Avalanches and clusters in planar crack front propagation

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(Received 12 November 2009; revised manuscript received 25 January 2010; published 27 April 2010)

We study avalanches in a model for a planar crack propagating in a disordered medium. Due to long-range interactions, avalanches are formed by a set of spatially disconnected local clusters, the sizes of which are distributed according to a power law with an exponent \( \gamma_s = 1.5 \). We derive a scaling relation \( \tau_c = 2 \gamma - 1 \) between the local cluster exponent \( \tau_c \) and the global avalanche exponent \( \tau \). For length scales longer than a crossover length proportional to the Larkin length, the aspect ratio of the local clusters scales with the roughness exponent of the line model. Our analysis provides an explanation for experimental results on planar crack avalanches in Plexiglas plates, but the results are applicable also to other systems with long-range interactions.

DOI: 10.1103/PhysRevE.81.046116 PACS number(s): 62.20.mt, 45.70.Ht, 64.60.av

I. INTRODUCTION

Driven elastic manifolds in disordered media have been used to model a number of physical systems ranging from crack propagation in solids [1–8] to dynamics of magnetic domain walls in ferromagnets [9] and vortices in type-II superconductors [10], and to charge density waves [11]. Such systems exhibit nonequilibrium phase transitions as an external driving force \( f \) is varied, so that below the depinning threshold \( f_c \), the system is pinned by the disorder, while for \( f > f_c \), the system moves at a finite average velocity. For \( f \) close to but above \( f_c \), such motion typically occurs as a sequence of avalanches with a broad distribution of sizes. A large class of such systems—including advancing crack fronts in solids [1,4–8], contact lines in wetting [12], magnetic domain walls with dipolar interactions [9], and plastically deforming crystals [13]—is characterized by a long-range interaction kernel. The peculiarity of these systems is that due to the long-range interactions governing the avalanche dynamics, the area swept over by an avalanche is not that due to the long-range interactions governing the avalanche dynamics, the area swept over by an avalanche is not necessity simply connected in space as it would be the case if the interactions were only short ranged. Instead, global avalanches are formed by a set of spatially disconnected local clusters.

The propagation of a planar crack confined in a weak plane represents an ideal system to realize experimentally the depinning transition predicted by the crack line model [1–3]. Yet, the first experimental results for the crack front roughness [4,5] and for the avalanche size distribution [6] disagreed with the theory. The roughness exponent of the crack front was found in the range \( \xi = 0.55–0.6 \) [4,5], instead of the theoretical value \( \xi = 0.39 \) [14], while the avalanche sizes defined by the waiting time matrix (WTM) [6], measuring only spatially connected parts of the avalanches, were found experimentally to be power-law distributed with an exponent much higher than the value predicted by the long-range line model for avalanches extracted by considering the global velocity signal of the crack front. These earlier claims have been revised recently: by testing a larger range of length scales it was shown that the roughness exponent would cross over to the theoretical value at sufficiently large scales [15,16]. Furthermore, in Ref. [7] it was shown that the crack line model can reproduce the avalanche statistics of the experiment, if also the numerical data are analyzed using the waiting time matrix introduced in [6], while a smaller exponent value is obtained when considering avalanches extracted from the global velocity signal of the crack front model. Still, a complete understanding of the origin of the different scaling exponents is lacking.

Here, we clarify these issues by studying the statistics of global avalanches and their spatially connected parts, the local avalanches (or clusters), for a long-range elastic string moving in a disordered medium under quasistatic external driving. Both the avalanche and cluster size distributions are found to exhibit scaling, but with different power-law exponents. In particular, by studying the model with a larger range of length scales, we obtain a more accurate description of the scaling behavior, and therefore a better estimate of the related exponents than in previous numerical studies [7]. These numerical results are compared with data from experiments on Plexiglas plates [6], and excellent agreement is found. Moreover, we propose a scaling relation between the power-law exponents of the avalanche and cluster size distributions by considering the avalanche breakup process, supported by our numerical results. We finally study the morphology of the local clusters and find that their aspect ratio scales with the roughness exponent of the line model above a crossover length proportional to the Larkin length of the crack front. Previous studies [7] using a continuous time version of the crack line model have presumably been sampling the short length scale regime. Thus, our results explain and clarify a number of open issues related in particular to planar crack front propagation, such as the different scaling of global avalanches and local clusters, as well as the connection between roughness and avalanche morphology. However, the implications of our study extend beyond crack front propagation, as the results can be immediately extended to a number of other systems described by the same model, such as contact line dynamics in wetting [12] and low angle grain boundaries in crystalline solids [17]. Furthermore, a number of other driven systems in which the avalanche dynamics is governed by long-range interactions, ranging from domain walls in ferromagnets with dipolar interactions [9] to plastic...
cally deforming crystalline solids [13], are expected to obey similar scaling laws. The paper is organized as follows. In the next section, the simulation model and the methods used to extract the local clusters from the simulations and the experiments are presented. These are then followed by the results of the numerical simulations along with a direct comparison with experimental results from Plexiglas plates in Sec. III. Section IV finishes the paper with conclusions and discussion.

II. MODEL AND METHODS

A. Quasistatic model of crack front propagation

The model of the propagating planar crack front, introduced in Ref. [2], is represented by a vector of integer heights \( h_i \), \( i=1, \ldots, L \), with \( L \) as the system size. Crack propagation is driven by the local stress intensity factor (SIF) \( K_i \), which represents the asymptotic prefactor of the \( 1/\sqrt{r} \) divergence of the stress field near the crack tip. To model the different contributions of the local SIF \( K_i \) acting on a front element \( i \), it is taken to be of the form \( K_i = K_{i,\text{elastic}} + K_{i,\text{inst}} \), where

\[
K_{i,\text{elastic}} = \frac{1}{\gamma} \sum_{j \neq i} b_{j-i} (h_j - h_i) \tag{1}
\]

represents the first-order variation of the stress intensity factor due to a first-order perturbation of the front position; \( b \) is the front segment spacing and \( \gamma \) tunes the strength of the elastic interactions [1]; \( K_{i,\text{random}} \) is a time-independent Gaussian random variable of zero mean and unit variance, with

\[
\langle K_{i,\text{random}} \rangle = \delta(i-j) \delta(h_i - h_j),
\tag{2}
\]

representing random toughness of the material; and \( K_{i,\text{inst}} \) is the contribution of the external load. Notice that the long-range kernel in Eq. (1) is identical in many other problems, such as contact lines [12] and low angle grain boundaries [17]. Periodic boundary conditions are imposed. The dynamics is defined in discrete time \( t \) by setting

\[
v_i(t) = h_i(t+1) - h_i(t) = \theta(K_i),
\tag{3}
\]

where \( v_i \) is the local velocity of the front element \( i \), and \( \theta \) is the Heaviside step function. During a single time step, the front elements with \( v_i > 0 \) are advanced by a unit step, \( h_i(t+1) = h_i(t) + 1 \), and new random forces are generated for those elements. The local forces are then computed again for each element, and the process is repeated until \( v_i = 0 \) for all \( i \) and the avalanche stops. The size \( s \) of the avalanche is the total number of elementary moves during such an avalanche. Then the external stress is increased, so that exactly one of the elements becomes unstable, and a new avalanche is initiated. While such a discretization of the dynamics neglects the fact that the local velocity of the crack front should be proportional to the local SIF, this is a standard technique employed in models of the depinning transition, going back to the work of Leschhorn [18], and is known to have no influence on the scaling behavior we study here. Moreover, this procedure is essential to be able to drive the system quasistatically. Such quasistatic driving has the advantage over continuous time models with a finite driving velocity [7] that avalanches and clusters can be defined without ambiguity also for a zero threshold. As the crack front advances, the applied SIF \( K_{i,\text{inst}} \) decreases at a rate proportional to the instantaneous average velocity \( v(t) = 1/L \sum_i v_i(t) \) of the front, with a proportionality constant \( k \). Thus, the cutoff of the avalanche size distributions is expected to scale with \( k^{-1} \).

B. Methods to extract the local clusters

To extract the clusters from an avalanche, we use two methods. The first is the avalanche map (AM), based on a two-dimensional array \( A(x,y) \), for \( k = 0.05 \) and \( \gamma = 1 \), with the area swept over by different avalanches denoted by different shades of gray (time is running from black to white). (b) The same data as in (a) represented by the velocity matrix \( V(x,y) = 1/W(x,y) \) with a zero threshold. The crack front is moving to the positive \( y \) direction. Black corresponds to zero velocity, while regions of finite velocity are shown in white. (c) An example of a structure of a single avalanche as identified by the AM method. (d) The same avalanche as in (c) extracted by using the WTM method.

FIG. 1. (a) An example of the avalanche map \( A(x,y) \), for \( k = 0.05 \) and \( \gamma = 1 \), with the area swept over by different avalanches denoted by different shades of gray (time is running from black to white). (b) The same data as in (a) represented by the velocity matrix \( V(x,y) = 1/W(x,y) \) with a zero threshold. The crack front is moving to the positive \( y \) direction. Black corresponds to zero velocity, while regions of finite velocity are shown in white. (c) An example of a structure of a single avalanche as identified by the AM method. (d) The same avalanche as in (c) extracted by using the WTM method.
FIG. 2. (Color online) The main figure shows the size distributions of the avalanches for $\Gamma_0=1$ and various values of $k$ (open symbols), as well as the size distributions of the avalanches for a fixed $k=0.0125$ and various values of $\Gamma_0$ ranging from $\Gamma_0=0.4$ to $\Gamma_0=1.0$ (filled symbols, with the latter set of distributions displaced vertically for clarity). The solid line is a guide to the eye and corresponds to $\tau=1.25$. The inset shows a data collapse of the distributions with $\Gamma_0=1$ and various values of $k$, with $\tau=1.25$ and $1/\sigma_k=0.725$.

a line element moves $n$ steps, as both the initial and final pinned configurations are classified as immobile. Thus, the avalanche in Fig. 1(d) is a bit thinner in the y direction as compared to the one in Fig. 1(e), and parts of the avalanche involving motion of a single step forward only are lost completely. However, our results suggest that this does not affect the statistical properties of the clusters. In the case of the discrete line model, another issue is that for soft lines involving motion of a single step forward only are lost compared to the one in Fig. 1.

In the bottom panel of Fig. 2, open symbols $a_{WTM}$ and filled symbols, with the latter set of distributions displaced, as well as the size distributions of the avalanches for $\Gamma_0=1$ and various values of $k$, with $\tau=1.25$ and $1/\sigma_k=0.725$.

III. RESULTS

A. Avalanche and cluster size distributions

The probability distributions of avalanches and cluster sizes for $L=2048$, $\Gamma_0=1$, and various values of $k$ are shown in Figs. 2 and 3, respectively. The sizes $s$ of the avalanches (Fig. 2, open symbols) are distributed according to a power law with a $k$-dependent cutoff,

$$P(s) \approx s^{-k f_2(s/k^{1/\sigma_k})},$$

with $f_2(x)$ as a scaling function describing the shape of the cutoff, $\tau=1.25 \pm 0.05$, and $1/\sigma_k=0.725 \pm 0.08$. The value of $\tau$ is in good agreement with results from simulations [7] and unpublished experiments [19]. However, the scaling of the cutoff was not considered in Ref. [7]. Notice that the scaling function $f_2(x)$ displays a bump around the cutoff of the distribution [20], and thus the true power-law exponent can be seen directly only for $s$ well below the cutoff scale. The exponent values we quote are obtained by optimizing the data collapse of the distributions. To demonstrate the robustness of our results with respect to changing the values of various parameters of the model, we also show the avalanche size distributions for a fixed $k$ but varying $\Gamma_0$ in Fig. 2 (filled symbols): the distributions turn out to be virtually independent of $\Gamma_0$ in the range $0.4 \leq \Gamma_0 \leq 1.0$ considered in Fig. 2. Analogously, the cluster sizes or areas $\sigma_{\text{AM}}$ (top panel of Fig. 3) and $\sigma_{\text{WTM}}$ (bottom panel of Fig. 3) extracted by the AM and WTM methods, respectively, are both observed to scale according to

$$P(a) \approx a^{-\tau_{\text{AM}}(a/k^{1/\sigma_k})},$$

but with an exponent $\tau_{\text{AM}}$ significantly different from $\tau$, i.e., $\tau_{\text{AM}}=1.52 \pm 0.05$ and $\tau_{\text{WTM}}=1.53 \pm 0.05$. Thus, the two methods to define the clusters give the same $\tau_{\text{WTM}}$ exponent within error bars. In this case the scaling function $g_2(x)$ does not exhibit a bump, and the scaling of the distributions extends all the way to the cutoff scale. Also for the cluster size distributions similar results are obtained for different values of $\Gamma_0$ (not shown). The observed exponent value is somewhat lower than that quoted in Refs. [6,7] for the WTM approach, but our analysis below indicates that also the data presented...
FIG. 4. (Color online) The scaled size distributions of the WTM clusters for various threshold values \(v_{th}\) with \(\tau_a=1.5\) and \(1/\sigma_k=1.8\). Filled symbols correspond to the simulations, while open symbols are from an experiment on planar crack propagation in Plexiglas plates [6], with \(v\approx11\ \mu m/s\) and a pixel size of 1.7 \(\mu m^2\). The two data collapses (simulations and experiments) have been shifted on top of each other to facilitate comparison.

in [6,7] are perfectly consistent with \(v_{th}^{WTM}=1.5\). This is because in Refs. [6,7] also the region of the cutoffs of the distributions was included in the fits, thus overestimating the true exponent value. Notice also that the cutoff of the cluster size distributions scales with \(k\) as \(k^{-1/\sigma_k}\), with \(1/\sigma_k=0.7\pm0.08\) for the AM approach, while we obtain \(1/\sigma_k=0.8\pm0.08\) for the WTM. The different \(1/\sigma_k\) exponents have the same value within error bars, and we suspect that the small variation observed might be related to a finite-size effect.

B. Comparison with experimental results

In order to compare directly our results to the experiments [6], we study WTM clusters as a function of a finite velocity threshold \(v_{th}\). The threshold introduces a cutoff in the distribution that scales as

\[
P(a) = a^{-\tau_a}g_k(a^{1/\sigma_k}),
\]

with \(\tau_a=1.53 \pm 0.05\) and \(1/\sigma_k=1.8 \pm 0.1\) (see Fig. 4). There, we compare the simulation results with experimental data on planar crack propagation in Plexiglas plates [6]. Both sets of data can be collapsed by using exponent values \(\tau_a=1.5\) and \(1/\sigma_k=1.8\). By shifting the two data collapses on top of each other, we observe perfect agreement between the simulations and the experiment. Notice again that in Refs. [6,7], the cutoffs of the distributions were included in the power-law fits, thus overestimating the true exponent value.

C. Link between global avalanches and local clusters

To explain these results, we consider the breakup process of an avalanche into a number of clusters. When an avalanche grows by 1 unit (i.e., when a single element of the line moves forward by 1 unit), there are three possibilities of how this can affect the number \(N(s)\) of distinct clusters within an avalanche: (i) a new cluster is initiated \((N\rightarrow N+1)\), (ii) two existing clusters are merged into a single cluster \((N\rightarrow N-1)\), or (iii) an existing cluster grows by 1 unit \((N\rightarrow N)\). We associate probabilities \(p_1, p_2,\) and \(1-p_1-p_2\), respectively, to these events. These probabilities are dynamic variables as they depend on the instantaneous structure of the avalanche. On the average, however, \(p_1=p_2\) as \(p_1>p_2\) would clearly indicate an increase in \(p_2\) and a subsequent tendency for \(p_1\) to decrease again: when the number of clusters grows, there is less and less space to create more clusters and the probability to merge existing clusters should increase. Similarly, \(p_1<p_2\) would imply that sooner or later most of the clusters would have merged, and consequently the probability \(p_2\) for further cluster merging decreases. Thus, for large enough avalanches, the number \(N\geq1\) of clusters during the growth of an avalanche follows a random walk up to \(s\) steps, with a reflective boundary at \(N=1\) (as each avalanche consists of at least one cluster). Assuming for simplicity that the process is not correlated indicates that for \(s\gg1\), \(\langle N(s) \rangle \sim s^{\alpha}\), with \(\alpha=1/2\). Our numerical results confirm this relation, and a scaling form

\[
\langle N(s) \rangle = s^\alpha \tilde{N}(s/k^{-1/\sigma_k}),
\]

with \(\tilde{N}(s)\) as a scaling function, \(\alpha=0.47 \pm 0.05\), and \(\sigma=0.76 \pm 0.08\) (see Fig. 5, i.e., the cutoff scales with \(k\) again with the same exponent \(1/\sigma_k=0.75\)). Thus, the cluster sizes scale as \(a \sim /N(s)\) or \(s \sim a^2\). \(P(a)da=P(s)ds\) then implies that \(P(a) \sim a^{\tau_a-2}\), with

\[
\tau_a = 2 - \alpha - 1.
\]
where $a = \frac{L}{s}$ (as evidenced by the crossover to a plateau in Fig. 5), consistent with the observation that the cutoff of the avalanche and cluster size distributions scale with the same exponent $1/\sigma$. Notice also that while deriving Eq. (8) we did not make assumptions about the precise form of the long-range kernel nor the spatial dimension of the system, and we thus expect Eq. (8) to be valid for a wide class of avalanching systems where the avalanche dynamics is governed by long-range interactions. Notice that if the avalanche breakup process is correlated such that $\langle N(s) \rangle \sim s^{\alpha}$ with $\alpha \neq 1/2$, Eq. (8) can be generalized to read $\tau_c = (\tau - \alpha)/(1 - \alpha)$.

D. Cluster morphology and roughness of the crack front

In experiments [6] and simulations [7] it was found that the aspect ratio of the clusters was scaling as $l_y = \xi L_c$, with $\xi = 0.6$ and where $l_x$ and $l_y$ are the linear sizes of the cluster in the $x$ and $y$ directions, respectively. This result is in agreement with the roughness of the crack front measured in early experiments, but disagrees with the known value of the roughness exponent of the crack line model $\xi = 0.39$ [14,21]. Moreover, by investigating a larger range of length scales, a recent experimental work on roughness of the crack front has shown a crossover from a roughness exponent of around 0.6 to a lower value corresponding to the line model prediction of around 0.39 [15,16]. This is surprising since we have just shown that the experimental avalanche statistics is in perfect agreement with the predictions of the model. This puzzle can be resolved, noticing that $\xi = 0.39$ is expected to hold only for length scales larger than the Larkin length $L_c$ [22], which for our model scales like $L_c = \Gamma_0^{\zeta_0}/R^2$, where $\xi$ and $R$ are the correlation length and the amplitude of the disorder (here, $\xi = R = 1$), respectively [17]. Below $L_c$, the crack line roughness should scale with an exponent $\zeta_x = 1/2$.

In the top panel of Fig. 6 we report the scaling of $l_x$ with $L_c$ for different values of $\Gamma_0$. By rescaling $l_x$ by $L_c = \Gamma_0^{\zeta_0}$, we observe a crossover for the scaling exponent going from $\zeta_x = 0.55 \pm 0.05$ at small scales to $\zeta_x = 0.39 \pm 0.03$ at large scales. In the top panel of Fig. 6 we also report the scaling of the root-mean-square height fluctuations $\Delta h / \langle \delta \rangle^{1/2}$ [35] with $\Delta h = h_{\text{max}} - h_{\text{avg}}$ of the crack front, expected to scale like $\Delta h / \langle \delta \rangle^{1/2} \sim \delta$. Again, by rescaling $\delta$ by $\Gamma_0$, we observe a crossover from a small-scale regime with $\zeta_x = 0.48 \pm 0.05$ to the large-scale value $\zeta_x = 0.37 \pm 0.03$. The bottom panel of Fig. 6 shows that similar conclusion can be made by considering the power spectrum $S(q)$ of the line profile, expected to scale as $S(q) \sim q^{-2\zeta_0 + 1}$. Rescaling the data by $L_x = \Gamma_0^{\zeta_0}$ according to $S(q) = \Gamma_0^{2\zeta_0} \Gamma_0^{-2\zeta_0} \delta^2$, we observe a crossover for the scaling exponent going from $2\zeta_0 + 1 = 1.96 \pm 0.04$ for large $q$ to $2\zeta_0 + 1 = 1.76 \pm 0.04$ for small $q$, in agreement with the analysis of root-mean-square height fluctuations. We have verified that in the present model, where the crack front is constrained to move along the $y$ direction, there is no multiscaling [23]. Notice that while we observe different crossover scales for the roughness of the line and the scaling of the morphology of the clusters, both sets of data scale with the Larkin length $L_c$, and thus the crossover length is proportional to $L_c$, in both cases. However, the fact that these two length scales are not the same implies that there is a range of scales for which the roughness of the line appears to scale with the asymptotic exponent $\zeta \approx 0.39$, while the cluster aspect ratio is still characterized by the larger $\zeta_x \approx 0.55$. It is probably this range of scales that has been investigated in earlier numerical studies [7].

IV. CONCLUSIONS

To summarize, we have demonstrated how avalanches are broken into spatially disconnected clusters in systems with long-range interactions. Both the avalanches as a whole and their localized parts or clusters exhibit scaling, but with different power-law exponents, related by a scaling relation de-
rived from a simple random-walk argument for the avalanche breakup process. We also showed that large enough clusters exhibit aspect ratio scaling with an exponent consistent with the roughness exponent of the crack line model [14,21]. Due to the general nature of both the model considered as well as our arguments, we expect these results to be applicable in a large class of avalanching systems with long-range interactions, ranging from crack propagation to contact lines in wetting to domain walls in ferromagnets.

Of particular interest here is our explanation of the different scaling exponents observed for global avalanches and local clusters in the crack line model [7], and its potential implications for further studies of avalanche phenomena in systems with long-range interactions. An interesting prospect would be to explore the possibilities to explain the observed acoustic emission exponents in peeling of paper sheets [8] with similar arguments. The observation that the avalanche breakup process in the line model can be described by a simple random walk will be interesting to test in experiments not only on crack front propagation in Plexiglas plates, but also in other avalanching systems with long-range interactions, such as for domain-wall dynamics in ferromagnetic thin films [9], as well as for imbibition [24].

An important point regarding the roughness of the crack front and the related cluster morphology is that their scaling properties depend on the range of length scales considered: the asymptotic scaling is observed only for length scales larger than a crossover scale proportional to the Larkin length of the crack line. Earlier simulations [7] as well as experiments [6] have most likely been probing the regime below this crossover scale, possibly explaining the different exponent values observed in those cases. However, for these small length scales additional complications arise due to the overhangs observed in the experimental crack profiles, which are excluded by construction in the present line model. Such overhangs might be responsible for the multiscaling observed for the experimental crack profiles for short length scales [16]. We are working to extend the crack front line model to include the possibility to form overhangs and are planning to check if such an extension is sufficient to reproduce the experimentally observed short length scale scaling properties of the crack fronts.

The connection between avalanches and clusters in avalanching systems with long-range interactions opens up also interesting possibilities to understand the observed spatiotemporal correlations between avalanches in a number of systems, such as the space-time coupling of avalanches observed in plastically deforming crystals [25], and even the foreshock and aftershock sequences of earthquakes: one possibility that worth further studies would be to check if such correlated avalanches could in fact be local clusters, with the observed correlations arising naturally from the fact that the localized clusters are parts of the same global avalanche.

**ACKNOWLEDGMENTS**

M. J. Alava and X. Illa are thanked for interesting discussions, and D. Bonamy for sending us the data of the avalanche size distributions of Ref. [7]. K. J. Mályö is thanked for the possibility to publish the experimental data in Fig. 4. L.L. wishes to thank the Academy of Finland for financial support.

[23] Due to artifacts related to the discretization of the front y position $h(x)$, an effective multiscaling is observed at very small scales, but we have checked that this spurious effect disappears if we employ a continuous variable for $h(x)$.