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# Series expansion for a stochastic sandpile 

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#### Abstract

Using operator algebra, we extend the series for the activity density in a onedimensional stochastic sandpile with fixed particle density $p$, the first terms of which were obtained via perturbation theory (Dickman and Vidigal 2002 J. Phys. A: Math. Gen. 35 7269). The expansion is in powers of the time; the coefficients are polynomials in $p$. We devise an algorithm for evaluating expectations of operator products and extend the series to $\mathcal{O}\left(t^{16}\right)$. Constructing Padé approximants to a suitably transformed series, we obtain predictions for the activity that compare well against simulations, in the supercritical regime.


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## 1. Introduction

Sandpiles with a strictly conserved particle density (so-called fixed-energy sandpiles or FES [1]), exhibit absorbing-state phase transitions [2-4], and have recently attracted much interest. Until now, most quantitative results for FES have been based on simulations [5-8], an important exception being the solution by Priezzhev et al [9] of a directed, fixed-energy version of the Maslov-Zhang model [10], via the Bethe ansatz. It is therefore of great interest to develop theoretical approaches for FES.

Series analysis has proved to be one of the most accurate and reliable approaches to critical phenomena, in both equilibrium and nonequilibrium contexts [11-14]. Series expansion typically functions best in low-dimensional systems (because longer series can be derived), that is, for just those systems in which the renormalization group and expansion about an upper critical dimension $d_{\mathrm{c}}$ are less reliable. In the case of sandpiles, systematic epsilon expansions are as yet unavailable, and the value of $d_{\mathrm{c}}$ in fact remains controversial [15-17]. Simulation results suggest novel critical behaviour in the one-dimensional FES, although conflicting critical exponent values have been reported [7, 8, 18, 19], which may reflect finitesize effects. Series expansions, on the other hand, implicitly treat the infinite-size limit, and
so provide important information, complementary to that afforded by simulations. In light of these observations, we believe it highly desirable to apply series methods to sandpile models.

This paper is one of a series analysing a stochastic sandpile using operator methods. In an earlier work [20], a path-integral representation was developed and an expansion derived for the order parameter (activity density) in powers of time. While the path-integral formalism reveals interesting features of the model, and may be applied in any number of dimensions, the complexity of the diagrammatic expansion limits the number of terms that can be obtained. (In [20] terms up to $\mathcal{O}\left(t^{5}\right)$ are reported.) In this paper, we employ a different approach, which permits us to extend the series for the one-dimensional case considerably. After casting the master equation for the sandpile in terms of an operator formalism, we analyse the direct expansion of its (formal) solution, leading to an algorithm for generating the series coefficients.

We consider Manna's stochastic sandpile in its fixed-energy (particle-conserving) version [7, 20-22]. The configuration is specified by the occupation number $n$ at each site; sites with $n \geqslant 2$ are said to be active, and have a positive rate of toppling. When a site topples, it loses exactly two particles ('grains of sand'), which move randomly and independently to nearest-neighbour (NN) sites. (Any configuration devoid of active sites is absorbing, i.e., no further evolution of the system is possible once such a configuration is reached.) In this work, as in [20], we adopt a toppling rate of $n(n-1)$ at a site having $n$ particles, which leads us to define the order parameter as $\rho=\langle n(n-1)\rangle$. While this choice of rate represents a slight departure from the usual definition (in which all active sites have the same toppling rate), it leads to a much simpler evolution operator, and should yield the same scaling properties [20]. Preliminary simulation results [23] indicate that in one dimension the model exhibits a continuous phase transition at $p_{\mathrm{c}}=0.9493$.

In the following section, we define the model and review the operator formalism introduced in [20]. This is followed in section 3 by an analysis leading to an expansion in terms of socalled reduced commutators. Implementation of the expansion in a computational algorithm is described in section 4 . Then in section 5 we report numerical results of the series analysis. A summary and discussion is provided in section 6.

## 2. Model

As discussed in [20], the master equation for this model may be written in the form

$$
\begin{equation*}
\frac{\mathrm{d}|\Psi\rangle}{\mathrm{d} t}=L|\Psi\rangle \tag{1}
\end{equation*}
$$

where

$$
|\Psi\rangle=\sum_{\left\{n_{i}\right\}} p\left(\left\{n_{i}\right\}, t\right)\left|\left\{n_{i}\right\}\right\rangle
$$

is the probability distribution, and the evolution operator takes the form

$$
\begin{equation*}
L=\sum_{i}\left[\frac{1}{4}\left(\pi_{i-1}+\pi_{i+1}\right)^{2}-\pi_{i}^{2}\right] a_{i}^{2} \equiv \sum_{i} L_{i} \tag{2}
\end{equation*}
$$

Here $a_{i}$ and $\pi_{i}$ are, respectively, annihilation and creation operators associated with site $i$, defined via

$$
a_{i}\left|n_{i}\right\rangle=n_{i}\left|n_{i}-1\right\rangle
$$

and

$$
\pi_{i}\left|n_{i}\right\rangle=\left|n_{i}+1\right\rangle .
$$

The formal solution of the master equation is $|\Psi(t)\rangle=\mathrm{e}^{t L}|\Psi(0)\rangle$; that for the activity density is

$$
\begin{equation*}
\rho(t)=\langle | a_{0}^{2} \mathrm{e}^{t L}|\Psi(0)\rangle \tag{3}
\end{equation*}
$$

Here we have introduced the notation:

$$
\begin{equation*}
\langle | \equiv \sum_{\{n\}}\langle\{n\}| \tag{4}
\end{equation*}
$$

for the projection onto all possible states; thus normalization reads: $\langle\mid \Psi\rangle=1$. We consider a uniform Poisson-product initial distribution. Letting $p_{n}=\mathrm{e}^{-p} p^{n} / n!$, and using $|P\rangle_{i}=\sum_{n_{i}} p_{n_{i}}|n\rangle_{i}$ to denote a Poisson distribution at site $i$, we have

$$
\begin{equation*}
|\Psi(0)\rangle=\prod_{j}|P\rangle_{j} \tag{5}
\end{equation*}
$$

We shall expand equation (3) for the activity density in powers of $t$.

## 3. Operator algebra

To begin we note some basic properties of operators $a_{j}, \pi_{j}$ and $L_{j}$ :

$$
\begin{align*}
& a_{j}^{n}|P\rangle_{j}=p^{n}|P\rangle_{j}  \tag{6}\\
& \langle | \pi_{j}=\langle |  \tag{7}\\
& \langle | L_{i}=0  \tag{8}\\
& {\left[a_{i}, \pi_{j}\right]=\delta_{i, j}}  \tag{9}\\
& {\left[L_{i}, L_{j}\right]=\left[a_{i}, L_{j}\right]=0 \quad \text { for } \quad|i-j|>1} \tag{10}
\end{align*}
$$

The second relation expresses the fact that the creation operator conserves the normalization of any state, while the third shows that $L_{i}$ conserves probability, as it must.

The coefficient of $t^{n} / n!$ in the expansion of the activity is

$$
\begin{equation*}
\sum_{\mathcal{S}}\langle | a_{0}^{2} L_{s_{1}} L_{s_{2}} \cdots L_{s_{n}}|P\rangle \equiv \sum_{\mathcal{S}} \rho_{\mathcal{S}} \tag{11}
\end{equation*}
$$

where the sum is over all sequences $\mathcal{S}$ of sites $s_{0} \equiv 0, s_{1}, \ldots, s_{n}$ with $\left|s_{1}\right| \leqslant 1$, and $s_{j+1} \in\left\{s_{j, \min }-1, \ldots, s_{j, \max }+1\right\}$, for $j \geqslant 1$, where $s_{j, \min }=\min \left\{s_{0}, \ldots, s_{j}\right\}$, and $s_{j, \max }$ is the maximum of this set. The restriction on sequences follows from equations (8) and (10); if the condition were violated, it would be possible to move one of the $L_{j}$ to the left of all other operators, yielding a result of zero.

Our strategy for evaluating $\rho_{\mathcal{S}}$ is to commute each $L_{j}$ to the left of $a_{0}^{2}$. The first step replaces $a_{0}^{2} L_{s_{1}}$ by its commutator due to equation (8). If we write this commutator in normal order, that is, with all creation operators $\pi_{j}$ to the left of all annihilation operators, then the $\pi$ may be replaced by 1 by equation (7). Thus,

$$
\begin{equation*}
\langle | a_{0}^{2} L_{j}=\langle |\left[a_{0}^{2}, L_{j}\right]_{R} \tag{12}
\end{equation*}
$$

where the subscript $R$ denotes a reduced commutator, that is, the commutator in normal order, with all $\pi$ replaced by unity. Evidently $\left[a_{0}^{2}, L_{j}\right]_{R}$ involves only annihilation operators. The two nontrivial expressions of this kind are

$$
\begin{equation*}
\left[a_{0}^{2}, L_{0}\right]_{R}=-2 a_{0}^{2}-4 a_{0}^{3} \tag{13}
\end{equation*}
$$

and

$$
\begin{equation*}
\left[a_{0}^{2}, L_{1}\right]_{R}=\frac{1}{2} a_{1}^{2}+2 a_{0} a_{1}^{2} \tag{14}
\end{equation*}
$$

In the computational algorithm, discussed in some detail below, it is not necessary to generate the tree structure of sequences explicitly, since each monomial is processed separately and both translation and reflection symmetries may be used in the calculations of the contributions $\rho_{\mathcal{S}}$. Evaluating the expectation of each term in $\rho_{\mathcal{S}}$ is trivial, because

$$
\begin{equation*}
\langle | a_{s_{1}}^{m_{1}} \cdots a_{s_{n}}^{m_{n}}|P\rangle=p^{M} \tag{15}
\end{equation*}
$$

where $M=\sum_{j} m_{j}$ is the number of annihilation operators, irrespective of which sites are involved.

It remains to find a general expression for the reduced commutator $\left[F, L_{k}\right]_{R}$, where $F$ is a linear combination of products of annihilation operators. Recalling that $a_{i}$ and $L_{k}$ commute if $|i-k|>1$, we see that the problem reduces to evaluating

$$
\begin{equation*}
C(p, q, r) \equiv\left[a_{-1}^{p} a_{0}^{q} a_{1}^{r}, L_{0}\right]_{R} . \tag{16}
\end{equation*}
$$

(Commutators involving $L_{j}$ with $j \neq 0$ are obtained using translation invariance.) It is straightforward to evaluate $C(p, q, r)$ using the following identities. First we note that

$$
\begin{equation*}
\left[a_{j}^{p}, \pi_{j}\right]=p a_{j}^{p-1} \tag{17}
\end{equation*}
$$

as is readily shown by induction. Using this it is simple to show

$$
\begin{equation*}
\left[a_{j}^{p}, \pi_{j}^{2}\right]_{R}=p(p-1) a_{j}^{p-2}+2 p a_{j}^{p-1} \tag{18}
\end{equation*}
$$

and

$$
\begin{equation*}
\left[a_{j}^{p}, \pi_{j}^{2} a_{j}^{2}\right]_{R}=p(p-1) a_{j}^{p}+2 p a_{j}^{p+1} \tag{19}
\end{equation*}
$$

Finally, we may use equation (17) to show that for $i \neq j$,

$$
\begin{equation*}
\left[a_{i}^{p} a_{j}^{r}, \pi_{i} \pi_{j}\right]_{R}=p a_{i}^{p-1} a_{j}^{r}+r a_{i}^{p} a_{j}^{r-1}+p r a_{i}^{p-1} a_{j}^{r-1} . \tag{20}
\end{equation*}
$$

Applying these relations one readily finds

$$
\begin{align*}
C(p, q, r)= & a_{0}^{q+2}\left[\frac{1}{4} p(p-1) a_{-1}^{p-2} a_{1}^{r}+\frac{1}{4} r(r-1) a_{-1}^{p} a_{1}^{r-2}+\frac{1}{2} p r a_{-1}^{p-1} a_{1}^{r-1}\right. \\
& \left.+p a_{-1}^{p-1} a_{1}^{r}+r a_{-1}^{p} a_{1}^{r-1}\right]-q a_{-1}^{p} a_{1}^{r}\left[2 a_{0}^{q+1}+(q-1) a_{0}^{q}\right] \tag{21}
\end{align*}
$$

Using this result, we can evaluate the reduced commutators in a computer algorithm.

## 4. Computational algorithm

Let us discuss some details of the computer algorithm used to generate the series for the activity. We employ a recursive procedure to generate the contributions of order $n+1$ on the basis of those of order $n$. From equation (3) and with a Poisson-product initial distribution defined in equation (5) we note that
$\frac{\mathrm{d} \rho}{\mathrm{d} t}=\langle | a_{0}^{2} L \mathrm{e}^{t L}|P\rangle=\langle |\left[a_{0}^{2}, L\right]_{R} \mathrm{e}^{t L}|P\rangle=-\rho+\langle |\left(4 a_{0} a_{1}^{2}-4 a_{0}^{3}\right) \mathrm{e}^{t L}|P\rangle$.
The last equality above may be understood using the reduced commutators (13) and (14). Using reflection symmetry, we have $\langle |\left[a_{0}^{2}, L\right]_{R} \mathrm{e}^{t L}|P\rangle=\langle |\left\{\left[a_{0}^{2}, L_{0}\right]_{R}+2\left[a_{0}^{2}, L_{1}\right]_{R}\right\} \mathrm{e}^{t L}|P\rangle$ and further simplification is provided by translation symmetry. The coefficient of $t^{n} / n!$ in the expansion of the activity may be identified with $C_{n}$, where

$$
\begin{equation*}
C_{n}=\left.\frac{\mathrm{d}^{n} \rho}{\mathrm{~d} t^{n}}\right|_{t=0} \tag{23}
\end{equation*}
$$

Using the procedure described above we obtain a recursion relation

$$
\begin{equation*}
C_{n+1}=-C_{n}+\langle | F_{n+1}|P\rangle \tag{24}
\end{equation*}
$$

where $F_{n+1}=\left[F_{n}, L\right]$ and $F_{1}=4 a_{0} a_{1}^{2}-4 a_{0}^{3}$. To complete the algorithm, we have that $C_{0}(0)=\rho(0)=\langle | a_{0}^{2}| \rangle=p^{2}$, where relation (15) is used. An immediate consequence of these recursion relations is that the coefficient of $p^{2}$ in the term of order $n$ in $t$ is given by $(-1)^{n}$, since each monomial in the functions $F$ has at least three annihilation operators. This was already shown in [20].

The calculations were done in two steps. Initially, the functions $F$ were calculated up to order 12. Each monomial was represented by three integer variables: an eight-byte integer for the numerator of the coefficient, a four-byte integer for the denominator (which is always a power of 2 ), and another eight-byte integer to store the number of factors of each annihilation operator. Since all calculations are done in integer arithmetic, there are no roundoff errors. Using translation invariance, each monomial was put in a form such that the the annihilation operator of lowest spatial index is $a_{0}$. The power $m_{i}$ associated with each annihilation operator $a_{s_{i}}$ (as in equation (15)) is stored in four bits of the eight-byte integer variable mentioned above. As each new monomial is generated, a search is performed for any existing term with the same set of powers; storing all powers in a single integer facilitates the search. As a consequence of equation (10), when the reduced commutator of a monomial with $L$ is calculated, nonzero contributions may arise only from the commutator of the monomial with $L_{-1}, L_{0}, L_{1}, \ldots, L_{i+1}$, where $i$ is the largest index in the monomial. As is clear from equation (21), each of these commutators can give rise to up to seven new monomials. Thus, it is apparent that the number of monomials grows very rapidly as the order is increased; the function $F_{12}$ involves 519115 monomials. To go beyond order 12, it is necessary to handle integers larger than can be represented using eight bytes and to process monomials with more than 16 exponents, which can no longer be stored in a single eight-byte integer variable. In fact, at order 12 most of the processing time is used in the search procedure. Therefore, our results from order 13 to 16 were obtained processing the monomials in $F_{12}$ one-by-one, generating all contributions from it at orders 13-16. In these calculations, the numerators were represented by two eight-byte integers. The limiting order (16) is determined by the large number of new monomials generated; a single monomial in $F_{15}$ may generate on the order of 40 monomials in $F_{16}$. The results presented here required about 170 h of cpu time on an Athlon K7 1800 MHz computer.

It is convenient to write the expansion in the form

$$
\begin{equation*}
\bar{\rho}(t) \equiv \frac{\rho(t)}{p^{2}}=\sum_{n} \frac{(-t)^{n}}{n!} \sum_{m=0}^{n-1} b_{n, m} p^{m} \tag{25}
\end{equation*}
$$

The series coefficients $b_{n, m}$ are listed in table 1. In [20], it was shown that $m \leqslant n-1$, with

$$
\begin{equation*}
b_{n, n-1}=2^{4 n-1} \frac{(2 n-1)!!}{(2 n+2)!!} \tag{26}
\end{equation*}
$$

The coefficients reported in table 1 satisfy this relation at each order, and agree with those derived (for $n \leqslant 5$ ) using the path-integral formalism [20].

## 5. Analysis of series

The coefficients $b_{n, m} / n$ ! in the time series, equation (25), grow rapidly with $n$; the rate of growth appears to be faster than exponential. This is evident from an analysis of

$$
\begin{equation*}
h_{n, m} \equiv \ln \left(\frac{b_{n, m}}{n!}\right) \tag{27}
\end{equation*}
$$

Table 1. Series coefficients in the expansion of the activity.

| $n$ | $m$ | $b_{n, m}$ | $n$ | $m$ | $b_{n, m}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 0 | 0 | 1 | 10 | 4 | 3584915570625 |
| 1 | 0 | 1 |  | 5 | 7999349570432 |
| 2 | 0 | 1 |  | 6 | 6656488808368 |
|  | 1 | 8 |  | 7 | 2121777710528 |
| 3 | 0 | 1 |  | 8 | 227436059136 |
|  | 1 | 66 |  | 9 | 4402970624 |
|  | 2 | 80 | 11 | 0 | 1 |
| 4 | 0 | 1 |  | 1 | $\underline{12064410263}$ |
|  | 1 | 442 |  | 2 | 3142928518289 ${ }^{64}$ |
|  | 2 | 2076 |  | 3 | $\frac{85419503179415}{8}$ |
|  | 3 | 896 |  | 4 | 116020128091449 |
| 5 | 0 | 1 |  | 5 | 394806480115048 |
|  | 1 | 2842 |  | 6 | 514548057479072 |
|  | 2 | 35396 |  | 7 | 278154455793952 |
|  | 3 | 52240 |  | 8 | 61313513593600 |
|  | 4 | 10752 |  | 9 | 4683285856256 |
| 6 | 0 | 1 |  | 10 | 61641588736 |
|  | 1 | 18118 | 12 | 0 | 1 |
|  | 2 | 516880 |  | 1 | 76711895439 |
|  | 3 | 1737952 |  | 2 | $\frac{80070040225479}{32}$ |
|  | 4 | 1187968 |  | 3 | $\frac{1770456755814995}{8}$ |
|  | 5 | 135168 |  | 4 | $\underline{14610068149248089}$ |
| 7 | 0 | 1 |  | 5 | 18396700126638476 |
|  | 1 | $\underline{922265}$ |  | 6 | 35650110284461928 |
|  | 2 | 7040282 |  | 7 | 29745976515005712 |
|  | 3 | 45847512 |  | 8 | 11054665928232448 |
|  | 4 | 67368480 |  | 9 | 1747506609502464 |
|  | 5 | 25614368 |  | 10 | 97252577107968 |
|  | 6 | 1757184 |  | 11 | 872465563648 |
| 8 | 0 | 1 | 13 | 0 | 1 |
|  | 1 | $\underline{5865473}$ |  | 1 | $\underline{1951093993893}$ |
|  | 2 |  |  | 2 | $\underline{2037418656354491}$ |
|  | 3 | $\frac{4}{4} 078168434$ |  | 3 | $\begin{array}{r} 64 \\ 72926486692093419 \\ \hline \end{array}$ |
|  | 4 | 1078168434 |  | 3 | $\begin{array}{r} 16 \\ 905058014398112835 \\ \hline \end{array}$ |
|  | 4 | 2871388040 |  | 4 | $\frac{8}{8653914608555433}$ |
|  | 5 | 2283464832 |  | 5 | $\frac{1655391460208555433}{2}$ |
|  | 6 | 536472640 |  | 6 | 2309179626832648726 |
|  | 7 | 23429120 |  | 7 | 2816714002502804952 |
| 9 | 0 | 1 |  | 8 | 1601275099838022656 |
|  | 1 | $\frac{74596747}{16}$ |  | 9 | 426223203786122496 |
|  | 2 | 4797745191 |  | 10 | 49655626778919936 |
|  | 3 | 23841662132 |  | 11 | 2046635410882560 |
|  | 4 | 105679404154 |  | 12 | 12463793766400 |
|  | 5 | 147137780760 | 14 | 0 | 1 |
|  | 6 | 71353965088 |  | 1 | $\frac{2613736799297}{64}$ |
|  | 7 | 11072770560 |  | 2 | $\underline{11934019637184639}$ |
|  | 8 | 318636032 |  | 3 | 711799150376749517 |
| 10 | 0 | 1 |  | 4 | $\underline{53725847102644113051}$ |
|  | 1 | 474336627 |  | 5 | $\xrightarrow{142451880202934178839}$ |
|  | 1 | $\frac{16}{16}$ |  | 5 | $\frac{14245802}{4}$ |
|  | 2 | $\frac{123077063429}{8}$ |  | 6 | 140971191603315396510 |
|  | 3 | $\frac{1018938641745}{2}$ |  | 7 | 243993717303561711492 |

Table 1. (Continued.)

| $n$ | $m$ | $b_{n, m}$ | $n$ | $m$ | $b_{n, m}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 14 | 8 | 201186302850192322944 | 15 | 12 | 40856912571394580480 |
|  | 9 | 81746928038823569408 |  | 13 | 957525442027462656 |
|  | 10 | 16113035142846829824 |  | 14 | 2602440138424320 |
|  | 11 | 1415730263534155776 | 16 | 0 | 1 |
|  | 12 | 43811063460921344 |  | 1 | 422699161810361 |
|  | 13 | 179478630236160 |  | 2 | 61687281835997869817 |
| 15 | 0 | 1 |  | 3 |  |
|  | 1 | 531822407449409 |  | 4 | $\underline{12465559773385531628949}$ |
|  |  | $\begin{array}{r} 2048 \\ 606748047325325193 \\ \hline \end{array}$ |  | 5 | 133184916649036025384179 |
|  | 2 |  |  | 5 | $\frac{2}{2}$ |
|  | 3 | $\frac{116603968592784196927}{64}$ |  | 6 | 502866857598243404511546 |
|  | 4 | 819559563865455675379 |  | 7 | 1627359034988855536002199 |
|  | 5 | 12371190775573187034899 |  | 8 | 2537319446210036202445148 |
|  | 6 | 8499377166784889887638 |  | 9 | 2040132355769944077918400 |
|  | 7 | 20286352375993324998496 |  | 10 | 872811268389569306302976 |
|  | 8 | 23298464567721533566328 |  | 11 | 198169235678101485620992 |
|  | 9 | 13579307980015469945184 |  | 12 | 22853161358227259040768 |
|  | 10 | 4068154005082098401408 |  | 13 | 1195547596367062589440 |
|  | 11 | 607110051667479652352 |  | 14 | 21401594847260721152 |
|  |  |  |  | 15 | 37965009078190080 |

To see if the $h_{n, m}$ follow a systematic trend we analyse these quantities for a given $q \equiv(m-1) /(n-2)$. (For a fixed value of $n$, the $h_{n, m}$ appear to trace out a smooth curve, so that $h_{n}(q)$ for intermediate values of $q$ can be estimated via interpolation.) As shown in figure $1, h_{n}(q)$ appears to grow faster than exponentially with $n$, away from the limits $q=0$ and $q=1$. (Observe that for $m=0, h_{m, n} \rightarrow-\infty$ as $n \rightarrow \infty$ since $b_{n, 0}=1$, and similarly for $m=n-1$, since equation (26) implies that $b_{n, n-1}$ grows more slowly than $n!$ ) A reasonable description of the dominant growth in the series coefficientes is $h_{n}(q) \sim n^{\alpha}(q)$, with the exponent $\alpha$ (see the inset of figure 1) taking its maximum value of about 1.2 for $q \simeq 0.4$. This of course implies faster-than-exponential growth for the coefficients $b_{n, m} / n$ !.

Next we examine the behaviour of the coefficients in the time series for specific values of the particle density $p$. Let

$$
\begin{equation*}
\bar{\rho}(t)=\sum_{n} c_{n}(-t)^{n} \tag{28}
\end{equation*}
$$

where

$$
\begin{equation*}
c_{n} \equiv \frac{1}{n!} \sum_{m=0}^{n-1} b_{n, m} p^{m} \tag{29}
\end{equation*}
$$

For $p=1$ (slightly above the critical value of 0.9493 ), $c_{n}$ is simply the sum of all coefficients at order $n$, divided by $n!$. The coefficients $c_{n}$ again grow faster than exponentially, with $\ln c_{n} \sim n^{1.15}$ for $p=1$ and $\sim n^{1.10}$ for $p=2$. (Given the limited number of coefficients, we cannot make very precise estimates of the exponent. The key point is that the growth appears to be faster than exponential.) These results imply that equation (28) is a divergent series with zero radius of convergence.

We turn now to an analysis of the series for $\bar{\rho}(t)$. As is well known, it is often possible to obtain useful results from divergent series by means of a resummation technique. In the


Figure 1. Function $h_{n}(q)$ as defined in text, for $q=0.2(\square) ; q=0.4(\bullet) ; q=0.6$ (o); $q=$ $0.8(\square)$. Observe that $h$ grows faster than linearly for $q=0.2,0.4$. Inset: growth exponent $\alpha(q)$ defined via $h_{n}(q) \sim n^{\alpha(q)}$.
present case, we construct Padé approximants to the time series or to the series obtained via a transformation of variables [13, 25-27]. We have examined many transformations, for example

$$
\begin{align*}
& y=\frac{1-\mathrm{e}^{-b t}}{b}  \tag{30}\\
& x=\frac{t}{1+b t}  \tag{31}\\
& z=1-\frac{1}{(1+b t)^{\gamma}}  \tag{32}\\
& w=1-\frac{1}{1+\ln (1+b t)} \tag{33}
\end{align*}
$$

and

$$
\begin{equation*}
v=1-\exp \left\{b\left[1-(1+t)^{\gamma}\right]\right\} . \tag{34}
\end{equation*}
$$

Each transformation maps the interval $t \geqslant 0$ to a finite interval, and can be expanded as a power series in $t$ about $t=0$, with the lowest-order term $\propto t$. Each is readily inverted permitting one to express the time $t$ in powers of the new variable. The slow convergence associated with the power-law or logarithmic forms in the last four expressions is motivated by the numerical finding of slow relaxation in the sandpile model, even far from the critical point [24]. We analyse the transformed series for $\bar{\rho}$ (or for $\ln \bar{\rho}$ ) using Padé approximants. The degree of success depends greatly on the range of $p$ under consideration. (Each transformation includes a free parameter $b$, which can be adjusted to optimize the regularity of the result, or to obtain consistency between different approximants. Except where noted, the results do not exhibit much sensitivity to the choice of this parameter.)


Figure 2. Normalized activity $\bar{\rho} \equiv \rho(t) / p^{2}$ versus time for $p=1 / 2$ from simulation and various Padé approximants to the times series as indicated.

For small values of $p$, the best results are obtained via Padé approximants to the $t$-series without any transformation of variable. Figure 2 compares series predictions for $p=0.5$ (obtained using the $[6,7],[7,8]$ and $[7,9]$ approximants to the series for $\bar{\rho}(t)$ ) against the result of a Monte Carlo simulation for a system of 500 sites (for $p=0.5$ finite-size effects are negligible at this system size). The [7,8] approximant is reliable for $t \leqslant 10$. (Various other approximants, such as $[8,8]$ and $[7,7]$, are ill-behaved and provide reasonable predictions only for quite short times, typically $t \leqslant 2$.) We have not been able to improve the series prediction for longer times, either by a change of variable or through analysis of $\ln \bar{\rho}(t)$ or its time derivative. Although some improvement could be expected with longer series, it appears unlikely that the asymptotic decay of $\bar{\rho}(t)$ in the subcritical regime will be accessible through analysis of an expansion in powers of time.

For larger values of $p$ the transformation defined in equation (32), using $\gamma=1 / 2$, is the most useful of those studied. In figure 3 , we compare the $[8,8]$ approximant (obtained using $b=0.57$ ) with simulation data for $p=1$. The situation is markedly better than for $p=1 / 2$ : the series prediction accompanies the simulation result up to around $t=1000$. It must be noted, however, that the good agreement seen here depends on the choice of the transformation parameter $b$. For other values, the agreement with simulation is not as good. (A more suitable criterion for choosing $b$ would be by seeking agreement among various approximants [13]. In the present case, this is not possible because the off-diagonal approximants to the $z$ series are ill-behaved, while the $[7,7]$ approximant behaves very similarly to the $[8,8]$ used here.) Despite the good agreement up to times of the order of 1000 , the present series seems incapable of capturing the asymptotic long-time relaxation of $\bar{\rho}(t)$, which is non-monotonic, as shown in figure 3.

Remarkably, the reliability of the series improves dramatically at larger values of the particle density $p$. Series and simulation results for $\bar{\rho}(t)$ at $p=2$ are compared in figure 4 ; the maximum relative error is about $0.1 \%$. (The series prediction is generated as for $p=1$, but using $b=1.5$ in this case.) The good agreement, moreover, persists at long times, motivating a study of $\bar{\rho}_{\infty} \equiv \lim _{t \rightarrow \infty} \bar{\rho}(t)$, corresponding to the transformed series with $z=1$ in equation (32). (We again use the $[8,8]$ approximant to the $z$-series.)

The series prediction for $\bar{\rho}_{\infty}$ is compared with simulation in figure 5 , using parameters $b=0.57$ and $b=5$. Excellent agreement is found for $p \geqslant 2$, the relative error being $\leqslant 0.2 \%$.


Figure 3. Normalized activity for $p=1$. Symbols: simulation result; curve: series prediction as described in text.


Figure 4. As in figure 3 for but for $p=2$.

The smaller $b$ value yields better results for $p \simeq 1$, whereas slightly better results are obtained for larger $p$, using the larger $b$ value. For $p \geqslant 2$ we may claim quantitative accuracy for the series prediction. Nearer the critical point, the agreement appears reasonable (at least on the scale of figure 5), but it is clear that the 16 -term series cannot be used to study critical properties. For example, the prediction using $b=0.57$ yields a critical value of about 0.906 , that is, the extrapolated activity density goes to zero at this $p$ value. The critical value found in simulations is 0.9493 .

In summary, the present series seems quite reliable in the supercritical regime, both at short and at asymptotically long times, whereas its utility in the critical and subcritical regimes is restricted to rather short times. Just above the critical point, rather good predictions are possible for short and intermediate times, but this depends on a judicious choice of the transformation parameter $b$.


Figure 5. Main graph: limiting activity $\bar{\rho}_{\infty}$ versus particle density $p$. Points: simulation; solid curve: series prediction using transformation (32) with $b=0.57,[8,8]$ Padé approximant; dashed line: same approximant and transformation but using $b=5$. Inset: difference $\Delta=\bar{\rho}_{\infty, \text { series }}-\bar{\rho}_{\infty, \text { sim }}$ for $b=0.57$ (■) and $b=5$ ( $\square$ ).

## 6. Discussion

We develop an algebraic method leading to a time series for the activity density of the stochastic sandpile model introduced in [20]. Determination of the series coefficients depends on evaluation of certain commutators, an algebraic task readily codified in a computational algorithm. We extend the series for the one-dimensional case to 16 terms.

Analysis of the series yields disappointing results for the subcritical and critical regimes, but very good predictions in the supercritical region, as judged by comparison with Monte Carlo simulation. At first glance this is surprising, since in the subcritical regime the stationary state is inactive and might be regarded as trivial. Relaxation to this inactive state (and to the active state at or near the critical point $p_{c}$ ) is however nontrivial, characterized by stretched exponential, power-law or other slowly-converging forms [24]. It appears to be very difficult to capture such behaviour in the kind of temporal series developed here, which employs a Poissonian initial distribution. The reason is that for smaller values of the particle density ( $p<2$, say), the one-site stationary occupation distribution $P(n)$ is far from Poissonian. As $p$ increases, the second factorial moment $\langle n(n-1)\rangle=\rho$ approaches $p^{2}$, as expected for a Poisson distribution. Figure 6 shows that the stationary one-site distribution observed in simulations approaches the corresponding Poisson distribution with the same density $p$. (Even for $p=8$ there are significant differences between the distributions; but analysis of the third and fourth factorial moments suggests convergence to a Poisson distribution as $p \rightarrow \infty$.)

An important open question is whether simply increasing the number of terms would permit one to analyse the small- $p$ regime. The present results suggest that even with 20 or 30 terms this region would remain inaccessible. It appears to be more promising to approach the critical region from above, since for larger particle densities we find good agreement with numerical results. In this context, it is interesting that the quality of predictions near the critical point improves greatly on going from 12 to 16 terms. This suggests that further extension of the series, to 20 or more terms, would yield quantitative results for critical properties, through study of relaxational properties at $p_{\mathrm{c}}$, or of stationary properties (as in figure 5) as


Figure 6. Single-site occupancy distributions $P(n)$ obtained in simulation ( $\square$ ) compared with the corresponding Poisson distribution ( $\square$ ). Upper panel: $p=1.2$; middle: $p=3$; lower: $p=8$.
$p_{\mathrm{c}}$ is approached from above. It would also be of great interest to develop an expansion for a stationary property such as $\bar{\rho}_{\infty}$ directly in powers of the particle density $p$, but this appears to be much more difficult than deriving an expansion in powers of the time. We leave such investigations, using modified or extended series, as subjects for future work.

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