^{\dagger}] + [L_k, \rho L_k^{\dagger}]).$$
(4)

We use eight Lindblad operators acting locally on the first and last sites of each leg, injecting or absorbing fermions (spinons) with certain probability:

$$L_{1,2} = \sqrt{\Gamma(1 \mp \mu)} \sigma_1^{\pm}, \quad L_{3,4} = \sqrt{\Gamma(1 \pm \mu)} \sigma_L^{\pm}$$
 (5)

for the first and

$$L_{5,6} = \sqrt{\Gamma(1 \mp \mu)} \, \tau_1^{\pm}, \quad L_{7,8} = \sqrt{\Gamma(1 \pm \mu)} \, \tau_L^{\pm}$$
 (6)

for the second leg. Γ is the strength of the coupling to the baths, while μ is a driving strength playing the role of a chemical potential bias. As demonstrated in previous studies of 1D spin chains¹³ the precise form of Lindblad operators does not influence the bulk properties. Because of dissipative terms the time-dependent solution $\rho(t)$ of the Lindblad equation converges after a long time to a time-independent state called a nonequilibrium steady state (NESS), $\rho_{\infty} = \lim_{t \to \infty} \rho(t)$, which is unique.¹⁶ Once the steady state is reached, expressed in terms of a matrix product operator of a given bond dimension (following the method described in detail in Ref. 13 straightforwardly adapted for the spin ladder), expectation values of arbitrary observables in the NESS can be efficiently evaluated. All expectation values considered in this paper are taken with respect to the NESS, that is $\langle A \rangle = \operatorname{tr}(\rho_{\infty}A)$, which we will-when it is clear from the context, and to simplify notation-denote just by A. In each NESS calculation, we have carefully checked that the convergence is reached, i.e., we evolve the Lindblad equation (3) until a time-independent state is obtained, and that the results are stable with respect to increasing bond dimension.¹⁸ For $\mu = 0$, i.e., no driving, one has an equilibrium setting, resulting in a trivial NESS $\rho_{\infty} \propto 1$. This means that for small driving μ we are studying nonequilibrium behavior at an infinite temperature. Note that such an infinite temperature state is separable in the operator space. Since the efficiency of the numerical method crucially depends on the entanglement infinite-temperature nonequilibrium states are the easiest ones to calculate because the entanglement is expected to be smaller than at finite temperatures.

Expectation values of fermionic observables are obtained from the corresponding ones in the ladder formulation, for instance, particle densities are $n_{i\uparrow} = (\sigma_i^z + 1)/2$ and $n_{i\downarrow} =$ $(\tau_i^z + 1)/2$. Magnetization currents of the two spin species, defined through the continuity equations $d(\sigma_i^z/2)/dt = j_i^{(\sigma)}$ $j_{i-1}^{(\sigma)}$, $d(\tau_i^z/2)/dt = j_i^{(\tau)} - j_{i-1}^{(\tau)}$, are $j_i^{(\sigma)} = -\frac{t}{2}(\sigma_i^x \sigma_{i+1}^y \sigma_i^y \sigma_{i+1}^x)$, $j_i^{(\tau)} = -\frac{t}{2}(\tau_i^x \tau_{i+1}^y - \tau_i^y \tau_{i+1}^x)$. In fermionic picture, the particle (charge) current is $j_i^{(c)} = j_i^{(\sigma)} + j_i^{(\tau)}$, while the spin current is $j_i^{(S)} = (j_i^{(\sigma)} - j_i^{(\tau)})/2$. Particle density is $n_i =$ $n_{i\uparrow} + n_{i\downarrow}$, while spin density is $s_i = (n_{i\uparrow} - n_{i\downarrow})/2$. Because of the same driving at both ladder legs the currents $j^{(\sigma)}$ and $j^{(\tau)}$ are the same. Therefore NESS is such that it has a nonzero charge current and zero spin current. We have also performed simulations with Lindblad operators on the τ -chain driving the transport in the opposite direction, that is with

$$L_{5,6} = \sqrt{\Gamma(1 \pm \mu)} \tau_1^{\pm}, \quad L_{7,8} = \sqrt{\Gamma(1 \mp \mu)} \tau_L^{\pm}.$$
 (7)

In such a case of spin driving, the NESS has a nonzero spin current and zero charge current because $j^{(\tau)} = -j^{(\sigma)}$ holds. Furthermore, we stress that spin and charge transports are interchanged under the particle-hole transformation for the *down* spin fermions only and simultaneously changing the sign of *U*. Namely, taking $R := \prod_{i=1}^{L} \tau_i^{x} = R^{\dagger}$, one finds $Rj_i^{(S)}R^{\dagger} = j_i^{(C)}/2$ and $RH(U)R^{\dagger} = H(-U)$, provided one takes a symmetric interaction term $(n_{i\uparrow} - \frac{1}{2})(n_{i\downarrow} - \frac{1}{2})$ in Eq. (1) or, equivalently, adds a chemical term -UN/2 to *H* with $N = \sum_{i,s} n_{i,s}$. Even though our master equation evolution (3) does not strictly conserve *N*, we have checked explicitly that the results based on Hamiltonians *H* and H - UN/2 are identical.

III. RESULTS

A. Evidence of diffusion: density profiles and scaling of currents

We set t = 1, $\Gamma = 1$ and $\mu = 0.2$, except in Fig. 7 where $\mu = 1$. Driving $\mu = 0.2$ corresponds to equilibrium density in the reservoirs of $n_{L,s} = 0.4$ at the left end and $n_{R,s} = 0.6$ at the right end. The average filling ratio is therefore n = 1/2, $\sum_{i=1}^{L} n_{i\uparrow} = \sum_{i=1}^{L} n_{i\downarrow} = L/2.$ The value $\mu = 0.2$ is at the upper end of a linear response regime. For large drivings $\mu\gtrsim 0.6$, one gets a negative differential conductance effect, ¹⁹ where the current decreases with increasing driving. The main goal of this paper is to classify spin and charge transport, whether it is ballistic, diffusive, or anomalous. For NESS, different transport regimes are reflected in the scaling of the current on the system size. Fixing the driving strength μ , in a ballistic conductor the current is independent of the system length, $j \sim L^0$, for a diffusive conductor it scales as $j \sim 1/L$, whereas in the anomalous case the current is proportional to a fractional power of L. We therefore calculated NESSs for different sizes L. Typical density profile is shown in Fig. 1. One can see that the densities of spin-up and spin-down fermions are linear in the bulk. Jumps in the density at the boundaries are due to over-simplified Lindblad operators that are not



FIG. 1. (Color online) Density profile $n_{i,s}$ along the chain for L = 100, U = 1. Apart from jumps at the boundary, density is linear, which is typical for diffusive conductors. Solid black line, overlapping with the numerical points, is a best-fitting linear function.



FIG. 2. (Color online) Scaling of charge current $j^{(c)}$ divided by the extrapolated density drop $L\nabla n$ with the system size L for different interactions U. Thick full (red) line, overlapping with U = 1.0 data, is ~30.4/L, indicating a diffusive transport. Two dashed lines also suggest ~1/L scaling.

"matched" to the bulk dynamics, i.e., there are boundary resistances. Because these jumps are rather large, we have fitted a linear function to the density profile in the bulk, thereby obtaining the density gradient $\nabla n_{i\uparrow} = \nabla n_{i\downarrow} = \nabla n_i/2$. In Fig. 2, we then plot the scaling of the charge current (which is in the NESS independent of the site) with the gradient of the charge density. At interaction strength U = 1, one can see a nice scaling $i^{(c)} \sim 1/L$. Together with a linear density profiles this is a clear indication of diffusive charge transport. As mentioned, for spin transport virtually the same behavior is obtained (data not shown). For U = 2, the scaling is not quite as good. It seems that for shorter chains $j^{(c)}$ decreases with L slower than 1/L, however, for two largest sizes that we managed to calculate, a crossover to $\sim 1/L$ scaling is clearly suggested. For smaller interaction U = 0.5, the convergence seems better, but contrary to the U = 2 case, the current approaches the asymptotic scaling $(\sim 1/L)$ from above, i.e., it decays a bit faster for short chains. It is not clear whether U = 1 corresponds to a crossover point between the two behaviors since details of convergence in the thermodynamic limit might depend on a particular choice of boundary Lindblad operators. In Fig. 4(a), we show density profiles for U = 0.5which, apart from considerable boundary jumps, again look linear.

Looking at the density profiles at U = 0.5, 1, and 2 which are linear in the bulk already for rather small sizes L, it seems natural to conjecture that the transport is diffusive in TL $L \to \infty$ for all nonzero finite values of U. Different transient scaling of the current with L for shorter chains is likely due to rather strong boundary effects. That the boundary effects are notable can also be seen in the inset of Fig. 3(a), where we show the jump in the density between the reservoir and the first site $n_{1\uparrow} - n_{L\uparrow}$, as well as between the first two sites in the system $n_{2\uparrow} - n_{1\uparrow}$. While the boundary effects show a tendency to disappear in TL, at U = 2 and largest length L = 64 they are still non negligible. One can try to optimize the coupling constant Γ in order to minimize the boundary effects, however, we found that $\Gamma \approx 1$ is usually close to the optimal value which does not seem to depend on L. For U = 0.5, the



FIG. 3. (Color online) (a) Density profiles $n_{i,s}$ at U = 2 and (b) the corresponding connected density-density correlation function $\langle n_{i\uparrow}n_{i\downarrow}\rangle_c$. Data are shown for L = 16 (blue squares), L = 32 (green circles), and L = 64 (red triangles), all for charge driving. Inset in (a): scaling of the jump between the reservoir and the 1st particle and between the 1st and 2nd particles, with size L. Black dashed lines indicate $\sim 1/L$ and $\sim 1/L^{0.7}$. Note that at L = 64 boundary jumps still account for around 25% of the total density difference between the chain ends.

boundary effects are larger than for larger U cases studied, U = 1 and U = 2, see Fig. 4(a). This is probably due to a smaller bulk resistivity that makes the effects of the contacts (contact resistivity) relatively larger.

Our finite size simulations of course cannot exclude existence, in thermodynamic limit $L \to \infty$, of a very small ballistic contribution $j_{\infty} = \lim_{L\to\infty} j(L)$ to the current characterized by a small nonvanishing value of the Drude weight. Nonvanishing value of j_{∞} would be visible, for sufficiently larger *L*, as a plateau in j(L). This means that one should see a departure from a linear scaling for large values of *L* as shown in Fig. 2. Assuming the scaling $j(L) = \text{const}/L + j_{\infty}$, we can estimate $j_{\infty} \lesssim \delta[j(L_{\text{max}})]$ where $\delta[j(L_{\text{max}})]$ is the absolute error of data point at the largest size L_{max} . From the data of Fig. 2, we estimate that $j_{\infty} \lesssim 0.01$ (for the case of U = 1 and $L_{\text{max}} = 100$).

B. Density-density correlations

In Figs. 3(b), 4(b), and also 5(b), we show the connected spin-up/spin-down correlation function $\langle n_{i\uparrow}n_{i\downarrow}\rangle_c = \langle n_{i\uparrow}n_{i\downarrow}\rangle - \langle n_{i\uparrow}\rangle\langle n_{i\downarrow}\rangle$, that gives on-site correlations between two fermion species. If we extrapolate our finite-*L* data to TL, we find

$$\langle n_{i\uparrow}n_{i\downarrow}\rangle_{\rm c} \propto [\mu^2 - (2\langle n_{i\uparrow}\rangle - 1)(2\langle n_{i\downarrow}\rangle - 1)],$$
 (8)



FIG. 4. (Color online) Density profiles (a) and connected correlations (b) for U = 0.5. Other parameters are the same as in Fig. 3.

with a proportionality prefactor depending on U only, yielding a *parabolic correlation profile* for our *linear* density profiles. Interestingly, in the middle of the chain the connected correlations become independent of the system size, while they are going to zero at the boundaries.²¹ In Fig. 5, we in addition show also, for U = 1 case, the nonconnected correlations $\langle n_{i\uparrow}n_{i\downarrow}\rangle$ [top frame (a), red squares], or centered nonconnected correlations $\langle (n_{i\uparrow} - 1/2)(n_{i\downarrow} - 1/2)\rangle$ [bottom frame (b), red dotted line]. We can see that the nonconnected correlations have similar shapes as density profiles.

Considering spin driving (7), we find numerically the same diffusion constant as for charge driving (6), while the connected correlations $\langle n_{i\uparrow}n_{i\downarrow}\rangle_c$ change sign [see Fig. 5(b), blue circles]. Note that this implies that conductivities at infinite temperature do not depend on the sign of interaction U. We have also checked explicitly, by comparing data for U = -1 with U = 1, that density profiles, currents and correlations are insensitive to the sign of U. One can see that in the presence of spin current without charge current, nonconnected correlations, shown in Fig. 5(a) with blue circles, are practically independent of the site and are slightly smaller than 1/4.

Note that in both cases of charge and spin driving the connected correlations scale as $\sim \mu^2$ and are of purely nonequilibrium origin, i.e., they vanish in the equilibrium limit $(\mu = 0)$.

C. Large interaction U

In the limit $U \rightarrow \infty$, the low-energy excitations of the half-filled Hubbard model can be effectively described by the 1D isotropic Heisenberg model. In our open system formulation, this mapping cannot be strictly implemented, due



FIG. 5. (Color online) Density-density correlation functions for the case of charge driving (red squares) and for spin driving (blue circles). (b) shows the connected correlations (symbols) and (a) the nonconnected ones. In (b), we also show $\langle (n_{i\uparrow} - \frac{1}{2})(n_{i\downarrow} - \frac{1}{2}) \rangle$ (dashed curves). All is for U = 1 and L = 32.

to the presence of high-temperature baths that drive the system locally away from half-filling. It is therefore an interesting question whether the transport properties of the Hubbard model in the limit of large U are qualitatively the same as for the Heisenberg model.

In Fig. 6, we plot density profiles in our open Hubbard model for increasing values of U, keeping L fixed, and find increasingly arcsin like shape, similar as in the isotropic Heisenberg model that displays an anomalous transport²² with the magnetization current scaling as $\sim 1/\sqrt{L}$. This perhaps explains slower decay of the current with L in the Hubbard model for small L's and larger U, seen for instance in Fig. 2 at U = 2. Note that the limits $U \rightarrow \infty$ and $L \rightarrow \infty$ do not



FIG. 6. (Color online) As one increases U at fixed length L = 32 the rescaled density profiles $(2n_{i,s} - 1)/\mu$, shown for U = 1 and U = 5, become similar to $\frac{2}{\pi} \arcsin[(2i - 1)/L - 1]$ (full red curve), found in the isotropic Heisenberg model (see Ref. 22).



FIG. 7. (Color online) At maximal driving, $\mu = 1$, density profiles have a cosine shape (red squares, for L = 50 and U = 1), while the current scales as $\sim 1/L^2$ (data not shown), exactly as in the Heisenberg model at maximal driving.²³ In the inset, we show convergence of $\Delta n = n_{i,s} - \sin^2(\pi x/2)$ with L.

commute. In order to recover the Heisenberg behavior in TL, one has to first let $U \to \infty$ and only then $L \to \infty$.

D. Strong driving, $\mu = 1$

In sections III A, III B, and III C, we have shown that the weakly driven Hubbard model displays diffusive behavior. Here we show that for strong driving, where the system is far away from equilibrium, the behavior of physical observables can be dramatically different. For example, we briefly discuss



FIG. 8. (Color online) Density profiles (a) and currents (b) for a non-half-filled Hubbard system. We show data for L = 16 (squares) and L = 32 (circles), all for U = 1. Density profiles in the bulk are flat and currents do not depend on L, indicating ballistic transport. In both frames, upper two sets of symbols are for one fermion species, lower two for the other.

the case of maximal driving $\mu = 1$ and find that the current scales subdiffusively as $j^c \propto 1/L^2$. The corresponding density profile is shown in Fig. 7. We can see that the profile is in TL given by a simple cosine shape $n_{i,s} = \sin^2[\pi(2i - 1)/4L]$, exactly the same as has been found analytically in the isotropic Heisenberg model at strong driving.²³ This suggests that similar exact solution for NESS at maximum driving $\mu = 1$ as for the Heisenberg spin chain is also achievable for the Hubbard model and points to wider applicability of the algebraic method proposed in Refs. 23 and 24.

E. Non-half-filled case

So far, all the results shown were for symmetric driving, producing on average half-filled bands. However, if the filling of the two fermion species is not 1/2, we expect the transport to be ballistic at high temperatures. This follows from an existence of a nontrivial constant of motion, which in the non-half-filled case possesses nonvanishing overlap with the spin/charge currents.²⁰ In order to numerically verify the consistency of our nonequilibrium setup with this expectation, we choose a nonsymmetric driving with $L_{1,2} = \sqrt{\Gamma(1 \mp \mu_{L\uparrow})} \sigma_1^{\pm}$ and $L_{3,4} = \sqrt{\Gamma(1 \pm \mu_{R\uparrow})} \sigma_L^{\pm}$ for the first chain, where $\mu_{L\uparrow} = 0.5$ and $\mu_{R\uparrow} = 0.9$, while $L_{5,6} = \sqrt{\Gamma(1 \mp \mu_{L\downarrow})} \tau_1^{\pm}$ and $L_{7,8} =$ $\sqrt{\Gamma(1\pm\mu_{R\downarrow})} \tau_L^{\pm}$ with $\mu_{L\downarrow} = 0.4$ and $\mu_{R\downarrow} = 0.8$ for the second chain. Density profiles can be seen in Fig. 8. One can see that the gradient is very small (or zero); what is more, currents are almost independent of system size L. Namely, in Fig. 8(b) currents are within numerical errors the same for L = 16 and 32 (small inhomogeneities visible in the figure are due to truncation errors). If the transport were diffusive, as is the case for the half-filled system, the current for L = 32should be half as large as for L = 16. We therefore confirm that the nonequilibrium transport is clearly ballistic for a non-half-filled Hubbard model.

IV. CONCLUSION

Summarizing our findings about the transport in halffilled zero-magnetization 1D Hubbard model at an infinite temperature, we have shown that at finite interaction U both charge and spin transport are diffusive. This conclusion is based on the scaling of the currents with the system size for up to 100 lattice sites as well as on perfectly linear density profiles away from the boundaries. This complements ballistic spin transport at U = 0 and arbitrary temperature and anomalous transport at high temperature and $U = \infty$. Such diffusive nonequilibrium states are in turn characterized by nonvanishing quantum correlations between spin-up and spin-down particles at the same sites. For the case of extremal driving, we find subdiffusive (insulating) transport behavior with cosine shaped density profiles.

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- ¹⁶NESS is in our case unique and is therefore reached irrespective of the initial state $\rho(0)$. Evans theorem¹⁷ states that NESS [fixed point of a Liouvillian dynamical semigroup (3)] is unique iff the set of operators $\{H, L_1, L_2, \ldots\}$, in our case $\{H, \sigma_{1,L}^{\pm}, \tau_{1,L}^{\pm}\}$, generates, under multiplication and addition, the entire algebra of operators, in our case, over the spin ladder on *L* sites. Indeed, this can be verified by operator identities giving a recursive generation of σ_j^{\pm} : $\sigma_2^{+} = (1/4)\sigma_1^z[\sigma_1^+, [H, \sigma_1^z]]$, and for $j \ge 2$, $\sigma_{j+1}^+ = -\sigma_{j-1}^+ - (1/2)\sigma_j^z[\sigma_j^-, \sigma_j^+ H \sigma_j^+]$. Operators for the second leg τ_j^{\pm} are generated through a similar recursion, replacing σ_j^{α} by τ_j^{α} . Finally, all possible products of $\{\sigma_j^{\pm}\}$ and $\{\tau_j^{\pm}\}$ generate the entire operator algebra of the Hubbard chain.
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