Smooth or shock: Universality in closed inhomogeneous driven single file motions

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We study the nonequilibrium steady states in a unidirectional or driven single file motion (DSFM) of a collection of particles with hard-core repulsion in a closed system. For driven propulsion that is spatially smoothly varying with a few discontinuities, we show that the steady states are broadly classified into two classes, independently of any system detail: (i) when the steady state current depends explicitly on the conserved number density n, and (ii) when it is independent of n. This manifests itself in the universal topology of the phase diagrams and fundamental diagrams (i.e., the current versus density curves) for DSFM, which are determined solely by the interplay between two control parameters n and the minimum propulsion speed along the chain. Our theory can be tested in laboratory experiments on driven particles in a closed geometry.

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I. INTRODUCTION

Driven single file motion (DSFM) implies unidirectional particle movement along narrow channels where the particles cannot cross each other due to hard-core repulsion. It is an inherently nonequilibrium process that consumes energy for propulsion. We are particularly interested in DSFM with spatially nonuniform propulsion and finite resources, i.e., fixed available number of particles. This should be relevant in wideranging systems, e.g., vehicular or pedestrian movement along closed networks of roads with bottlenecks having varying strength, closed urban transport networks with enforced speed variations [1,2], and spatially varying electric fields in closed arrays of quantum dots [3]. This study could form the basis for further research on weakly number-conserving quasi-onedimensional (1D) transport models where particle number conservation approximately holds at timescales shorter than any nonconserving processes, e.g., ribosome translocations along closed mRNA loops with pause sites (for which ribosomes are typically reinitiated in translocation and breaking of ribosome number conservation is likely to have an effect only at relatively large timescales) [4–6]. It should also be useful in studies on the effects of quenched disorder on asymmetric exclusion processes with finite resources [7].

The general goal of this work is to theoretically understand the classes of steady states in spatially nonuniform systems with restricted one-dimensional (1D) motion with finite resources, and to elucidate their universal nature. For this, we construct a minimal theory for DSFM with positiondependent propulsion speed and hard-core repulsion in closed geometries, with the total number of particles $N_{\rm tot}$ being conserved. This theory adequately describes the interplay between inhomogeneity and conservation laws, and reveals the generic universal nature of the nonequilibrium steady states. It applies to all in vitro or in vivo systems where individual particles are nonactive or weakly active, i.e., do not actively push or pull the neighbors strongly and are undergoing quasi-1D motion, without mutual passage and having number conservation. It can also be useful and serve as a theoretical benchmark for quasi-1D systems with weak particle nonconservation, e.g., binding-factor-mediated enhancement of the probability of loop formation in mRNA in eukaryotes [8]. The results can be tested in carefully designed *in vitro* experiments on the collective motion of driven particles along a nonuniform closed track.

We focus on the steady state densities in DSFM and their dependencies on N_{tot} and position-dependent propulsion. In order to extract generic results from a minimal description without losing the essential physics, we model DSFM by the well-known 1D totally asymmetric simple exclusion process (TASEP), where each site can accommodate at most one particle that can hop only in one direction if the neighboring site is empty. TASEP with open boundaries is a simple model for nonequilibrium phase transitions in 1D open systems [4,9,10].

The rest of the paper is organized in the following manner. In Sec. II, we present our principal results. Then in Sec. III, we introduce the model. We evaluate the steady state densities using standard mean-field methods (along with corresponding Monte Carlo simulation results) in Sec. IV. Finally, in Sec. V we summarize and conclude.

II. PRINCIPAL RESULTS

In this article, we study closed TASEP with N sites as a model for DSFM. Space-dependent propulsion is described by quenched hopping rates that are spatially smoothly varying with a finite number of discontinuities having single or

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multiple point minima. The main results are as follows: (i) independently of the details of the heterogeneous hopping rates, there are generically two classes of steady states delineated by the steady state current J: (a) when J depends on mean density $n = N_{tot}/N$ (0 < n < 1) explicitly (hereafter *smooth* phase), and (b) when J is independent of n, characterized by a phase separation with localized (LDW) or delocalized (DDW) domain walls (hereafter shock phase); (ii) the phases and the reentrant transitions between them are controlled by the interplay between n and the global minima q_{\min} of the positiondependent propulsion speed; (iii) moving shocks appear only for multiple global minima in propulsion speed; multiple local minima with only one global minimum only produce a localized shock; and (iv) while accumulation of particles where the hopping rate is low is naively expected, we show below that the position of the peak of the density in the shock phase can actually be anywhere in the system, being controlled by n.

This article shows how the general concept of universality, well developed for equilibrium systems, applies for spatially varying steady state density profiles in driven inhomogeneous systems with number conservation. This remains hitherto unexplored. More specifically, the topology of the phase diagrams plotted as functions of n and q_{\min} and the associated fundamental diagrams (i.e., the J vs n plots) is argued to be *universal*, independently of the precise hopping rate functions; see Fig. 1.

III. MODEL

The model consists of a closed 1D lattice, and the particles execute unidirectional motion from one site to its neighbor, subject to exclusion. The model has a hopping rate $q_i \leq 1$ at a site *i*. The dynamics clearly conserves the total particle number $N_{\text{tot}} = \sum_{i=1}^{N} n_i$, where n_i is the occupation of site *i*. The dynamics of TASEP is formally given by rate equations for every site, which are not closed [11]. In mean-field theory (MFT), we write down the dynamical equations for TASEP in closed forms, amenable to analytical treatments. We label the sites by x = i/N; in the thermodynamic limit $N \to \infty$, x effectively becomes a continuous variable confined between 0 and 1. In this parametrization, the hopping rate function is given by $0 < q(x) \leq 1$; we assume q(x) to be piecewise continuous, smooth, slowly varying functions of x with a few point minima. Further, we define $\rho(x) = \langle n_i \rangle$ as the density at x; here $\langle \ldots \rangle$ refers to temporal averages in the steady states. Studies on quenched heterogeneous TASEP have a long history; see, e.g., Refs. [12–15] for some studies on different aspects of heterogeneous TASEP. Our model complements these existing works, and primarily investigates the notion of universality not discussed elsewhere.

IV. STEADY STATE DENSITIES

In the steady states

$$\frac{\partial \rho}{\partial t} = -\frac{\partial}{\partial x} [q(x)\rho(1-\rho)] = 0, \tag{1}$$

in MFT over a range of x in which q(x) is smooth [16]. This gives



FIG. 1. (a) Illustrations of universal topology of the phase diagrams in the q_{\min} -n plane: (i) $q(x) = (x - 0.5)^2 + q_{\min} = q_1(x)$, and (ii) (inset) $q(x) = 0.5x^2 + q_{\min} = q_2(x)$ (with n = [0, 1] and $q_{\min} = [0, 0.75]$). The range of q_{\min} is chosen in such a way that q(x)does not exceed unity anywhere. Magenta lines and blue points are overlapping MFT and MCS results, respectively. (b) Fundamental diagrams (J vs n) for $q_1(x) = (0.5 - x)^2 + 0.5$ (dashed curves) and $q_2(x) = 0.5x^2 + 0.25$ (solid curves) as chosen above for (a). The different phases that exist on the fundamental diagrams are marked. Both curves have the same form (see text). Note the saturation of current J with respect to the density n in the shock phase. Curves and points represent MFT and MCS results, respectively.

where *J*, a constant (yet unknown), is the steady state current. Equation (2) has two spatially nonuniform solutions $\rho_+(x)$ and $\rho_-(x)$:

$$\rho_+(x) = \frac{1}{2} [1 + \sqrt{1 - 4J/q(x)}] > \frac{1}{2}, \tag{3}$$

$$\rho_{-}(x) = \frac{1}{2} [1 - \sqrt{1 - 4J/q(x)}] < \frac{1}{2}, \tag{4}$$

for all *x*. Clearly, both $\rho_+(x)$ and $\rho_-(x)$ are smooth functions for all *x*, except where q(x) itself is discontinuous. Since $\rho(x) > 0$ for all x, $1 - \frac{4J}{q(x)} \ge 0$. Thus

$$J \leqslant q(x)/4,\tag{5}$$

everywhere. The maximum allowed value of J is thus independent of n:

$$J_{\max} = q_{\min}/4,\tag{6}$$

whence $\rho_+(x_0) = \rho_-(x_0)$; $x = x_0$ is the location of q; see Ref. [17] for analogous result in a disordered exclusion model.

We now outline the derivation of the results announced at the beginning of this article and the principles for the construction of Fig. 1, followed by some illustrations of the phases for some representative q(x). Since $\rho_+(x) \left[\rho_-(x)\right] >$ [<] 1/2 for all x except at isolated points x_0 where q(x) = q_{\min} , for a given q(x) if n is sufficiently close to 1/2, $\rho(x)$ must be a combination of $\rho_+(x)$ $(>\frac{1}{2})$ and $\rho_-(x)$ $(<\frac{1}{2})$. This is the shock phase mentioned above. In contrast, when n approaches 0, there are only a few particles in the system and $\rho(x) =$ $\rho_{-}(x)$ in the steady state throughout the system. Analogously, for *n* approaching unity, $\rho(x) = \rho_+(x)$ in the steady state. The smooth phases with $\rho(x) = \rho_{-}(x) < 1/2$ and $\rho(x) =$ $\rho_{+}(x) > 1/2$ are spatially nonuniform, and hence generalize the spatially uniform low-density (LD) and high-density (HD) phases, respectively, of TASEP with open boundaries [16]. Thus, a reentrant smooth-shock-smooth nonequilibrium phase transition is expected as *n* rises from 0 to 1 for any given q(x). The precise boundaries between these phases for a given nand q(x)—which will tell us for a given q(x) how close n must be to 1/2 for the smooth-shock transition—are obtained by imposing particle number conservation on (3) and (4) and using (6); see below.

Current J is fixed by the particle number conservation:

$$\int_0^1 \rho_a(x) = n, \ a = +, -.$$
(7)

From (3) and (4), the maximum (minimum) of $\rho_{-}(x) [\rho_{+}(x)]$ coincides with the minimum of q(x), a fact borne out by our Monte Carlo simulation (MCS) studies: q_{\min} effectively acts as a bottleneck, and as a result, particles tend to accumulate behind it (see below).

As *n* increases from zero, *J* rises and eventually reaches J_{max} . For $J = J_{\text{max}}$, $\rho_+(x_0) = \rho_-(x_0)$, where x_0 is the location of q_{\min} . On increasing *n* further, the additional particles are accommodated by representing $\rho(x)$ as a combination of $\rho_-(x)$ and $\rho_+(x)$ which meet smoothly at x_0 . Since particles should accumulate behind the *bottleneck* at x_0 , we expect that additional particles will go over to the high-density solution represented by $\rho_+(x)$. Since we have a closed system, the two solutions must meet at another point x_w , such that $\rho_+(x_w) > \rho_-(x_w)$ [since $\rho_+(x) = \rho_-(x)$ only at $x = x_0$], leading to a *discontinuous jump* in the form of a localized domain wall (LDW) in $\rho(x)$ at x_w , thus giving rise to the *shock phase*, with a jump $\overline{\rho}$ given by

$$\rho_+(x_w) - \rho_-(x_w) = \overline{\rho},\tag{8}$$

controlled by *n* and the functional form of q(x). As more particles are added, x_w shifts to make the region of existence for $\rho_+(x)$ larger and $\rho_-(x)$ smaller. This indeed leaves $J = J_{\text{max}} = q_{\text{min}}/4$ unchanged. Thus the current in the shock phase saturates to its maximum value J_{max} . This continues till $\rho_+(x)$ spans the full system. Thus, as *n* rises from the low to moderate values, a smooth-to-shock transition is encountered.

Interestingly, independently of the form of q(x) this transition is *reentrant*—since, as *n* rises further, the system moves from shock phase to smooth phase again, with ρ_+ now being the only valid solution. This reentrant transition can also be understood from the particle-hole symmetry of the model. Since particle density $\rho_+(x)$ can be interpreted as the hole density $1 - \rho_+(x) = \rho_-(x)$, if $\rho_-(x)$ is a steady state solution for overall particle density n, $\rho_+(x)$ is a steady state solution for particle density 1 - n. This picture remains valid even when there are additional local minima (but only one global minimum q_{\min} at x_0): J_{\max} is still controlled solely by q_{\min} [Eq. (6) above], with $\rho_+(x_0) = \rho_-(x_0)$; the other local minima having no effect on J_{max} are effectively screened in any steady state current measurements in the shock phase. However, the form of the LDW, i.e., the functional form of $\rho(x)$, depends on any local minima through its dependence on the full form

Assuming only one global minimum for q(x), at the phase boundary between smooth and shock phases, $J = J_{\text{max}} = q_{\min}/4$ and $\rho(x) = \rho_{-}(x)$ (for n < 1/2) for all *x*, or for n > 1/2, $\rho_{+}(x)$ for all *x*. Thus

of q(x).

$$\int_{0}^{1} dx \,\rho_{\pm}(x, J_{\max}) = \int_{0}^{1} dx \,\frac{1}{2} \left[1 \pm \sqrt{1 - \frac{q_{\min}}{q(x)}} \right] = n \quad (9)$$

gives the quantitative dependence of *n* on q_{\min} for the reentrant transition, or equivalently, the boundaries between smooth and shock phases in Fig. 1(a). As our arguments above are independent of the precise form of q(x), the topology of the phase diagrams in Fig. 1 (top) should remain independent of the precise forms of q(x) having the same q_{\min} . This is the universality in DSFM mentioned at the beginning that is also manifest in the fundamental diagrams in Fig. 1 (bottom), obtained from (2) and (7). The results in Fig. 1 are also obtained from the MCS studies, which corroborate the MFT predictions closely. This holds true even if there are multiple global minima with value q_{\min} ; in this case, shock phases correspond to moving shocks (see below). Notice that the topology of the phase diagram and the corresponding fundamental diagram are the same as those obtained in Ref. [18], where a single slow site controls the current, establishing an equivalence between the model of Ref. [18] with a single slow site and our model here. The strength of the single slow site in the model of Ref. [18] corresponds to q_{\min} here.

We now illustrate the phases with a few representative choices for q(x). See Fig. 2 for a plot of $\rho(x)$ vs x in the smooth phase; see also Figs. 7–9 in Appendix C for plots of $\rho(x)$ with different q(x) in the smooth phase; good agreements between MFT and MCS results are evident. Further see Fig. 3 for a plot of the density in the shock phase for a choice of q(x), again showing strong agreements between MCS and MFT results; see also Fig. 10 in Appendix C.

Notice that the peak of the density in the smooth phase in Fig. 2 coincides with the location of q_{\min} . As argued above, this can be understood from the form of $\rho_{-}(x)$ as given in (4)—clearly, the maximum of $\rho_{-}(x)$ must coincide with the minimum of q(x). In contrast, the location of the extrema of the density profiles in the shock phase have no such relation with the minimum of q(x). Instead, as we note in Fig. 3, $\rho(x)$ is continuous at the location x_0 of q_{\min} , so long as q(x) itself



FIG. 2. Plot of $\rho(x)$ vs x in the LD phase (n = 0.2). Magenta line indicates the hopping rate function $q(x) = (x - 0.5)^2 + 0.5$. Red continuous line and blue points denote $\rho_-(x)$, respectively, from MFT and MCS studies (see text). Excellent agreement between MFT and MCS results is clearly visible. We have used N = 2000.

is continuous at x_0 (since $\rho_+ = \rho_-$ at x_0). If however, q(x) itself is discontinuous at x_0 , the density is also discontinuous there; see Fig. 10 (right) in Appendix C. In both these cases, the density has a discontinuity *elsewhere* whose location is controlled by *n* for a given q(x) with q(x) being *continuous* there. This is one of the principal results in this article. Furthermore, given that $\rho_+(x)$ reduces and $\rho_-(x)$ grows when q(x) grows in *x* [see Eqs. (3) and (4) above], the density in the shock phase, being a linear combination of $\rho_+(x)$ and $\rho_-(x)$, can have a peak anywhere in the system that is controlled by *n*.

We further illustrate the occurrence of a single LDW in the shock phase with q(x) having multiple local but one global minimum in Fig. 4, with $\rho_+(x) = \rho_-(x)$ only at the location of the global minimum of q(x), in agreement with the theoretical predictions. Thus, experimental measurements of the steady state densities will not reveal the actual existence of other local minima.

When there are multiple global minima, the abovementioned theory of LDW breaks down and very different physics emerges in the shock phase. For instance, consider



FIG. 3. Plot of $\rho(x)$ vs *x* in the shock phase with n = 0.7, N = 2000. Magenta line indicates q(x) = 1 - x, $0 \le x < 0.5$, and q(x) = x, $0.5 \le x < 1$. Red line and blue points, respectively, represent MFT and MCS results for $\rho(x)$.



FIG. 4. One LDW with n = 0.6, N = 2000 and q(x) = 1 - x, x = [0, 0.25]; q(x) = 0.5 + x, x = [0.25, 0.5]; q(x) = 2 - 2x, x = [0.5, 0.75]; q(x) = 2x - 1, x = [0.75, 1] with two local minima of q(x) at x = 0.25 and x = 0.75. Excellent agreement between MFT and MCS is observed.

q(x) to have only two global minima at diametrically opposite points x_1 and x_2 : $q(x_1) = q(x_2) = q_{\min}$; see Fig. 5 for such a choice of q(x). Thus there are now two effective bottlenecks at x_1 and x_2 which split the ring into two identical segments, say T_A and T_B [19,20], of equal length. Furthermore, from (3) and (4) we have

$$\rho_{-}(x_1) = \rho_{+}(x_1) = \rho_{-}(x_2) = \rho_{+}(x_2).$$
(10)

In each of T_A and T_B , MFT for the shock phase applies. Therefore, two domain walls are expected. Particle number conservation then can only yield a single relation between the two domain wall positions and *n*, and cannot determine both x_{w1} and x_{w2} separately. This means that a shift in x_{w1} can be balanced by an equivalent reverse shift in x_{w2} that still satisfies particle number conservation. Due to the inherent stochasticity of the system, all possible solutions of x_{w1} and x_{w2} that are consistent with particle number conservation are visited by the system, if we wait long enough. This leads to *two* DDWs. Under long-time averages, envelopes of the



FIG. 5. Plot of $\rho(x)$ vs *x* showing two DDWs; n = 0.5. Magenta lines indicate q(x) = 1 - x, $0 \le x < 0.5$, and q(x) = 1.5 - x, $0.5 \le x < 1$; red line and blue points, respectively, are MFT and MCS results for DDW, which agree with each other.



FIG. 6. Kymograph for n = 0.5, N = 800, q(x) = 1 - x for $0 \le x < 0.5$; q(x) = 1.5 - x for $0.5 \le x < 1$. See [21] for technical details of obtaining kymographs of the same nature in closed TASEP systems.

moving DDWs will be observed; see Fig. 5 and Appendix A for detailed calculations.

As the DDWs in Fig. 5 move, T_A and T_B exchange particles. Particle number conservation ensures that when a particle enters (leaves) T_A , it must necessarily leave (enter) T_B , leading to perfect synchronization of the DDWs in T_A and T_B ; see Appendix B for additional discussions and an associated kymograph in Fig. 6.

If there are more than two global minima of q(x), then there will be as many subchannels (T_A , T_B , T_C , ...), and thus, as many DDWs. However, the argument for synchronization breaks down in such cases; thus there should be no synchrony in movement between any two of the DDWs; see Fig. 11 (left) [and the corresponding kymograph Fig. 11 (right)] and related discussions in Appendix B.

The shocks in open TASEPs, which are always delocalized, appear when the entry (α) and exit (β) are equal and not exceeding 1/2. This is equivalent to $J_{in} = J_{out}$, where J_{in} and $J_{\rm out}$ are, respectively, the currents determined by the entry side and exit side boundaries. The shock phase of the present model is analogous to the shocks in open TASEPs. To see this we note that our ring model can be viewed as an open TASEP with effective entry $[\alpha_e = \rho_-(x_0)]$ and exit $[\beta_e = 1 - \rho_+(x_0)]$ rates that are joined at x_0 , the location of q_{\min} (assuming a single global minimum) [19–21]. From the condition of the shock phase, $J_{in} = J_{out} = J_{max} = q_{min}/4$ holds automatically. That we obtain an LDW here as opposed to a DDW in an open TASEP under equivalent conditions is due to the strict particle number conservation here; see Refs. [7,18,19,21]. When there are two or more global minima of q(x), the system can be thought to consist of those many TASEPs joined together to form a ring; see Refs. [19,21]. The conditions for the shock phase are satisfied in all these TASEP segments simultaneously, leading to formation of a domain wall in each of them. As explained in Refs. [19,21], in this situation number conservation cannot pin the domain walls completely giving rise to DDWs; see also Appendices A and B. We lastly note one point of dissimilarity between the shocks here and the DDWs in an open TASEP: for the latter, the current is less than its maximum value, where as in the shock phase here, the current necessarily saturates at its maximum value of J_{max} .

V. SUMMARY AND OUTLOOK

We have thus developed the theory for DSFM in a closed system with position-dependent propulsion by studying a closed TASEP with quenched hopping rates q(x). This theory reveals the universal form of the phase diagrams and the fundamental diagrams of the model, for generic smooth q(x)with a finite number of discontinuities and global minimum, but independent of the precise forms for q(x). Our theory is sufficiently general and applies to any smoothly varying q(x) with a finite number of discontinuities; it generalizes the analyses in Ref. [12]. From the perspectives of nonequilibrium systems, these results generalize the studies in Refs. [13,18,19,21,22]. In the more complex situation where individual particles can actively push (or pull) its neighbor, thus only dynamically modulating the effective hopping rate, we expect our results to remain valid for weak activity. For strong activity, competition between background heterogeneous hopping and the active processes will determine the steady state, whose full analysis is beyond the scope of the present work. We have here studied smooth q(x) with a small number of discontinuities. There are in vivo situations where q(x) could be rapidly fluctuating, e.g., the DNA strands [23]. It would be interesting to study to what degree our results remain valid for rapidly fluctuating q(x).

Lastly, MFT developed here neglects spatial density correlations, which is an approximation. This may be improved by systematically including two-point correlations; see, e.g., Ref. [15]. It has been shown that extended particles instead of point particles significantly affect the steady state densities of an open TASEP with site-dependent hopping rate; see Refs. [15,24]. It will be interesting to study this effect in a closed TASEP.

This theory may be verified in model experiments on the collective motion of driven particles with light-induced activity [25] in a closed narrow circular channel [26,27]. Unidirectionality of the motion can be ensured by suppressing rotational diffusion, e.g., by choosing ellipsoidal particles with the channel width shorter than the long axis of the particle everywhere, or by using dimer particles. Propulsion speed can be tuned by applying patterned or spatially varying illumination [25]. Steady state densities can be measured by microscopy with image processing. While technical challenges are anticipated in setting up appropriate experimental arrangements, we hope this will be realized in the near future.

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APPENDIX A: DENSITY PROFILES FOR DELOCALIZED DOMAIN WALLS

We calculate here the steady state density profiles when q(x) has two symmetrically placed global minima of same value. The system then can be considered to consist of two TASEP chains of equal size [19], say T_A (with $0 \le x \le 1/2$) and T_B (with $1/2 \le x \le 1$), each spanning from one global minimum of q(x) to the other. While the total particle number in the ring is conserved, the number of particles in each of T_A and T_B can fluctuate. We closely follow Ref. [28] in our analysis below.

Now consider one delocalized domain wall (DDW) in each of T_A and T_B . Let x_w^A and x_w^B be the instantaneous positions of the DDWs in T_A and T_B , respectively, and let the respective heights be $\Delta_A(x_w^A)$ and $\Delta_B(x_w^B)$. We note here that the DDW heights are explicit functions of their positions, since the steady state density is not uniform for an arbitrary q(x).

Now, increasing the number of particles in T_A by 1 would imply shifting x_w^A by an amount $\delta x_w^A = \frac{-1}{L\Delta_A}$. Similarly, decrease of a particle would mean $\delta x_w^A = \frac{1}{L\Delta_A}$. In order to understand why this is so, we note that the "height Δ_A of the domain wall (DW) at x_w^A " means that the Δ_A number of excess particles are needed to fill up one lattice spacing $(=\frac{1}{L})$, or to cause one lattice spacing leftward or rightward movement of the DW (and thus, the above values of δx_w^A). Let us now note that there are two basic processes which can alter the number of particles individually in T_A and T_B , i.e., if a particle enters T_A through its left boundary (equivalent to a particle leaving T_B through its right boundary) and vice versa.

For the following analysis, we will focus on T_A . Let $P(x_w^A, t)$ be the probability of finding a DW at x_w^A at time t. For a given x_w^A , one can evaluate x_w^B at time t, uniquely, using total particle number conservation. The transition rate for a particle entering T_A through the left boundary can be written as $W_L = J_{in} = q(x = 0)\alpha_e^A(1 - \alpha_e^A)$, $\delta x_w^A = \frac{-1}{L\Delta_A}$. Similarly, the transition rate for the particle leaving through the right boundary is given by $W_R = J_{out} = q(x = 1/2)\beta_e^A(1 - \beta_e^A)$, $\delta x_w^A = \frac{1}{L\Delta_A}$. Here, α_e are β_e are the densities at x = 0 and x = 1/2, respectively, in T_A , e.g., $\alpha_e = \rho(x = 0)$, etc.

With these transition rates, we can calculate the average shift or the expectation value of the change, $\langle \delta x_w^A \rangle$, which is given by the product of the increment (with sign) and the sum of the different transition rates:

$$\left\langle \delta x_w^A \right\rangle = \frac{1}{L \Delta_A \left(x_w^A \right)} [J_{\text{out}} - J_{\text{in}}]. \tag{A1}$$

It should be noted here that the domain wall itself performs a random walk about its mean position, $\overline{x_w^A}$. For the fixed point of the random walk, i.e., the value of x_w^A for which $\langle \delta x_w^A \rangle = 0$, we obtain

$$J_{\rm out} - J_{\rm in} = 0, \tag{A2}$$

the well known condition for formation of domain walls. In order to calculate the steady state profiles of the DDWs, we need to study the fluctuations in the DW positions that we do below.

Using the expressions for the transition rates defined above, we can write down the master equation for $P(x_m^A, t)$, the

probability of finding the DW at x_w^A at time t:

$$\frac{dP(x_w^A, t)}{dt} = \Sigma_{\delta x_w^A} \Big[P(x_w^A + \delta x_w^A, t) W(x_w^A + \delta x_w^A \to x_w^A) \\ - P(x_w^A, t) W(x_w^A \to x_w^A + \delta x_w^A) \Big].$$
(A3)

To proceed further, we employ a Kramers-Moyal expansion [29] of the master equation above around $\overline{x_w^A}$, up to second order in δx_w^A . This gives

$$\frac{dP(x_w^A, t)}{dt} = -\frac{\partial}{\partial y}[a(y)P(y, t)] + \frac{1}{2}\frac{\partial^2}{\partial y^2}[b(y)P(y, t)], \quad (A4)$$

where $y = \delta x_w^A$, $a(y) = \sum_y y W(x_w^A + \delta x_w^A \to x_w^A)$, and $b(y) = \sum_y y^2 W(x_w^A + \delta x_w^A \to x_w^A)$. Using the already known values for W and δx_w^A , and Eq. (A2), we arrive at the following results for a and b:

$$a(x_{w}^{A}) = \frac{1}{L\Delta_{A}(x_{w}^{A})} \left[-\alpha_{e}^{A} (1 - \alpha_{e}^{A}) + \beta_{e}^{A} (1 - \beta_{e}^{A}) \right] = 0$$
(A5)

and

$$b(x_w^A) = \frac{1}{L^2 \Delta_A^2(x_w^A)} \Big[\alpha_e^A (1 - \alpha_e^A) + \beta_e^A (1 - \beta_e^A) \Big] > 0.$$
(A6)

Thus up to this order

$$\frac{dP(x,t)}{dt} = \frac{1}{2} \frac{\partial^2}{\partial x^2} [b(x)P(x)].$$
(A7)

Since $J_{in} = J_{out}$, the DW position effectively follows the detailed balance condition. This means the fluctuations in the DW position should follow an equilibrium distribution in the steady state. Hence, the probability current, given by $J_{DW}(x) = \frac{\partial}{\partial x} [b(x)P(x)] = 0$. This yields

$$P(x) = \frac{C}{b(x)},\tag{A8}$$

where *C* is a constant which can be evaluated by the normalization condition on P(x).

Construction of the density profiles

We can now construct the density profile $\rho(x)$ with the knowledge about P(x). Since the long-time-averaged steady state density involves averaging over P(x), we argue that

$$\frac{\partial \rho}{\partial x} = AP(x),\tag{A9}$$

where *A* is a constant of proportionality. Clearly from Eq. (A9), we can see that if P(x) = constant, as in the case for a DDW in an open TASEP, $\rho(x)$ varies linearly with *x*, a known result. The constant *A* in this example can be evaluated by the boundary conditions. In yet another example, for an LDW as $P(x) \propto \delta(x - x_w)$, $\rho(x)$ is a Heaviside Θ function according to Eq. (A9), whose height can be determined using the boundary conditions. We now obtain the DDW steady state density profiles.



FIG. 7. (a) Plot of $\rho(x)$ vs x in the smooth phase for q(x) = 1 - x, $0 \le x < 0.5$; q(x) = x, $0.5 \le x < 1$ (purple continuous line), n = 0.8. (b) Plot of $\rho(x)$ vs x in the smooth phase for $q(x) = (x - 0.5)^2 + 0.5$ (purple continuous line), n = 0.8. (c) Plot of $\rho(x)$ vs x in the smooth phase for $q(x) = 0.5x^2 + 0.5$ (purple continuous line), n = 0.8. Continuous magenta line and overlapping blue points represent, respectively, MFT and MCS data in each plot.

By using Eq. (A9) we write

$$\rho(x) = \tilde{A} \int \frac{dx}{b(x)} + D = A_1 \int dx [\rho_+(x) - \rho_-(x)]^2 + D,$$
(A10)

where, \tilde{A} , A_1 , and D are constants, and we have substituted the value of the DW height $\Delta_A(x) = \rho_+(x) - \rho_-(x)$ in b(x). Using already derived expressions for $\rho_-(x)$ and $\rho_+(x)$ above, we finally arrive at the following expression for the DDW profile,

$$\rho(x) = A_1 \int dx \left(1 - \frac{q_{\min}}{q(x)} \right) + D.$$
 (A11)

The value of the constants can be fixed using the boundary conditions on $\rho(x)$. But there is one more undetermined quantity that we are yet to address. The DDW in general has a certain extent of wandering in T_A that is less than the length of the channel, depending on the number density. We can have two situations: one where $\rho(x)$ shows a mix of



FIG. 8. (a) Plot of $\rho(x)$ vs *x* in the smooth phase for q(x) = 1 - x, $0 \le x < 0.5$; q(x) = x, $0.5 \le x < 1$ (purple continuous line), n = 0.2. (b) Plot of $\rho(x)$ vs *x* in the smooth phase for q(x) = 1 - x, $0 \le x < 0.5$; q(x) = 1.5 - x, $0.5 \le x < 1$ (purple), n = 0.1. Continuous magenta line and overlapping blue points represent, respectively, MFT and MCS data in each plot.



FIG. 9. Plot of $\rho(x)$ vs x in the smooth phase for q(x) = 1 - x, $0 \le x < 0.5$; q(x) = 1.5 - x, $0.5 \le x < 1$ (purple), n = 0.9. Continuous magenta line and overlapping blue points represent, respectively, MFT and MCS data.

LD and DDW profiles, or one where it is a mix of DDW and HD profiles. Therefore, within x = [0, 1/2], we can have either a situation with an LD phase from x = 0 to, say, $x = \overline{x}$, followed by the DDW from $x = \overline{x}$ to x = 1/2, or the situation with the DDW from x = 0 to $x = \overline{x}$, followed by an HD phase from $x = \overline{x}$ to x = 1/2. Now, what determines the value of \overline{x} is the condition $\int_0^{1/2} \rho_A(x) dx = n$, where $\rho_A(x)$ is the complete density profile for T_A and n is the number density (notice that T_A and T_B are identical, and both must have the same average number density). If the DDW does not span the entire T_A , an additional unknown parameter \overline{x} must be determined, which we fix numerically by using the particle number conservation. This in turn yields the complete density profile. We use this scheme to obtain $\rho(x)$ for q(x) = 1 - x for x = [0, 0.5] and q(x) = 1.5 - x for x = [0.5, 1]; see Fig. 5 in the main text. Good agreement with the MCS result is clearly visible, establishing our analytical framework.

APPENDIX B: SYNCHRONIZATION OF DDW MOVEMENT

We now show that the two DDWs formed when q(x) has two global minima placed at diametrically opposite points in the ring move with perfect synchrony. To do that we first consider the two basic microscopic processes in the dynamics that lead to the movement of individual domain walls:

- (i) A particle leaves T_A and enters into T_B .
- (ii) A particle leaves T_B and enters into T_A .

Let δx_w^A and δx_w^B be the shifts in the instantaneous positions x_w^A and x_w^B of the domain walls in T_A and T_B , respectively, due to the processes mentioned above. Jumps in the densities at x_w^A and x_w^B are given by

$$\rho_+(x_w^A) - \rho_-(x_w^A) = \overline{\rho}_A, \quad \rho_+(x_w^B) - \rho_-(x_w^B) = \overline{\rho}_B.$$
(B1)

In general, $\overline{\rho}_A \neq \overline{\rho}_B$.

Since $\alpha_e = \beta_e$ for both T_A and T_B , both processes (i) and (ii) take place with equal rate $W = \alpha_e(1 - \alpha_e)$. By process (i) above, $\delta x_w^A = -WL^{-1}\overline{\rho}_A^{-1}$, $\delta x_w^B = WL^{-1}\overline{\rho}_B^{-1}$. Similarly, by process (ii) above, $\delta x_w^A = WL^{-1}\overline{\rho}_A^{-1}$, $\delta x_w^B = -WL^{-1}\overline{\rho}_B^{-1}$. Thus,

$$\langle \delta x_w^A \rangle + \langle \delta x_w^B \rangle = 0$$
 (B2)

identically. This means $\langle \delta x_w^A \rangle = -\langle \delta x_w^B \rangle$, which is the essence of synchronization of the movements of the two DDWs in T_A and T_B . This synchronization manifests pictorially in a kymograph given in Fig. 6.

The above argument for synchronization breaks down when the number of global minima of q(x), which is same as the number of TASEP segments that make up the ring system, exceeds two. Now imagine q(x) to have N > 2 global minima of value q_{\min} , placed at equal spacing. Thus the ring system can now be considered to be composed of N TASEPs. The microscopic dynamical process for each TASEP consists of receiving a particle from the previous TASEP and releasing a particle to the next one. Clearly, the general analog of (B2) would be

$$\langle \delta x_w^A \rangle + \langle \delta x_w^B \rangle + \langle \delta x_w^C \rangle + \dots = 0.$$
 (B3)

This ensures that for any two successive TASEP segments the sum of the shifts of the corresponding DDW is no longer zero, leading to the loss of synchronization in the movement of the DDWs in any two successive TASEP segments.



FIG. 10. (a) Plot of $\rho(x)$ vs x in the shock phase for $q(x) = (x - 0.5)^2 + 0.5$ (purple continuous line), n = 0.6. (b) Plot of $\rho(x)$ vs x in the shock phase for $q(x) = 1 - 0.5x^2$ (purple continuous line), n = 0.5. Continuous magenta line and overlapping blue points represent, respectively, MFT and MCS data in each plot.



FIG. 11. (a) MCS plot (blue dashed lines) of $\rho(x)$ vs x for $q(x) = 1 - 0.5 \cos^2(4\pi x)$, n = 0.5, N = 1200. The red solid lines and the blue dashed lines represent q(x) and MCS data for $\rho(x)$, respectively. (b) Kymograph for $q(x) = 1 - 0.5 \cos^2(4\pi x)$, n = 0.5, N = 600. Existence of four DDWs is clearly visible which move without any synchronization between any two of them.

APPENDIX C: DENSITY PROFILES

We show below some representative plots of $\rho(x)$ vs x from MCS studies along with MFT predictions. MCS studies have generally been performed with N = 2000 with random sequential updates (except for Fig. 5, where random updates have been used for reasons of limitations on computational resources).

1. Density profiles in the smooth phase

Here we show plots of $\rho(x)$ vs x in the smooth phase with various choices for q(x); see Figs. 7–9.

2. Density profiles in the shock phase

Here we present a few illustrative examples of the steady state density profiles in the shock phase; see Fig. 10.

3. Density profile in the shock phase with four DDWs

Here we show results for the density profile when there are four DDWs; see Fig. 11.

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