

When are cellular automata random?

J. B. Coe¹³, S. E. Ahnert⁴, and T. M. A. Fink¹²³

¹INSERM U900 and ²CNRS UMR144, Curie Institute, Paris F-75248, France

³Ecole des Mines de Paris, ParisTech, Fontainebleau, F-77300 France and

⁴Theory of Condensed Matter, Cavendish Laboratory, Cambridge CB3 0HE, UK

A random cellular automaton is one in which a cell's behaviour is independent of its previous states. Analytical conditions for the existence of random cellular automata are derived and we find that a multitude of non-trivial random cellular automata exist. We develop an indicator variable formalism to further investigate these random automata and confirm analytical results with simulation.

Introduction

A cellular automaton is a collection of cells, each of which is in one of a finite number of states. The cells' states are updated according to some local rules which take into account the state of the cell and the states of its neighbours. For instance: a two dimensional grid with two states per cell is used in Conway's Game of Life in which each cell is alive or dead and transition rules are intended to mimick the effects of birth, death from isolation and death from overcrowding [1]. Cellular automata have attracted attention from the physics community as they are an appealing modelling tool for any process driven by local interactions.

Cellular automata have been used to model a wide range of physical phenomena including: traffic flow [2–8]; disease epidemics [9, 10]; stochastic growth [11]; predator-prey dynamics [12, 13]; invasion of populations [14]; earthquakes [15] and dynamics of stock markets [16].

Analytical work has been fruitful in classifying behaviour and studying properties of specific forms. Wolfram exhaustively catalogued and classified behaviour of simple deterministic automata [17, 18]. Phase diagrams and critical exponents have been evaluated for automata with absorbing states [19–23]. Certain probabilistic automata have been shown to fall into the same universality class as directed percolation [24, 25]. Coarse-graining has been shown to predict the emergence of large scale properties even for computationally irreducible systems [26]. Fuks et. al. identified conservative automata in which the number of coloured cells remains constant [27–29].

Classification of rules is of fundamental importance in any broad analytical treatment of cellular automata. Random behaviour is among the most elementary forms of behaviour possible and must be firmly understood if a general theory of structure — deviation from randomness — is to be developed.

In deterministic cellular automata the update rules have no probabilistic component: for a given configuration of cell states the updated cell state is always the same. In probabilistic cellular automata, local rules may have a probabilistic element to them: rather than dictating the state of an updated cell, the rule gives the prob-

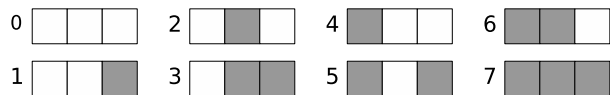


FIG. 1: The eight possible neighbourhoods for a simple cellular automaton are numbered as illustrated. The probability that the neighbourhood numbered i produces a coloured cell is denoted by a_i .

ability that an updated cell will be in each of a number of states.

The simplest cellular automaton is a line of cells, each of which is in one of two states: coloured or uncoloured. The cells are updated simultaneously; the updated state of a cell depends on its state and the states of its nearest neighbours. There are eight possible configurations for the states of a cell and its nearest neighbours; this gives eight different update rules which we label as a_i . The configurations and rules are numbered according to Fig. 1.

We denote the state of the central cell by x_c : when the central cell is coloured $x_c = 1$, when it is uncoloured $x_c = 0$. The left neighbour is denoted by x_l and the right neighbour by x_r . Uncoloured cells are denoted by \bar{x}_l , \bar{x}_c and \bar{x}_r ($\bar{x}_i = 1 - x_i$). The state of the central cell after an update, x'_c is given by

$$\begin{aligned}
 P(x'_c = 1) &= a_0 \bar{x}_l \bar{x}_c \bar{x}_r + a_1 \bar{x}_l \bar{x}_c x_r \\
 &+ a_2 \bar{x}_l x_c \bar{x}_r + a_3 \bar{x}_l x_c x_r \\
 &+ a_4 x_l \bar{x}_c \bar{x}_r + a_5 x_l \bar{x}_c x_r \\
 &+ a_6 x_l x_c \bar{x}_r + a_7 x_l x_c x_r.
 \end{aligned}
 \tag{1}$$

In this paper we present the conditions for random cellular automata (which we derive in a later section).

We first solve these conditions over the space of deterministic cellular automata. We find that of the 256 simple cellular automata, 28 of them exhibit random behaviour. In all but two of them, cells are coloured with probability $\frac{1}{2}$. Ten of these random automata are not related by symmetry (reflection or inversion), all of which we illustrate in fig. 2.

Second, we solve the conditions for random behaviour for probabilistic cellular automata. Within the eight-

cube — the corners of which are the 256 deterministic cellular automata — we find an expression for the density of random cellular automata and find two disjoint five-volumes both of which have volume 0.12. We give examples of random probabilistic cellular automata through illustrations.

In the third section we derive the conditions for randomness by applying an indicator variable formalism to nearest neighbour one-dimensional cellular automata.

We employ indicator variables in the fourth section to study *correlation current* which manifests itself as a visually apparent flow in many of the random automata. We develop an analytical expression for correlation current and compare predictions with simulation.

In the fifth section we show that the total randomness in random cellular automata, must come from either of two sources: spatial randomness in the initial conditions which is converted to temporal randomness; or randomness injected directly through probabilistic update rules. We derive explicit expressions to calculate these quantities for any random cellular automaton.

Random cellular automata

In a random cellular automaton the probability of a cell being coloured is independent of all of its previous states: all cells are taken to be coloured with probability p . There are two sets of criteria on the rules; satisfying either of these produces random behaviour:

$$\begin{aligned} (1-p)a_0 + pa_4 = p, & \quad (1-p)a_1 + pa_5 = p, & (2) \\ (1-p)a_2 + pa_6 = p, & \quad (1-p)a_3 + pa_7 = p; \end{aligned}$$

$$\begin{aligned} (1-p)a_0 + pa_1 = p, & \quad (1-p)a_2 + pa_3 = p, & (3) \\ (1-p)a_4 + pa_5 = p, & \quad (1-p)a_6 + pa_7 = p. \end{aligned}$$

Note that these conditions have certain symmetry properties. Switching over left- and right-handed rules turns one set of conditions into the other. State-inversion (where coloured cells and uncoloured cells are switched round) leaves both sets of conditions unchanged.

We examine the behaviour of random automata for both deterministic and probabilistic rules. Since conditions have left-right reflective symmetry, we only look at examples for the second set of conditions, (3).

For a random cellular automaton which satisfies (3), the state of the central cell after an update is given by

$$\begin{aligned} P(x'_c = 1) = & x_r + (\bar{x}_r - \left(\frac{1-p}{p}\right)x_r) & (4) \\ & \times (a_0\bar{x}_l\bar{x}_c + a_2\bar{x}_l x_c + a_4x_l\bar{x}_c + a_6x_l x_c). \end{aligned}$$

Wolf.	$a_7 \dots a_0$	x'_c	R	I	$R\&I$
85	01010101	\bar{r}	15	85	15
86	01010110	$\bar{l}\bar{c}r + l\bar{r} + c\bar{r}$	30	149	135
89	01011001	$\bar{l}rc + \bar{c}\bar{r} + l\bar{r}$	75	101	45
90	01011010	$l\bar{r} + \bar{r}l$	90	165	165
102	01100110	$c\bar{r} + \bar{c}r$	60	153	195
105	01101001	$\bar{l}\bar{c}\bar{r} + lc\bar{r} + l\bar{c}r + \bar{l}cr$	105	105	105
106	01101010	$lc\bar{r} + \bar{c}r + l\bar{r}$	120	169	225
150	10010110	$l\bar{c}\bar{r} + \bar{l}c\bar{r} + \bar{l}\bar{c}r + lcr$	150	150	150
154	10011010	$l\bar{c}\bar{r} + \bar{l}r + lr$	210	166	180
170	10101010	r	240	170	240

TABLE I: The 10 sets of rules which give random behaviour along with their symmetric partners. The rules themselves are shown along with a simplified boolean expression for the updated state of the central cell. Wolfram numbers and Wolfram numbers of symmetric partners are given. The symmetry operations considered are reflection (R), state-inversion (I) and both (R&I). The value of p for all rule sets is 0.5 except for rule set 170, for which p can take any value.

Random deterministic automata

For deterministic cellular automata all the rules must be either zero or one. We are not interested by trivial solutions in which all cells are coloured ($p = 1$) or uncoloured ($p = 0$); this imposes limits on possible choices of p . For (3) this means that if $a_{2j} = 0$ then $a_{2j+1} = 1$ and if $a_{2j} = 1$ then $a_{2j+1} = 0$. One choice of rules, where all $a_{2j} = 0$, gives random behaviour for arbitrary p , all other choices give $p = \frac{1}{2}$. There are 16 possible choices of free rules which give random behaviour, though some are related by symmetry.

The left-handed conditions provide an equal number of possible rule sets leading to random behaviour some of which are already given by the right-handed conditions. There are 4 possible choices which satisfy both left- and right-handed conditions for randomness giving a total of 28 distinct random deterministic cellular automata.

Wolfram assigns deterministic automata numbers according to the values of the rules [17]. These Wolfram numbers can be obtained by evaluating $\sum_0^7 a_i 2^i$. Of the 28 rule sets which give random behaviour when symmetric partners (either left-right or inversion symmetries) are discounted, there are 10 distinct rule sets. We list these rule sets and their symmetries in table I. Several of these automata belong to Wolfram's classes 3 and 4 which exhibit the most complex kinds of behaviour [18].

Of the 10 rule sets, some can be qualitatively explained. 170 copies the right site to itself, propagating the initial sequence along space and time. 85 copies the conjugate of the right site to itself. 90 is a NAND gate on the left and right sites, and 102 is a NAND gate in the left and centre sites. 105 gives a coloured cell if its neighbourhood contains one or three zeros. 150 gives a coloured cell if its neighbourhood contains none or two zeros.



FIG. 2: Simulation-generated pictures for the 10 right-handed random deterministic automata not related to one another by symmetry. Space runs from left to right, time from top to bottom. From left to right, starting at the top, the Wolfram numbers are: 85, 86, 89, 90, 102, 105, 106, 150, 154 and 170.

Random probabilistic automata

When the rules are no longer constrained to be zero or one we cannot count configurations so instead we consider the density of random states. As pairs of rules are related to one another by the conditions for random behaviour, there are four free rules and a probability of a cell being coloured, p , to choose. If these five free parameters were each randomly picked from the interval zero to one it would not necessarily result in random behaviour. The non-free rules must also lie between zero and one which may not be compatible with (2) and (3), the conditions for random behaviour.

Within the five-volume marked out by the free rules and p , there is a density of random configurations given by

$$\begin{aligned} dV_5 &= \left(\frac{p}{1-p}\right)^4 dp, & p \leq \frac{1}{2}, \\ dV_5 &= \left(\frac{1-p}{p}\right)^4 dp, & p > \frac{1}{2}. \end{aligned} \quad (5)$$

Integrating over all possible values of p , we find

$$V_5 = 17/3 - 8 \ln 2 \simeq 0.12. \quad (6)$$

Within the five-volume, 12% of possible selections of free rules and colouring probabilities will lead to random behaviour. A second five-volume exists when the other set of conditions is satisfied. Note that the five-volumes that correspond to different sets of conditions are not the same (different rules are free) and have zero intersection in five dimensions.

Satisfying both sets of conditions gives a three-volume within which 23% of randomly picked points correspond to random behaviour.

Imposing left-right symmetry by satisfying both sets of conditions, (2) and (3), yields

$$\begin{aligned} P(x'_c = 1) &= 1 - \bar{x}_l \bar{x}_r - r x_l x_r \\ &+ (a_0 \bar{x}_c + a_2 x_c)(\bar{x}_r - r x_r)(\bar{x}_l - r x_l), \end{aligned} \quad (7)$$

which is symmetric under interchange of x_l and x_r as expected. In this expression the state of the central cell and its left and right neighbours are all important.

Some illustrative examples, all with $p = 0.5$, of random probabilistic cellular automata are shown in Fig. 3. A Java program [31] can be freely downloaded and used to generate figures for probabilistic cellular automata.

Indicator variables

Indicator variables are taken from the theory of random processes and have been used to solve the asymmetric exclusion process [32]. In a powerful formalism in population genetics, indicator variables are used to keep track of genes and correlations between genes [33]. Here we make use of indicator variables to represent the state of a cell and use cancellation properties to simplify expressions for high-order correlations.

Every cell in a one-dimensional cellular automaton can be assigned a unique integer index increasing from left to right. We represent the state of each cell by an indicator variable τ_i , where i is the index of the site. The indicator variable takes the value 1 if a site is coloured and 0 if it is not. For notational brevity we define an anti-indicator $\bar{\tau}_i$ which shows if a cell is uncoloured: $\bar{\tau}_i = 1 - \tau_i$.

If p_i is the probability that cell i is coloured, then the expected value of the indicator variable τ_i is p_i . We represent this as

$$\mathbb{E}[\tau_i] = p_i. \quad (8)$$

Several identities follow from the above definitions and can be used to simplify complicated expressions in which indicator variables appear:

$$\mathbb{E}[\tau_i \tau_i] = \mathbb{E}[\tau_i]; \quad \mathbb{E}[\bar{\tau}_i \bar{\tau}_i] = \mathbb{E}[\bar{\tau}_i]; \quad \mathbb{E}[\tau_i \bar{\tau}_i] = 0. \quad (9)$$

Correlations between indicators and anti-indicators can always be expressed as correlations between indicators by exploiting the identity $\bar{\tau}_i = 1 - \tau_i$:

$$\mathbb{E}[\bar{\tau}_i \bar{\tau}_j] = 1 - \mathbb{E}[\tau_i] - \mathbb{E}[\tau_j] + \mathbb{E}[\tau_i \tau_j] \quad (10)$$

$$\mathbb{E}[\tau_i \bar{\tau}_j] = \mathbb{E}[\tau_i] - \mathbb{E}[\tau_i \tau_j]. \quad (11)$$



FIG. 3: Simulation-generated pictures for some representative samples of random probabilistic cellular automata. In all of these automata the probability that a cell is coloured is $\frac{1}{2}$. The rules, from left to right, from a_7 to a_0 are given at the end of this manuscript [30].

For an unknown configuration of cells, the probability that the cell with index 1 is coloured after an update follows from the definitions of the update rules: each possible neighbourhood can be described by a three-fold product of indicator variables; each neighbourhood has a probability a_i of producing a coloured site after an update. We write the indicator variable for cell i after an update as τ'_i .

$$\begin{aligned} \mathbb{E}[\tau'_1] &= a_0\mathbb{E}[\bar{\tau}_0\bar{\tau}_1\bar{\tau}_2] + a_1\mathbb{E}[\bar{\tau}_0\bar{\tau}_1\tau_2] \\ &+ a_2\mathbb{E}[\bar{\tau}_0\tau_1\bar{\tau}_2] + a_3\mathbb{E}[\bar{\tau}_0\tau_1\tau_2] \\ &+ a_4\mathbb{E}[\tau_0\bar{\tau}_1\bar{\tau}_2] + a_5\mathbb{E}[\tau_0\bar{\tau}_1\tau_2] \\ &+ a_6\mathbb{E}[\tau_0\tau_1\bar{\tau}_2] + a_7\mathbb{E}[\tau_0\tau_1\tau_2]. \end{aligned} \quad (12)$$

Without assumptions this cannot be simplified.

Conditions for random behaviour

A random configuration is one in which the probability of any given cell being coloured is uniformly and independently distributed through time. As the probability of a cell being coloured can be expressed in terms of neighbouring cells being coloured in the previous time step (12), for a cell to be independent of its previous states requires inter-cell independence at any time step. Single-cell temporal independence requires system wide spatial independence at any timestep. If the cells are coloured with probability p , a random configuration can be expressed in terms of the indicator algebra as

$$\mathbb{E}[\tau'_i] = p, \quad (13)$$

$$\mathbb{E}\left[\prod_{i \in \mathcal{S}} \tau'_i\right] = \prod_{i \in \mathcal{S}} \mathbb{E}[\tau'_i]. \quad (14)$$

The first condition, (13), can be written in terms of indicator variables and simplified using (14). This results in a polynomial in p .

$$\begin{aligned} p &= a_0q^3 + a_1pq^2 + a_2pq^2 + a_3p^2q \\ &+ a_4pq^2 + a_5p^2q + a_6p^2q + a_7p^3, \end{aligned} \quad (15)$$

where $q = 1 - p$. While this can be satisfied for many choices of a_i and p , the second condition (14) is much stricter and any set of rules satisfying (14) will satisfy this polynomial.

To satisfy (14) no inter-cell correlations can exist for any possible set of cells. For certain choices of a_i , larger correlation terms can be expressed in terms of shorter ones: repeating this procedure reduces all correlation terms to products of single cell expectations. This reduction of correlation terms can be performed in one of two ways:

$$\mathbb{E}\left[\prod_{i \in \mathcal{S}} \tau_i\right] = \mathbb{E}[\tau_0]\mathbb{E}\left[\prod_{i \in \mathcal{S} \setminus \{0\}} \tau_i\right], \quad (16)$$

$$\mathbb{E}\left[\prod_{i \in \mathcal{S}} \tau_i\right] = \mathbb{E}\left[\prod_{i \in \mathcal{S} \setminus \{N\}} \tau_i\right]\mathbb{E}[\tau_N]. \quad (17)$$

where N and 0 have been chosen, without loss of generality, to be the largest and smallest indices in the set \mathcal{S} . We label \mathcal{S} without the index 0 as $\mathcal{S} \setminus \{0\}$ and \mathcal{S} without the index N as $\mathcal{S} \setminus \{N\}$. We refer to removal of the smallest index as left-handed reduction and removal of the largest index as right-handed reduction as the removed index corresponds to the right-most or left-most cell in the set.

For reduction of a correlation, the most restrictive set of conditions exists when the index to be removed from the set is among three neighbouring cells. Adding further cells to the set will not introduce more conditions as there is no overlap between the neighbourhoods of the cell to be removed from the set and any other cells that may be added. As reduction can be performed from the left or right, there are two possible choices

$$\mathbb{E}[\tau'_0\tau'_1\tau'_2] = \mathbb{E}[\tau'_0]\mathbb{E}[\tau'_1\tau'_2], \quad (18)$$

or

$$\mathbb{E}[\tau'_0\tau'_1\tau'_2] = \mathbb{E}[\tau'_0\tau'_1]\mathbb{E}[\tau'_2]. \quad (19)$$

Substituting in indicator variable expressions for all of the τ'_i into these two expressions leads directly to (2) and (3), the conditions for random behaviour.

Correlation current

Some of the random automata we have seen show a flow of some sort. Rather than injecting randomness through probabilistic rules, rule 170 copies the state of an adjacent cell which is guaranteed to be random and independent

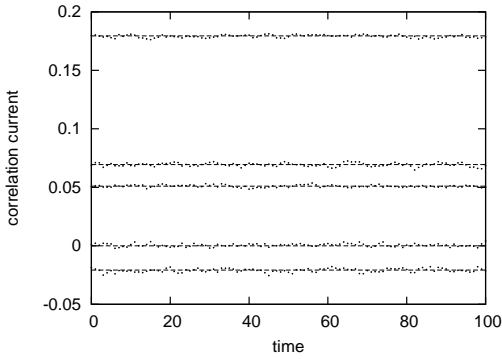


FIG. 4: Measurements of correlation current each taken from 100 iterations of a 10^5 cell cellular automata. The rule sets used are the same as those used to generate Fig. 3. Simulation results are represented by points and predictions by the dotted lines.

of the history of the updated cell. Randomness requires that a cell's state is independent of its previous states so this copying is allowed. The flow seen in many of the automata can be quantified by defining a spatio-temporal correlation between an updated cell and its left- or right-sided neighbour before the update. We define this as *correlation current*.

The magnitude of right-handed flow is given by

$$\mathbb{E}[\tau_0 \tau'_1] - p^2 = (a_4 - p)p(1-p)^2 + (a_5 - p)p^2(1-p) + (a_6 - p)p^2(1-p) + (a_7 - p)p^3, \quad (20)$$

and left-handed by

$$\mathbb{E}[\tau_1 \tau'_0] - p^2 = (a_1 - p)p(1-p)^2 + (a_3 - p)p^2(1-p) + (a_5 - p)p^2(1-p) + (a_7 - p)p^3. \quad (21)$$

Correlation current can be left- or right-handed depending on which neighbour is considered, and can be positive or negative. The sign and direction are unconnected, a negative right-handed correlation current and a positive left-handed correlation current are not the same thing. A negative correlation current is anti-correlation a positive one is correlation.

Random cellular automata permit right-handed correlation current or left-handed correlation current, but not both. Satisfying both sets of conditions, (2) and (3), means that no correlation current exists. Analytical predictions for correlation current are compared with simulation results in Fig. 4.

Information flow

The information flow through a channel with input Z and output Y is given by the mutual information $I(Z; Y)$:

$$I(Z; Y) = H(Y) - H(Y|Z). \quad (22)$$

Where $H(Z)$ is the Shannon entropy of the random variable Z and the conditional entropy $H(Y|Z)$ is given by $H(Y|Z) = H(Z, Y) - H(Z)$ [34].

We consider the eight possible cellular automata neighbourhoods as our Z . We label them in a similar manner to the rules so that $z_0 = 000, z_1 = 001, \dots, z_7 = 111$. In random cellular automata, the probability of one of these neighbourhoods occurring is given by $P(z_i) = p^{w_i} q^{3-w_i}$ where w_i is the number of coloured cells z_i and $q = 1 - p$.

In neighbourhood i , a coloured cell is produced with probability a_i and an uncoloured cell with probability $1 - a_i$, which we denote by \bar{a}_i . The information flow $F(p, \mathbf{a})$ for a random automaton is given by:

$$F(p, \mathbf{a}) = \sum_i P(z_i) \left(a_i \log_2 \frac{a_i}{p} + \bar{a}_i \log_2 \frac{\bar{a}_i}{q} \right). \quad (23)$$

We can write the total entropy of the system as

$$H_{total} = H_{flow} + H_{rules}, \quad (24)$$

where the various entropies are given by:

$$H_{total} = -p \log_2 p - q \log_2 q, \quad (25)$$

$$H_{flow} = F(p, \mathbf{a}), \quad (26)$$

$$H_{rules} = \sum_i -p(z_i)(a_i \log_2 a_i + \bar{a}_i \log_2 \bar{a}_i). \quad (27)$$

This relationship tells us how much of the total entropy (H_{total}) at a particular cellular automata site arises from the probabilistic nature of the set of rules (H_{rules}) and how much is due to information flow from the neighbouring cells (H_{flow}).

Conclusion

We have presented analytical conditions for the existence of random cellular automata with both probabilistic and deterministic rules. The techniques developed, in addition to giving a deeper understanding of random behaviour, can be more broadly applied to analytical solution of cellular automata models.

Randomness is equivalent to the mean-field approximation employed by [17] and [19] which was extended to a broader approximation known as local structure theory [35]. Within limits, both approximations have had considerable success. Analytical limits on the validity of approximation will allow more sophisticated approximation techniques to be developed and applied appropriately.

The indicator variable approach used is readily generalisable to larger neighbourhoods and dimensionalities allowing the techniques employed in this paper to be used to find conditions for random behaviour in larger, more complex automata.

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- [1] M. Gardner, *Scientific American*, **223** 120 (1970).
- [2] Xiao-Bai Li, Rui Jiang, and Qing-Song Wu, *Phys. Rev. E*, **68**, 016117 (2003).
- [3] Xiaobai Li, Qingsong Wu, and Rui Jiang, *Phys. Rev. E*, **64**, 066128 (2001).
- [4] E. Brockfeld et al, *Phys. Rev. E*, **64**, 056132 (2001).
- [5] Rui Jiang, Qing-Song Wu and Bing-Hong Wang, *Phys. Rev. E*, **66**, 036104 (2002).
- [6] V. Belitsky et al, *J. Phys. A*, **40** 11221 (2007).
- [7] Satoshi Sakai et al, *J. Phys. A*, **39** 15327 (2006).
- [8] Wei Zhang et al, *J. Phys. A*, **39** 9127 (2006).
- [9] Quan-Xing Liu, Zhen Jin, and Mao-Xing Liu, *Phys. Rev. E*, **74**, 031110 (2006).
- [10] R. M. Zorzenon dos Santos and S. Coutinho, *Phys. Rev. Lett.*, **87**, 168102 (2001).
- [11] T. G. Mattos et al, *J. Phys. A*, **40** 13245 (2007).
- [12] T. Tome et al, *J. Phys. A*, **40** 12901 (2007).
- [13] E. Arashiro et al, *J. Phys. A*, **40** 887 (2007).
- [14] M. J. Simpson et al, *Phys. Rev. E*, **76**, 021918 (2007).
- [15] C. J. Boulter and G. Miller, *Phys. Rev. E*, **71**, 016119 (2005).
- [16] G. Qiu, D. Kandhai, and P. M. A. Sloom, *Phys. Rev. E*, **75**, 046116 (2007).
- [17] S. Wolfram, *Rev. Mod. Phys.*, **55** 601 (1983).
- [18] S. Wolfram, *Physica D*, **10** 1 (1984).
- [19] L. S. Schulman and P. E. Seiden *J. Stat. Phys.*, **19**, 293 (1978).
- [20] W. Just, *Phys. Rev. E*, **74**, 046209 (2006).
- [21] F. Bagnoli, F. Franci and R. Rechtman, *Phys. Rev. E*, **71**, 046108 (2005).
- [22] F. Bagnoli and N. Boccara, *Phys. Rev. E*, **63**, 046116 (2001).
- [23] A. P. F. Atman, R. Dickman, and J. G. Moreira, *Phys. Rev. E*, **67**, 016107 (2003).
- [24] P. Grassberger *J. Phys. A*, **22**, L1103, (1989).
- [25] Gyrgy Szab and Istvn Borsos 2002 *J. Phys. A*, **35**, L189 (2002).
- [26] N. Israeli and N. Goldenfeld, *Phys. Rev. E*, **73**, 026203 (2006).
- [27] H. Fuks, *Phys. Rev. E*, **66**, 066106 (2002).
- [28] H. Fuks *Nonlinearity*, **17**, 159 (2004).
- [29] J. V. Kujala and T. J. Lukka, *Phys. Rev. E*, **65**, 026115 (2002).
- [30] The eight rule sets used to generate the random probabilistic cellular automata, shown in Fig. 3, listed from a_7 to a_0 are:
 0.833, 0.167, 0.698, 0.302, 0.937, 0.063, 0.968, 0.032;
 0.018, 0.982, 0.749, 0.251, 0.982, 0.018, 0.251, 0.749;
 0.639, 0.361, 0.639, 0.361, 0.639, 0.361, 0.639, 0.361;
 0.886, 0.114, 0.886, 0.114, 0.031, 0.969, 0.031, 0.969;
 0.240, 0.760, 0.964, 0.036, 0.240, 0.760, 0.964, 0.036.
- [31] <http://www.jbcoe.net/#software>
- [32] B. Derrida et al *J. Phys. A*, **26**, 1493 (1993).
- [33] M. Kirkpatrick, T. Johnson and N. H. Barton *Genetics*, **161**, 1727 (2002).
- [34] D. J. C. MacKay *Information Theory, Inference and Learning Algorithms*, Cambridge University Press (2003).
- [35] H. A. Gutowitz et al *Physica D*, **28**, 18 (1987).