Critical behaviour of annihilating random walk of two species with exclusion in one dimension

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The $A + A \to 0$, $B + B \to 0$ process with exclusion between the different kinds is investigated here numerically. Before treating this model explicitly, we study the generalized Domany-Kinzel cellular automaton model of Hinrichsen on the line of the parameter space where only compact clusters can grow. The simplest version is treated with two absorbing phases in addition to the active one. The two kinds of kinks which arise in this case do not react, leading to kinetics differing from standard annihilating random walk of two species. Time dependent simulations are presented here to illustrate the differences caused by exclusion in the scaling properties of usually discussed characteristic quantities. The dependence on the density and composition of the initial state is most apparent. Making use of the parallelism between this process and directed percolation limited by a reflecting parabolic surface we argue that the two kinds of kinks exert marginal perturbation on each other leading to deviations from standard annihilating random walk behavior.

I. INTRODUCTION

Non-universal dynamical behaviour seems to be a controversial issue in non-equilibrium models. An outstanding example is the debated behaviour of systems exhibiting infinitely many absorbing states [1–4]. There is no analytic treatment up to now; argumentation of various authors, in most of the cases, is based on simulation results. Despite intensive study, the critical behaviour of such systems is poorly understood, non-universality remains an unresolved problem and even scaling behaviour is questioned. Roughly speaking, in these coupled processes the ‘primary’ particles follow a branching and annihilating random walk while the other species just provide a slowly changing environment that effects the branching rates of the primaries. The spreading exponents of the primaries depend on the initial conditions of the environment.

A possible way which might lead to a deeper understanding of the mechanism behind non-universal spreading could be the study of simpler coupled systems. Perhaps the simplest case is the coupled annihilating random walk of two species ($A + A \to 0$, $B + B \to 0$). Naively one would expect that this could be described by the exactly solved field theory of the $A + A \to 0$ process (ARW). In one dimension, however, the situation is more subtle than in higher dimensions. Particles of different type can block the motion of each other. The difference between one and two dimensions has been found to give rise to different phase diagrams in the case of the general epidemic model [2]. The question now arises how relevant the exclusion perturbation caused by this blocking mechanism is to a fixed point of the kind determined in [3].

Another motivation of this study originates from the investigations of Hinrichsen [9], who found, by simulations, a strange scaling behaviour in some special case of his model [21], for which, however, an explanation is still lacking. In section II Hinrichsen’s model will be introduced. It is easy to see, that the kinks in this model at the symmetry point corresponding to the compact directed percolation point of the Domany-Kinzel automaton, exhibit the process described above. In sections III and IV we present our high precision time dependent simulation results from random and seed initial conditions. In section V these results are compared with those obtained by rigid (i.e. parabolic) boundaries. We further investigate this analogy on the mean-field level in section VI, while section VII is devoted to results in the explicit two-species Annihilating Random Walk model with exclusion (ARW2e). We summarise our numerical results in section VIII and give an outlook toward $N$-species generalisation in IX. A qualitative description of the behaviour of Hinrichsen’s model outside the symmetry point on the line of compactness is presented in section X and finally in section XI we summarise and discuss our results.

II. THE GENERALISED DOMANY-KINZEL SCA

The Domany-Kinzel (DK) stochastic cellular automaton (SCA) [7] is one of the simplest models which show a non-equilibrium phase transition into an absorbing state. This one dimensional SCA is defined on a ring with two states ‘1’ and ‘0’ with the following rule of update:

$t$: 0 0 0 0 0 0 1 1 1 1 1 0 0 1 1
$t+1$: 0 p p p q
where at \( t + 1 \) the probability of '1'-s is shown.

In the plane of the parameters \((p, q)\), the phase diagram of the DK SCA is as follows. A line of critical points separates the active phase (with a finite concentration of '1'-s) from the absorbing (vacuum) phase (with zero steady state density of '1'-s). This continuous transition belongs to the universality class of directed percolation (DP) \([8]\). The endpoint of this line \((q = 1, p = 1/2)\) describes a transition, however, outside the DP class; it corresponds to compact directed percolation. Here the model exhibits Ising symmetry and can be solved exactly \([7]\).

In 1997 Hinrichsen \([9]\) introduced a generalized version of the DK SCA including more than one symmetric inactive states \((I_1, I_2, \ldots)\) and one active state \((A)\). The motivation for this study was to look for a possible change of the critical behaviour of the above described system, on the line of compactness, at \( q = 1, p = 0.5 \). Red: \( I_1 \), blue: \( I_2 \), white: \( A \)

The phase diagram exhibited on Fig1.b) shows that the line of PC transitions ends at \( q = 1, p = 1/2 \), a point which corresponds the Ising symmetry point of the DK automaton. The primary aim of the present work is to investigate the scaling properties of GDK at this point, which will be called CDP2 transition point. A typical time evolution of the GDK model at this special point when starting from a random initial arrangement of \( I_1\)-s, \( I_2\)-s and \( A\)-s is shown on Fig. 2. Here active islands can be spatially extended, thus three kinds of compact clusters can grow. Nevertheless only the \( I_1 \) and \( I_2 \) phases are \( Z_2 \)-symmetric while the active phase plays a special role. (The situation is different from a 3-states Potts model with Glauber kinetics).

It is well known that the CDP process in 1d is equivalent to an annihilating random walk process of kinks separating compact domains of 0-s and 1-s. In the model investigated here two types of kinks can be defined, namely kink \( K_1 \) between domains \( A - I_1 \) (and \( I_1 - A \)) and kink \( K_2 \) between neighbouring \( A - I_2\)-s (and \( I_2 - A\)-s). The rules of the model inhibit occurrence of kinks between domains of absorbing phases, i.e. between \( I_1-I_1 \) and \( I_2-I_2 \).

Kinks \( K_1 \) and \( K_2 \) perform annihilating random walks: \( K_1 + K_1 \rightarrow \emptyset, K_2 + K_2 \rightarrow \emptyset \), while the processes \( K_1 + K_2 \rightarrow \emptyset, K_2 + K_1 \rightarrow \emptyset \) are, however, forbidden. In other words, upon meeting, a \( K_1 \) and a \( K_2 \) “block” each other (do not annihilate and do not exchange sites) \([21]\). To our knowledge such kind of kinetics has not been studied before. Motivated by this fact we have decided to explore the critical behaviour of the above described system, on the line of compactness \((q = 1)\), by computer simulation. In this study special attention will be paid to the \( p = 1/2 \) symmetry point CDP2.
III. SIMULATIONS FROM RANDOM INITIAL STATE

A. Kink decay simulations

We have performed time dependent simulations starting from states with uniformly distributed species $A$, $I_1$, and $I_2$, with respective densities: $\rho_0(A)$, $\rho_0(I_1)$ and $\rho_0(I_2)$. At the CDP2 point unusual scaling behaviour of the density of kinks has been observed previously [21]: a deviation from the ordinary annihilation-diffusion process with kink-density decay $\rho(t) \sim \frac{t}{\sqrt{t}}$. Instead, $\rho(t) \sim t^{-\alpha}$ with $\alpha \approx 0.55$ has resulted from the first simulations.

To check whether the observed deviation from standard ARW behaviour is only a cross-over effect, or it heralds some basic feature of altered kinetics we have performed very long-time ($t_{\text{max}} = 10^6$ MCS) simulations on systems with $L = 24000$ (Fig. 3, 22). Throughout the whole paper $t$ is measured in units of Monte-Carlo sweeps.

![Figure 3](image)

**FIG. 3.** Total kink number as the function of time started from symmetrical homogeneous random initial states: $\rho_0(I_1) = \rho_0(I_2) = 0.3$ (solid line) and $\rho_0(I_1) = \rho_0(I_2) = 0.1$ (dotted line). The dashed line corresponds to the single species annihilating random walk ($\rho_0(I_1) = 0$, $\rho_0(I_2) = 1/2$) exhibiting $\rho(t) \propto t^{-0.5}$.

Figure 3 shows the results of simulations. It is seen that the deviation from the standard ARW value of the decay exponent remains present asymptotically as well: the local slopes of the decay curves

$$\alpha(t) = \frac{\ln[\rho(t)/\rho(t/m)]}{\ln(m)}$$

(1)

(where we use $m = 8$ usually) go to constant values. Moreover, another interesting feature has become apparent: the kink-decay exponent depends on the initial concentrations of the components $\rho_0(I_1) = \rho_0(I_2)$ and in such a way that for higher initial kink density (lower average distance between the kinks) the decay is faster.

Asymptotically, as $\rho_0 \to 0$, the average distance of dissimilar kinks goes to zero and the decay exponent tends to the ARW value: $\alpha \to 0.5$.

In the case of asymmetric initial condition ($\rho_0(I_1) \neq \rho_0(I_2)$) $K_1$-s and $K_2$-s decay with different rates. The type that has smaller initial density decays faster. Example: in case of $\rho_0(I_1) = 1/9$, $\rho_0(I_2) = 1/3$, $K_2$ decays roughly like $t^{-0.5}$ ( unperturbed by $K_1$-s) but the local slopes of the log-log $\rho(K_1)-t$ dependence decrease from $-0.5$ strongly.

IV. SIMULATIONS FROM AN ACTIVE SEED

The cluster simulations [12] were started from a state with uniformly distributed $A$-s and $I_1$-s except a single $I_2$ pair in the middle and the following characteristic quantities for the $I_2$-s were followed:

- the average number of $I_2$-s, $N_{I_2}(t)$,
- their survival probability $P_{I_2}(t)$,
- and the average mean square distance of spreading of $I_2$’s from the center $R_{I_2}^2(t)$

The above quantities were averaged over $N_s$ independent runs at the CDP2 point ( in case of $R_{I_2}^2(t)$ only for surviving samples ). At the critical point we expect these quantities to behave for $t \to \infty$, as

$$N_{I_2}(t) \propto t^\eta,$$

(2)

$$P_{I_2}(t) \propto t^{-\delta},$$

(3)

$$R_{I_2}^2(t) \propto t^z.$$

(4)
Upon varying the initial density $\rho_0(I1)$, for the exponents $\delta$ and $\eta$ (defined similarly to eq.1), the local slopes of $N_{I2}(t)$ and $P_{I2}(t)$ continuously changing values have been observed (Figs. 5, 6). The deviation of these exponents from those of the single-species annihilation random walk process: $1/2$ and 0, respectively is remarkable. The spreading exponent, $z$, on the other hand, seems to be constant within numerical accuracy and equals that of the single species annihilating random walk: $z = 2/Z = 1$ such that the generalised hyper-scaling law of the compact directed percolation \cite{26} is satisfied.

\begin{equation}
\eta + \delta = z/2
\end{equation}

is satisfied. In this respect it is important that $\eta$ has been found to be negative.

FIG. 5. Local slopes of the number of $I2$-s. The initial state is uniformly distributed with initial densities: $\rho_0(I1) = 0.1$ (solid line), 0.25 (dotted line), 0.5 (dashed line), 0.75 (long-dashed line).

FIG. 6. The same as on Fig. 5 for the cluster survival probability.

V. CLUSTER SIMULATIONS OF COMPACT DIRECTED PERCOLATION CONFINED IN A PARABOLA.

To understand the physics of our numerical results up to now we set up a parallelism with an other case where DP process is bounded by parabolic space-time boundary conditions. We perform simulations on the compact cluster version of this and compare the results with those of the GDK model in section VII.

Kaiser and Turban have investigated \cite{23,24} the 1+1 dimensional DP process limited by a special, parabolic boundary condition in space and time directions:

\begin{equation}
y = \pm C t^k
\end{equation}

where $C$ changes under uniform length rescaling (by $b$) to:

\begin{equation}
C' = b^{2k-1} C
\end{equation}

Here $Z$ is the dynamical critical exponent. By referring to conformal mapping of the parabola to straight lines and showing it in the mean-field approximation Kaiser and Turban claim that for $k < 1/Z$ this surface gives relevant, for $k > 1/Z$ irrelevant and for $k = 1/Z$ marginal perturbation to the DP process. The marginal case results in $C$ dependent non-universal power-law decay, (for details see next section), while for the relevant case stretched exponential functions have been obtained. The above authors have given support to this claim by numerical simulations.

We have investigated the effect of parabolic and reflecting boundary conditions for the CDP2 process numerically. Time-dependent cluster spreading simulations have been performed in the GDK model with parabolic boundaries such that at each time step the simulation region is bounded by two $I1$-s at $y_{min}$ and $y_{max}$, where

\begin{equation}
y_{min} = L/2 - 2 + C t^k
\end{equation}

\begin{equation}
y_{max} = L/2 + 2 + C t^k
\end{equation}

Two $I2$-s have been put initially at the centre ($L/2, L/2 + 1$) and some initial space (two A-s to the left and right) between $K1$-s and $K2$-s has been added. Therefore the role of $I1$-s now is purely the formation of parabolic boundaries around $I2$-s and in fact we investigate the plain CDP process with reflective boundary conditions. A typical 1+1 dimensional run looks like as shown below (1, 2 and 0 stand for $I1$, $I2$ and $A$, respectively):
When we fixed the exponent at $k = 1/2$, to make the situation marginal we found continuously changing exponents for the exponents of the survival probability $P_{I_2}$ and the number of $I_2$-s, $N_{I_2}$, by varying the shape $C$ (Fig. 7). One can see that the exponent-slopes of $N_{I_2}(t)$ (Fig. 8) and those of $P_{I_2}(t)$ (Fig. 9) change by varying $C$. The spreading exponent of the $I_2$-s, $z$, seems to be constant: equal to unity. (Fig. 10). These results are very similar to those of the seed simulations in GDK of section V.

The analysis based on local slopes (Figs. 8, 9, 10) shows again plateaus for high values of $t$, indicating true power-law behaviours. The magnitude of the exponent characterizing the decay of the density of $I_2$-s decreases as $C$ is increased reminiscent of a similar situation in [23].

\[ z/2 = \eta + \delta = 1/2 \]

is fulfilled. In this case it is important, again, that $\eta$ takes negative values.
VI. THEORETICAL CONSIDERATIONS FOR CDP CONFINED IN A PARABOLA

A. Anisotropic scaling

In 1+1 dimensional anisotropic systems the correlation length diverges as \( \xi_{\parallel} \sim t^{-\nu_{\parallel}} \) in time and as \( \xi_{\perp} \sim t^{-\nu} \) in the space direction with a dynamical exponent \( Z = \nu_{\parallel}/\nu \) (\( \nu \) is also denoted as \( \nu_{\perp} \) in the literature). Covariance under a change of the length scales then requires two different scaling factors, \( b_{\parallel} = b^Z \) and \( b_{\perp} = b \).

We consider now a system displaying anisotropic critical behaviour and limited by a free surface in the \((t,y)\) plane as given in eq.(6). Under rescaling, with \( t' = t/b^Z \) and \( y' = y/b \), \( C \) transforms according to eq.(7), as discussed in the previous section.

In the marginal case, which we will consider now, \( Z = 1/k \), the scaling dimension \( x_m \) of the tip order parameter becomes \( C \)-dependent \( x_m(C) \).

The order parameter correlation function between the origin and a point at \((t,y)\) transforms as

\[
G\left(\Delta, t, y, \frac{1}{C}\right) = b^{-2x_m} G\left(\frac{b^{1/\nu}}{\nu}, \frac{t}{b^Z}, \frac{y}{b^{1/\nu}}\right)
\]

when \( L \) is infinite. With \( b = t^{1/Z} \), equation (10) leads to:

\[
G\left(\Delta, t, y, \frac{1}{C}\right) = t^{-2x_m/Z} G\left(\frac{t}{\Delta^{\nu_{\parallel}}}, \frac{y^{\nu_{\parallel}}}{t^{1/\nu_{\parallel}}}, \frac{1}{C}\right).
\]

(11)

Here \( l_C = C^{-1/2\nu_{\parallel}} \), \( \Delta = \left(\frac{p-p_c}{p_c}\right) \) and \( x_m \) is the scaling dimension of the order parameter. The latter is connected to \( \beta \), the critical exponent of the order parameter via \( \beta = \nu x_m \). We will use this scaling form in the following. \( \beta \) is the usual order-parameter exponent, defined, for the DKCA, through \( \rho_1 \propto (p-p_c)^\beta \), for \( p > p_c \); \( \rho_1 \) is the stationary density of \( 1 \)'s. In case of a first order transition as is the case with compact directed percolation the following considerations hold.

As already mentioned,

\[
P(t) \propto t^{-\delta},
\]

(12)

is the survival probability of \( 1 \)'s for spread of particles (\( 1 \)'s, in our notation) about the origin. Away from the critical point \( \beta' \) governs the ultimate survival probability (starting from a localized source): \( P_\infty \equiv \lim_{t \to \infty} P(t) \propto (p-p_c)^{\beta'} \). It is known that \( \beta' = 1 \) in CDP. The order-parameter exponent, \( \beta \), however, is zero. This is because \( p = 1/2 \) marks a discontinuous transition, by symmetry. \( \rho_1 = 0 \) for \( p < 1/2 \) and \( \rho_1 = 1 \) for \( p > 1/2 \); strictly speaking, \( \beta \) is not defined here but it is natural to associate the value \( \beta = 0 \) with the discontinuous transition.

This problem with the ill-defined exponent \( \beta \) can be avoided following the lines of Grassberger and de la Torre’s scaling argument [12] for discontinuous transitions as it has been presented by Dickman and Tetryakov [20]. Consider a model with a transition from an absorbing to an active state at \( \Delta = 0 \), with exponents \( \delta, \eta, \gamma, \) and \( \beta' \) defined as above. Suppose, however, that the order parameter \( \rho \) is discontinuous, being zero for \( \Delta < 0 \), and

\[
\rho = \rho_0 + f(\Delta),
\]

(13)

for \( \Delta > 0 \), where \( \rho_0 > 0 \), and \( f \) is continuous and vanishes at \( \Delta = 0 \). According to the scaling hypothesis for spreading from a source there exist two scaling functions, defined via [12]

\[
\rho(y,t) \sim t^{\eta - \delta + \gamma/2} G(y^2/t^{\gamma}, \Delta t^{1/\nu}),
\]

(14)

and

\[
P(t) \sim t^{-\delta} \Phi(\Delta t^{1/\nu}).
\]

(15)

(Here \( \rho(y,t) \) is the local order-parameter density. \( \tau \sim \Delta^{-\nu} \).) Existence of the limit \( P_\infty \) implies that \( \Phi(x) \sim x^{\beta'} \) as \( x \to \infty \), with \( \beta' = \delta \nu \). In a surviving trial, the local density must approach the stationary density \( \rho \) as \( t \to \infty \), so \( \rho(y,t) \sim \Delta^{\beta' \nu} \rho_0 \) for \( t \to \infty \) with fixed \( y \) and \( \Delta \) small but positive. It follows that \( G(0,x) \sim x^{\beta' \nu} \) for large \( x \).

An important consequence is that we can use as scaling dimension of the order parameter for CDP the value \( \beta' \) in the relation \( x_m = \frac{\beta}{\nu} \) instead of \( \beta \). Via scaling relations \( \beta' = \delta \nu \), the values obtained by computer simulations for \( \delta \) will be compared with results for CDP+parabolic boundary conditions. In this context the connection, again via scaling relation, between \( \delta \) and the decay exponent of the density of kinks when starting from a random initial state \( \alpha \) will also be made use of.
B. Mean field analysis for CDP confined in a parabola

In this section we will follow the lines of the mean-field analysis of the $1+1$ dimensional DP process confined by a parabola as given in ref. [24], but now applied to compact directed percolation.

The basic processes are:

![Site update rules for compact directed percolation](image)

FIG. 11. Site update rules for compact directed percolation.

The order parameter correlation function is the probability density $P(t, y)$ for a site at $(t, y)$ to be connected to the origin.

First we consider the case without confinement. In mean-field approximation one can set up an equation for the connectedness at $(t + 1, y)$:

$$P(t + 1, y) = p[P(t, y + 1)[1 - P(t, y - 1)] + P(t, y - 1) [1 - P(t, y + 1)] + P(t, y + 1)P(t, y - 1)$$

Going to the continuum limit the following differential equation is obtained:

$$\frac{\partial P}{\partial t} = p\frac{\partial^2 P}{\partial y^2} + (2p - 1)P + (1 - 2p)P^2$$  \hspace{1cm} (17)

The homogeneous, stationary solution of eq.(17) is:

$$P_0 = \begin{cases} 1 & \text{for } p > 1/2 \\ 0 & \text{for } p \leq 1/2 \end{cases}$$  \hspace{1cm} (18)

This describes a first order transition for CDP at $p_c = 1/2$, as it is the case. At the transition, $p = p_c$, eq.(17) reduces to

$$\frac{\partial P}{\partial t} = \frac{1}{2} \frac{\partial^2 P}{\partial y^2}$$  \hspace{1cm} (19)

which is the ordinary diffusion equation with RW solution

$$P(t, y) = \frac{\exp\left(-\frac{y^2}{4t}\right)}{\sqrt{2\pi t}}$$  \hspace{1cm} (20)

which is exact in the CDP case. From comparison with the scaling form in the previous subsection, the following (well-known) exponents for CDP arise:

$$\nu_\parallel = 1 \quad \nu = 1/2 \quad Z = 2 \quad x_m = \frac{1}{2}$$  \hspace{1cm} (21)

On a parabolic system, we use the new variables $t$ and $\zeta(t, y) = y/t^k$ for which the free surface is shifted to $\zeta = \pm C$ and equation (13) is changed into

$$\frac{\partial P}{\partial t} = \frac{1}{2\kappa^2} \frac{\partial^2 P}{\partial \zeta^2} + k \frac{\zeta}{t} \frac{\partial P}{\partial \zeta}$$  \hspace{1cm} (22)

with the boundary condition $P(t, \zeta = \pm C) = 0$. Through the change of function

$$P(t, \zeta) = \exp\left[-\frac{k}{2}\zeta^2 t^{2k-1}\right]Q(t, \zeta)$$  \hspace{1cm} (23)

equation (22) leads to

$$\frac{\partial Q}{\partial t} = \frac{1}{2\kappa^2} \frac{\partial^2 Q}{\partial \zeta^2} + \frac{k}{2} \left[(k - 1)\zeta^2 t^{2k-2} - \frac{1}{t}\right]Q$$  \hspace{1cm} (24)

for which the variables separate when $k = 1, 1/2$ or 0. These values of $k$ just correspond to irrelevant, marginal and relevant perturbations.

For $k = 1$ the critical behaviour is the same as for unconfinement percolation as expected for an irrelevant perturbation.

For the true parabola which is the marginal geometry, one may use equation (22) with $k = 1/2$ to obtain

$$\kappa \frac{\partial P}{\partial \zeta} = \frac{1}{2} \frac{\partial^2 P}{\partial \zeta^2} + \frac{\zeta}{t} \frac{\partial P}{\partial \zeta} \quad \zeta = \frac{y}{t^{1/2}}$$  \hspace{1cm} (25)

which is of the form studied in [23] for the directed walk problem. Writing $Q(t, \zeta) = \phi(t)\psi(\zeta)$ leads to the following eigenvalue problem for $\psi(\zeta)$

$$\frac{1}{2} \frac{d^2 \psi}{d \zeta^2} + \frac{\zeta}{t^{1/2}} \frac{d \psi}{d \zeta} = -\lambda^2 \psi$$  \hspace{1cm} (26)

with $\phi(t) \sim t^{-\lambda^2}$. The solution is obtained as the eigenvalue expansion

$$P(t, y) = \sum_{n=0}^{\infty} B_n \ t^{-\lambda_n^2} \ F_1\left[\lambda_n^2, \frac{1}{2}; \frac{y^2}{2t}\right].$$  \hspace{1cm} (27)

The behaviour at large $t$ is governed by the first term in this expansion which decays as $t^{-\lambda_n^2}$, i.e. with a $C$-dependent exponent as expected for a marginal perturbation. The dimension of the tip-to-bulk correlation function is the sum of the tip and bulk order parameter dimensions, the first one being variable. Comparing with the form of the decay in eq.(14) gives $\lambda_n^2 = |x_m^f(C) + x_m|/Z$ and, using (21), the tip order parameter dimension is given by

$$x_m^f(C) = 2\lambda_n^2 - \frac{1}{2}.$$  \hspace{1cm} (28)

Its dependence on $C$ is shown in figure 2. of [24]
Analytical results can be obtained only in limiting cases which have already been discussed in [23]. When $C$ is infinite, $\lambda^2 = 1/2$, only the first term in the expansion remains, which satisfies the initial and boundary conditions, giving back the free solution in equation (24). For large C-values the tip exponent is $x_m = (C) = 1/2 + \sqrt{2\pi C \exp\left(\frac{-C^2}{2}\right)} \left[1 + O(\varepsilon)\right]$, where $\varepsilon$ is the correction term itself. For narrow systems, the hyper-geometric function gives a cosine to leading order in $C^2$. One obtains the following asymptotic behaviour in $t$

$$P(t,y) \sim t^{-\pi y/8C^2} \cos \left(\frac{\pi y}{2C\sqrt{t}}\right)$$

(29)

and the tip exponent diverges as $\pi^2/4C^2$.

For $0 < k < 1/2$ the dependence on $t$ is expected to be a stretched exponential function. For details see [23,25].

VII. ANNihilating RANDOM WALK OF TWO SPECIES WITH EXCLUSION

To check our results concerning the scaling properties of kinks in the GDK model at the CDP2 point we have carried out an explicit simulation of the annihilating random walk of two species ($A, B$) with exclusion. The model we have investigated has been suggested by Hinrichsen [21] and is as follows. A(B) will hop to a neighbouring empty site with probability $p_A$ ($p_B$) or annihilate with a neighbour $A$ ($B$) with probability $p_{2A}$ ($p_{2B}$) while $A$ and $B$ do not react when getting into neighbouring positions. The initial configuration was chosen in such a way that allows pairs of the same kind to annihilate always within some finite time interval (i.e. the system evolves into an empty state), namely:

$$...A.A.B.B.A.A.A.A.B.B...$$

That means that $AA$ and $BB$ pairs have been put in a 1d ring with initial probability $\rho(0)$. Had we not chosen the initial state like this the system would have ended up in some finite particle configuration where $A$-s and $B$-s follow each other alternatingly, separated by arbitrary empty regions. (This initial configuration is in agreement with the arrangement of the two kinds of kinks in some random initial state of the GDK model, too.) The probabilities $p_A$, $p_B$, $p_{2A}$ and $p_{2B}$ have been chosen to be unity, to achieve maximum simulation efficiency; no qualitative difference in the results have been found upon lowering them.

Clearly this process is different from the simple annihilating random walk of two species $A + B \rightarrow \emptyset$ [6], therefore we may expect that a field theory describing this model (which, however, is still missing) would result in a different fixed point with different critical exponents as well. Furthermore one can argue that when comparing the simple random walk and the random walk+exclusion (SEP) processes one also observes different dynamical behaviours. This latter case is nothing else but the $T = 0$ dynamics of the 1d Ising Model with Kawasaki exchange, where we have different domain growth properties than in case of a simple random walk.

An extensive numerical simulation with look-up table algorithm seems to confirm this expectation. As Figure 12 shows the slopes of the density decay started from the special pairwise random states described above depend on the initial density $\rho(0)$.

![Figure 12](image.png)

**FIG. 12.** Local slopes of particle decay in the annihilating random walk + exclusion process of two species results ($L = 24000$). The initial conditions are: $\rho_0 = 0.1, 0.25, 0.5$ from bottom to top curve.

The local slopes tend to constant values greater than $\alpha = 0.5$ in agreement with the GDK kink results. The level-off in case of $\rho_0 = 0.5$ happens only for $t > 1.5 \times 10^6$ MCS. The average $AA$ and $BB$ distances confining an other type of particle have also been measured during the simulations, that enables us to extract the amplitude ($C$) of the confinement in the function fitted $(C \times t^{-\alpha})$. These values will be used to compare the results with those of the GDK (see next section).

VIII. SUMMARY OF TIME DEPENDENT RESULTS

Since in all of the previously shown cases we found non-universal scaling depending on the initial conditions and the generic model to account for such behavior seems to be the CDP2 with parabolic boundary condition, we have decided to measure the region of confinement in all cases and plotted the survival probability exponents $\delta$ and the kink decay exponents $\alpha$ as a function of the shape of the measured parabola.

In the present case the final survival probability of a cluster plays the role of the order parameter exponent $\beta$ as explained at the end of Sec.VI and for the characteris-
tic exponents we have: $\delta = \beta / \nu_\perp$. Thus we have plotted the results for $\delta(C)$. In a common graph the fitted values for $\alpha(C)$ are also shown; on the level of kinks the order parameter $\beta$ is connected to $\alpha$ in the same way as $\beta'$ to $\delta$ for 'spins' (see eg. [25]).

For random initial conditions in the GDK model the characteristic distance between two neighbouring kinks of a given type has also been measured. The average neighbour distance $l_{K2-K2}$ shown on Fig. 13 have been obtained for initial densities: $\rho_0(I1) = \rho_0(I2) = 1/3$. The power law increase for large $t$ (see the plateau for $t > 30.000$) with the same scaling exponent as the decay exponent is not very surprising because $\rho_{kink}(t) \propto 1/l_{K2-K2}$.

Since the $K2-K2$ and the $K1-K1$ pairs confine the motion of each other ( a $K1-K2$ pair can not exchange to $K2-K1$) this power-law increasing length scale imposes a 'stochastic' boundary condition (pressure on kinks) with a mean value of a parabola similarly that was investigated by Kaiser and Turban [23] in case of 1 + 1 d DP processes [24] and adapted for the case of a CDP-like first order transition in section VI. As discussed before, the scaling dimension of the order parameter changes continuously with the amplitude of the parabolically growing confining box size if it grows with the same exponent as the cluster inside.

![FIG. 13. Local slopes of the $K2-K2$ neighbour distances in GDK model of size $L = 24000$. The initial state is uniform with $\rho_0(I1) = \rho_0(I2) = 1/3$](image)

In our case we encounter similar situation. The kink density decay exponent $\alpha$ seems to vary continuously in case of symmetrical initial conditions. The initial conditions effect the amplitudes of the density decays (as was shown to be valid by field theory for pure the reaction diffusion of A,B particles [5]) and therefore the amplitudes of the confinement region sizes ($C$) (see Fig. 14). To compare our results with those of [23,24] the form $A + C t^\alpha$ has been fitted to the $l_{K2-K2}(t)$ distances determined from the density decay simulations (assuming $l_{K2-K2}(t) = 2/\rho_{kink}(t)$). The following table summarises the results for GDK with random initial conditions:

| $\rho_0$ | $C$ | $\alpha$ |
|---------|-----|---------|
| 0.0     | $\infty$ | 0.5000(3) |
| 0.05    | 14.09 | 0.517(2)  |
| 0.10    | 7.62  | 0.528(2)  |
| 0.15    | 6.11  | 0.534(2)  |
| 0.2     | 5.85  | 0.537(2)  |
| 0.3     | 4.18  | 0.540(1)  |

This is in agreement with Fig. 8 of [24], where an increasing $C$ causes a decreasing exponent. Note that the first line in the table corresponds to the simple ARW process, therefore there is no confinement (amplitude $C$ is $\infty$).

The main difference between the rigid parabola boundary case and the "stochastic" confinement is that in the latter case the boundary generates an additional noise to the motion of confined particles. Therefore we don’t simple have "free" particles confined in a parabola, but they are also perturbed by the noise in such a K1 ↔ K2 symmetrical way that the outcome perturbation is marginal.

![FIG. 14. Possible explanation for the non-universal scaling. The symmetric I1 an I2 clusters exert marginal exclusion perturbation on each other.](image)

In ARW simulations again the form $A + C t^\alpha$ has been fitted for the measured AA, BB distances.

In cluster simulations we fitted the form: $y = C \times t^\delta$ for the region of confinements and determined the respective C-s in all cases.
As Figure 15 shows we obtained similar monotonically decreasing curves in all cases that also agrees with the results of [23]. The GDK-uniform and the ARW2 results seem to lie on the same curve. The spreading simulation results of GDK are different from those of the CDP+parabolic boundary condition case.

This can be understood, however, since in the former case the confined particles ($I_2$-s) have a back effect on the bulk ($I_1$-s) particles, while this is not the case when the boundary is fixed.

We also show the $\pi^2/8C^2$ curve, determined as the asymptotic solution $C \to 0$ of the mean-field approximation, see section VI eq.(29). This seems to be in fair agreement with the case of CDP+fixed parabolic boundary condition.

**IX. GENERALISATION FOR $N > 2$: SYMMETRIC ANNIHILATING EXCLUSION PROCESS OF $N$ SPECIES**

We have carried out preliminary simulations in the generalized version of the model introduced in section VII. The system was started from configurations like:

$$\ldots A\ldots A\ldots B\ldots B\ldots C\ldots C\ldots DD\ldots E\ldots EFF \ldots$$

where species of the same type can annihilate each other but different types can not exchange. Our results show that concerning the time dependence of the density the deviations from the square root decay persist for $N > 2$ and this property seems to remain valid also for $N \to \infty$.

**X. GDK ON THE LINE OF COMPACTNESS**

On the line $q = 1$ the role of the absorbing states ($I_1$ and $I_2$) is symmetric. The above reported simulation results for the GDK model refer to $p = 1/2$, the CDP2 point. Now we will discuss the situation for $p \neq 1/2$. For $p > 1/2$ the creation of new $A$-s happens with probability greater than 0.5, the active domain size grows exponentially and the inactive regions die out quickly (and symmetrically); the all-$A$- phase plays the same role as the all- 0- phase of the original DK automaton [7]. The deviation from the DK picture is quite apparent, however, for $p < 1/2$, as instead of the all-1-phase of DK, for all values of $p$ with the exception of $p = 0$, Glauber-Ising -like kinetics governs the motion of kinks. The kinks here are extended objects ($A$-s), somewhat similarly to those in Grassberger’s CA models [10] (for $p < p_c$, where $p_c$ is the critical point of a parity conserving phase transition ) where also kinks of different extension (and there even of different structure) separate absorbing-phase clusters. (At $p = 0$ diffusion stops and a striped space-time picture of $I_1$ and $I_2$ domains freezes in, again like in Grassberger’s model A cited above.) The average size of $A$-s goes to zero between two domains of the same type quickly, the kink ‘particles’ of same type perform a *biased* random walk toward each other. On the other hand between domains of different types remains a film of $A$-s of average size 1, since a collision of an $I_1$ and $I_2$ domain always creates a new $A$ at the next time step. That means that kinks of different types still block the motion of each
other. Therefore the role of A-s is similar to the kinks of the $T = 0$ Glauber Ising model. On the whole line of $1/2 > p > 0$, in the long time limit, therefore one can expect the number of such kinks to decrease as $(t^{-1/2})$. On the other hand this is a line of compactness, as all clusters growing from a seed are compact, the characteristic exponents ($\eta$, $\delta$ and $z$) though strongly dependent on $p$ and the composition of the initial state, satisfy the hyper-scaling law valid for compact cluster [26]. This statement involves that this line is a line of first order transition points with order parameter exponent $\beta_s = 0$. This first order transition occurs in a symmetry-breaking ('magnetic') field coupled to the $I_1$ and $I_2$ spins. For the detailed description of a similar situation see [15]. All these features have been supported by simulations. As an example we can give some results obtained at $p = 0.4$ starting with a single $I_2$ in the sea of $I_1$. A-s (25% of $I_1$, 75% of $A$); $\delta = 0.45$, $z/2 = 0.47$, $\eta_{I_2} = -0.02$. It is worth mentioning, that for hyper-scaling to hold it is again important that $\eta_{I_2}$ is negative.

XI. DISCUSSION

We have investigated numerically the one dimensional generalized Domany-Kinzel cellular automaton on a line in the plane of its parameters where only compact clusters grow. The two types of kinks in the simplest version of this compact GDK model (two absorbing phases) follow annihilating random walk with exclusion (no reaction) between different types. The equivalence with an explicit two-species ARW model with exclusion is shown provided the initial state is prepared in such a way that the kinks are arranged in pairs with some density. High precision simulations revealed that this system relaxes in a non-trivial way: the decay exponent of the kink density depends on the initial density of kinks. We argue that this is a kind of (internal) surface effect; similar to the ARW process confined by a rigid space-time parabola provided the power of the parabola is chosen to be marginal. This case has been explicitly investigated with the result that the spreading exponents behave qualitatively the same way as expected from the corresponding mean-field approximation. We have no proof of the marginality for the theory including fluctuations, but rely on symmetry arguments. If we assumed that particles would exert relevant perturbation ($k < 1/2$) on each other, the corresponding parabola picture would predict stretched exponential decay (a behaviour that is very difficult to differentiate from power-laws by simulations) and the local slopes should go to some higher value as a function of time (meaning faster that any power-law decay). However our high precision data show just the opposite, the local slopes decrease as a function of time tending to a value somewhat greater than 1/2.

Nevertheless, the possibility of pure square-root decay masked by some tremendously long crossover function can not be ruled out. One could still expect a non-universal scaling of the survival probability of particles in the same way as was observed in [23][24] or in another similar situation [28] where a diffusing "prisoner" confined by marginally growing cage was investigated. In the latter case the the boundary condition was absorbing and an exact solution was possible giving an exponent for the survival probability which is a continuous function of the amplitude of the marginal parabola. Furthermore the survival of a diffusing prisoner (with diffusivity D) inside a cage where both walls diffuse (with diffusivity A) has been solved exactly and the decay exponent was found to be $\pi/2 \cos^{-1}(D/(D + A))$ [29].

Non-standard scaling in a 1d ARW model was also observed by Frachebourg et al. [10]. They have shown that the survival probability of particles in an ARW with one free boundary depends on the location of the particles. If we count the particles from the free boundary the survival probability of odd particles decay with exponent 0.225, while those of even numbered decay with exponent 0.865. The explanation for this is based on the fact that even numbered particles always have left and right neighbours during the process, while odd numbered particles lack one of the neighbours and since the ARW in 1d is diffusion limited they can escape. One can notice a similarity of this mechanism to the one in ARW2 models we investigated. Namely in our case there are infinitely many internal boundaries (generated by particles of different types which can not exchange sites).

Recently Bray [31] has shown that the relaxation towards the critical state in the 2d XY model depends on the initial state. This is very different from what is expected from field-theoretical RG predictions that can not take into account low-dimensional topological effects. Moreover, Bray has shown that the non-universal behavior of the persistence exponent in this case can be described by the random walk of a particle moving under an attractive central power-law force that creates marginal perturbation as compared to free random walk [21]. This scenario is similar to ours since we also have particles with RW exhibiting pressure of marginal strength on each other.

Right after our submission two other preprints appeared on cond-mat [32][33], dealing with models very similar to those presented here and reporting results which are in accord with ours for those quantities they also investigated.

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[1] I. Jensen and R. Dickman, Phys. Rev. E 48, 1710 (1993); I. Jensen, Phys. Rev. Lett. 70, 1465 (1993).
[2] P. Grassberger, H. Chaté and G. Rousseau, Rev. E 55, 2488 (1997).
[3] M. A. Muñoz, G. Grinstein, R. Dickman and R. Livi, Phys. Rev. Lett. 76, 451 (1996) and Physica D 103, 485 (1997).
[4] J. F. F. Mendes, R. Dickman, M. Henkel and M. C. Marques, J. Phys. A 27, 3019 (1994).
[5] B.P. Lee, J.Phys.A 27, 2633 (1994).
[6] B.P. Lee and J. Cardy, J.Stat.Phys 80, 971 (1995).
[7] E. Domany , W. Kinzel, Phys. Rev. Lett. 53, 311 (1984).
[8] Various papers on directed percolation are to be found in *Percolation Structures and Processes*, edited by G. Deutsch, R. Zallen and J. Adler, Annals of the Israel Physical Society, 5 (Hilger, Bristol, 1983).
[9] H. Hinrichsen, Phys. Rev. E 55, 219 (1997).
[10] P. Grassberger, F. Krause and T. von der Twer, J. Phys. A:Math.Gen., L105 17 (1984).
[11] P. Grassberger, J.Phys. A:Math.Gen. 22, L1103 (1989).
[12] P. Grassberger and A. de la Torre, Ann.Phys.(NY) 122 122, 373 (1979).
[13] H. Takayasu and A. Yu. Tretyakov, Phys. Rev. Lett. 68, 306 0, (1992).
[14] I. Jensen, Phys.Rev.E 50, 3623 (1994).
[15] N. Menyhárd, J.Phys.A:Math.Gen., 27, 6139(1994).
[16] N. Menyhárd and G. Ódor, J. Phys. A 28, 4505 (1995).
[17] N. Menyhárd and G. Ódor, J.Phys.A:Math.Gen. 29, 7739 (1996).
[18] N. Menyhárd N. and G. Ódor, J.Phys.A:Math.Gen. 31, 6771(1998).
[19] J. Cardy and U. Täuber, Phys. Rev. Lett. 77,4780 (1996).
[20] K. E. Bassler and D. A. Browne, Phys. Rev. Lett. 77, 4094 (1996); K. E. Bassler and D. A. Browne, Phys. Rev. E 55 (1997 May 1).
[21] H. Hinrichsen, private comm. (1997).
[22] G. Ódor, A. Krikelis, G. Vesztergombi and F. Rohrbach: Proceedings of the 7-th Euromicro workshop on parallel and distributed processing, Funchal (Portugal) Feb. 3-5 1999, IEEE Computer society press, Los Alamitos, ed.: B. Werner; e-print: physics/990905.
[23] C. Kaiser and L. Turban, J. Phys. A:Math.Gen 27, L579 (1994).
[24] C. Kaiser and L. Turban, J. Phys. A:Math.Gen 28, 351 (1995).
[25] L. Turban, J. Phys. A: Math. Gen 25, L127 (1992).
[26] R. Dickman and A. Yu. Tretyakov, Phys. Rev. E 52, 3218 (1995).
[27] N. Menyhárd N. and G. Ódor, cond-mat/0001101

[28] P. L. Krapivsky and S. Redner, Am. J. Phys.64, 546(1996)
[29] D. ben-Avraham J. Chem. Phys. 88, 941 (1988); M.E.Fisher and M.P.Gelfand J.Stat.Phys. 53,175 (1988).
[30] L. Frachebourg, P. L. Krapivsky and S. Redner, J. Phys. A: Math. Gen 31, 2791 (1998).
[31] A.J.Bray cond-mat/99-10-135 Random Walks in Logarithmic and Power-Law Potentials....
[32] S. N. Majumdar, D. S. Dean, cond-mat/0002217, Novel Coarsening in the Zero Temperature Dynamics of an Ising Chain with Kinetically Induced Field.
[33] S. Kwon, J. Lee, H. Park, cond-mat/0002214, Does hard-core interaction change absorbing type critical phenomena?