Exact Solution of the One-Dimensional Deterministic Fixed-Energy Sandpile

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> By reason of the strongly nonergodic dynamical behavior, universality properties of deterministic fixedenergy sandpiles are still an open and debated issue. We investigate the one-dimensional model, whose microscopical dynamics can be solved exactly, and provide a deeper understanding of the origin of the nonergodicity. By means of exact arguments, we prove the occurrence of orbits of well-defined periods and their dependence on the conserved energy density. Further statistical estimates of the size of the attraction's basins of the different periodic orbits lead to a complete characterization of the activity vs energy density phase diagram in the limit of large system's size.

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Sandpile models have been introduced by Bak, Tang, and Wiesenfeld (BTW) as simple nonequilibrium systems exhibiting self-organized criticality (SOC), providing a theoretical framework for the emergence of power laws in a large number of natural phenomena [1]. More recently, the occurrence of SOC has been related to the presence of underlying absorbing state phase transitions (APTs) governing the critical behavior of the system [2]. The relationship between sandpiles and nonequilibrium APTs is emphasized by a class of models with locally conserved dynamics, the fixed-energy sandpiles (FESs) [3]. Finite size scaling techniques led to the complete characterization of the universality class for driven-dissipative sandpile models with stochastic dynamics [4], but failed in determining the critical properties of the deterministic BTW model, that have been partially recovered by means of multiscaling analysis [5]. Stochastic FESs seem to belong to a different universality class with respect to directed percolation [3,6-8], while the deterministic BTW dynamics shows very strong nonergodic effects, that cannot be understood using a purely statistical mechanics approach [6,9]. In dissipative systems, the nonergodicity of the BTW rule is related to the group structure of the configuration space [10], and emerges when the process of grain addition and dissipation is not completely random [11]. Without dissipation, the group structure breaks down, but the nonergodic behavior is recovered as the result of a selfsustained activity that eventually enters a periodic orbit.

In this Letter, we report the exact solution of the onedimensional deterministic fixed-energy sandpile (1D-DFES), providing an exhaustive description of the microscopic dynamics and its steady states, and determining the system's macroscopic behavior in the limit of large size. Moreover, our explanation of the microscopic origin for nonergodicity in BTW FESs elucidates the relation, already suggested in Ref. [9], with mode-locking phenomena in nonlinear automata [12].

Let us consider a ring of N sites, each one endowed with an *energy* z_i , assuming non-negative integer values. Whenever the energy of a site equals or exceeds the threshold value $z^{\text{th}} = 2$, the site becomes *active* and undergoes a toppling event; i.e., it loses 2 units of energy, that are equally redistributed among its 2 neighbors. On the contrary, if $z_i < 2$, the site *i* is *stable*. The system evolves with parallel dynamics, therefore the update rule can be written as follows:

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$$z_i(t+1) = z_i(t) - 2\theta(z_i(t) - 2) + \sum_{j \in V_i} \theta(z_j(t) - 2) \quad (1)$$

for i = 1, 2, ..., N, where $\theta(x) = 1$ (0) if $x \ge 0$ (x < 0) and V_i is the neighborhood of the site *i*. When the total energy $E = \sum_{i} z_{i}$ is sufficiently large, periodic boundary conditions prevent the system from relaxing in the absorbing state $(z_i < 2 \forall i)$. Since the configuration space is finite and the dynamics is deterministic, after a transient, the system enters a periodic orbit (steady active state). We will conventionally assume sites to have energy between 0 and 3 (relaxing this condition does not modify the physical properties of the system), initial configurations being any possible sequence of N symbols randomly chosen in the alphabet {0, 1, 2, 3}. In addition to the invariance under the action of the finite group of cyclic permutations and reflections on the ring, the system admits another internal symmetry: the configurations are dynamically equivalent under the transformation $z_i \rightarrow z'_i = 3 - z_i$. This means that the whole orbit traced starting from a configuration $Z' = \{z_i\}$ is known if we know that traced by the configuration $Z = \{z_i\}$, and the two limit cycles have the same period. In the mean-field description [3,6], the critical behavior of the system around the absorbing-active transition point is deduced studying the activity phase diagram, the plot of the activity $\rho = \frac{1}{N} \sum_{i} \theta(z_i - 2)$ (the density of active sites) vs the energy density $\zeta = E/N$. The symmetry of the problem allows us to restrict the analysis of the 1D-DFES to the energy density range $\zeta \in$ [0, 3/2] (or in $\zeta \in [3/2, 3]$).

Numerical results show that large systems converge to periodic orbits of well-defined periods *T* in the whole range of the energy density. Figure 1 reports the diagrams $\rho(\zeta)$ and $T(\zeta)$ for a system of size N = 20 and 100 random initial conditions. To emphasize the nonergodicity, the



FIG. 1. The activity diagram $\rho(\zeta)$ (a) and the behavior of the period *T* as a function of the energy density ζ (b) for a 1D-DFES of size N = 20. Figures display all results, without averaging, obtained starting from 100 randomly chosen initial configurations at each energy value.

results of all runs are displayed, without computing any average. In particular, for $\zeta < 1$ the system converges to a completely frozen configuration (period T = 1) with no active sites, while in the range $1 < \zeta < 2$ only orbits of period T = 2 are observed. As imposed by symmetries, $\zeta > 2$ corresponds to fixed-point configurations with all sites active. At the critical points $\zeta = 1, 2$, orbits of period N seem to be statistically dominant for large systems, but other periods (e.g., T = 1, 2) are observed with lower frequency.

The reasons for such a particular phase diagram can be understood exploiting a result obtained in the mathematical literature, in which BTW sandpiles have been studied on generic graphs called parallel chip firing games (CFGs) [13]. When the underlying graph is undirected, with neither sinks nor sources, the CFG conserves the total energy. To our knowledge, the only result determining the properties of the periodic orbits of CFGs has been proved by Bitar and Goles [14] in the case of trees. Their theorem states that on a tree the steady states of the BTW dynamics are fixed points or cycles of period two. If the graph contains some loops, the theorem does not hold.

Notwithstanding, the method used in Ref. [14] can be exploited to study the 1D-DFES. Let us consider a system in a periodic steady state at a time t_0 , and a temporal window $P_{t_0} = [t_0, t_0 + T - 1]$ that corresponds to the first period from t_0 . If the system entered the periodic orbit O at time t_{in} , the whole temporal support $[t_{in}, \infty)$ of \mathcal{O} will be indicated with supp(\mathcal{O}). We define $s_i(t)$ as a binary variable that assumes value 1 (0) if the site *i* is active (stable) at the time t. The sequence of all these binary values forms the activity vector $S_i = \{s_i(t) | t \in \text{supp}(\mathcal{O})\}$ of a site *i*. Moreover, a set $[t, t + p - 1] \subseteq \text{supp}(\mathcal{O})$ is a maximal active set of length $p \ge 1$ if $s_i(t+r) = 1$ for r =0, 1, ..., p and s(t-1) = s(t + p + 1) = 0. Similarly, a set $[t, t + q - 1] \subseteq \text{supp}(\mathcal{O})$ is a maximal stable set of length $q \ge 1$ if s(t-1) = s(t+q+1) = 1 and $s_i(t+q) = 1$ r = 0 for $r = 0, 1, \dots, q$. According to Lemma 1 of Ref. [14], the following statement holds on a generic graph: if $[s - k, s] \subseteq \text{supp}(\mathcal{O})$ is a maximal active (stable) set for a site *i*, then there exists a neighbor *j* of *i* such that $[s - k - 1, s - 1] \subseteq \operatorname{supp}(\mathcal{O})$ is a maximal active (stable) set for j. We call V(W) the maximum number of consecutive 1's (0's) in all the activity vectors along the period, i.e., the length p(q) of the largest maximal active (stable) set over all sites. The analysis can be restricted to 0 < V < T(0 < W < T), the cases V = 0 (W = T) and V = T (W =0) corresponding to fixed-point configurations in which all sites are active (stable) [14]. Moreover, because of the internal symmetry with respect to $\zeta = 3/2$, *W*'s properties in the interval $\zeta \in [0, 3/2]$ are equivalent to those of *V* in the interval $\zeta \in [3/2, 3]$. As a consequence, we limit our study to *W* and corresponding maximal stable sets in the energy range $\zeta \in [0, 3/2]$. Two cases, W = 1 and W > 1, should be distinguished.

If W = 1, the activity vectors of all sites in the system are 2 periodic. Suppose, indeed, that a site *i* topples at a time *t* together with m < 2 of its neighbors: at the following time step t + 1 the site *i* does not topple, but the remaining 2 - m neighbors topple. Thus, having lost 2 - menergy units during the update at time *t* and gained the same quantity at the following time step, the value of site *i* has period 2. This argument holds for all sites in the periodic state (W = 1 for all sites), that consequently has global period T = 2.

In order to study the case W > 1, suppose the largest maximal stable set [t, t + W - 1] is at a site i_0 . For the above Lemma there is a neighbor i_1 of i_0 whose largest maximal stable set is [t - 1, t + W - 2]. In turn, site i_1 has a neighbor with largest maximal stable set [t-2, t+W-3]. However, W > 1 imposes that $i_2 \neq i_0$, because the Lemma implies $s_{i_2}(t-1) = 0$ while from the definition of maximal stable set we need $s_{i_0}(t-1) = 1$. Proceeding step by step along the ring in the same direction, we reach the site $i_N = i_0$, whose maximal stable set is [t - N, t +W - N - 1]. The properties of the system in i_0 (and in every other site) at time t - N result in the same as those at time t; thus we can conclude that during this process, the system performs one or more periodic cycles and returns in the starting configuration after exactly N temporal steps. In other words, for W > 1 (V > 1) the period T divides N.

The rest of this Letter provides a deeper insight into the structure of the periodic orbits using a simple method that is sketched in Fig. 2 for W = 2. Suppose that the interval [t, t + W - 1] of length 0 < W < N is the largest maximal stable set for a site *i*, then $s_i(t + s) = 0$ for s = 0, ..., W - 1, but also $s_i(t - 1) = 1$ and $s_i(t + W) = 1$. In addition, $s_{i-1}(t - 1 + r) = 0$ for r = 0, ..., W - 1, with $s_{i-1}(t - 2) = 1$ and $s_{i-1}(t + W - 1) = 1$; and similarly $s_{i+1}(t + 1 + r) = 0$ for r = 0, ..., W - 1, with $s_{i+1}(t) = 1$ and $s_{i+1}(t + W + 1) = 1$. These binary values can be drawn on a spatiotemporal grid as in Fig. 2. In order to determine the real values assumed by *i* in the interval [t, t + W - 1], we need to discuss separately the cases W = 1, W = 2 and W > 2.

The case W = 1 is simpler because all maximal stable sets have length 1, thus $s_i(t) = 0$ implies that at least one of the two neighbors is active. Drawing a spatiotemporal grid as in Fig. 2, it follows that spatial configurations can be



FIG. 2 (color online). Graphical sketch of the method used to find out the structure of periodic configurations. The case W = 2 is considered. Top figure shows the correspondence between the binary representation of the spatiotemporal dynamics and the evolution of the real configuration. An example of the procedure used to recover the real values assumed by a given site during the period is reported in the bottom series of draws. Site values can be univocally determined in W + 1 steps. At each step, the value of the site at a certain time is computed from a number of constraints on its (spatiotemporal) neighbors.

composed of all possible combinations of two-site blocks of the type 20, 21, 30 and 31 (and those with opposite order 02, 12, 03 and 13). Note that the set of all possible combinations of these blocks completely fills the energy interval $\zeta \in [1, 2]$ of 2-periodic orbits; while at the energy density $\zeta = 1$ ($\zeta = 2$), the configurations of the unique 2-periodic orbit are $Z_{20} = \{\dots 202020 \dots\}$ and $Z_{02} = \{\dots 020202 \dots\}$ ($Z_{31} = \{\dots 313131 \dots\}$ and $Z_{13} = \{\dots 131313 \dots\}$).

When W > 1 (see Fig. 2), the site *i* does not topple at time t + W - 1 but it has to topple at time t + W, after having received a single energy unit (from the site i - 1) at time t + W - 1, then $z_i(t + W - 1) = 1$ and, consequently, $z_i(t + W) = 2$. In particular, if W = 2, the sites *i* and i - 1 do not topple at time t + W - 2 = t, but i + 1topples, providing *i* of the unique unit of energy that we find at time t + W - 1 = t + 1. Hence, $z_i(t) = 0$, from which follows that $z_i(t - 1) = 2$.

In the case W > 2, the same arguments show that $z_i(t + W - 2) = 1$ (none of the neighbors provides any grain). If we go backward along the maximal complementary set, a site *i* assumes value 1 up to a time t + 1, then $z_i(t) = 0$ (one of its neighbors topples at that time). In summary, during the time interval [t - 1, t + W] with W > 1, the site under study assumes a well-determined sequence of values $2011 \dots 12$ with exactly W - 1 values equal to 1.

Figure 2 shows the existing relation between the values assumed by a site during these temporal intervals and those admitted in the spatial arrangement of the corresponding periodic configurations. If the maximal stable set of length W for i starts at time t, the uniqueness of the previous construction allows us to completely determine the spatial

block $\{z_{i-1}(t+W-1)z_i(t+W-1)\dots z_{i+W}(t+W-1)\}$ that will be of the form $\{211...102\}$ with W-1 sites at 1. Applying a spatiotemporal translation to our construction, the spatial block $\{z_{i-W}(t) \dots z_i(t) z_{i+1}(t)\} =$ $\{211...102\}$ is determined as well. The remaining structure of the configuration at time t can be studied with a similar technique, provided that we erase the spatial block $\{z_{i-w}(t) \dots z_i(t)\}$ and consider a reduced system composed of N - W - 1 sites. Since the role of the erased block was actually only that of transporting an activity "soliton" from the site i - W to the site i + 1, this operation does not alter the dynamics that maintains a periodic behavior. After the reduction, the above methods are applied on the system, with, in principle, a largest maximal stable set of different length $W' \leq W$. At the end of this process, the system (at a time t in the periodic regime) will be decomposed in blocks of decreasing length (from W + 1 to 2) and structure of the type 211...0. The only allowed blocks of length 2 are those of the form 20. The fronts direction of motion is determined by the initial conditions, but the dynamics is invariant under spatial reflections. Moreover, the analysis of maximal active sets for 0 < V < N leads to identical results (in the range $3/2 \le \zeta \le 3$) with spatial blocks of the type 122...31 (and 13...221) instead of 211...02 (and 20...112). We have proved by construction that the case W > 1 (V > 1) corresponds exactly to systems with an energy density $\zeta = 1$ ($\zeta = 2$). For all other values of ζ , only periods of length 1 and 2 are allowed. In the region $\zeta < 1$ (and $\zeta > 2$) the limit cycles are fixed-point configurations with all stable (active) sites, but in the interval $1 < \zeta < 2$ fixed points are forbidden and W = 1. Then, in this region the period of the limit cycles is T = 2, in agreement with Fig. 1. At the critical points $\zeta =$ 1 ($\zeta = 2$), the knowledge of the dynamics of building blocks also allows us to compute the exact number of periodic configuration of period T as a combinatorial enumeration problem. In particular, the number $\Pi_{all}(N)$ of configurations belonging to orbits with T > 1 equals the number of ways a numbered ring of N sites can be filled with an ordered set of blocks of length k comprised between 2 and N [15]. The generating function for the combinations with blocks of length $2 \le k \le N$ is O(z) = $1/(1 - \sum_{2 \le k \le N} z^k)$. Hence, $\prod_{all}(N)$ is obtained as function of the coefficients of the derivatives in z = 0 of Q(z); i.e., $\Pi_{\text{all}}(N) = \sum_{2 \le k \le N} \frac{2k}{(N-k)!} \frac{d^{(N-k)}}{dz^{(N-k)}} Q(z=0) - A_N, \quad \text{where}$ $A_N = 0$ if N is odd and $A_N = 2$ if N is even. The last term is introduced to compensate for the double counting (due to the factor 2k) of configurations composed of only 20 blocks. The number $\Pi_T(N)$ of configurations of period T < N is given by the number of different ways of filling the ring with identical blocks of length and period equal to T, i.e., $\Pi_T(N) = \Pi_T(T)$ for $2 \le T < N$. Formally, $\Pi_N(N) = \Pi_{\text{all}}(N) - \sum_{2 \le T < N} \Pi_T(N) \text{ (with } T \text{ divisor of } N).$ Starting from $\Pi_2(2) = \Pi_{\text{all}}(2) = 2$ and $\Pi_3(3) =$ $\Pi_{\text{all}}(3) = 6$, a recursive relation allows us to compute all the other terms $\Pi_T(N)$ ($2 \le T \le N$). The explicit compu-



FIG. 3. In panel (a), the number $\mu_T(N)$ of configurations belonging to the attraction's basin of orbits of period T = 1(crosses), 2 (squares), and N (circles) is plotted as function of N. The exponential rate of growth is estimated by curve fitting [dashed line for $\mu_1(N)$ and dot-dashed line for $\mu_N(N)$]. The inset shows that the ratio $\mu_1(N)/\mu_N(N)$ decreases with N. Panel (b) reports the total number $\Omega(N)$ of configurations at $\zeta =$ 1 (or $\zeta = 2$) as function of the system size N (circles). The fit (dashed line) shows an exponential growth as $\Omega(N) \propto \alpha^N$ with $\alpha \simeq 3.5$.

tation shows that, in the large N limit, the mass of orbits of period T = N grows faster than that of orbits of periods T < N.

However, in order to establish which period occurs with higher frequency at $\zeta = 1$ ($\zeta = 2$), the mass of the whole attraction's basin of an orbit is necessary, not only the number of periodic configurations. We have measured numerically the exact number of configurations in the basins of attraction of orbits of different periods. Figure 3(a) shows results up to N = 16. The size $\mu_1(N)$ of the basin of attraction of the fixed point grows exponentially with a rate $\mu_1(N)/\mu_1(N-1) \simeq \xi_1$ with $\xi_1 \simeq 3.25$. A similar growing rate is observed for orbits of period T =2 (N even), while the other orbits of period T < N present smaller attractions' basins. The largest growing rate is that of orbits of period N, for which $\mu_N(N)/\mu_N(N-1) \simeq \xi_N$ with $\xi_N \simeq 3.5$. This means that for N = 100 the probability of observing T = 1 or T = 2 is about 10^{-4} smaller than that of observing T = N. The conjecture that N-periodic orbits have probability 1 in the large N limit is corroborated by the observation of very similar scaling laws for $\mu_N(N)$ and the total number $\Omega(N)$ of possible configurations at energy E = N, whose behavior is displayed in Fig. 3(b) (data are computed analytically using simple combinatorics similar to that used in Ref. [16]). Consequently, at the critical energies $\zeta = 1, 2$, with probability 1, the system enters very long orbits characterized by a steady current of active solitons transported along the system. The activity can assume all (rational) values between 0 and 1/2.

In conclusion, our solution provides the full understanding of the mode-locking phenomena and the sharp activity transitions observed in the phase diagram of 1D-DFES. The arguments can be partially extended to higher dimensions, but the analysis is messed up by the complexity of the solitonic motion. In the case of two dimensional lattices, the present analysis together with symmetry arguments allow us to obtain an approximated picture of the correct phase diagram [17]. We hope that these results could also represent a kind of benchmark for the study of other automata with conserved dynamics, in which spatiotemporal patterns are governed by the motion of solitonic fronts.

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