Absorbing-state phase transitions with extremal dynamics

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Abstract

Extremal dynamics represents a path to self-organized criticality in which the order parameter is tuned to a value of zero. The order parameter is associated with a phase transition to an absorbing state. Given a process that exhibits a phase transition to an absorbing state, we define an “extremal absorbing” process, providing the link to the associated extremal (nonabsorbing) process. Stationary properties of the latter correspond to those at the absorbing-state phase transition in the former. Studying the absorbing version of an extremal dynamics model allows to determine certain critical exponents that are not otherwise accessible. In the case of the Bak-Sneppen (BS) model, the absorbing version is closely related to the “f-avalanche” introduced by Paczuski, Maslov and Bak [Phys. Rev. E 53, 414 (1996)], or, in spreading simulations to the “BS branching process” also studied by these authors. The corresponding nonextremal process belongs to the directed percolation universality class. We revisit the absorbing BS model, obtaining refined estimates for the threshold and critical exponents in one dimension. We also study an extremal version of the usual contact process, using mean-field theory and simulation. The extremal condition slows the spread of activity and modifies the critical behavior radically, defining an “extremal directed percolation” universality class of absorbing-state phase transitions. Asymmetric updating is a relevant perturbation for this class, even though it is irrelevant for the corresponding nonextremal class.

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I. INTRODUCTION

Extremal dynamics has been employed extensively in modelling far from equilibrium systems such as biological evolution [1] and driven interfaces [2,3]. Although processes with extremal dynamics do not have a phase transition (there is no control parameter) they exhibit scaling properties reminiscent of those observed at continuous phase transitions [4,5]. Indeed, it was suggested some time ago that the appearance of “self-organized” scaling properties in extremal dynamics and in sandpile models corresponds to forcing the order parameter (associated with an underlying phase transition) to zero from above [6]. The purpose of this work is to explore the connection between these scaling properties and those observed at a phase transition to an absorbing state. The connection between extremal dynamics and directed percolation (DP), the prime example of an absorbing-state phase transition, was first suggested by Paczuski, Maslov and Bak [4] and investigated in detail by these authors in the context of the Bak-Sneppen (BS) model and related processes [5]. (The latter work, as well as Ref. [7], clearly demonstrated that the critical exponents of the BS model are not those of DP.) Sornette and Dornic [8] and Grassberger and Zhang [9] have shown how a variant of directed percolation may be transformed via extremal dynamics to display SOC. These studies indicate that self-organized criticality (SOC) [10] under extremal dynamics arises because the system is driven to a critical point associated with a phase transition to an absorbing state [8], as is also the case for sandpiles [11].

In the present work we are particularly interested in the modifications needed to transform a (non-extremal, non-SOC) model having an absorbing state to one exhibiting SOC under extremal dynamics. We develop a general scheme relating the two classes of models via an intermediate, “extremal-absorbing” process whose absorbing-state critical point corresponds exactly to the critical behavior observed in the corresponding SOC model. Two examples (the BS model and an extremal contact process) are studied in detail, yielding refined estimates for critical properties, and evidence of a new universality class associated with absorbing phase transitions under extremal dynamics.

The prime example of extremal dynamics is the BS model [1,12], proposed to explain mass extinctions observed in the fossil record. While its application in the evolutionary context is debated [13], it remains an intriguing and incompletely understood example of scaling behavior far from equilibrium. The contact process (CP) [14] is the most familiar example of a Markov process exhibiting a phase transition to an absorbing state. We focus on the absorbing version of the BS model, and the extremal version of the CP, to illustrate the relations between extremal dynamics and absorbing phase transitions. Our analysis of the spread of activity leads to new or refined values for the exponents $\delta$, $\eta$, $\nu_{\parallel}$, $\beta'$ and $z_{sp}$, and for the critical threshold of the BS model. (For extremal dynamics, the avalanche exponent is $\tau = 1 + \delta$.) Finite-size scaling analysis of stationary properties at the critical point yields estimates of the exponent ratios $\beta/\nu_{\perp}$ and $\nu_{\parallel}/\nu_{\perp}$. 
Studies of modified BS models have shown that scaling properties are insensitive to changes that preserve its basic symmetries (that is, invariance under translation and reflection) [15–17], pointing to the existence of a BS universality class. Nonextremal models that exhibit a phase transition to an absorbing state, and that possess these same symmetries, and no additional ones, belong generically to the directed percolation (DP) universality class [18,19]. Here we show that such models fall in a new “ extremal-DP” universality class when modified to follow extremal dynamics. The absorbing phase transition corresponding to the scaling behavior of the BS and other extremal models belongs to the extremal-DP class, not that of ordinary directed percolation. In other words, extremal dynamics is a relevant perturbation for absorbing-state transition, just as was shown by Sneppen in the context of interface depinning [2].

The balance of this paper is organized as follows. Section II presents a general scheme linking ordinary absorbing-state phase transitions and extremal dynamics via an intermediary “extremal-absorbing” model. In Sec. III we describe our simulation method and results. Our findings regarding scaling and universality are discussed in Sec. IV. Mean-field analyses are presented in the Appendix.

II. ABSORBING STATE MODELS AND EXTREMAL DYNAMICS

In this section we examine how a stochastic model with an absorbing-state phase transition may be transformed to exhibit scale invariance under extremal dynamics. We begin, for generality, by defining a rather abstract scheme, and then discuss specific examples. A large class of models exhibiting an absorbing-state phase transition may be formulated as follows [20–23]. Consider a stochastic process $S$ defined on a connected graph $G$ of $N$ sites. ($G$ consists of a set of sites with links between certain pairs of sites. Typical examples are a ring of $N$ sites, and the $d$-dimensional hypercubic lattice $Z^d$, with links between nearest neighbors.) The state $\sigma(i)$ of site $i$ is 0 or 1, the latter value denoting an active site, the former an inactive one. For each site $i$ in $G$ we define a neighborhood $v(i) \subset G$, or, more generally, a set of neighborhoods $v_1(i), v_2(i),...,v_n(i)$.

The dynamics of $S$ proceeds in steps. Each step involves choosing an active site $i$ (the central site for this step), at random, and changing the states of the sites in $v(i)$ according to a certain rule $f$. In case there are two or more neighborhoods, one of them, $v_r(i)$ say, is chosen at random from the collection, with probability $p_r$, and a rule $f_r$ is applied to the site or sites in $v_r$. In general $f$ (or $f_r$) is a probabilistic rule. At each step the number of active sites may change, and if at any moment there are no active sites ($\sigma(i) = 0, \forall i \in G$), the process has fallen into an absorbing state and there is no further evolution. Otherwise the dynamics proceeds to the next step.

Using $\sigma_n$ to denote the entire set of activity variables $\sigma(1), \sigma(2),...,\sigma(N)$ at step $n$, the dynamics generates a sequence $\sigma_1, \sigma_2,...$ starting from the initial
configuration \(\sigma_0\). It is frequently of interest to associate a continuous time variable \(t\) with the process. This is usually done by associating a time increment \(\Delta t = 1/N_a\) with each step, where \(N_a\) is the number of active sites just before the step is realized. We define the order parameter as \(\rho(t) = \text{Prob}[\sigma_t(i) = 1]\), i.e., the fraction of active sites at time \(t\). (The event space here is the set of all realizations of the process up to time \(t\), starting from a given initial probability distribution on configuration space.) If the stationary order parameter, defined by \(\lim_{t \to \infty} \lim_{N \to \infty} \rho\), vanishes, the process is said to be in the absorbing phase; otherwise it is in the active phase.

A simple model exhibiting a phase transition to an absorbing state is the contact process (CP) [14]. Here we consider the one-dimensional version. There are three sets \(v_r(i)\), conveniently denoted as \(v_0(i) = i\) and \(v_\pm(i) = i \pm 1\). The associated probabilities are \(p_0\) and \(p_\pm = (1 - p_0)/2\). In terms of the usual parametrisation [14,23], \(p_0 = 1/(1 + \lambda)\), where \(\lambda \geq 0\) represents the rate of spread of activity. (In the “epidemic” interpretation of the CP, active sites represent infected organisms, inactive sites susceptibles, and \(\lambda\) is the infection rate.) The updating rules are: \(f_0 = 0, f_\pm = 1\). In other words, an active site has a probability per unit time of \(1/(1 + \lambda)\) to become inactive, while an inactive site \(j\) becomes active at rate \(\lambda n_a(j)/[2(1 + \lambda)]\), where \(n_a(j)\) is the number of active neighbors of site \(j\). As is well known, the one-dimensional CP exhibits a continuous phase transition between an absorbing phase and an active one at \(\lambda_c \approx 3.29785\) [20,22,23].

It is convenient to associate the control parameter with the updating rule \(f\) rather than with the probabilities \(p_r\). We therefore reformulate the CP as follows. With the sets \(v_0(i)\) and \(v_\pm(i)\) defined above, we take \(p_0 = 1/2\) and \(p_\pm = 1/4\), and define \(q = \lambda/(1 + \lambda)\). The updating functions are:

\[
f_0 = \begin{cases} 1, \text{ w.p. } q \\ 0, \text{ w.p. } 1-q \end{cases}
\]

(‘w.p.’ denotes “with probability”), and

\[
f_\pm = \begin{cases} 1, \text{ w.p. } q & \text{if } \sigma(i \pm 1) = 1 \\ 1, \text{ w.p. } q & \text{if } \sigma(i \pm 1) = 0 \\ 0, \text{ w.p. } 1-q \end{cases}
\]

It is easy to verify that the transition rates satisfy \(w(0 \to 1)/w(1 \to 0) = n_a\lambda/2\), just as in the original formulation. The critical value \(q_c \approx 0.76733\).

The following three-site contact process (CP3) will play an important role in our analysis [24]. For each site we define the set \(v(i) = \{i-1, i, i+1\}\) (the central site and its nearest neighbors). The updating function \(f\) takes values of 1 and 0 with probabilities \(q\) and \(1-q\) respectively, independently at each of the three sites in \(v(i)\). (In Ref. [24] this is called 'model 3'.) Simulations of the CP3 show that it exhibits a continuous phase transition at \(q = q_c \approx 0.63523(3)\). Further interacting particle systems, such as the pair contact process [25] and the diffusive CP [26], can be accommodated within the scheme set out above.
We shall assume that the process $S$ is defined so that the control parameters (such as $q$) are associated with the updating rule $f$. Each time a site is updated, the value of $f$ may be determined by comparing a random number $x$ with the parameter in question. (This is of course the usual procedure in simulations.) In the CP3, for example, we take $f = 1$ if $x < q$, and zero otherwise, where $x$ is uniformly distributed on the interval $[0,1]$. Call the random number associated with the most recent updating of site $i$, $x_i$, so that $\sigma(i) = \Theta(q - x_i)$ with $\Theta$ the unit step function. (The initial values of the $x_i$ are assigned according to the state variables $\sigma(i)$. For example, if all sites are initially active, we draw the initial $x_i$ from the distribution uniform on $[0, q]$. For the CP, Eq. (1) requires that we update the central site $i$ with a number chosen uniformly from $[0,1]$. According to Eq. (2), the same applies when updating an inactive neighbor ($i \pm 1$), but when updating an active neighbor, the random number $x$ is drawn from the interval $[0, q]$, since an active neighbor remains active.

Summarizing, we have shown how a particle system $S$ may be formulated using a set of random variables $x_i$, such that site $i$ is active if $x_i$ is smaller than a certain parameter $q$. $S$ suffers an absorbing-state phase transition at $q = q_c$.

We now define two related processes, $S_{EA}$ and $S_E$. The former, extremal absorbing process, is obtained by modifying how the central site is selected. Instead of choosing it at random from among the currently active sites, it is taken to be the active site having the smallest $x_i$. As in the original process $S$, if there are no active sites (i.e., $x_j > q, \forall j \in G$), the process has reached an absorbing configuration and the evolution ceases. Thus $S_{EA}$ possesses an absorbing state, and since the original process $S$ exhibits a phase transition between an active and an absorbing phase as the control parameter $q$ is varied, we expect $S_{EA}$ to do so as well, at some value $q_{c,E}$. (The reason is that the relative likelihood of generating and destroying active sites varies with $q$, just as in $S$.) Mean-field theory (see Appendix) yields $q_{c,E} = q_c$. Due to the different correlations generated under extremal dynamics, however, the critical value $q_{c,E}$ of $S_{EA}$ is in general different from $q_c$.

We define the extremal process $S_E$ by relaxing the condition in $S_{EA}$, that the extremal site $i$ must be active (i.e., have $x_i < q$) for the dynamics to proceed. If $S$ is the original contact process, then $S_E$ is a process in which either the minimal site or one of its nearest neighbors is updated at each step. If $S$ is the CP3, $S_E$ is the familiar Bak-Sneppen model. Note that $S_E$ has no absorbing state, hence no phase transition to such a state. Its stationary properties are nevertheless intimately connected with the critical-point properties of $S_{EA}$, as we now explain.

Of particular interest is the stationary probability density $p(x)$ of site variables under extremal dynamics. As is well known, $p(x) = C\Theta(x - q_{c,E})\Theta(1 - x)$ in the Bak-Sneppen model, in the infinite-size limit. ($C = 1/(1 - q_{c,E})$ is the normalization factor.) We expect $p(x)$ to exhibit a step-function singularity in any extremal model $S_E$ [15]. This feature is in fact already present in the original absorbing-state model $S$ at its critical point, because at the critical point $q = q_c$, the stationary density of active sites (having $x_j < q$) tends to
zero as the system size $N$ goes to infinity. The distribution on the “allowed” region $x > q$ is uniform, since the $x_j$ are drawn from a uniform distribution. Thus $\overline{p}(x)$ jumps from zero to a finite value at $x = q_c$. In the supercritical regime ($q > q_c$), $\overline{p}(x)$ is equal to a constant $p_1$ for $x < q$, such that $qp_1 = \rho$ (the order parameter), and takes a different constant value, $p_2$, on the interval $[q, 1]$. Once again, the stationary density is discontinuous at $x = q$.

What holds for $S$ also holds qualitatively for $S_{EA}$. The critical value $q_{c,E}$ may, as noted, differ from $q_c$, but since $S_{EA}$ exhibits an absorbing-state phase transition, its stationary distribution $\overline{p}(x)$ also has a step-function singularity. Just at $q = q_{c,E}$, the order parameter $\rho = 0$, but as $N \to \infty$ the survival time of the process tends to infinity. This means that the process can survive indefinitely, with the choice of the central site restricted to the set having $x \leq q_{c,E}$. The presence of active sites in the range $q_{c,E} < x < q$ is then irrelevant, since a site with $x \leq q_{c,E}$ is always available. Thus for $q \geq q_{c,E}$, the distribution $\overline{p}(x)$ exhibits a step-function singularity at $q_{c,E}$. Extremal dynamics effectively “pins” the singularity at $q_{c,E}$. The foregoing remarks on $S_{EA}$ obtain in the infinite-size limit; for finite $N$ there is a nonzero probability (for $q < 1$), that all sites have $x > q$ so that the system eventually becomes trapped in the absorbing state. (The mean lifetime, however, is expected to grow exponentially with $N$, for $q > q_{c,E}$.)

In summary, if $S$ exhibits an absorbing-state phase transition, then $S_{EA}$ should as well, although not necessarily at the same value of $q$. The distribution $\overline{p}(x)$ possesses a step function singularity in both cases. Near the critical point ($q \gtrsim q_c$) of the original process $S$ we expect $\rho \sim (q - q_c)^\beta$, with $\beta$ the critical exponent associated with the order parameter. Below the upper critical dimension ($d_c = 4$ for DP [18,19]), $\beta < 1$. In the supercritical regime of $S_{EA}$, on the other hand, $\rho = \int_0^q \overline{p}(x)dx \propto (q - q_{c,E})$ since $\overline{p}(x)$ jumps from zero to a finite value at $x = q_{c,E}$. Thus the order parameter exponent $\beta$ is unity in $S_{EA}$. This is but one example showing that the extremal condition changes the value of a critical exponent. Another example, established some time ago by Paczuski and coworkers [5,27], is that the spreading exponent $\eta$ (defined Sec. III.C) is generically zero for extremal dynamics, whereas $\eta > 0$ for directed percolation. Further evidence that extremal dynamics modifies critical exponent values will be given below.

In $S_E$, the central site is always chosen (in the $N \to \infty$ limit) from the set $\{i : x_i \leq q_{c,E}\}$, just as in $S_{EA}$ at its critical point. Since the order parameter is zero in the latter case, we may assert that extremal dynamics effectively tunes the order parameter to zero. ($\rho$ approaches zero from above as $N \to \infty$.) As we have seen, the existence of sites with $q_{c,E} < x < q$ becomes irrelevant in $S_{EA}$, in the infinite-size limit. In other words, $S_E$ is identical to the critical process $S_{EA}$ in this limit. (Starting from the same initial configuration, and using the same set of random numbers, the same sequence of sites will be updated in the two processes.)

Since the extremal version of the CP3 is the familiar Bak-Sneppen model, we shall refer to the CP3$_{EA}$ as the absorbing Bak-Sneppen (ABS) model. Our
objective is to characterize the behavior of the ABS model and of the extremal
and EA versions of the contact process. The ABS model is closely related to
the f-avalanche process studied in Ref. [5]. An f-avalanche (in the present
notation, q-avalanche), begins when the minimal site variable $x_{\min} < q$, the
minimum having been greater than $q$ at the preceding step or steps, and con-
tinues until the minimum is once again $> q$. (As $q$ approaches $q_{c,E}$ from below,
the mean avalanche duration diverges.) The dynamics of the BS model contin-
ues, regardless of whether a given avalanche has terminated or not. But in the
ABS model $x_{\min} > q$ represents an absorbing state and the dynamics ceases.
In the BS model, it is common to analyze the properties of q-avalanches in the
stationary state. It is similarly of interest to study stationary properties of the
ABS model, attained once the system has relaxed, after an initial transient pe-
riod. We may also study the mean lifetime of the active state as a function of
system size. Another approach to studying absorbing-state phase transitions
consists in following the spread of activity starting from a single active site.
This spreading phenomenon in the BS model was studied in Refs. [4,5], where
it is called the BS branching process, and in Ref. [7] under the name of the
BS($\tilde{p}$) model.

The assertions regarding extremal and extremal-absorbing models are sup-
ported by mean-field theory (MFT), which is presented in the Appendix. In
particular, for the ABS model $\overline{p} = \frac{2}{3} \Theta(x - \frac{1}{3})$, just as in the MFT of the orig-
inal BS model. In the extremal and the extremal-absorbing contact process,
the stationary probability densities exhibit (for $q > q_{c,E} = 1/2$), two disconti-
nuities, one at $x = \kappa \equiv q^2/(3q - 1)$, the other at $x = q > \kappa$. [These coalesce
at $q = 1/2$. Note that the parameter $q$ continues to influence the form of $\overline{p}$ in
the extremal contact process, due to the nature of the updating rule, Eqs. (1)
and (2).] These predictions are in qualitative agreement with simulation.

III. SIMULATION RESULTS

A. Absorbing Bak-Sneppen Model: Simulation Method

Before discussing our results we insert an observation on simulations of the
BS model. Since the site with the smallest variable, $x_{\min}$, must be identified
at each step, it becomes important to devise an effective search strategy. An
efficient general-purpose search algorithm uses a binary tree structure to iden-
tify $x_{\min}$. One approach [7] utilizes a lattice of $2^n$ sites. At the first level of
selection, each site is compared with one of its neighbors and the minimum of
the pair selected. At the next level the minimum between each neighboring
pair is selected, and so on, so that at the $n$-th level the global minimum is
identified.

A second binary scheme [28] is formulated as follows. Site 0 is placed
at the apex of the tree. Site 1 is placed on the level below the apex, to the left
of 0 if $x_1 < x_0$, to the right if $x_1 > x_0$. A site $i$ is added to the tree in the
following way: we go down the tree comparing $x_i$ with the variables $x_1, ..., x_{i-1}$,
turning left or right depending on whether \( x_i \) is smaller or larger than \( x_j \), until we find an empty site. Building the tree in this way, \( x_{\text{min}} \) will occupy the leftmost position in the tree. In these schemes, maintaining the tree structure, once constructed from the initial set of variables \( x_i \), requires a small number of operations at each step, and is many times more efficient that a repeated global search for the minimum. We find, nonetheless, that a suitably restricted search requires less cpu time in the stationary state.

A special property of the BS model (shared by its absorbing version, and by other extremal models), is that the minimal site falls, with a probability approaching unity as the system size grows, in the interval \([0, q_{c,E}]\). At the same time the density of sites with values in this interval approaches zero as \( N \to \infty \). This suggests maintaining a list of sites having \( x < q_{c,E} \) \([29]\). Then the search for \( x_{\text{min}} \) may be restricted to the list, except for the rare instances in which the latter is empty. (For the ABS we must in any case restrict the search to sites with \( x \leq q \).) If the system is large, so that the typical number of sites with \( x < q \) is not small, it becomes advantageous to introduce a second list, of sites having \( x < q^{**} < q_{c,E} \). When this relatively short list is nonempty (as is usually the case) the search for \( x_{\text{min}} \) is restricted to it. In studies of the BS model, we obtained the greatest efficiency using \( q^{**} = 0.54 \), while the criterion for the first list was \( x < 0.65 \), that is, slightly below \( q_{c,E} \). (The occasional need to perform a global search, in the rare instances when both lists are empty, is more than compensated by their reduced sizes when using these values.) Compared with the binary tree method, our approach results in threefold reduction in CPU time, in the stationary state, for a system of 1000 sites. (The binary tree approach may prove more efficient for studying transients, since initially the lists will not be short, if the \( x_i \) values are chosen uniformly on \([0,1]\).)

B. Absorbing Bak-Sneppen Model: Stationary Properties

Using the simulation method described above, we conducted extensive studies of the one-dimensional ABS model. We initialize the system with all sites active and allow it to relax until mean properties fluctuate about stationary values. The stationary properties are then obtained from temporal averages over the set of surviving realizations. Each step corresponds to a time interval of \( \Delta t = 1/N_a \), with \( N_a \) the number of active sites just before the updating is performed. Results for the stationary order parameter (i.e., the density of sites with \( x < q \)), on lattices confirm that in the supercritical regime (\( q > q_{c,E} \)), the order parameter grows linearly with \( q - q_{c,E} \), as anticipated in Sec. II.

We study the finite-size scaling behavior of the stationary order parameter \( \overline{\rho} \) and of the lifetime \( \tau \) at the critical point. (\( \tau \) is obtained from an exponential fit to the survival probability \( P_s(t) \).) The expected finite-size scaling behaviors at the critical point are: \( \overline{\rho} \sim L^{-\beta/\nu_\perp} \) and \( \tau \sim L^{\nu_1/\nu_\perp} \). We performed simulations at \( q = 0.66701 \) and \( 0.66702 \), the latter being the preferred literature value for the threshold in the one-dimensional BS model, while the former is favored by
the results discussed in the following subsection. We studied systems of 1000, 2000, 4000,...,32000 sites in simulations of $2 \times 10^7$ to $3 \times 10^8$ time steps.

For $L = 4000 - 32000$, the results for the order parameter follow a power law with $\beta/\nu_\perp = 0.755(5)$. The data for smaller system sizes, however, show systematic deviations from a pure power law, leading us to seek a correction to scaling term; a correction decaying $\propto L^{-1/2}$ leads to a good fit. We fit the expression

$$\ln \rho = -\frac{\beta}{\nu_\perp} \ln L - \frac{b}{L^{1/2}}$$

(3)

to the data for $L \geq 1000$, allowing $\beta/\nu_\perp$ and $b$ as adjustable parameters; the best-fit values are $\beta/\nu_\perp = 0.769(7)$ and $b = 3.69(20)$. The simulation data, and the difference from best fit of Eq. (3) are plotted in Fig. 1, showing the high quality of fit. The data for the lifetime, using system sizes of 125, 250,...8000 yield $\nu_\parallel/\nu_\perp = 2.12(1)$, with no obvious correction term (see Fig. 2). We also determined the stationary moment ratio $m = \rho^2/\rho^2$ at the critical point. The estimates for $m$ decrease slowly with $L$, and appear, when plotted versus $L^{-0.25}$, to approach a limiting ($L \to \infty$) value of 1.030(5) (see Fig. 3). (For the one-dimensional CP, $m = 1.1737$ at the critical point.) Essentially the same results are obtained regardless of whether we use $q_{c,E} = 0.66701$ or 0.66702 in these simulations.

C. Absorbing Bak-Sneppen Model: Spread of activity

Scaling properties at an absorbing state phase transition are also reflected in the spread of activity from an initially localized region [30]. In spreading simulations of the ABS model we start the system with a single site $x_0 < q$ and all others above this value. (This is completely equivalent to the BS branching process studied in [4,5,7].) At $q = q_{c,E}$, the process generates a scale-invariant pattern of activity that may be characterized by power-laws for the survival probability $P_s(t)$, the mean number of active sites $n(t)$ and the mean-square distance $R^2(t) = [n(t)]^{-1}(\sum_j r_j^2(t))$. ($r_j(t)$ denotes the position of the $j$–th active site at time $t$. Note that $n(t)$ is taken over all realizations, including those that have become trapped in the absorbing state at or before time $t$.) The scaling laws are typically written in the form

$$P_s \sim t^{-\delta}, \quad n \sim t^{\eta}, \quad R^2 \sim t^{z_{sp}},$$

(4)

relations that have been verified to high precision for various examples [21,23]. (We use $z_{sp}$ to denote the spreading exponent, to avoid confusion with the dynamical exponent $z$.) The appearance of power laws is commonly used to locate the critical point [23].

The spreading exponent $\delta$ is related to the avalanche size exponent $\tau$, defined (in the BS model) via $P_D(s) \sim s^{-\tau}$, where $P_D(s)ds$ is the probability of an avalanche having a duration between $s$ and $s + ds$. Thus the survival probability $P_s(t) = \int_t^\infty P_D(s)ds$, implying $\tau = 1 + \delta$. 

9
We performed spreading simulations of the ABS model at \( q = 0.66699, 0.66700, 0.66701, 0.66702 \) and \( 0.66703 \). Each realization was followed up to a maximum time of about \( 2.7 \times 10^5 \); the total number of realizations ranged from \( 4 \times 10^5 \) to \( 1.6 \times 10^6 \), depending on the value of \( q \). To locate the critical point we study the local slope \( \delta(t) = d \ln P/d \ln t \), plotted versus \( t^{-1} \). For \( q < q_c \), the local slope is expected to veer downward at large times, and vice-versa. (Numerically, \( \delta(t) \) is given by the slope of a least-squares linear fit to the data in an interval \([t_0, 20t_0]\), with geometric mean \( t \).) On the basis of the local slope data (see Fig. 4) we conclude that \( q_c = 0.66701(1) \). This is consistent with previous estimates, which place the threshold at \( 0.66702(8) \) \[7\]. (We did not find analyses of the local slopes \( \eta(t) \) or \( z_{sp}(t) \), defined analogously to \( \delta(t) \), useful in locating the critical point.)

Analyzing the data at the critical point, we are unable to obtain good fits to \( P_s, n \) and \( R^2 \) using simple power-law expressions. Including a subdominant power-law correction in the relations of Eq. (4) greatly improves the quality of fit. In particular, the survival probability can be fit quite accurately using

\[
\ln P_s \simeq -\delta \ln t + \phi_pt^{-1/4} + C
\]  

where \( C \) is a constant and the best-fit values are \( \delta = 0.084(1) \) and \( \phi_p = 0.115 \). The same value for \( \delta \) is found using the data for \( q = 0.66702 \). (The choice of a correction term decaying as \( t^{-1/4} \) is motivated by the fact that the local slopes \( \delta(t) \) and \( z_{sp}(t) \) are essentially linear when plotted versus \( t^{-1/4} \), as seen in the inset of Fig. 4.) In Fig. 5 we plot \( P_s \) and the ratio of \( P_s \) to the fitting function, Eq. (5); the ratio is seen to be essentially constant for \( t > 50 \). The mean-square displacement may also be fit using an asymptotic power law and correction term. We find

\[
\ln R^2 \simeq z_{sp} \ln t - \phi_Rt^{-1/4} + C'
\]

with \( z_{sp} = 0.921(10) \) and \( \phi_R = 1.703 \).

It has been argued that \( \eta = 0 \) quite generally for extremal dynamics \[5,27\]. Our data for the one-dimensional ABS model support this conclusion: on a double-logarithmic plot, \( n(t) \) clearly grows more slowly than a power law. While \( \eta = 0 \) is compatible with \( n(t) \) growing without limit as \( t \to \infty \) (for example, \( \propto (\ln t)^{\phi} \), as suggested in Ref. \[7\]), our results support the conclusion that \( n(t) \) saturates at a finite value \( n_\infty \) at long times. Specifically, we are unable to fit the long-time behavior using an expression of the form \( n \sim (\ln t)^{\phi} \). On the other hand, we find \( d \ln n/d \ln t \propto t^{-\omega} \), with \( \omega \simeq 0.149 \), suggesting that \( n(t) \approx n_\infty \exp[-ct^{-\omega}] \). In fact an excellent fit is obtained using \( c = 1.92 \) and \( n_\infty = 14.574 \), as can be seen in Fig. 6. (Saturation of \( n(t) \) does not occur on the time scale of the simulation; for the anistropic case, shown in the inset of Fig. 6, saturation is in fact evident.)

In the absorbing phase (\( q < q_{c,E} \)), the survival probability must vanish as \( t \to \infty \). Our data follow \( P_s \sim t^{-\delta}e^{-t/\tau} \), where \( \tau \sim |q - q_{c,E}|^{-\nu} \), with \( \nu = 2.54(2) \). On the other hand, for \( q > q_{c,E} \), the survival probability tends to a finite value as \( t \to \infty \). We obtain \( \lim_{t \to \infty} P_s \equiv P_\infty \sim (q - q_{c,E})^{\beta} \), with
\( \beta' = 0.20(1) \). (In DP and allied models \( \beta' = \beta \) [30], but this need not hold for models in other universality classes.)

In the CP and other nonextremal models, the spread of activity in the supercritical regime follows a simple pattern: the size \( R \) of the active region grows linearly with time, and the number of active sites \( n \) grows \( \propto t^d \). Our observation of subdiffusive spreading (and saturation in \( n \)) at the critical point lead us to investigate supercritical spreading in the ABS model. We find that spreading is indeed sublinear. For example, using \( q = 0.7 \) in a study extending to \( t \approx 9 \times 10^6 \) to avoid transient effects, we obtain \( R^2 \sim t^\chi \) with \( \chi = 1.20(4) \) and \( n \sim t^\lambda \) with \( \lambda = 0.61(2) \). (The exponent governing \( R^2 \) should be twice that for \( n \), since active regions have a finite activity density in the supercritical regime.) Once again, extremal dynamics is seen to limit the growth of activity.

### D. CP3 Model

We performed spreading simulations of the CP3 model, using the approach described in the previous subsection. Each realization is followed up to a maximum time of \( 6 \times 10^4 \). Using the power-law behavior of \( P_s(t) \) and \( n(t) \) as the criterion for criticality, we find \( q_c = 0.63525(3) \) for the CP3. (Note that this is some 5\% smaller than the critical value of the corresponding extremal model.) Analyzing the local slopes, we obtain \( \delta = 0.162(2), \eta = 0.312(2) \) and \( z_{sp} = 1.265(4) \). These values are fully consistent with those for directed percolation (see Table I), confirming that the CP3 model belongs to the same universality class as the original contact process.

A striking difference between extremal and nonextremal models with an absorbing state is that the spread of activity in the critical process is much slower in the former. This is of course reflected in the value \( \eta = 0 \) for extremal models, (while for example \( \eta = 0.314 \) for DP in one dimension), and in the subdiffusive growth in \( R^2 \) in the ABS model. In Figs. 7 and 8 we compare typical evolutions in the ABS model and its nonextremal analog, the CP3, at their respective critical points. It is evident that the activity spreads much more slowly in the ABS than in the CP3. A further notable difference is that in the ABS a site can remain active for a very long time, i.e., while the site is not the minimum site or a neighbor of it. Thus the rates of both addition and loss of active sites are much smaller in the critical extremal process than in the corresponding nonextremal one.

### E. Extremal CP

In light of the discussion of Sec. II, it is of interest to study the behavior of other absorbing-state models under extremal dynamics. As a first step we report simulation results for the extremal-absorbing contact process (CP\(_{EA} \)). We performed spreading simulations to determine \( q_E \) and the exponents \( \delta, \eta \) and \( z_{sp} \), using simulations running to a maximum time of \( 6 \times 10^4 \) in \( 5 \times 10^5 \)
independent realizations. We find \( q_{c,E} = 0.79415(5) \) for the extremal CP, compared with 0.76733 for the original (non-extremal) process. (Note that, as in the comparison between the CP3 and ABS models, \( q_{c,E} > q_c \). This may reflect the slower spread of activity under extremal dynamics.)

As in the case of the ABS model, the decay of the survival probability at the critical point follows an expression of the form of Eq. (5), here with best-fit parameters \( \delta = 0.0855 \) and \( \phi = 0.226 \). The exponent \( \delta \) is essentially the same as found for the ABS model, while the correction term is about twice as large. At the critical point the derivative \( d \ln n/d \ln t \sim t^{-0.1} \), again indicating a behavior of the form \( n(t) \simeq n_{\infty} \exp[-ct^{-\omega}] \), here with \( \omega = 0.1 \). The mean-square distance of active sites from the origin grows in a manner similar to that in the ABS model. We are again able to fit the data for \( R^2 \) using an expression of the form of Eq. (6), with \( z_{sp} = 0.932 \) and \( \phi_{R^2} = 2.026 \). These results strongly suggest that the CP\(_{EA}\) belongs to the same universality class as the ABS model.

We turn now to the rather surprising behavior of the stationary probability density \( p(x) \) in the extremal CP. Recall that mean-field theory predicts \( p(x) = 2\Theta(x - 1/2) \) for \( q < q_{c,E} = 1/2 \), while for \( q > 1/2 \) there are two steps, one at \( x = \kappa = q^2/(3q - 1) \), the other at \( x = q \). In simulations of the CP\(_E\) on a ring we find a single step discontinuity for \( q < q_{c,E} = 0.79415 \), and, for \( q > q_{c,E} \), a pair of steps, one at \( x = q \), the other at \( x = q_s < q \). The positions of the singularities as obtained in simulation (using data for system sizes \( L = 100, 200, \ldots, 1600 \) to extrapolate the position in the infinite-size limit), are shown in Fig. 9. The lower singularity \( q_s \) is seen to bifurcate from \( q = q_{c,E} \) just at the critical point, in qualitative agreement with MFT. Note however that the position of the singularity is not constant for \( q < q_{c,E} \), as predicted by MFT.

The density \( p(x) \) is shown for \( q = 0.794 \simeq q_{c,E} \) and \( q = 0.85 \) in Fig. 10. In the latter case it is evident that the step at \( x = q \) is sharp (this is true even for small systems) and derives from the singular nature of the updating rule. The step at \( x = q_s \), by contrast, is subject to finite-size rounding, and becomes sharper with increasing system size, as is characteristic of a critical singularity. The finite width of the peak at \( x = q_c \) (in the process with \( q = q_{c,E} \)), appears to be a finite size effect as well: it becomes sharper with increasing \( L \), suggesting that the singularities merge in the limit \( L \to \infty \).

F. Anisotropic ABS Model

The scaling behavior of the Bak-Sneppen model changes when the updating rule is asymmetric [31]. The same critical exponents are found for a highly anisotropic version in which at each step, only the minimal site and its neighbor on the right are updated [32], and for weak anisotropy [16,33], so that one may identify an anisotropic BS universality class. In this section we report results of spreading simulations of the anisotropic absorbing BS model. To obtain these results, we simulated the anisotropic ABS (in the highly anisotropic version) in studies extending to a maximum time of \( 1.6 \times 10^5 \), using \( 3 \times 10^5 \) realizations.
Analysing the local slope $\delta(t) = d\ln P/d\ln t$, we determined the threshold of this model to be $q_{c,E} = 0.72370(2)$. (The best previous estimate is $0.7240(1)$ [34].) A typical evolution of the critical spreading process is shown in Fig. 11. The local slope $\delta(t)$ yields the estimate $\delta = 0.234(5)$. For anisotropic models we define $R^2(t)$ as the mean-square radius of gyration, i.e., the distance is measured relative to the current center of mass of the set of active sites, rather than to a fixed origin. This is done to eliminate a spurious contribution due simply to the overall drift in the active region. For the anisotropic ABS model $R^2(t)$ may again be fit with an expression of the form of Eq. (6), with $z_{sp} = 1.425(10)$ and $\phi_R = 2.3(2)$. The exponents $\delta$ and $z_{sp}$ are quite different from those of the isotropic model. Despite these differences, we again find $\eta = 0$ for the anisotropic model. As before, the mean number of active sites $n(t)$ saturates at long times, more rapidly in fact than in the isotropic ABS model (see Fig. 6, inset). We are able to fit the data well using $n(t) = n_{\infty}[1 - e^{-c(t/\varepsilon)^{\phi}}]$ with parameters $n_{\infty} = 5.206(3)$ and $c = 0.348$.

The nonextremal model corresponding to the anisotropic ABS model is a two-site contact process, CP2, which is simply the CP3 with updating restricted to the central site and its neighbor on the right. We have verified that the spreading exponents of the CP2 model are those of directed percolation. (Here again, we define $R^2$ as the mean-square radius of gyration.) This leads to the interesting conclusion that a perturbation (asymmetric updating) that is irrelevant for a nonextremal model is relevant for the corresponding extremal system. (We note that, because the two sites in the CP2 are updated in the same manner, the model does not fall in the so-called anisotropic-DP class, for which bonds along different axes are present with different probabilities [35].)

**IV. CONCLUSIONS**

We investigate the relation between extremal dynamics, exemplified by the Bak-Sneppen model, and nonextremal models exhibiting a phase transition to an absorbing state, using general arguments, mean-field theory and simulation. The relation between the BS model and directed percolation was already suggested some time ago [4,5]. Here we clarify this connection by showing how a generic absorbing-state model can be transformed to an extremal one via the associated extremal-absorbing model. The nonextremal ‘precursor’ of the BS model is a three-site contact process [24], CP3, which, like the original CP, belongs to the directed percolation universality class. The BS model and the extremal version of the CP belong to a common universality class that may be dubbed ‘extremal DP’ (EDP). A number of extremal dynamics classes distinct from EDP are discussed in [5]; another example is the anisotropic BS model. We expect that further extremal dynamics universality classes exist, for example an extremal parity-conserving class [36], although examples of the latter have yet to be studied.

Our results for the critical exponents of the EDP class, which includes the BS model and the extremal CP, are compared against those of ordinary DP (in
one spatial dimension) in Table I. (Here we have taken the values $\eta = 0$ and $\beta = 1$ to be exact for EDP.) The differences between the two sets of exponent values are evident. Our results $\tau = 1 + \delta = 1.084(1)$, and $z = \nu_\parallel/\nu_\perp = 2.12(1)$ are in agreement with the earlier estimates [5] of 1.07(1) and 2.10(5), respectively. Our result is however somewhat higher than Grassberger’s result $\tau = 1.073(3)$ [7].

Certain scaling relations are expected to hold among the critical exponents [21,23,30]. In spreading processes one expects $z_{sp} = 2\nu_\perp/\nu_\parallel$; our data are nearly consistent with this, yielding $2\nu_\perp/\nu_\parallel - z_{sp} = 0.022(14)$. The relation $\beta' = \delta\nu_\parallel$ is also satisfied: our data yield $\beta' - \delta\nu_\parallel = -0.013(14)$. Finally, we consider the generalized hyperscaling relation [37]

$$2 \left(1 + \frac{\beta}{\beta'}\right)\delta + 2\eta = dz_{sp},$$  \hspace{1cm} (7)

in $d$ dimensions. Using our data, we find the difference between the two sides of this relation to be 0.09(6). Our results are marred by another inconsistency that may reflect corrections to scaling or finite size effects: the product $\left(\beta/\nu_\perp\right)^{-1}(\nu_\parallel/\nu_\perp)\nu^{-1}$, with the first two factors determined from finite-size scaling at the critical point, and the final factor obtained from the decay of the survival probability in the subcritical regime, should equal $\beta = 1$; our data yield 1.08(3). These minor inconsistencies suggest that one or more of the exponents may be in error by 5% or so. Refining their values will require accumulating larger data sets in simulations of larger systems, a task we leave for future work. (The studies reported here were quite demanding computationally, representing approximately 6 months’ cpu time on an alpha workstation.)

In the course of our study we revisit a three-site contact process (CP3) that is the nonextremal analog of the BS model [24]. We verify that the CP3 belongs to the universality class of directed percolation, as expected [7]. Similarly, we define extremal and extremal-absorbing versions of the original contact process (CP$_E$ and CP$_{EA}$, respectively) and verify that their scaling properties are the same as those of the BS model. The stationary probability density for the CP$_E$ follows, in general terms, the predictions of mean-field theory, but certain interesting differences exist, as detailed in Sec. III.E.

It is clear that when an absorbing-state model is modified to follow extremal dynamics, its critical exponents are altered. In general extremal dynamics tends to slow the spread of activity in the critical and supercritical regimes. One may nevertheless inquire whether any more general features of the original model are preserved under ‘extremalization’. A candidate for such a conserved property is the critical dimension $d_c$. In critical phenomena, various universality classes (differing in the symmetry group of the order parameter, or the presence of conserved quantities) may share the same $d_c$ if the algebraic structure of their continuum description (in particular, the power of the lowest-order nonlinear term in the order parameter, in a Landau-Ginzburg-Wilson effective Hamiltonian) is the same. Thus $d_c = 4$ for all models in the $n$-vector family. Extending this idea to extremal models is questionable, since there is no continuum description at hand. (At first glance, the notion of extremal dynamics
in a description using a continuous activity density seems problematic, since there is always one and only one extremal site.) Be that as it may, it seems plausible that if the field theory for DP \([18,19,38]\) were somehow modified to reflect extremal dynamics, the dominant nonlinearity would not change, so that \(d_c\) would retain its value of four, as in DP. The upper critical dimension \(d_c = 4\) for the BS model was established some time ago by Boettcher and Paczuski \([39]\). Our argument suggests that extremal versions of other absorbing-state models have the same upper critical dimension as the corresponding nonextremal model. We hope to test this prediction in future work.

Studying the anisotropic ABS model and its nonextremal counterpart, the CP2 model, we find that anisotropy is a relevant perturbation for extremal DP, while it is irrelevant for the corresponding nonextremal class. We suspect that other perturbations, such as diffusion, may exhibit a similar pattern of relevance.

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APPENDIX: MEAN-FIELD THEORY

A. Extremal dynamics as a zero-temperature limit

There are several ways of formulating a mean-field theory (MFT) for extremal models. First we consider an approach [11,33] in which the probability of a site \( i \) being chosen as the central site is proportional to \( e^{-\beta x_i} \); extremal dynamics is recovered in the limit \( \beta \to \infty \). (In the present discussion the parameter \( \beta \) bears no relation to the critical exponent denoted by the same symbol in the main text.) Applied to the BS model, this approach yields the distribution \( p(x) = \left(\frac{3}{2}\right) \Theta(x - 1/3) \Theta(1 - x) \) when \( \beta \to \infty \).[11,33]

In the ABS model the distribution \( p(x) \) evolves via

\[
\frac{\partial p(x, t)}{\partial t} = -e^{-\beta x} p(x) \Theta(q - x) + 3 \int_0^q e^{-\beta y} p(y, t) dy - 2p(x, t) \int_0^q e^{-\beta y} p(y, t) dy
\]

(8)

The first term represents a site with value \( x \) being selected as the central site, which is only possible if \( x < q \). The second term reflects updating three sites with new variables uniform on \([0, 1]\), with the integral representing the overall rate of events. The final term represents updating of the two neighboring sites, and is obtained using the mean-field factorization of the nearest-neighbor joint probability density: \( p(x, y, t) \simeq p(x, t)p(y, t) \). Note that in writing Eq. (8) we have associated a time increment \( dt = 1/N \), with \( N \) the number of sites, with each event.

Eq. (8) admits an infinite set of stationary solutions for which \( p(x) = 0 \) on the interval \( 0 \leq x \leq q \). These represent absorbing states. To seek an active stationary solution we let \( I = \int_0^q e^{-\beta y} p(y, t) dy \), and equate the time derivative to zero, yielding

\[
\overline{p}(x) = \frac{3I}{2I + \Theta(q - x)e^{-\beta x}}
\]

(9)

To determine \( I \) we multiply Eq. (9) by \( e^{-\beta x} \) and integrate from \( x = 0 \) to \( x = q \), leading to \( I = (e^{-\beta/3} - e^{-\beta q})/[2(1 - e^{-\beta/3})] \), so that

\[
\overline{p}(x) = \frac{\frac{3}{2}(e^{-\beta/3} - e^{-\beta q})}{(e^{-\beta/3} - e^{-\beta q}) + \Theta(q - x)e^{-\beta x}(1 - e^{-\beta/3})}
\]

(10)

In the limit \( \beta \to \infty \), we find, for \( q > 1/3 \), the singular density \( \overline{p}(x) = (3/2)\Theta(x - 1/3) \). This is precisely the MF result for the original BS model. [When we take \( \beta \to \infty \), the above expression reduces to \( \overline{p}(x) = (3/2)\Theta(x - q) \) for \( q < 1/3 \). But this density is not normalized on \([0, 1]\) and so must be rejected. We are left with only absorbing stationary solutions for \( q < 1/3 \).] Thus \( q_{c,E} = 1/3 \) in the MFT of the absorbing Bak-Sneppen model. Note that the parameter \( q \) is irrelevant for \( q > q_{c,E} = 1/3 \), as was argued in Sec. II.
A moment’s reflection shows that the evolution of \( p(x) \) in the (nonextremal) CP3 model is given by Eq. (8) with \( \beta = 0 \), since all active sites are then equally likely to be chosen as the central site. Taking the limit \( \beta \to 0 \) of the stationary solution, Eq. (10), one finds, for \( q \geq 1/3 \), the stationary density

\[
\overline{p}(x) = \begin{cases} 
\frac{1}{2}(3 - q^{-1}), & x < q \\
\frac{3}{2}, & q < x \leq 1 
\end{cases}
\] (11)

Eq. (11) confirms that the stationary density of a (nonextremal) model exhibiting an absorbing state phase transition is characterized by a steplike singularity, as asserted in Sec. II. For \( q < 1/3 \), Eq. (11) yields an unphysical, negative density, showing that \( q_c = 1/3 \) for the CP3, in the mean-field approximation.

The foregoing analysis is readily extended to the extremal-absorbing contact process (CP_{EA}) defined in Sec. II. The rate of events is again given by \( I = \int_0^q e^{-\beta x} p(x) dx \). At each event, there is a probability of 1/2 that the central site (which must have \( x < q \)) is replaced, while with probability 1/2 a neighbor is updated. Thus the loss terms in the equation for \( p(x, t) \) are: \( -(1/2)[e^{-\beta x} \Theta(q - x) + I]p(x) \). The gain term corresponding to updating of the central site is simply \( I/2 \), but for updating a neighbor it is \( (I/2)[1 - P(q) + (1/q)\Theta(q - x)P(q)] \), where \( P(x) = \int_0^x p(y) dy \) is the probability that a given site \( i \) has \( x_i < x \). (The reason is that when updating an active neighbor, the new variable is chosen from the distribution uniform on \([0, q]\)). Thus the MF equation of motion is

\[
\frac{\partial p}{\partial t} = -\frac{1}{2}e^{-\beta x}p(x)\Theta(q - x) + \frac{I}{2} \left[ 2 - P(q) + \frac{\Theta(q - x)}{q} P(q) - p(x) \right] \] (12)

To find the stationary solution \( \overline{p}(x) \) we first note that for \( 0 \leq x < q \), setting \( \partial p/\partial t \) to zero yields

\[
\overline{p}(x) = \frac{AI}{(I + e^{-\beta x})} \] (13)

where \( A = 2 + (q^{-1} - 1)P(q) \). Integrating Eq. (13) from \( x = 0 \) to \( x = q \), we find

\[
P(q) = Aq - \frac{A}{\beta} \int_{e^{-\beta q}}^1 \frac{du}{I + u} \] (14)

If we now multiply Eq. (13) by \( e^{-\beta x} \) and integrate over the same interval, we obtain

\[
\frac{A}{\beta} \int_{e^{-\beta q}}^1 \frac{du}{I + u} = 1 \] (15)

leading to \( P(q) = 2 - q^{-1} \) in the stationary state. We see that \( q_{c,E} = 1/2 \) as in the MFT of the original contact process. Evaluating the integral in Eq. (15)
one finds \( I = (e^{-\beta \kappa} - e^{-\beta q})/(1 - e^{-\beta \kappa}) \), where \( \kappa = q^2/(3q - 1) \). The stationary density is

\[
\overline{p}(x) = \frac{1}{q} \frac{(e^{-\beta \kappa} - e^{-\beta q})[1 + (2 - q^{-1})\Theta(q - x)]}{(e^{-\beta \kappa} - e^{-\beta q}) + e^{-\beta x}(1 - e^{-\beta \kappa})\Theta(q - x)}
\]  

(16)

For \( q > 1/2 \), we have \( \kappa < q \), and in the limit \( \beta \to \infty \),

\[
\overline{p}(x) = \begin{cases} 
0, & x < \kappa \\
\frac{1}{\kappa}, & \kappa < x < q \\
\frac{1}{q}, & x > q
\end{cases}
\]  

(17)

which is normalized and exhibits step-function singularities at \( x = \kappa \) and \( x = q \). For \( q < 1/2 \) on the other hand, \( \kappa > q \) and Eq.(16) does not yield an acceptable probability density, and we conclude that the only stationary state is the absorbing one. The critical point of the \( CP_{EA} \) thus falls at \( q_{c,E} = 1/2 \) in MF approximation.

Taking \( \beta \to 0 \) in Eq. (16), we obtain the probability density for the original CP:

\[
\overline{p}(x) = \begin{cases} 
\frac{2q - 1}{q^2}, & x < q \\
\frac{1}{q}, & x > q
\end{cases}
\]  

(18)

Finally, for the extremal CP, the equation of motion is

\[
\frac{\partial p}{\partial t} = -\frac{1}{2} e^{-\beta x} p(x) + \frac{I}{2} \left[ 2 - P(q) + \frac{\Theta(q - x)}{q} P(q) - p(x) \right]
\]  

(19)

with \( I = \int_0^1 e^{-\beta x} p(x) dx \). To find the stationary solution we write

\[
\overline{p}(x) = \frac{2 + [q^{-1}\Theta(q - x) - 1]P(q)}{1 + I^{-1}e^{-\beta x}}
\]  

(20)

Integrating from 0 to \( q \) and solving for \( P(q) \) we find

\[
P(q) = \frac{2(q - \gamma)}{q + \gamma(q^{-1} - 1)}
\]  

(21)

where

\[
\gamma = \frac{1}{\beta} \ln \frac{I + 1}{I + e^{-\beta q}}
\]  

(22)

Now multiply Eq. (20) by \( e^{-\beta x} \) and integrate from 0 to 1 to obtain

\[
1 = A \int_0^q \frac{e^{-\beta x}dx}{I + e^{-\beta x}} + A' \int_q^1 \frac{e^{-\beta x}dx}{I + e^{-\beta x}}
\]  

(23)
where $A = 2/[q + \gamma(q^{-1} - 1)]$ and $A' = \gamma A/q$. If $P(q) > 0$, the first term on the r.h.s. of Eq. (23) is nonzero and dominates as $\beta \to \infty$. Equating the first term to unity then leads to $\gamma = q^2/(3q - 1) = \kappa$, and thence to $P(q) = 2 - q^{-1}$ which is positive for $q > 1/2$. A simple calculation then yields the distribution of Eq. (17) in the limit $\beta \to \infty$.

If $q < 1/2$ the above solution is not valid since it implies $P(q) < 0$. We therefore take $P(q) = 0$, implying $\gamma = q$, and so $A = A' = 2$. Eq. (23) now reads

$$1 = 2q + 2\beta \ln \frac{I + e^{-\beta q}}{I + e^{-\beta}}$$

Solving for $I$ and inserting the result in Eq. (20), we find in this case

$$\lim_{\beta \to \infty} \bar{p}(x) = 2\Theta(x - 1/2).$$

These results have been verified via numerical integration.

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**B. Extremal dynamics on a complete graph**

Another approach to formulating MFT for the BS model considers extremal dynamics on an $N$-site complete graph or random-neighbor model (two neighbors are selected at random each time a site is updated); the stationary density $\bar{p}(x)$ becomes a step function in the infinite-size limit [12,16,40,41]. We now extend this approach to the ABS model. Let $P(x) = \text{Prob}[x_i < x] = \int_0^x p(y)dy$ be the distribution function and let $Q(x) = 1 - P(x)$. By definition $Q$ is a non-increasing function with $Q(0) = 1$ and $Q(1) = 0$, since $p(x) = 0$ outside the interval $[0,1]$.

Activity in the ABS model is predicated on the minimal site $x_{min}$ being smaller than $q$; the probability of this event, under the MF factorization, is $1 - [Q(q)]^N$. Given $x_{min} < q$, updating the extremal site and two neighbors results, on the average, in the increment: $dP(x) = (1/N)[-1 - Q(x_{<})^N] - 2P(x) + 3x$, where $x_{<} \equiv \min\{x, q\}$, so that the first term represents loss of the minimal site, the second removal of two neighbors, and the third random replacement of the three site variables with numbers uniform on $[0,1]$. If we adopt a time increment $dt = 1/N$ for each such event, the equation of motion for $P$ is

$$\frac{\partial P(x,t)}{\partial t} = -[1 - Q(x_{<},t)^N] + [1 - Q(q,t)^N][3x - 2P(x,t)]$$

(25)

Note that the evolution ceases if $Q(q,t) = 1$, i.e., if there are no active sites. (Since $Q$ is nonincreasing $Q(q) = 1 \Rightarrow Q(x_{<}) = 1$.)

For $q > 1/3$, the stationary solution to Eq. (25) corresponds to a density $\bar{\eta}(x)$ that approaches a step function, $(3/2)\Theta(x - 1/3)$, as $N \to \infty$. A simple calculation yields the dominant contribution for large $N$:

$$\bar{\eta} \simeq \begin{cases} 
(1 - 3x)^{1/N}, & x < \frac{1}{3} \\
\frac{2}{3}(1 - x) + O(e^{-\text{const.}N}) & x > \frac{1}{3}
\end{cases}$$

(26)
(One should note however that the convergence is nonuniform in $x$, being slower the closer $x$ is to the critical value of $1/3$.) For $q < 1/3$ we are unable to find an acceptable stationary solution with $Q < 1$ (i.e., $\bar{p} > 0$), for $x < q$, and conclude that only absorbing solutions exist.

The analysis of the ABS model on a complete graph confirms that in the infinite-size limit, the model enjoys the usual properties of the BS model for $q > q_{c,E} = 1/3$, and falls into the absorbing state for $q < 1/3$.

The evolution of $P(x, t)$ in the extremal CP follows, in MF approximation, the equation

$$\frac{\partial P(x, t)}{\partial t} = -\frac{1}{2}[1 - Q(x, t)^N] + \frac{x}{2} - \frac{1}{2}P(x, t) + \frac{1}{2}[x^*P(q, t) + xQ(p, t)] \quad (27)$$

where $x^* = \min\{x/q, 1\}$. Numerical integration shows that the solution converges, for large $N$, to a stationary distribution consistent with the singular density found above in the limit $\beta \to \infty$. 

22
TABLES

TABLE I. Critical exponents for the one-dimensional absorbing Bak-Sneppen model (ABS) and contact process (CP). CP exponents from Refs. [23] and [41].

| Exponent   | ABS       | CP         |
|------------|-----------|------------|
| $\beta$    | 1         | 0.27649(4) |
| $\beta'$   | 0.20(1)   | ( = $\beta$) |
| $\nu_\parallel$ | 2.54(2)   | 1.73383(3) |
| $\beta/\nu_\perp$ | 0.77(1)   | 0.25208(5) |
| $\nu_\parallel/\nu_\perp$ | 2.12(1)   | 1.58071(11) |
| $\delta$   | 0.084(1)  | 0.15947(3) |
| $\eta$     | 0         | 0.31368(4) |
| $z_{sp}$   | 0.921(10) | 1.26523(3) |

TABLE II. Spreading exponents for the CP3 and CP$_{EA}$ models and the anisotropic absorbing Bak-Sneppen (AABS) model in one dimension.

| Exponent   | CP3       | CP$_{EA}$  | AABS       |
|------------|-----------|------------|------------|
| $\delta$   | 0.162(2)  | 0.0855(20) | 0.234(5)   |
| $\eta$     | 0.312(2)  | 0          | 0          |
| $z_{sp}$   | 1.265(4)  | 0.932(20)  | 1.425(10)  |
FIGURE CAPTIONS

FIG. 1. Stationary activity density (filled symbols) $\bar{\rho}$ versus system size $L$ in the one-dimensional ABS model at the critical point. Open symbols: difference between $\bar{\rho}$ and the fitting function, Eq. (3), shifted vertically for visibility.

FIG. 2. Mean lifetime $\tau$ (filled symbols) versus system size $L$ in the one-dimensional ABS model at the critical point. Open symbols: $\tau/L^{\nu_\parallel/\nu_\perp}$ (shifted vertically for visibility).

FIG. 3. Moment ratio $m$ for the ABS model versus system size $L^{-0.25}$. Points: simulation data; line: best linear fit, $m = 1.0295 + 0.268L^{-0.25}$.

FIG. 4. Local slope $\delta(t)$ versus $1/t$ in the ABS model. $q$ values (bottom to top) 0.66700, 0.66701, 0.66702 and 0.66703. Inset: data for $q = 0.66701$ plotted versus $1/t^{0.25}$.

FIG. 5. Survival probability $P_s(t)$ in the ABS model at the critical point, $q = 0.66701$. The nearly constant function represents the ratio of $P_s$ to the fitting function, Eq. (5).

FIG. 6. Mean number of active sites $n(t)$ in the ABS model at the critical point, $q = 0.66701$. The solid curve represents the fitting function described in the text. Inset: a similar plot, for the critical anisotropic ABS model.

FIG. 7. Spread of activity in a typical realization of the critical CP3 model ($q = 0.63525$).

FIG. 8. Spread of activity in a typical realization of the critical ABS model ($q = 0.66701$).

FIG. 9. Position $q_s$ of the singularity in the stationary probability density of the extremal CP. The upper line of singularities, $x = q$, bifurcates from $q_s$ at the critical value $q_{c,E}$.

FIG. 10. Stationary probability density $\overline{\rho}(x)$ in the CP$_E$ for $q = 0.794$ (left) and $q = 0.85$ (right); system size $L = 1600$.

FIG. 11. Spread of activity in a typical realization of the critical anisotropic ABS model ($q = 0.72370$).
