Self-organized model of cascade spreading

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(Dated: June 16, 2010)

We study simultaneous price drops of real stocks and show that for high drop thresholds they follow a power-law distribution. To reproduce these collective downturns, we propose a self-organized model of cascade spreading based on a probabilistic response of the system elements to stress conditions. This model is solvable using the theory of branching processes and the mean-field approximation. For a wide range of parameters, the system is in a critical state and displays a power-law cascade-size distribution similar to the empirically observed one. We further generalize the model to reproduce volatility clustering and other observed properties of real stocks.

I. INTRODUCTION

Cascade spreading is an important emergent property of various complex systems. Real life examples of cascades are numerous and range from infrastructure failures and epidemics to traffic jams and cultural fads [1, 2]. Theoretical models of cascades usually assume that agents can be in one of two states (healthy or failed) and an agent's failure puts some stress on its neighbors which may consequently fail too. See [3] for a recent survey of this field offering a novel unifying view.

In this paper we focus on cascades in economic systems which can be identified with stock prices suddenly dropping in a major market crash [4] or with companies going bankrupt simultaneously and leading to global recession [5]. Corresponding theoretical models are based on shortage and bankruptcy propagation in production networks [6], default propagation in credit networks [7, 8], interaction of firms through one monopolistic bank [9] or in a complex credit network economy [10], and herding behavior of traders [11, 12]. While these models help us to understand cascade processes in economic systems, they are mostly too involved to allow for analytical solutions—their study relies on numeric simulations and agent-based modeling [13].

A simpler point of view on cascade phenomena is offered by the concept of self-organized criticality (SOC) which has had a deep impact on the science of complexity. First introduced more than twenty years ago to explain the ubiquitous 1/f noise [14], it caused a blossoming of toy models, computer simulations, and real life experiments [15]. The analytical techniques employed include scaling arguments [16], mean-field theories [17], renormalization methods [18, 19], and rigorous algebraical techniques [20].

SOC is a mechanism which explains the emergence of complex behavior in many diverse real world systems [21, 22]. The generic behavior of SOC models is: (a) they evolve so that they always stay close to the critical point, (b) long periods of robustness and moderate activity are interrupted by sudden breakdowns. This indeed qualitatively resembles "stock markets which expand and grow on relatively long time scales but contract in stockmarket crashes on relatively short time scales" [15] and "stock crashes caused by the slow buildup of long-range correlation leading to a global cooperative behavior of the market eventually ending into a collapse in a short time interval" [4]. This similarity provides motivation for the present study.

We begin our work with an empirical investigation of simultaneous price drops of real stocks and show that the size distribution of observed events is broad (for high drop thresholds it follows a power-law distribution). This observation confirms that simultaneous stock downturns are a collective phenomenon. We propose a dynamical model which for a wide range of parameters selforganizes into a critical state. Unlike most SOC models, our model assumes a probabilistic response mechanism where a node has only a certain probability of reacting to the current stress conditions. The basic idea behind modeling simultaneous stock downturns with cascades is that decline of a single stock may provoke investors' reactions which consequently may cause other stocks to decline and the "infection" to spread. The key premise is that while failed nodes become significantly more resistant in the next time step, healthy nodes become slightly less resistant. This close parallel with the slow growth/fast decay picture described above is further supported by our analysis of empirical data which shows that majority of stocks behave in this way.

The basic proposed model has the advantage of being simple, analytically solvable in some cases, and easily generalizable to more complicated settings. We analyze it using the formalism of branching processes, the mean-field approximation and, for complex topologies of nodes' interactions, using numerical simulations. Obtained cascade-size distributions exhibit a close similarity to our empirical observations. Introduction of memory within the model allows us to reproduce other empirically observed features, such as volatility clustering, though at the cost of analytical tractability. We conclude our study with a discussion of further model's generalizations and possible areas of application.

II. EMPIRICAL DATA

Here we investigate co-occurring price movements of real stocks. Adopting the vocabulary of cascade models,



FIG. 1. The empirical size distribution observed with real stock prices for thresholds H = 0% and H = 10%.

we say that a stock fails when the relative loss of its price over a given time interval Δt exceeds a certain threshold H. Denoting the price of stock i at time t as $p_i(t)$, its failure occurs when $[p_i(t) - p_i(t + \Delta t)]/p_i(t) > H$. The number of stocks failing at time t, $n_F(t)$, is a direct analog of the cascade size in a model of cascade spreading. As the input data we use daily prices (hence $\Delta t = 1$ day) of 500 stocks from the standard U.S. index S&P 500 (this data is freely available at, for example, finance.yahoo.com). To achieve a fixed system size, we consider only those 332 companies which are in the stock market since the beginning of 1992 and use their prices during the 18-years long period ending in May 2010 for our analysis.

The empirical distribution of failure sizes is shown in Fig. 1 for H = 0% and H = 10%. We see that for this large value of H (which is in line with the notion of stock *failures*), the observed size distribution has a powerlaw shape. Using the methodology described in [23], we obtained the power-law exponent 2.19 ± 0.05 with the lower bound for the power-law behavior $n_{\min} = 3$. The corresponding p-value (obtained using the standard Kolmogorov-Smirnov statistic) is 0.92, which confirms that the data is consistent with the hypothesis of a power-law distribution. Similar results are obtained also for other threshold values so long as $H \gtrsim 8\%$. When $H \lesssim 8\%$, the resulting size distributions are broad but probably not power laws. Finally, when H = 0% (*i.e.*, any price drop is interpreted as a failure), the size distribution is roughly symmetric around the value corresponding to one half of the system size (see Fig. 1).

The power-law shape itself suggests that big stock downturns occurring on the same day are rather a collective phenomenon than independent events. This hypothesis is further supported by the average correlation of simultaneously failing stocks, 0.35 (again including only events with at least three simultaneously failing stocks), which is significantly higher than the overall average stock correlation, 0.25. Another sign of a strong connection among simultaneously failing stocks comes from their known division to ten different industrial sectors. The effective number of sectors participating in a cascade is defined as

$$e = \left(\sum_{i=1}^{10} r_i^2\right)^{-1}$$

where r_i is the share of sector *i* in the cascade and $\sum_{i=1}^{10} r_i = 1$. By averaging this quantity over all cascades of a given size *S*, we obtain e(S). This number can be compared with the effective number of sectors corresponding to selecting failed stocks at random, e'(S). The analysis of stock prices shows that for any S > 3, e(S) is significantly smaller than e'(S) which implies that simultaneous stock failures preferentially affect strongly connected stocks in one sector or in a small number of sectors.

The second part of our analysis of empirical data is devoted to time correlations of consecutive failures. The autocorrelation of the number of failing stocks with the time lag one day $C(n_F(t), n_F(t+1)) \approx 0.15$ is comparable with the autocorrelation of absolute returns $C(|r(t)|, |r(t+1)|) \approx 0.25$. Both values are signs of volatility clustering which is known to appear in financial data [24]. We further estimate conditional failure probabilities for individual stocks: for example, P(F|N)denotes stock's failure probability given that this stock didn't fail in the previous time step. When the results are averaged over all stocks, we obtain P(F|F) = 0.039which is much higher than the overall failure probability P(F) = 0.003—this is another sign of volatility clustering in our data. On the level of individual stocks, however, 62% of all stocks with at least three failures strongly satisfy the inequality P(N|F) > P(N) which is equivalent to P(F|F) < P(F) (because P(F|F) + P(N|F) = 1). (By strong satisfying we mean that the difference of the two probabilities is greater than the sum of their uncertainties.) We see that despite volatility clustering in the data, most stocks are more "resistant" to failures after they have just undergone one. For the remaining stocks, probabilities P(N|F) and P(N) either differ less than the sum of values' uncertainties (for 14% of stocks) or even strongly satisfy the opposite inequality P(N|F) < P(N), with corresponding values of P(F|F) often as high as $0.30 \ (24\% \text{ of stocks})$. We can conclude that while after a failure, most stocks become more resistant to another failure, there is a fraction of stocks which are prone to consecutive failures.

III. BASIC MODEL AND ITS MEAN-FIELD SOLUTION

In this section we present a simple model which is amenable for analytical treatment and qualitatively reproduces some of the features observed in empirical data. In its basic formulation, this model is particularly suitable for stocks that, as discussed in the previous section, after a failure become more robust. Generalization of the model aiming at reproducing other observed features (volatility clustering, for example) is discussed in Section IV.

Consider a system of N nodes where node i (i = 1, ..., N) has two possible states: failed $(i \in \mathcal{F})$ and healthy $(i \notin \mathcal{F})$. With each node i we further associate fragility $f_i \in [0, 1]$ which measures how this node reacts to failures of its neighbors (the higher the fragility, the more likely is the node to follow a neighbor's failure). The dynamics of the model is governed by the following simple rules. (i) In each time step, the first failed node ("trigger") is chosen at random and may induce failures of other nodes. (ii) If a neighbor of node i fails, node ifollows it with the probability f_i . The cascade of failures propagates until all remaining nodes resist the damage. (iii) At the end of the time step, fragilities of all nodes are updated according to

$$f_i(t+1) = \begin{cases} \lambda f_i(t) & i \in \mathcal{F} \\ (1+\beta)f_i(t) & i \notin \mathcal{F} \end{cases}$$
(1)

where $0 < \beta \ll 1$ and $\lambda \in (0,1)$ are parameters of the model (in effect, failed nodes become less fragile and healthy nodes become slightly more fragile). All values $f_i(t+1) > 1$ are truncated to 1 (this may occur when the parameter β is large). After this update is finished, all nodes are again marked as healthy, the current time step ends and the new one begins with point (i). Note that unlike some other models of cascade spreading, failed nodes are not removed from the system in our case. If a long enough equilibration period is applied before measuring the system behavior, the initial fragility values $f_i(0)$ are of little importance (see Section III E for a detailed discussion); we set them randomly in the range (0; 1) in all our simulations (unless stated otherwise).

According to the rules above, when n neighbors of node i fail, node i resists with the probability $(1-f_i)^n$ and fails with the complementary probability

$$P_F(f_i, n) = 1 - (1 - f_i)^n.$$
(2)

This response to failures is "path-independent" in some sense: the probability that a node resists failure of n of its neighbors, $(1 - f_i)^n$, is the same as the probability of resisting two consequent waves of failures of x and n - x neighboring nodes, $(1 - f_i)^x (1 - f_i)^{n-x}$.

We simplify the system by assuming that interactions of all nodes are equally strong (the general case will be studied in Section IIID). This renders the notion of "node's neighbors" superfluous because each node's failure affects all remaining healthy nodes in the system. Now assume that after the initial failed node is chosen, n_1 nodes respond to this failure and fail too. Each of the remaining $N - n_0 - n_1$ nodes (here $n_0 = 1$ is the initial number of failed nodes) then has some n_1 -dependent failure probability which results in n_2 new failures, and so on, until in iteration m, $n_m = 0$ is achieved. The cascade size is then defined as the total number of failures,



0.010

 $P_{F} 0.005$

0

 β FIG. 2. Average failure probability P_F : analytical (solid black line), approximate analytical (dashed blue line) and numerical (symbols) results for $N = 10^3$, $\lambda = 0.1$, averaged over 10^6 time steps. The vertical dotted line indicates β_0 obtained with Eq. (5).

0.01

 $S = n_0 + \cdots + n_m$, and node fragilities are consequently updated according to Eq. (1). Since cascade sizes are limited by the system size, $S \leq N$.

The dynamics of the system, based on a failure propagation process and fragility updating, is fully contained in the three described simple rules. In the following paragraphs we shall study when these rules drive the system to the critical point and how this self-organized criticality manifests itself in the distribution of cascade sizes P(S).

A. Failure probability

Let P_F be the average failure probability of a given node in one time step (or, equivalently, the average fraction of failed nodes in one time step). Assuming that t_{eq} is some sufficiently long time after which the system is equilibrated (we use $t_{eq} = 10^4$ for all our simulations), fragility values averaged over time, $\langle f_i \rangle$, do not evolve anymore. All nodes interact equally strongly, hence $\langle f_i \rangle$ is independent of *i* and it can be replaced with $\langle f \rangle$. Since in a large number of time steps *T* each node undergoes $P_F T$ failures and $(1 - P_F)T$ non-failures, Eq. (1) implies

$$\langle f(t_{\rm eq} + T) \rangle = \langle f(t_{\rm eq}) \rangle \lambda^{P_F T} (1 + \beta)^{(1 - P_F)T}.$$
 (3)

Using the equilibrium condition $\langle f(t_{eq} + T) \rangle = \langle f(t_{eq}) \rangle$, we can solve this equation with respect to P_F to get

$$P_F(\beta,\lambda) = -\frac{\ln(1+\beta)}{\ln\frac{\lambda}{1+\beta}} \approx -\frac{\beta}{\ln\lambda}$$
(4)

where the approximate result is valid for $\beta \ll 1$ (see Fig. 2 for a comparison of these results with numerical simulations).

A node may fail because it is selected as the first failed node (with probability 1/N) or due to failure propagation

0.02

(with probability P_P); P_F thus can be written as $P_F = 1/N + P_P$. Since the value of P_F depends solely on β and λ , $P_P = P_F - 1/N$ may be negative for a small system which is, of course, impossible in practice. This situation occurs when for given λ , N, the value of β is smaller than a certain threshold β_0 and hence it does not suffice to compensate for the fragility decay due to λ . Eq. (3) then has only the trivial solution $\langle f \rangle = 0$ and hence $P_F(\beta, \lambda) = 1/N$ (failures do not spread). When β is small, the approximate form of P_F can be used to solve this equation with respect to β and we get

$$\beta_0 \approx -\frac{\ln \lambda}{N} \tag{5}$$

which agrees with numerical simulations (see Fig. 2). Note that if the number of initial failed nodes grows with the system size as wN ($w \ll 1$), we get $\beta_0 \approx -w \ln \lambda$ which is independent of N.

When model parameters are set to extreme values (for example, $N = 10^3$, $\beta = 10^3$, $\lambda = 10^{-3}$), the system exhibits unusual modes of behavior where active turns (with nearly all nodes failed) alternate with calm turns (with nearly all nodes healthy). While Eq. (4) holds also in such conditions, our further analysis focuses on $\beta \ll 1$ which renders more realistic behavior.

B. Average fragility

When $nf_i \ll 1$, Eq. (2) can be approximated with $P_F(f_i, n) = nf_i$ which can be interpreted as independence of stress inflicted by n individual failed nodes. This further means that each failed node has its failing descendants independently of other failed nodes and hence one can use the theory of branching processes [25] to describe the cascade spreading. Note that by use of this theory we implicitly assume that the system size is infinite. For a discussion of the finite-size effects on the size of an epidemic outbreak see [26].

As already mentioned, when interactions of all nodes are equal, $\langle f_i \rangle$ is independent of *i*. If we further neglect fluctuations of f_i , then all nodes have identical fragility $\langle f \rangle$. This is a mean-field-like approximation which replaces the exact cascade spreading with cascade spreading in a homogeneous averaged medium. Since the number of direct descendants now follows a simple binomial distribution with mean $N\langle f \rangle$, we can use elementary results of branching process theory to express the average cascade size (the total progeny) as $\langle S \rangle = 1/(1 - N\langle f \rangle)$. Further, using $\langle S \rangle = NP_F(\beta, \lambda)$ we obtain the average fragility

$$\langle f \rangle = \frac{1}{N} \left(1 - \frac{\ln\left[(1+\beta)/\lambda\right]}{N\ln(1+\beta)} \right). \tag{6}$$

Since $\beta > 0$ and $\lambda < 1$, $\langle f \rangle$ is always less than 1/N. Comparison with numerical simulations (not shown) confirms that Eq. (6) is valid only for $\beta \ll 1$.

C. Cascade size distribution

The size distribution for a critical branching process is known to have the power-law form $P(S) \sim S^{-3/2}$ (see, for example, [27]). However, critical cascades represent only one of possible modes of model's behavior and hence we briefly discuss the general size distribution $P(S|\beta, \lambda)$. According to a theorem from [28], if the generating function for the number of direct descendants d is $\pi(x)$, the total progeny of the resulting branching process Y has the distribution

$$P(Y|n_0) = \frac{n_0}{Y} p_{Y-n_0}^{(Y)}$$
(7)

where $p_a^{(b)}$ is defined using

$$[\pi(x)]^b = p_0^{(b)} + p_1^{(b)}x + \dots$$
(8)

and n_0 is the number of ancestors (in our case, the number of initial failed nodes). Since *d* obeys a binomial distribution its generating function is simply $\pi(x) = (1 - \langle f \rangle + \langle f \rangle x)^N$ and we get

$$P(S|\beta,\lambda) = \frac{1}{S} {NS \choose S-1} \langle f \rangle^{S-1} (1 - \langle f \rangle)^{NS-S+1}$$
(9)

where we used $n_0 = 1$ and $\langle f \rangle$ is given by Eq. (6). Note that the resulting probability is positive for S > N which contradicts the model assumptions (each node fails at most once in a given turn). This is a direct consequence of using the theory of branching processes which assumes that the system size is infinite. This problem is of little importance for small values of β when the obtained values of P(S) are negligible for S > N.

When $1 \ll S \ll N$, Eq. (9) can be approximated with

$$P(S|\beta,\lambda) = \frac{(N\langle f \rangle)^{S-1} \mathrm{e}^{S(1-N\langle f \rangle)}}{\sqrt{2\pi} S^{3/2}}.$$
 (10)

According to Eq. (6), $\lim_{N\to\infty} N\langle f \rangle = 1$ for any given β, λ and hence in the limit of large system size is $P(S|\beta,\lambda) \sim S^{-3/2}$ which corresponds to the critical branching process mentioned before. For any finite-size system, the smaller the value of β , the larger the value of $1-N\langle f \rangle$. Consequently, the power-law scaling holds only for $S \ll \beta N$ (this agrees with Fig. 3 where for $\beta = 10^{-3}$, the power-law behavior disappears at $S \approx 10$). On the other hand, the range of β and λ for which the system self-organizes to a critical state is wide and we can say that this is an SOC system.

A comparison of the obtained analytical results with numerical simulations is shown in Fig. 3. The agreement is good for small values of β ($\beta \leq 0.01$) and the initial slope of the distributions (before the finite size effects become apparent) is close to -3/2. Results obtained with $\beta = 0.001$ confirm that when β is small enough, P(S)decays faster than as a power law. When β is large, true P(S) deviates from the analytical prediction and exhibits



FIG. 3. Cascade size distribution: numerical results (color lines), analytical results according to Eq. (9) (dashed lines) and the power-law decay with exponent -3/2 (thick solid line) for $N = 10^4$, $\lambda = 0.1$, 10^7 time steps, and $\beta = 0.001$ (red line, fastest decay), $\beta = 0.01$ (green line, medium decay), $\beta = 0.1$ (blue line, slowest decay). The analytical solution is not plotted for $\beta = 0.1$ because it is very similar to that for $\beta = 0.01$.

a secondary maximum at a large size value—this effect is well visible in Fig. 3 for $\beta = 0.1$. This maximum, formally simply a super-critical phase of the model, resembles the so-called meaningful outliers discussed in [29]. To estimate the value of β at which the secondary maximum appears and Eq. (9) ceases to hold, we take the average number of failures computed both from Eq. (9) and from Eq. (4). By comparing the two results we obtain

$$NP_F(\beta, \lambda) = \sum_{S=1}^{N} SP(S|\beta, \lambda).$$
(11)

When β is small, both sides of this equation depend on β and the equality can hold. However, Eq. (10) shows that when β is sufficiently large, the size distribution is approximately power-law and it does not change with β anymore. As we increase β further, the power-law distribution does not suffice to provide enough failures and for Eq. (11) to hold, an additional contribution must appear on the rights side. The value β^* when this happens can be found by substituting $P(S) \sim S^{-3/2}$ on the right side and approximating the summation with integration. When N is large, we obtain

$$\beta^* \approx -\ln\lambda \sqrt{\frac{2}{\pi N}} \tag{12}$$

which complements the previously found threshold β_0 given by Eq. (5). For $N = 10^4$ and $\lambda = 0.1$, we obtain $\beta^* \approx 0.02$ which agrees with our empirical observation ($\beta \leq 0.01$ for Eq. (9) to hold) above.

Finally, by comparing the empirical observations presented in Fig. 1 with the obtained analytical results, we can conclude that the presented model exhibits qualitative agreement with the studied real system.

D. Generalizations

To test how robust are the obtained results, we consider simple generalizations of the proposed model. First of all, when the multiplicative fragility update rule Eq. (1) is replaced by an additive one, the behavior of the system does not change considerably. The second generalization relates to the assumed even influence of a node's failure on all the remaining nodes. Denoting the strength of failure propagation from node i to node j as $C_{i,j}$, the probability that node j fails as a result of i's failure can be generalized to $C_{i,j}f_j$. The probability that node j fails as a result of a group \mathcal{F} of failed nodes (given by Eq. (2) before) generalizes to the form

$$P_F(f_j, \mathcal{F}) = 1 - \prod_{i \in \mathcal{F}} (1 - C_{i,j} f_j).$$

$$(13)$$

Matrix ${\sf C}$ encodes the structure of the network of node interactions.

When the elements $C_{i,j}$ are drawn independently from a given distribution and the system size is large, the mean-field approximation is again appropriate to describe the system behavior and the power-law size distribution with exponent 3/2 results. Similarly when C contains a block structure with inter-block elements drawn from a different distribution than intra-block elements (this mimics the well-known sector structure of the stock correlation matrix), the original power-law size distribution remains largely unchanged (unless either the block division of C or one of the two probabilistic distributions are such that they do not allow to use the mean-field approximation). Analogous behavior results from the "random neighbor approximation" in which node's neighbors are chosen anew repeatedly (see [30] for this kind of analysis of a different model).

When all elements $C_{i,j}$ are either zero or one, matrix C can be represented by a network and a complex topology of node interactions can be introduced by using some of the standard network models [31]. We studied two different types of networks: the Erdös-Rényi network where $C_{i,j} = 1$ with probability p and $C_{i,j} = 0$ otherwise and the growing Barabási-Albert network where each new node is attached to I old nodes. (These two kinds of networks are structurally very distinct as the former consists of nodes of approximately identical degree and the latter exhibits a power-law degree distribution.) Numerical results for both cases are shown in Fig. 4. As expected, for the Erdös-Rényi network with p > 1/N, the size distribution exponent remains unchanged. When p < 1/N, the network consists of small unconnected components and hence big cascades cannot occur. The irregular size distribution P(S) observed for $\beta = 5 \cdot 10^{-5}$ is due to topological properties of the particular network realization where the model was simulated (*i.e.*, positions of respective ups and downs of the size distribution depend on the network realization). These results agree with a previous study of the sandpile dynamics [32] (see [33] for



FIG. 4. The cascade size distribution on complex networks: (a) sparse Erdös-Rényi networks with $p = 5 \cdot 10^{-5}, 10^{-4}, 10^{-3}$ (red, green, and blue line, respectively; the indicative thick line has slope 1.5), (b) Barabási-Albert networks with I = 1and I = 10 (red and blue line, respectively; the indicative thick line has slope 1.65). Parameters of the system: $N = 10^4$, $\beta = 0.005, \lambda = 0.1, 10^7$ time steps.

an extensive recent review of critical phenomena in complex networks). By contrast, Barabási-Albert networks yield cascade size distributions with significantly higher exponents (approximately 1.65) which is probably due to strong inhomogeneity of the network. When I = 1, the size distribution deviates from a power law, probably as a consequence of the scale-free network topology (the same shape of the distribution is observed for different realizations of the network).

E. Role of the initial fragility values

While it sounds plausible that due to model's stochasticity, the initial fragility values have no influence on the equilibrium fragility distribution, the situation is in fact more complicated. For example, a simple numerical simulation with $f_i(0) = 1/N$ for all *i* shows a case where: (i) no stationary fragility distribution arises, (ii) at any time step, only a small number of distinct fragility val-



FIG. 5. Fragility distributions at different time steps (the initial fragility values are set to 1/N, $N = 10^4$, $\beta = 0.001$, $\lambda = 0.1$).

ues is observed (see Fig. 5). What causes the discreteness of fragility values? Denoting the number of failing and healthy time steps of node i as F_i and H_i , respectively, it must hold that $F_i + H_i = t$ where t is the current time step. This node's fragility now can be written as

$$f_i(t) = f_i(0)(1+\beta)^t \left[\lambda/(1+\beta)\right]^{F_i}.$$
 (14)

When all $f_i(0)$ are identical, the possible values of $f_i(t)$ are discrete at any time step t and the ratio of neighboring possible values is $(1 + \beta)/\lambda$. If λ is small (as it is in our simulations), this ratio is large and hence the number of actually observed fragility values is small (because values much smaller or greater than the average fragility are unlikely). Eq. (14) implies that possible fragility values depend on t and hence there can be no stationary fragility distribution—this is confirmed by Fig. 5 where fragility peaks constantly shift to higher values and change their relative heights. Interestingly, even this peculiar setting of $f_i(0)$ does not alter the long-term model's behavior substantially and the aggregate quantities (such as the average failure probability or the cascade size distribution) are similar to those found for randomized initial fragility values before.

Differences between neighboring peaks are $\lambda/(1 + \beta)$, hence the time after which the fragility distribution pattern repeats can be estimated as $\ln [\lambda/(1+\beta)]/\ln(1+\beta)$. Since this is a typical time of fragility evolution, one can use it also as an estimate of the initial equilibration time T_{eq} . For the smallest value of β in our simulations $(\beta = 0.005)$ we obtain $T_{eq} \approx 4\,600$ which ex post confirms our setting of the equilibration time to 10^4 . Finally, note that while the random setting of $f_i(0)$ prevents discrete fragility values from appearing, some remnants of the initial fragility values can be preserved by Eq. (14). To obtain a fragility distribution truly independent of the initial values, one has to assume annealed dynamics, *i.e.* fragility updating by randomized values of β and λ .

IV. GENERALIZED MODEL WITH PARTIAL MEMORY

Fragility updating rules defined by Eq. (1) imply that nodes become more robust after a failure and hence autocorrelation of their failures as well as autocorrelation of the total number of failures are negative (their magnitudes depend on β and λ). As discussed in Section II, this is true for majority of stocks but certainly not for all of them. To allow for repeatedly failing stocks, we introduce the probability α with which a failed node stays failed also in the next time step (and consequently acts as an additional initial failed node). This probability has the role of *partial memory* in the system and, as we shall see, gives rise to volatility clustering and other effects observed in real financial data. Note that memory or delayed stress propagation are quite often part of cascade spreading models as in, for example, [34]. In our case it is assumed that fragilities of nodes which stay failed due to α are not updated in the given time step (as long as α is small, this assumption has little influence on the results).

Since $P(F) \ll 1$, the probability of a node's repeated failure is now $P(F|F) \approx \alpha$ which, in combination with the empirical results presented in Section II, allows us to set $\alpha = 0.04$. We further choose $\beta = 0.01$ and $\lambda = 0.1$ which best correspond to the critical regime in Fig. 3. Using this setting we numerically obtain conditional probabilities consistent with those observed in the empirical data: P(F|F) = 0.041 (empirical value is 0.039), P(F) = 0.004 (empirical value is 0.003) and P(F|N) = 0.004 (empirical value is 0.003). As long as we stay in the critical regime, these values depend on β and λ weakly. Presence of volatility clustering is confirmed by the significantly positive autocorrelation of the number of failures $C(n_F(t), n_F(t+1)) \approx 0.3$ (empirical value is 0.15). The precise value depends on α and λ (and much less on β) but strongly positive autocorrelation is always observed. For comparison, $\alpha = 0$ yields $P(F|F) \approx P(F)$ and $C(n_F(t), n_F(t+1)) \approx -0.04$. Finally, Fig. 6 shows P(S) for different values of α . We see that for small values of α , the size distribution remains power-law with exponent k gradually decreasing as α grows. Due to the additional complexity introduced by partial memory, the cascade size distribution for this generalized model has not been obtained yet.

V. DISCUSSION

We studied empirical data with stock prices and found that large simultaneous downturns follow a broad distribution consistent with a power law with exponent 2.19 ± 0.05 . To reproduce this behavior, we proposed a simple stochastic model of failure propagation. Using a mean field approximation and branching process theory we derived the general cascade size distribution and determined the range of parameters which give rise to the



FIG. 6. Cascade size distribution for the modified model: numerical results for $\beta = 0.01$, $\lambda = 0.1$, $N = 10^4$, 10^7 time steps, and $\alpha = 0$ (steepest green line), $\alpha = 0.04$ (medium blue line), $\alpha = 0.08$ (least steep red line).

critical regime. To reproduce other features observed in financial data, such as volatility clustering, partial memory was introduced in the basic model.

Our results still leave some open and challenging questions which deserve further investigation. Firstly, since the cascade sizes corresponding to the secondary maximum in Fig. 3 are comparable with system size, this behavior cannot be described within the formalism of branching processes where an infinite system size is assumed. While we found an approximate condition for the appearance of the secondary maximum, how to proceed further towards an exhaustive description of the resulting size distribution is still an open question. Secondly, it would be interesting to reach an analytical expression for the size distribution exponent in scale-free networks where it appears to differ from the mean-field value 3/2. Thirdly, generalized model with "partial memory", only studied numerically here, calls for analytical approaches. Fourthly, it would be interesting to know whether the model can be modified to produce power-law size distributions with exponents considerably higher than those reported here. One opportunity for such a generalization is to assume a dynamic network structure whose evolution depends on nodes' failures, similarly to the approach used in [35, 36] for different models. Alternatively, as a generalization of the current binary model where nodes are either healthy or failed, one could define a multi or continuous-state model in which the probability of following a neighbor's failure depends on the failure's magnitude.

We stress that the probabilistic spreading mechanism proposed here is a general one and its use is not limited to market crashes or firm bankruptcies. For example, on a two or three dimensional lattice, a similar mechanism might be employed to model earthquakes as the failure at one place of the Earth's crust exerts some stress on its neighborhood (the number of failed nodes would then represent the earthquake size). In summary, the proposed model, together with its generalizations, has proven to be simple yet rich in behavior. It poses a variety of new research questions and we are looking forward to its future development and applications.

ACKNOWLEDGMENTS

This work was partially supported by the Future and Emerging Technologies programme FP7-COSI-ICT of

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