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Critical properties of cluster size distribution in an asymmetric diffusion-aggregation model

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We consider a stochastic dynamics for a system of diffusing hard-core particles on a periodic chain with asymmetric diffusion rules. A given cluster of particles can diffuse to the right as a whole but the particle located on the left boundary of the cluster is allowed to break-off from it and diffuse to the left. Clusters of particles can eventually merge with other clusters. These rules allow for the creation of clusters of different sizes. We discuss the size distribution of the clusters in the long time or steady state limit, as a function of the particle concentration \( c \). We consider the general time dependent master equation based on Smoluchowski’s theory for local cluster merging or fragmentation and diffusion processes, and study the solutions using the generating function in the large size limit. We found that there exists a critical density \( c^* = \sqrt{2} - 1 \) for which the cluster distribution decays like a power law with exponent \( 5/2 \).

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I. INTRODUCTION

Statistical properties of coalescence and fragmentation processes have been studied in various fields of physics and chemistry [1–7] where for example chemical constituents react to form new molecules, or fragment into elementary units. Both aggregation and fragmentation
reactions are governed by stochastic processes, influenced eventually by two-body collision properties [8, 9]. The kinetic of such processes is described by a Smoluchowski’s master equation [10], which is a balance equation for determining the particle distribution in size or mass in the steady or equilibrium state [11, 12], and can be used for example to determine the size of interstellar dust grains [13] or even particle size in Saturn rings [14].

The dynamics is usually controlled by reaction rates depending on the size or mass of the particles, and it is assumed in most of Smoluchowski’s models that both aggregation and fragmentation rates follow a scaling law: they depend intrinsically on the sum or product of the individual mass or size of the merging components with some power exponent. Therefore a general description of the physical properties such as mass or size distribution depends indeed on the microscopic details of those kernels [5, 11, 15, 16]. For example, in the case of gravitational systems, the aggregation rate is proportional to the sum of the constituent masses, with eventually some power law, whereas in the case of branched polymerization they are proportional to the product of the molecule weights. In particular, the size distribution when the rate is constant and when a monodisperse source is introduced exhibits, in addition to an exponential damping factor, a power law with a decay exponent equal to $3/2 + \beta$ [11], $\beta$ being the weight exponent in the kernel.

In this paper we present a dynamical model of hard-core particles on a periodic chain of size $N$ and which can diffuse or fragment from or merge with other clusters. We can view the clusters of particles as polymers confined on a one-dimensional (1D) chain and which can react with the other neighboring clusters by chemical reaction. The dynamics we consider is defined more precisely by the following asymmetric and nonlocal rules: We assume that, at every time step, a particle located at the left boundary of a cluster can break-off from it at rate $1/\tau$, and/or eventually aggregate with another cluster. A whole cluster can move to the right by one elementary step with the same rate, and eventually fuses with a neighboring cluster if present. The schematic process is displayed in Fig. 1. From these two processes emerge clusters of small and large sizes. Diffusion of large clusters in one direction allows them to grow in size by merging with other clusters, whereas their partial fragmentation on the opposite direction increases the number of small clusters. Overall, we expect that the size distribution will display a tail that decays like a power law. Also, the structure of the distribution will depend crucially on the particle concentration. The periodic boundary conditions are important for the asymmetric dynamics we consider, as they allow for the
current of particles, due to the asymmetry of the dynamics, to flow along the chain and particles to be transferred to the other side once they reach a boundary.

The dynamics is microscopically irreversible and does not reach an equilibrium state [17], but there is a steady state as we will see below from which a well defined size distribution exists. This is due in particular to the fact that one-particle clusters are formed constantly by fragmentation of larger clusters, in absence of any source, and therefore there is no absorbing state. A total aggregation process would otherwise lead to a unique cluster, a phenomena known as gelation [18–20] where the largest cluster is the final state.

We can relate this dynamics to a traffic flow problem [21], where for example groups (clusters) of vehicles coherently move to the right, the last vehicle of the groups being allowed to slow down and detach from its cluster. A particle current is present and oriented to the right direction as only one-particle clusters can move to the left. This kind of dynamics can also be seen as a variation of the TASEP (Totally Asymmetric Exclusion Process) model [22], with non-local or coherent motion involved as clusters of particles can move collectively. The exclusion process comes from the fact that there is at most one particle per site (hard core interaction), and therefore individual particles cannot move if the neighboring sites are already occupied, like the ASEP model or Asymmetric Simple Exclusion Process.

The aim of this study is to show that this particular asymmetric fragmentation or merging process induces criticality in the cluster size distribution for a special concentration value. The master equation that we define in the next section is closely related to a non-linear differential-integral Smoluchowski’s equation but with the constraint of dimensionality and space: Distances and therefore concentration between clusters influence the rate of merging.

From a technical point of view, many models based on Smoluchowski’s equation can be solved exactly using generating functions [11, 23, 24] which are an adequate tool to obtain distribution moments. However the confinement effect in the present model renders the analysis more complicated as momentum interaction occurs in the Fourier space.

II. MASTER EQUATION FOR CLUSTER DISTRIBUTION

The stochastic process presented in the Introduction is characterized by a master equation. We first need to define the local probability for each \( n \)-particle cluster at a given time \( t \) and location on the lattice, which is a succession of particles surrounded by two empty
FIG. 1: Elementary processes with three particles. (a) Clusters can move to the right only. (b) The particle located on the left boundary can detach from the cluster and diffuse on the left.

sites:

\[ P_n(r) = \text{Prob}(\circ \bullet \cdots \bullet ^{r+n-1} \circ) \] (1)

where the first particle of the cluster is located at site \( r \in \{1, \cdots, N\} \). The system is periodic with \( P_n(r + N) = P_n(r) \), the number of clusters of size \( n \) is \( N_n = \sum_r P_n(r) \), and the total number of clusters \( N_c = \sum_n N_n \). The number of particles is conserved, and equal to

\[ N_p = \sum_r \sum_n nP_n(r) \] (2)

and their concentration is denoted by \( c = N_p/N \), with maximum cluster or cutoff size \( N_p \).

There are different processes which contribute to the change of probability \( P_n(r) \) after a small time interval \( \Delta t \). In the following, we note \( \partial P_n/\partial t \) the rate of change of the probability when \( \Delta t \to 0 \). The first process (a) is the decay or destruction of the cluster configuration by breaking-off and diffusion of the leftmost particle or diffusion of the whole cluster to the right

\[ \left( \frac{\partial P_n}{\partial t} \right)_a = -\text{Prob}(\circ \bullet \cdots \bullet ^{r+n-1} \circ) - \text{Prob}(\circ \bullet \cdots ^{r+n-1} \circ) \] (3)

Another decay is provided by the aggregation of neighboring clusters on the left, and particles on the right, processes (b) and (c)

\[ \left( \frac{\partial P_n}{\partial t} \right)_{b+c} = -\text{Prob}(\times \circ \bullet \cdots ^{r+n-1} \circ) - \text{Prob}(\circ \bullet \cdots ^{r+n-1} \circ) \] (4)

where the symbol (\( \times \)) represents a site occupied or not by a particle. The positive contributions to \( P_n \) come from a series of two-clusters aggregations by avalanches, process (d)

\[ \left( \frac{\partial P_n}{\partial t} \right)_d = \text{Prob}(\circ ^{r-1} \circ \bullet \cdots \bullet ^{r+n-1} \circ) + \text{Prob}(\circ ^{r} \circ \circ \circ \cdots \bullet ^{r+n-1} \circ) + \cdots + \text{Prob}(\circ ^{r-1} \circ \circ \circ \cdots \bullet ^{r+n-1} \circ) \] (5)
The cluster can also be generated by one-particle diffusion from the right, process \((e)\)
\[
\left( \frac{\partial P_n}{\partial t} \right)_e = \text{Prob}(\circ \bullet \ldots \circ \bullet \circ \circ \times) \quad (6)
\]
Process \((f)\) is the avalanche contribution of the same cluster originally located at \(r - 1\) and moving to the right
\[
\left( \frac{\partial P_n}{\partial t} \right)_f = \text{Prob}(\circ \bullet \ldots \circ \bullet \circ \circ) \quad (7)
\]
The last process \((g)\) consists in creating the cluster from a cluster of size \(n + 1\) from which the first particle separates by diffusing to the left hand side
\[
\left( \frac{\partial P_n}{\partial t} \right)_g = \text{Prob}(\circ \bullet \ldots \circ \bullet \circ \circ) \quad (8)
\]
The analytical expressions of the previous processes are given by products of cluster size distributions and their convolution, in the case where several clusters are contributing to the dynamics, and which are detailed in Appendix A. Overall, all the changes in \(P_n(r)\) can be summed up and we obtain for \(n \geq 2\) the time evolution:
\[
\tau \frac{\partial P_n}{\partial t} = -2P_n(r) - P_n(r) \sum_{m=1}^{N_p-n} P_m(r - m - 1) - P_n(r) \sum_{m=1}^{N_p-n} P_m(r + n + 1) \\
+ \sum_{m=1}^{n-1} P_m(r - 1)P_{n-m}(r + m) + P_{n-1}(r) \sum_{m=1}^{N_p-n+1} P_m(r + n) \\
+ P_n(r - 1) \left[ 1 - \sum_{m=1}^{N_p-n} P_m(r + n) \right] + P_{n+1}(r - 1) \quad (8)
\]
For the particular case \(n = 1\), there is no \((d)\) process, and the \((e)\) process is replaced by a new process \((e')\) which can be written as
\[
\left( \frac{\partial P_1}{\partial t} \right)_{e'} = \text{Prob}(\circ \bullet \circ \times) = \frac{1}{\tau} \Pr(\circ \bullet \circ \times) \\
= \frac{1}{\tau} \left[ \sum_{m=1}^{N_p} P_m(r + 1) - \sum_{m=1}^{N_p-1} P_m(r + 1) + \sum_{m'=1}^{N_p-m} P_{m'}(r - m') \right] \quad (9)
\]
It is convenient to introduce the generating function in the Fourier space for Eq.(8) and space average the probability distribution \(\bar{P}_n\),
\[
\hat{G}(x, k) = \sum_r e^{-ikr} \sum_{n \geq 1} x^n P_n(r) = \sum_{n \geq 1} x^n \hat{P}_n(k),
\]
\[
\bar{P}_n = \frac{N_n}{N} = N^{-1} \hat{P}_n(0)
\]
with momenta $k = 2\pi m/N$, $m = 0, \cdots, N-1$. From this definition, we have the relations

$$N_c = \hat{G}(1, 0), \quad N_p = \partial_x \hat{G}(1, 0)$$  \hspace{1cm} (11)

The generating function satisfies a master equation which incorporates all the previous processes, as well as the one-particle dynamics. Applying the transformation Eq.(10) to Eq.(8), we obtain

$$\tau \frac{\partial}{\partial t} \hat{G}(x, k) = [-2 + (1 + x^{-1})e^{-ik}] \hat{G}(x, k) - \frac{1}{N} \sum_{k'} e^{-i(k-k')} \hat{G}(x, k') \hat{G}(e^{-i(k-k')}, k - k')$$
$$+ (x - 1) \frac{1}{N} \sum_{k'} e^{i(k-k')} \hat{G}(xe^{i(k-k')}, k') \hat{G}(1, k - k')$$
$$+ \frac{1}{N} \sum_{k'} e^{-ik'} \hat{G}(xe^{i(k-k')}, k') \left[ \hat{G}(x, k - k') - \hat{G}(1, k - k') \right]$$
$$+ xe^{ik} \hat{G}(1, k) - x \frac{1}{N} \sum_{k'} e^{ik'} \hat{G}(1, k') \hat{G}(e^{-i(k-k')}, k - k') - e^{-ik} \hat{P}_1(k)$$  \hspace{1cm} (12)

with $\hat{P}_1(k) = \partial_x \hat{G}(x, k)|_{x=0}$. We should notice that we assume $N \gg 1$ and that the finite size cutoff at $n = N_p$ in Eq.(8) is not taken into account in the expression for $\hat{G}(x, k)$. It can be checked from the previous equation that the number of particles is conserved, by computing directly $\partial_t N_p = \partial_t \partial_x \hat{G}(1, 0) = 0$. The last term in Eq.(12) corresponds to the contribution of one-particle clusters by fragmentation of bigger clusters. Since there is a continuous production of one-particle clusters, we expect the dynamics to reach a steady state. In particular, the time evolution of the cluster number $N_c = \hat{G}(1, 0)$ is equal to

$$\tau \frac{\partial}{\partial t} \hat{G}(1, 0) = -2 \frac{1}{N} \sum_{k'} e^{ik} \hat{G}(1, k') \hat{G}(e^{ik}, -k') + \hat{G}(1, 0) - \hat{P}_1(0)$$  \hspace{1cm} (13)

The dynamics is non-local as $\hat{G}(1, 0)$ depends on non-zero momentum contributions. Let us consider the steady state of Eq.(12). If we assume the dynamics is governed by low momenta close to zero, we may try to study more precisely solutions given by $\hat{G}(x, k) = N \delta_{k,0} \hat{G}_0(x)$ which is a mean-field approximation. This approximation will be compared later with the numerical results. The functional equation Eq.(12) can be solved for this solution, which yields

$$\hat{G}_0(x) = \frac{1}{8x} \left[ 4(1 - x) \sqrt{1 - \alpha x + \beta x^2} - 4 + \left( 7 - 3 \sqrt{1 - 8N^{-1} \hat{P}_1(0)} \right) x \right]$$
$$+ \left( \sqrt{1 - 8N^{-1} \hat{P}_1(0)} - 1 \right) x^2 \right]$$
$$\alpha = \frac{3}{2} \left( 1 - \sqrt{1 - 8N^{-1} \hat{P}_1(0)} \right), \quad \beta = \frac{1}{8} \left( 1 - 8N^{-1} \hat{P}_1(0) - \sqrt{1 - 8N^{-1} \hat{P}_1(0)} \right)$$  \hspace{1cm} (14)
The value of $\hat{P}_1(0)$ is determined by the condition of particle conservation $\hat{G}_0(1) = c$. Solving this condition leads to the following expressions for $\hat{P}_1(0)$ and coefficients $\alpha$ and $\beta$ as function of $c$

$$\hat{P}_1(0) = \frac{c(1-c)(1-c+2c^2)}{(1+c)^2} N,$$

$$\alpha = \frac{6c(1-c)}{1+c}, \quad \beta = \frac{c^2(1-c)^2}{(1+c)^2}, \quad \frac{\alpha^2}{\beta} = 36$$

Therefore $\hat{G}_0(x)$ depends only on the concentration. The series expansion of the generating function can then be expressed using the series expansion of $(1 - \alpha x + \beta x)^{1/2} = \sum_{n \geq 0} C_n x^n$ which leads to

$$\hat{G}_0(x) = N^{-1} \hat{P}_1(0) x + \frac{1}{2} \sum_{n \geq 2} (C_{n+1} - C_n) x^n$$

from which we can identify the size distribution $\bar{P}_n = \frac{1}{2} (C_{n+1} - C_n)$ for $n \geq 2$. Expansion coefficients $C_n$ can be written and rearranged as a double sum (see Appendix B), and we obtain the formula given in Eq.(B3)

$$C_n = \frac{(-1)^{1+[n/2]} \alpha^n}{2^n (2([n+1]/2] - 1) \mathcal{P}^{(-n+1/2,-1)}(7/9)}$$

where $\mathcal{P}^{(-n+1/2,-1)}(z)$ is the Jacobi polynomial which depends on $n$ only and $[.]$ in the integer part. We should distinguish between odd and even $n$ values, and study the asymptotic limit when $n \to \infty$ (see Appendix B for details). In particular we find that

$$\bar{P}_{2m} = \frac{(-1)^m \alpha^{2m+1}}{2^{2m+1}} \left[ \mathcal{P}_m^{(-2m+1/2,-1)}(7/9) - \frac{\alpha \mathcal{P}_m^{(-2m-1/2,-1)}(7/9)}{4m+2} \right] \approx \frac{A_0}{\sqrt{\pi m}} \left( \frac{1 - A(c)}{2m+1} + \frac{2}{4m^2 - 1} \right) \exp(2m \ln A(c))$$

where

$$A(c) = \frac{c(1-c)}{(1+c)(\sqrt{2} - 1)^2}, \quad A_0 = \frac{\sqrt{2} - 1}{2^{1/4}}$$

and for odd values

$$\bar{P}_{2m+1} = \frac{(-1)^m \alpha^{2m+1}}{2^{2m+2}(2m+1)} \left[ \frac{\alpha}{2} \mathcal{P}_{m+1}^{(-2m-3/2,-1)}(7/9) - \mathcal{P}_m^{(-2m-1/2,-1)}(7/9) \right] \approx \frac{A_0}{2 \sqrt{\pi}} \left( \frac{1 - A(c)}{\sqrt{m}} + \frac{A(c)}{2m^{3/2}} \right) \exp[(2m + 1) \ln A(c)]$$
FIG. 2: Average number of clusters with their standard deviation as function of time for $N = 500$ and $10^3$ initial random configurations.

The quantity $A(c)$ is positive and less than unity. $\bar{P}_n$ is exponential with a corrective power law factor $n^{-3/2}$, except at the value where $A(c) = 1$, which is achieved at the critical point $c^* = \sqrt{2} - 1$. For this value, $\bar{P}_n$ follows a power law with decay exponent $5/2$ since the amplitude $1 - A(c)$ for the $n^{-3/2}$ contribution vanishes, as seen in the previous expressions Eq.(18) and (20). In general, $1 - A(c)$ remains a small quantity around $c^*$, so that the apparent exponent is close to $5/2$. For example $A(0.3) \simeq 0.94$, or $A(0.5) \simeq 0.97$ which is close to unity. This corresponds to $\beta = 1$ in the model [11] since $\eta = 3/2 + \beta$ if we identify the two models, and this is represented by an aggregation kernel equal to the size product of its constituents. Also, near $c^*$, ln $A(c) \simeq -(1-c/c^*)^2/\sqrt{2}$, which gives a diverging characteristic cluster size $n_c = \sqrt{2}(1-c/c^*)^{-2}$.

Numerically, we simulated the stochastic dynamics using a number of random initial configurations at fixed concentration and system size. A site is chosen randomly with a uniform probability and if it is occupied by a particle, this one is allowed to move either on the left or right by one step with equal probability, provided that the destination site is empty. If the particle moves to the left, it leaves the cluster where it was attached eventually. If the particle moves to the right, it moves with all the cluster it is attached to accordingly.
FIG. 3: Numerical evaluation of the cluster size distribution $\bar{P}_n$ for $N = 2000$ and different concentrations $c$. $10^5$ random initial realisations were generated, and averages were performed after $2 \times 10^5$ time steps. In inset, black dot symbols represent coefficient $-\ln A(c)$ of fitted function $\bar{P}_n \simeq A(c)^{-n} n^{-\eta}$, and square red symbols are the power law exponent $\eta$.

The process is repeated at each time iteration, and the size distribution is evaluated after the steady state is reached, see Fig. 2, typically after at least $10^5$ time steps for the sizes considered in simulations. For different concentration values, the averaged size distribution is displayed in Fig. 3 for $N = 2000$. The cutoff at $n \simeq cN$ induces a sharp decrease of the curves at large concentration beyond the power law region for finite $N$. As discussed before, this cutoff is not included in the computation of Eq.(14), since we consider the thermodynamical limit $N \gg 1$. In the inset of the figure we have plotted the parameters for the fitting function $\bar{P}_n \simeq A(c)^{-n} n^{-\eta}$ at different concentrations. Using a polynomial fit, $1 - A(c)$ appears to vanish for $c \simeq 0.443$, which is close to the theoretical value 0.414 from the approximation method Eq.(14). The exponent $\eta$ is close to $3/2$ for most of concentrations around $c^*$, which is the exponent extracted from the asymptotic values Eq.(18) and (20). We would normally expect an exponent close to $5/2$ in the critical region, due to the smallness
FIG. 4: Numerical average values of one-particle cluster number $N_1$ and total cluster number $N_c$. $10^5$ random initial realisations were generated, and averages were performed after $2 \times 10^5$ time steps. Dotted lines are the approximated values given by the generating function Eq.(14).

of factor $1 - A(c)$. However the mean-field approximation gives a qualitative explanation of the dynamics with the possible numerical evidence of a critical point near $c^*$. At higher concentrations $c > c^*$, we observe that large clusters tend to aggregate and dominate the dynamics since particles are getting closer to each other, and then merge more easily. Indeed, the finite size effect induces strong contributions to large clustering as $\bar{P}_n$ approaches the singular limit $\bar{P}_n \simeq \delta_{n,cN}$ when $c \simeq 1$, which is observed in the curves above $c = 0.78$ of Fig. 3 by the appearance of a broad peak in the distributions. At low concentrations, the distribution follows mostly an exponential law with a large amplitude $A(c)$ as the clusters tend to diffuse and break into small independent particles. In Fig. 4, we have plotted the average number of clusters $N_c$ and the average number of one-particle clusters $N_1$ as a function of the concentration. Dotted lines represent the values given by Eq.(14), i.e. $N_c = Nc(1 - c)/(1 + c)$ and $N_1 = \hat{P}_1(0)$; see Eq.(15). The agreement is correct for small values of concentrations, and the behavior of $N_1$ is modified near $c \simeq 0.4$ above which it
For comparison, let us consider the simplest symmetric and local dynamics where particles are allowed to diffuse only on one of the two neighboring sites if these are unoccupied. The averaged size distribution $\bar{P}_n$ of the $n$-particle clusters can be evaluated naively as function of the particle concentration $c$. Indeed a cluster of $n$ particles exists with a probability proportional to $(1-c)^2 c^n$, and the averaged distribution follows therefore a simple exponential law $\bar{P}_n = (1-c)^2 c^n$ with a decay rate equal to $-\ln c$, as exemplified in Fig. 5.

Moreover the average number of clusters is equal to the semi-circle law $N_c = Nc(1-c)$. Any cluster can only fragment by releasing one particle at its edges and the distribution is purely exponential with no power law contribution. Therefore the asymmetric and non-local dynamics studied in this paper induces a power law contribution in the distribution. The asymmetry also induces the presence of an internal current to the right direction, with an average particle velocity equal to $v = N^{-1} \sum_n (1-n) \bar{P}_n = 1 - N_c/N > 0$. The velocity is maximum for states with large clusters or at low concentration. Using the approximate value for $N_c$, we find that $v = (1+c^2)/(1+c)$. The velocity is minimum at the critical concentration $c^* = \sqrt{2} - 1$. 
III. DISCUSSION

Stochastic dynamics of clusters with a strong asymmetric kinetic rule displays a critical behavior at a finite concentration \( c^* \simeq 0.414 \), where the size-distribution becomes a power-law with a decay exponent equal to \( 5/2 \). This critical point can be simply viewed as the singular point of the square root term in the generating function Eq.(14) at \( x = 1 \), where the quantity \( 1 - \alpha x + \beta x^2 \), which is always positive for \( 0 \leq x < 1 \), vanishes. It is interesting to compare these results with a model of coagulation with single-particle fragmentation [18] and using the product form for the aggregation kernel, which leads to gelation after a finite time. The mass or size distribution displays, when the gelation time is approached from below, a power law with the same exponent \( 5/2 \) at the critical concentration. We can deduce that the present model might belong to a class of models defined by an aggregation kernel proportional to the product of individual masses or sizes. A quasi-gelation state exists if the chain is not periodic and boundary conditions are fixed at both ends, which modifies the stochastic rules at the boundaries. Indeed the current tends to aggregate all particle to the right end of the chain where a large cluster forms since no particle can be transferred to the left hand side of the chain. After a short time, almost all particles belong to the large cluster, with few small clusters surviving in the steady state. This may be in analogy with percolation problems in higher dimensions where small clusters persist after the spanning cluster is formed.

Corrections to the theory of the present model would incorporate finite size cutoff of the cluster sizes at \( n = cN \) in Eq.(14), which would explain the broad contribution of large clusters at higher concentrations in Fig. 3. Also, we need to take into account the momenta contributions \( k \neq 0 \) in Eq.(12) in order to give more accurate results, for example by considering fluctuations around the uniform solution: \( \hat{G}(x, k) = N\delta_{k,0}\hat{G}_0(x) + \delta G(x, k) \) with \( \delta G(x, 0) = 0 \). This latter substitution would however not modify the average value of the cluster number as non-linear effects should be important. We should also ask whether the critical point is robust when we modify the fragmentation rule. For example Eq.(8) can be generalized to any fragmentation distribution, for example we could impose that one particle can break-off from the cluster from the left. In this context, there should be a class of fragmentation rules that allows the existence of the critical point.

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APPENDIX A: BALANCE EQUATIONS FOR THE DIFFERENT PROCESSES

In this section, we derive the explicit expressions for the different processes that contribute to the master equation Eq.(8). For the process (a), one has simply

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ) = \text{Prob}(\circ \cdots \circ r^{-1} \circ) = \frac{1}{\tau} P_n(r) \quad (A1) \]

For the process (b) one has instead

\[ \text{Prob}(\times \circ \cdots \circ r^{-1} \circ) = \text{Prob}(\circ \circ \circ \cdots \circ r^{-1} \circ) + \text{Prob}(\circ \circ \circ \cdots \circ r^{-1} \circ) + \cdots \]
\[ = \frac{1}{\tau} P_n(r) \times [P_1(r - 2) + P_2(r - 3) + \cdots + P_{N_p-n}(r - N_p + n - 1)] \quad (A2) \]

Process (c) is given by

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ \times) = \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ) + \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ \circ) + \cdots \]
\[ = \frac{1}{\tau} P_n(r) \times [P_1(r + n + 1) + P_2(r + n + 1) + \cdots + P_{N_p-n}(r + n + 1)] \quad (A3) \]

Process (d) is given by

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ) + \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ) + \cdots + \text{Prob}(\circ \cdots \circ r^{-1} \circ \cdots \circ r^{-1} \circ) \]
\[ = \frac{1}{\tau} [P_1(r - 1) P_{n-1}(r + 1) + P_2(r - 1) P_{n-2}(r + 2) + \cdots + P_{n-1}(r - 1) P_1(r + n - 1)] \]

Process (e) is given by

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ \times) = \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ) + \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ \circ) + \cdots \]
\[ = \frac{1}{\tau} P_{n-1}(r) \times [P_1(r + n) + P_2(r + n) + \cdots + P_{N_p-n}(r + n)] \quad (A4) \]

Process (f) is given by

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ) = \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ \circ) - \text{Prob}(\circ \cdots \circ r^{-1} \circ \circ \circ \circ \circ) \]
\[ = \frac{1}{\tau} P_n(r - 1) \times [1 - P_1(r + n) - P_2(r + n) - \cdots - P_{N_p-n}(r + n)] \quad (A5) \]

And the last process (g) is given by

\[ \text{Prob}(\circ \cdots \circ r^{-1} \circ) = \frac{1}{\tau} P_{n+1}(r - 1) \quad (A6) \]
APPENDIX B: EXPANSION COEFFICIENTS $C_n$

We consider in this appendix the expression of coefficients $C_n$ in Eq.(16), and their behavior when $n$ is asymptotically large. From Eq.(14), we first have to expand the quantity

$$\sqrt{1 - \alpha x + \beta x^2} = \sum_{n \geq 0} a_n (-\alpha x + \beta x^2)^n = \sum_{n \geq 0} a_n x^n \sum_{k=0}^{n} \left( \begin{array}{c} n \\ k \end{array} \right) (-\alpha)^{n-k} \beta^k x^k$$

$$a_n = \frac{(-1)^{n+1}(2n)!}{2^{2n}n^2(2n-1)}$$

We can rearrange the terms in the double sum so that

$$\sum_{n \geq 0} a_n x^n \sum_{k=0}^{n} \left( \begin{array}{c} n \\ k \end{array} \right) (-\alpha)^{n-k} \beta^k x^k = \sum_{n \geq 0} x^n \sum_{k=0}^{\lfloor n/2 \rfloor} a_{n-k} \left( \begin{array}{c} n - k \\ k \end{array} \right) (-\alpha)^{n-2k} \beta^k$$

$$= \sum_{n \geq 0} x^n (-\alpha)^n \sum_{k=0}^{\lfloor n/2 \rfloor} a_{n-k} \left( \begin{array}{c} n - k \\ k \end{array} \right) 36^{-k} = \sum_{n \geq 0} C_n x^n$$

(B2)

For $n$ even or odd respectively, the sums over $k$ have a representation in term of Jacobi polynomials

$$C_{2m} = \frac{(-1)^{m+1} \alpha^{2m}}{4^m(2m-1)} P_{m}^{(-2m+1/2, -1)}(z_0),$$

$$C_{2m+1} = \frac{(-1)^{m+1} \alpha^{2m+1}}{4^m(4m+2)} P_{m}^{(-2m-1/2, -1)}(z_0)$$

(B3)

with $z_0 = 7/9$. For $m$ large, we would like to obtain an expansion of the polynomials $P_{m}^{(-2m+1/2, -1)}(z_0)$. Although there are asymptotic formulas in the literature for a large parameter $m \gg 1$ [25, 26], there is no result for the value $z_0 = 7/9$ which is outside the domain of convergence of the asymptotic series. We will use instead an integral representation of the polynomials in the complex plane and perform a simple saddle point analysis. If $w(z) = (1-z)^{-2m+1/2}(1+z)^{-1}$ is the weight of the Jacobi polynomials in Eq.(B3), we have

$$P_{m}^{(-2m+1/2, -1)}(z_0) = \frac{(-1)^m}{2^m w(z_0)} \int \frac{dz}{2i\pi} \frac{w(z)(1-z^2)^m}{(z-z_0)^{m+1}}$$

(B4)

where the contour is a small circle around $z_0$. A change of variable $z = z_0 - (1-z_0)s$ leads to

$$P_{m}^{(-2m+1/2, -1)}(z_0) = \frac{(1+z_0)^m}{2^m} \int \frac{ds}{2i\pi s^{m+1}} \frac{(1+s)^{\pm 1/2}}{1-rs} \left( \frac{1-rs}{1+s} \right)^m$$

(B5)

where $r = (1-z_0)/(1+z_0)$ and the contour is a small circle around the origin. For $m$ large, we apply the saddle point method by extremizing the function $\phi(s) = \ln[(1-rs)/s(1+s)]$. 
We obtain the relevant solution $s^* = (1 - \sqrt{1 + r})/r$ and expand $\phi(s)$ around $s^*$ up to the second order, then compute the resulting Gaussian integral. We obtain

$$\mathcal{P}_m^{(2m+\epsilon/2, -1)}(z_0) \simeq \frac{(1 + s^*)^{\pm 1/2} \exp(m\phi(s^*) + m \ln(1 + z_0) - m \ln 2)}{s^* (1 - rs^*)} \sqrt{2\pi m \phi''(s^*)}$$

with the values

$$\phi(s^*) = i\pi - \ln(1 + z_0) + 2 \ln(\sqrt{2} + \sqrt{1 + z_0}),$$

$$\phi''(s^*) = \frac{2^{1/2}(1 - z_0)^2 (3 + z_0 - 2^{3/2} \sqrt{1 + z_0})}{(1 + z_0)^{1/2} (\sqrt{2} - \sqrt{1 + z_0})^4}$$

Therefore, the asymptotic behavior of $\mathcal{P}_m^{(2m+\epsilon/2, -1)}(z_0)$, with $\epsilon = \pm 1$, and which is valid at least in the interval $0 < z_0 < 1$, is equal to

$$\mathcal{P}_m^{(2m+\epsilon/2, -1)}(z_0) \simeq \frac{(-1)^m (\sqrt{2} - \sqrt{1 + z_0})^{1+\epsilon/2} (1 + z_0)^{1/4} e^{m(2\ln(\sqrt{2} + \sqrt{1 + z_0}) - \ln 2)}}{2^{(5-\epsilon)/4} (1 - z_0)^{\epsilon/2} \sqrt{3 + z_0 - 2^{1/2} \sqrt{1 + z_0} \sqrt{\pi m}}}$$

where we have, in this limit

$$\mathcal{P}_m^{(2m+\epsilon/2, -1)}(z_0) \simeq (1 + s^*) \mathcal{P}_m^{(2m-1/2, -1)}(z_0) \mathcal{P}_m^{(2m-1/2, -1)}(z_0)$$

The asymptotic behavior Eq.(B8) has been checked with ©Maple2018 for $z_0 = 7/9$ with a relative error of about 0.4% up to $m = 80$.

REFERENCES


