Growth Law and Superuniversality in the Coarsening of Disordered Ferromagnets

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Abstract

We present comprehensive numerical results for domain growth in the two-dimensional Random Bond Ising Model (RBIM) with non-conserved Glauber kinetics. We characterize the evolution via the domain growth law, and two-time quantities like the autocorrelation function and autoresponse function. Our results show evidence for the crossover from a pre-asymptotic regime with “power-law growth with a disorder-dependent exponent” to an asymptotic regime with “logarithmic growth”. We compare this behavior with previous results on one-dimensional disordered systems and we propose a unifying picture in a renormalization group framework. We also study the corresponding crossover in the scaling functions for the two-time quantities. Super-universality is found not to hold. Clear evidence supporting the dimensionality dependence of the scaling exponent of the autoresponse function is obtained.
1 Introduction

The *coarsening dynamics* or *domain growth* of a thermodynamically unstable system, such as a ferromagnet or a binary mixture quenched below the critical point, is a well-established paradigm of slow relaxation. From the pioneering work of Marro et al. [1] up to the reviews of Bray [2] and Puri [3], the equal-time properties have been extensively investigated. Recently, the interest has shifted to the study of two-time quantities, since the ordering process offers a simple and non-trivial arena for the understanding of aging phenomena [4, 5]. There is also a long-standing interest in studying the robustness of the basic features of coarsening systems when quenched disorder (without frustration) is introduced [6, 7]. Taking the Ising model with ferromagnetic couplings as the basic system, the class of disordered systems investigated includes the *random field Ising model* (RFIM) [8] and the *random bond Ising model* (RBIM) [7].

The first effect of disorder is to change the relaxation mechanism from purely energy lowering to activated, due to the creation of energy barriers [9]. The question is to what extent this affects the growth law, namely the dependence on time of the linear size of domains $L(t)$. Intuitively, the presence of energy barriers is expected to slow down the coarsening process. However, the problem turns out not to be so straightforward, since the barrier size may depend on the domain size, thus determining the nature of the asymptotic growth law through a non-trivial feedback process. So far, this issue has been well understood in the case of one-dimensional systems. The $d = 1$ RFIM [10] and the $d = 1$ RBIM [11] offer instructive and clear instances of disorder forming either $L$-dependent barriers (in the RFIM) or $L$-independent barriers (in the RBIM). In both cases, there is a crossover. In the $d = 1$ RFIM, it occurs from pure-like power-law growth to disorder-dominated logarithmic growth. In the $d = 1$ RBIM, the crossover takes place in the reverse order: from a pre-asymptotic disorder-dependent power law to an asymptotic pure-like power-law growth.

In higher dimensions, a crossover from algebraic to logarithmic behavior has been found in the growth law of the RFIM [12, 13, 14, 15]. A crossover from algebraic to slower than power-law behavior, possibly logarithmic, has been reported also in the case of the random site Ising model (RSIM) [16, 9, 17]. According to the phenomenological theory of Huse and Henley (HH) [18], this means that in these systems energy barriers ought to scale like a power of $L$. 
Instead, in the $d = 2$ RBIM the role of barriers and the nature of the growth law has remained controversial, since the available numerical evidence shows a steady algebraic growth with a disorder-dependent dynamical exponent \cite{19, 21, 22}. The open issue is whether the data manage to probe the genuine late time behavior, in which case energy barriers ought to scale logarithmically with $L$ \cite{19}, or the HH paradigm is valid also for this system, in which case the data would be probing a long pre-asymptotic regime, eventually to be followed by a crossover to logarithmic growth \cite{23, 24}. That an HH-type scenario might yield an effective algebraic law at intermediate times had already been noticed by Bouchaud et al. \cite{25} in the context of spin glasses. Our contribution in this paper is precisely the clearing up of this point, producing the numerical evidence for the existence of the crossover to logarithmic growth in the $d = 2$ RBIM.

A related important question is how disorder affects the scaling functions, namely whether it enters only through the modified growth law or if there is also an explicit disorder dependence. In the first case, scaling functions ought to be the same for pure and disordered systems, once time is reparametrized through $L(t)$. This is referred to as the super-universality (SU) hypothesis \cite{26}, whose validity has been supported mainly from numerical results for the equal-time correlation function or structure factor \cite{6, 14, 15, 19}. More recently, the validity of SU has been extended by Sicilia et al. \cite{20} to the geometrical properties of domain structures and by Henkel and Pleimling (HP) \cite{21, 22} to two-time quantities, such as the order parameter autocorrelation and autoresponse functions in the $d = 2$ RBIM. However, next to the features supporting SU validity, there exists also some evidence showing SU violation, as reported by HP \cite{22} in the short distance behavior of the $d = 2$ RBIM and by Park and Pleimling \cite{17} in the space-time correlation function in the RSIM. What is missing, here, is the general framework where to fit coherently these conflicting pieces of information, with the end-result that the issue of SU validity is regarded as yet to be understood.

By contrast, a clear picture emerges from the $d = 1$ cases mentioned above \cite{10, 11}, where the crossover can be investigated with considerable precision and turns out not to be limited to the growth law, but to take place also in the correlation and response functions. Then, the issue of SU validity or lack of validity is reduced to deciding whether the asymptotics are or are not disorder-dominated. Adopting this point of view, our finding of the crossover in the growth law in $d = 2$ RBIM strongly suggests that in this system too there might be full violation of SU and that this might fit into the
pattern extracted from the $d = 1$ systems. Indeed, this is what we find from the autocorrelation and autoresponse function data and, putting together these new data with those for $d = 1$ systems, we have succeeded in constructing a new and coherent picture of general character for the coarsening dynamics of weakly disordered systems. This is most effectively formulated into a renormalization group framework in terms of the competition between a pure fixed point and a disordered fixed point. Then, the variety of different scenarios observed arises from the different possible choices in the relative stability of these fixed points.

Finally, the presence of disorder allows for an exhaustive investigation of the scaling properties of the dynamical susceptibility. The data in the pre-asymptotic regime of the $d = 2$ RBIM, where the growth law is algebraic with good precision, together with those for the $d = 1$ RBIM, produce conclusive evidence for the $d$-dependence of the exponent involved. Our results are in agreement with the conjecture put forward in studies of the pure system [5, 27, 28, 29]. A preliminary account of this result has been presented in Ref. [11].

This paper is organized as follows. In Sec. 2, we provide a general framework, based on renormalization group ideas, to understand the effects of disorder on domain growth. In Sec. 3, we summarize our understanding of coarsening in $d = 1$ disordered systems. The bulk of the paper is contained in Sec. 4, where we present comprehensive numerical results for domain growth in the $d = 2$ RBIM. We conclude the paper with a summary and discussion in Sec. 5.

2 General Framework

This section is devoted to an overview, using scaling arguments and renormalization group (RG) terminology, of the possible deviations from pure-like behavior due to disorder.

2.1 Growth Law

The linear size of domains in pure systems asymptotically grows according to the power law

$$L(t) = D_0 t^{1/z},$$

(1)
which holds when time $t$ is large enough for $L(t)$ to dominate all other lengths in the problem. Here, $D_0$ is a prefactor which is dependent on temperature, and $z$ is the dynamical exponent. It is straightforward to verify that the above power law satisfies the homogeneity relation

$$L(t) = lL \left( \frac{t}{l^z} \right)$$

where $l$ is arbitrary and can be regarded as the scale parameter in an RG transformation. In this paper we shall consider only non-conserved dynamics. In pure systems the dynamical exponent is $z = 2$, independent of the system dimensionality [2].

Assuming scaling, the growth law for disordered systems is obtained generalizing Eq. (2) to

$$L(t, \epsilon) = lL \left( \frac{t}{l^z}, \frac{\epsilon}{l^\phi} \right)$$

where $\epsilon$ is a parameter quantifying the amount of disorder and $\phi$ is the corresponding scaling exponent. Once the growth law is written in this form, the standard crossover scenario follows from the competition between the pure fixed point at $\epsilon = 0$ and the disordered fixed point at $\epsilon = \infty$. The relative stability of these two fixed points is regulated by the sign of $\phi$, with the pure fixed point being stable or unstable for $\phi > 0$ or $\phi < 0$, respectively. In the first case the asymptotic growth law would remain the pure one, while, in the second case, there would be the crossover to an asymptotic behavior different from that of Eq. (2) and controlled by the disordered fixed point.

In place of investigating directly the sign of $\phi$, it turns out to be more convenient to introduce the crossover length $\lambda(\epsilon) \sim \epsilon^{1/\phi}$ and to rewrite Eq. (3) as

$$L(t, \epsilon) = lL' \left( \frac{t}{l^z}, \frac{\lambda(\epsilon)}{l} \right).$$

At fixed points the crossover length values are either $\lambda^* = 0$ or $\lambda^* = \infty$, and since $\lambda'(\epsilon)$ flows from $\infty$ toward 0, the fixed point corresponding to $\lambda^* = 0$ is stable, while the one with $\lambda^* = \infty$ is unstable. Therefore, the relevancy or irrelevancy of disorder can be deduced from the behavior of $\lambda(\epsilon)$ as $\epsilon \to 0$. Specifically, since $\epsilon = 0$ corresponds to the pure fixed point and the two

\footnote{As we shall see, in the case of the $d = 1$ RBIM Eq. (4) applies also to the more general case where $\lambda(\epsilon)$ does not scale like a power of $\epsilon$.}
possible limits of $\lambda(\epsilon)$ are

$$\lim_{\epsilon \to 0} \lambda(\epsilon) = \begin{cases} 0 \\ \infty \end{cases}$$

if the first one applies, the pure fixed point is stable, while if the second limit applies, the disordered fixed point is stable.

As a final remark, notice that making the choice $l = t^{1/z}$ in Eq. (4), the growth law takes the standard scaling form

$$L(t, \epsilon) = t^{1/z} \mathcal{L} \left( \frac{\lambda(\epsilon)}{t^{1/z}} \right).$$

### 2.2 Two-time Quantities

Once the role of $\lambda(\epsilon)$ is understood in the growth law, the next step is to establish how it affects the other observables in the problem.

In this paper, we shall be concerned with spin systems which are quenched below the critical temperature at $t = 0$. For what follows, the basic two-time quantities are the autocorrelation function and the autoresponse function. The first of these is defined by

$$C(t, t_w) = \langle \sigma_i(t) \sigma_i(t_w) \rangle - \langle \sigma_i(t) \rangle \langle \sigma_i(t_w) \rangle.$$

Here, $\sigma_i = \pm 1$ is the spin variable at the lattice site $i$, $(t, t_w)$ are a pair of times after the quench, with the convention $t_w \leq t$, and angular brackets $\langle \cdot \rangle$ denote average over all sources of randomness in the problem: initial condition, thermal fluctuations and quenched disorder. After taking these averages $C(t, t_w)$ is space-translationally invariant and there is no dependence on the choice of $i$.

The autoresponse function is defined by

$$R(t, t_w) = \frac{\delta \langle \sigma_i(t) \rangle_h}{\delta h_i(t_w)} \bigg|_{h=0},$$

where $h_i(t)$ is a space and time-dependent external field conjugate to $\sigma_i$, and $\langle \cdot \rangle_h$ denotes the average in the presence of the field. This instantaneous response function is very difficult to measure, both numerically and experimentally. For this reason, time integration is used as a means to improve the signal to noise ratio [30]. We shall focus on the zero-field-cooled susceptibility
(ZFCS), obtained by switching on the external field from the time $t_w \geq 0$ up to the observation time $t$:

$$\chi(t, t_w) = \int_{t_w}^{t} dt' R(t, t').$$  \hspace{1cm} (9)

The motivations and the advantages for using this form of integrated response function, with respect to others, have been discussed at length in Ref. [28].

Let us denote by $O(t, t_w)$ either one of the above observables. In the pure case and for quench temperatures low enough to neglect thermal fluctuations, the aging-scaling relation is obeyed [4, 5, 31]

$$O(t, t_w) = L^{-\alpha}(t_w) F\left(\frac{L(t)}{L(t_w)}\right).$$  \hspace{1cm} (10)

where $\alpha$ is the scaling exponent and $F(x)$ is the scaling function.

The extension of Eq. (10) to the disordered case is a non-trivial problem. According to the SU hypothesis, disorder should affect only the growth law, leaving the form of Eq. (10) unaltered with the same scaling function, that is

$$O(t, t_w, \epsilon) = L^{-\alpha}(t_w, \epsilon) F\left(\frac{L(t)}{L(t_w)}\right).$$  \hspace{1cm} (11)

In other words, once time is re-parametrized by $L(t, \epsilon)$, there should be no difference between disordered and pure systems. This is a strong statement.

Alternatively, one can allow for an explicit disorder-dependence of the scaling function

$$O(t, t_w, \epsilon) = L_w^{-\alpha} \tilde{F}\left(\frac{L}{L_w}, \frac{\lambda}{L_w}\right),$$  \hspace{1cm} (12)

and extend the crossover scenario to two-time observables. Here $L(t, \epsilon)$ has been denoted by $L$, $L(t_w, \epsilon)$ by $L_w$ and $\lambda(\epsilon)$ by $\lambda$. Using the notation $x = L/L_w$ and $y = \lambda/L_w$, the $x \ll y$ region is controlled by the unstable $y^* = \infty$ fixed point, while the $x \gg y$ region is controlled by the stable $y^* = 0$ fixed point. Adopting the convention that hat-free symbols refer to the pure fixed point and hatted symbols to the disordered fixed point, if disorder is an irrelevant perturbation from the analysis presented in the previous subsection follows that the stable fixed point $y^* = 0$ is pure and that the general scaling relation in Eq. (12) yields the limiting behaviors

$$L_w^{-\alpha} \tilde{F}(x, y) \sim \begin{cases} L_w^{-\alpha} \tilde{F}(x), & \text{for } x \ll y, \\ L_w^{-\alpha} F(x), & \text{for } x \gg y, \end{cases}$$  \hspace{1cm} (13)
where $\tilde{F}(x) \neq F(x)$. The exponent $\tilde{\alpha}$ is different from or equal to $\alpha$, depending on whether $\lim_{y \to \infty} \tilde{F}(x, y)$ is singular or not. Conversely, if disorder is a relevant perturbation the stable fixed point $y^* = 0$ is disorder-controlled and the limiting behaviors are reversed

$$L^{-\alpha}_w \tilde{F}(x, y) \approx \begin{cases} L^{-\alpha}_w F(x), & x \ll y, \\ L^{-\tilde{\alpha}}_w \tilde{F}(x), & x \gg y \end{cases} \quad (14)$$

with $\tilde{\alpha}$ being different from or equal to $\alpha$, depending on whether $\lim_{y \to 0} \tilde{F}(x, y)$ is singular or not. For instance, in the case of the autocorrelation function there are no singularities in $\tilde{F}(x, y)$, since $\alpha_C = \tilde{\alpha}_C = 0$ as a consequence of the compact nature of domains, both for pure and disordered systems.

In the general framework outlined above, the SU hypothesis holds as an asymptotic statement when disorder is irrelevant, while SU is violated and the asymptotics of disordered and pure systems are different when disorder is relevant.

### 3 Domain Growth in $d = 1$ Disordered Ising Systems

Let us now discuss $d = 1$ disordered systems. These will provide clear realizations of the possible scenarios outlined in the previous section. As a general comment, it should be kept in mind that even if $d = 1$ systems are paramagnetic at any finite temperature, as long as the domain size $L(t)$ is smaller than the equilibrium correlation length $\xi(T)$, the coarsening dynamics is the same as in the quench to $T = 0$ where ferromagnetic ordering occurs. This is explained in detail in Refs. [10, 11, 32].

We consider the ferromagnetic Ising chain with Hamiltonian

$$\mathcal{H} = - \sum_{i=1}^{N-1} J_i \sigma_i \sigma_{i+1} - \sum_{i=1}^{N} h_i \sigma_i, \quad \sigma_i = \pm 1. \quad (15)$$

The system evolves with non-conserved Glauber spin-flip dynamics [3]. Interfaces in this case are point defects, separating domains of up spins from domains of down spins. In order to identify the length $\lambda$, and to keep the discussion simple, let us focus on the energetics of the particular configuration
[σ]_j, containing a single defect breaking the bond J_j:

\[ \sigma_i = \begin{cases} 1, & \text{for } i \leq j, \\ -1, & \text{for } i > j. \end{cases} \] (16)

The energy change when the defect moves from the bond j to another location r lattice spacings away with, say, r > 0, is given by

\[ \Delta E_{j,j+r} = 2[J_{j+r} - J_j - \sum_{i=j}^{j+r} h_i]. \] (17)

In the pure system, the ferromagnetic coupling is the same for all pairs of nearest neighbors (J_i = J > 0). In the absence of an external field (h_i = 0), \( \Delta E_{j,j+r} = 0 \) for any r. The potential landscape is flat and defects perform free diffusion, yielding the behavior of Eq. (11) in the large-time regime [32].

In the \( d = 1 \) ferromagnetic RBIM, the couplings J_i are non-negative independent random variables, and h_i = 0 as in the pure case. For a given configuration of the couplings, there are now local minima and maxima in the potential landscape. A local minimum occurs when a defect breaks a bond flanked, to the right and to the left, by stronger bonds. Similarly, a defect sits on a local maximum when the opposite bond configuration occurs. The energy difference in this case is given by

\[ \Delta E_{j,j+r} = 2(J_{j+r} - J_j), \] (18)

which depends on the actual values of the couplings involved, but not on their distance. For instance, taking the J_i to be uniformly distributed in the interval \((1 - \epsilon, 1 + \epsilon)\) with \(0 \leq \epsilon \leq 1\), the upper bound on \( \Delta E_{j,j+r} \) is given by \(4\epsilon\). Using the Arrhenius-type relation between escape time and barrier height \( t_B \sim e^{E_B/T} \), there remains defined a characteristic time

\[ \tau \sim e^{4\epsilon/T}, \] (19)

as the time needed to overcome the highest energy barrier, and a characteristic length scale

\[ \lambda \sim \left[ e^{4\epsilon/(\epsilon T)} - 1 \right]. \] (20)

We take the Boltzmann constant \( k_B = 1 \) throughout the paper.

In the \( d = 1 \) RFIM, the ferromagnetic coupling is the same for all pairs of nearest neighbors (J_i = J > 0), as in the pure case, while h_i = ±\( \epsilon \) is an
uncorrelated random field with expectations $\langle h_i \rangle = 0$ and $\langle h_i h_j \rangle = \epsilon^2 \delta_{ij}$. In this case, domain walls perform random walks in a random potential of the Sinai type [33]. The landscape contains minima and maxima, as in the $d = 1$ RBIM. However, the crucial difference is that the height of energy barriers is not bounded and scales with the distance traveled by the defect:

$$\Delta E_{j,j+r} = - \sum_{i=j}^{j+r} h_i \sim \epsilon r^{1/2}. \quad (21)$$

The characteristic disorder length is obtained by balancing thermal energy with the barrier height

$$\lambda = (T/\epsilon)^2. \quad (22)$$

Therefore, from Eqs. (20) and (22), it follows that when $\epsilon \to 0$ the two possible scenarios, originating from Eq. (5) and analyzed in detail in Sec. 2, find realization in the $d = 1$ RBIM and RFIM, respectively. Indeed, this is what has been found [10, 11]. In the $d = 1$ RBIM disorder is irrelevant since from Eq. (20) follows $\lim_{\epsilon \to 0} \lambda = 0$. Our numerical simulations [11] have shown very precisely that the growth law, the autocorrelation function and the ZFCS scale according to Eqs. (6), (12) and (13), with a crossover from pre-asymptotic disordered to asymptotic pure-like behavior. Actually, in the pre-asymptotic regime, the growth law is algebraic with a disorder-dependent growth exponent, indicating that $\lambda$ should more properly be regarded as a marginal scaling field. A similar behavior will be encountered in the $d = 2$ RBIM discussed in Sec. 4. Furthermore, the ZFCS exponent vanishes:

$$\alpha_\chi = 0 \quad (23)$$

for all disorder values, as in the pure case [32]. Hence, the ZFCS scaling function, like that of the autocorrelation function, does not have singularities for $y \to \infty$.

On the other hand, the $d = 1$ RFIM [10] is a realization of the opposite case, since Eq. (22) shows that $\lim_{\epsilon \to 0} \lambda = \infty$ and, therefore, that disorder is relevant. The growth law, the autocorrelation function and ZFCS scale according to Eqs. (6), (12) and (14) and the crossover occurs from pure-like to disordered behavior. Again, as in the $d = 1$ RBIM, the exponents of the autocorrelation function and ZFCS vanish in the pure and disordered

2Clearly, here the average is only over the quenched randomness.
systems, revealing the absence of singularities in the scaling functions as $y \to 0$.

Because of its importance in this paper, it is useful to recall the result for the growth law in the $d = 1$ RFIM. The limiting forms of the scaling function in Eq. (6) are given by [10]

$$L(y) \sim \begin{cases} y(\ln y)^2, & \text{for } y \ll 1, \\ D, & \text{for } y \gg 1. \end{cases}$$

(24)

Thus, in the pre-asymptotic regime (where $y \gg 1$ or $L(t) \ll \lambda$), the thermal energy exceeds the barrier height and growth takes place via free diffusion, as in the pure system (there is a possible dependence of the diffusion constant $D$ on disorder.) In the asymptotic regime (where $y \ll 1$ or $L(t) \gg \lambda$), diffusion is of the Sinai type with $L(t) \sim (\ln t)^2$ [34]. This is a consequence of the Arrhenius law with a barrier height which depends on $L$ according to Eq. (21). Then, the overall growth law, which interpolates between limiting behaviors and accounts for Eq. (24), can be written as

$$t \approx L^z e^{(L/\lambda)^{1/2}}.$$  

(25)

4 Domain Growth in the $d = 2$ Random Bond Ising Model

Let us next consider the $d = 2$ RBIM in the absence of an external field, with Hamiltonian

$$\mathcal{H} = -\sum_{\langle ij \rangle} J_{ij} \sigma_i \sigma_j,$$

(26)

where the sum runs over nearest-neighbor pairs. The ferromagnetic couplings $J_{ij}$ are independent random variables with the same statistics as in the $d = 1$ case, i.e., uniformly distributed in the interval $(1 - \epsilon, 1 + \epsilon)$ with $0 \leq \epsilon \leq 1$.

Let us first summarize the preexisting theoretical and numerical situation. According to the phenomenological HH [18] theory, energy barriers in this system (when $d > 1$) ought to scale as a power of the domain size:

$$E_B(L) = \kappa L^\psi,$$

(27)

where $\kappa$ is a disorder-dependent prefactor such that

$$\lim_{\epsilon \to 0} \kappa = 0$$

(28)
and $\psi$ is an exponent dependent on dimensionality with the value $\psi = 1/4$ for $d = 2$.

Let us, then, see what are the consequences on the basis of the general considerations made above. Identifying the characteristic disorder scale through the matching of thermal energy with barrier height, one has

$$\lambda = (T/\kappa)^{1/\psi}$$

which implies $\lim_{\epsilon \to 0} \lambda = \infty$. Therefore, in the HH theory disorder is relevant, which implies that the scenario found in the $d = 1$ RFIM should be replicated with a diffusive pre-asymptotic regime followed by asymptotic logarithmic growth

$$L(t) \sim \begin{cases} \frac{t^{1/z}}{\lambda}, & \text{for } L \ll \lambda, \\ (\ln t)^{1/\psi}, & \text{for } L \gg \lambda \end{cases}$$

and with an interpolating formula analogous to Eq. (25)

$$t \simeq L^z e^{(L/\lambda)^{\psi}}.$$ 

However, this HH scenario has never been actually observed. Rather, from a number of experiments on random magnets [35] and from extensive simulations of the $d = 2$ RBIM [19, 21, 22], the observed growth seems to be well-described by an algebraic law

$$L(t) \sim t^{1/\tau},$$

with a disorder-dependent exponent $\tau$. This growth law applies across the accessible time region, and there is no hint of a crossover.

A possible explanation of the numerical findings has been put forward by Paul et al. [19] on the basis of a logarithmic dependence of $E_B$ on $L$, in place of the power law in Eq. (27). However, Cugliandolo et al. [23, 24] have argued that the observation of a behavior consistent with Eq. (32) does not necessarily rule out a growth law obeying Eqs. (30) and (31), since a disorder-dependent algebraic growth is compatible with Eq. (31) as an intermediate regime, as it had been previously remarked by Bouchaud et al. [25] in the context of spin glasses. Thus, if we rewrite Eq. (31) as

$$L^\tau \simeq L^z e^{(L/\lambda)^{\psi}},$$

we obtain

$$\tau \simeq z + \frac{1}{\lambda^\psi \ln L}.$$
and, if $L^\psi/\ln L$ can be treated as a constant $c$ over the time interval of interest, the effective exponent is given by

$$z \simeq z + \frac{c}{\lambda^\psi} = z + \frac{cK}{T}. \tag{35}$$

The present understanding of the observed algebraic growth (32) is to a sort of standstill between these two mutually excluding explanations.

In order to settle the issue, we have performed the most comprehensive simulations, to date, of coarsening in the $d = 2$ Glauber-RBIM. Our simulations have been done on $N^2$ square lattices up to maximum time $t_m$ Monte Carlo steps (MCS). Taking a random configuration of up and down spins, which mimics the disordered $T = \infty$ state before the quench, as initial condition for a run, we have let the system evolve with the updating rule which aligns spins with probability 1 with the Weiss field, if there is a majority of nearest neighbours pointing in the same direction, while the usual Glauber rule is used when the Weiss field is zero. This is equivalent to take couplings of the form $J_{ij} = J + \delta_{ij}$, with $\delta_{ij}$ uniformly distributed in $(-\epsilon, \epsilon)$ and then letting $T \to 0$ and $\epsilon \to 0$, while keeping $\epsilon/T$ finite. Specifically, we have studied the coarsening dynamics with values $\epsilon/T = 0, 0.5, 1, 1.5, 2, 2.5$. Kinetics in this form is appropriate for deep quenches, where it provides a considerable gain in the speed of computation, since updating trials are restricted to spins not aligned with the Weiss field, whose number roughly scales like $1/L(t)$. In the above limit disorder and temperature enter the transition rates in the form $\epsilon/T$ and, therefore, also results depend on disorder and temperature only through the ratio $\epsilon/T$. Finally, we have checked that there is no difference with the growth law and the aging properties of two-time quantities computed with the standard Monte Carlo algorithm.

All statistical quantities have been obtained as an average over $N_{\text{run}}$ independent runs. For each run, the system has a different initial condition and disorder configuration. The values of $N, t_m$ and $N_{\text{run}}$ for different values of $\epsilon/T$ are listed in Table 1. Notice that we have taken huge system sizes for $\epsilon/T = 0$ and $\epsilon/T = 0.5$. This is necessary to obtain a reasonable time-window of domain growth before we encounter finite-size effects, as coarsening is more rapid for pure and near-pure systems with nonconserved dynamics. We have performed several checks to ensure that the data presented here are free of finite-size effects. This is crucial as the crossover phenomena we are investigating are subtle and are expected to occur at very late times. We also stress that the $t_m$-values for the present data sets are an order of magnitude

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Disorder values

| $\epsilon/T$ | System size $N^2$ | Maximum time $t_m$ (MCS) | Number of runs $N_{\text{run}}$ |
|--------------|-------------------|--------------------------|-------------------------------|
| 0.0          | $12000^2$         | $1.01027 \times 10^8$    | 20                            |
| 0.5          | $8192^2$          | $3.12973 \times 10^9$    | 20                            |
| 1.0          | $4096^2$          | $1.006735 \times 10^6$   | 75                            |
| 1.5          | $4096^2$          | $1.006735 \times 10^6$   | 75                            |
| 2.0          | $4096^2$          | $1.006735 \times 10^6$   | 75                            |
| 2.5          | $4096^2$          | $1.006735 \times 10^6$   | 75                            |

Table 1: Numerical parameters for $d = 2$ RBIM simulations for various disorder strengths.

(or more) larger than those of Paul et al. [19].

Our findings for the growth law, the autocorrelation function and the ZFCS are presented in the following subsections.

### 4.1 Growth Law

We have obtained $L(t)$ from the inverse density of defects, which is measured by dividing the number of sites with at least one oppositely-aligned neighbor by the total number of sites. The plot of $L$ vs. $t$ in Fig. 1 shows the existence of two time regimes, separated by a microscopic time $t_0$ of order 1. In the early time regime, for $t < t_0$, there is no dependence on disorder and growth is fast. The defects seeded by the initial condition execute rapid ballistic motion toward the nearby local minima. For $t > t_0$, the data sets in Fig. 1 appear to confirm previous observations [19, 21, 22] of a late-time regime with algebraic growth as in Eq. (32). Measuring slopes, the values of the disorder-dependent exponent $\tau$ are displayed in Table 2 and Fig. 2, with an $\epsilon/T$ dependence of the type observed in Ref. [22].

The improvement with respect to previous work comes from the quality of our data, which allow for a more refined analysis. In addition to estimating $\tau$ we have extracted the effective exponent $z_{\text{eff}}(t)$ defined by

$$ \frac{1}{z_{\text{eff}}(t)} = \frac{d(\ln L)}{d(\ln t)}. $$ (36)

3We have checked that this definition is in good agreement with the usual method of measuring $L(t)$ by the half-peak-width of the equal-time structure factor.
Figure 1: Growth law in the pure case (top curve) and for various disorder strengths.

Table 2: Exponent $\tau$ for various disorder strengths.
Figure 2: Disorder-dependent exponent $\overline{z}$ vs. $\epsilon/T$. 

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In Fig. 3, we have plotted $z_{\text{eff}}(t)$ vs. $L(t)$ and here comes the main point in the paper, since this plot allows to discriminate between algebraic and logarithmic asymptotic growth in favour of the latter choice. Indeed, if growth was described correctly by Eq. (32), then $z_{\text{eff}}(L)$ ought to be independent of $L$, for $L \geq L_0 = L(t_0)$, with $z_{\text{eff}}(L) \equiv \bar{z}$. It is clear from Fig. 3 that this is the case only in the pure system ($\epsilon/T = 0$). In all other cases, there is a slow time-increase which is not consistent with Eq. (32). In order to make this feature clear, in Fig. 3 we have drawn the straight horizontal broken lines corresponding to the values of $\bar{z}$ from Table 2. From the data follows i) that we can exclude the asymptotic validity of the algebraic growth law (32) (and with it of the logarithmic dependence on $L$ of energy barriers) and ii) that the rather smooth and near-linear behavior of $z_{\text{eff}}(L)$, for the smaller values of $\epsilon/T$ and $L \geq L_0$, is of the form

$$z_{\text{eff}} = \zeta + bL^\varphi,$$

where $\zeta, b, \varphi$ are fit parameters. Hence, inserting Eq. (37) into Eq. (36) and integrating over time from $t_0$ onward, we derive the growth law in the scaling form

$$\frac{t}{t_0} = \frac{h(L/\lambda)}{h(L_0/\lambda)},$$

where

$$h(x) = x^\zeta \exp(x^\varphi)$$

and where we have set

$$\lambda = (\varphi/b)^{1/\varphi}.\quad (40)$$

We stress that in the above derivation there are no assumptions, other than the form (37) of $z_{\text{eff}}$. Therefore, from the analysis of the numerical data we predict the existence of a crossover from algebraic to logarithmic growth, with the crossover length given by Eq. (40).

With the assumption $L_0 \ll \lambda$ (to be verified a posteriori), and using $h(x) \simeq x^\zeta$ for $x \ll 1$, Eq. (38) becomes

$$(L/\lambda)^\zeta e^{(L/\lambda)^\varphi} = (L_0/\lambda)^\zeta t/t_0$$

with the limiting behaviours

$$L(t) \simeq \begin{cases} \frac{Dt^{1/\zeta}}{\lambda}, & \text{for } L \ll \lambda, \\ \lambda[\ln(t/t_0)]^{1/\varphi}, & \text{for } L \gg \lambda, \end{cases}$$

(42)
Figure 3: Effective exponent $z_{\text{eff}}$ vs. $L$. Straight dashed lines indicate $\tau$. 
Table 3: Values of $\lambda$ and $\zeta$ for various disorder strengths.

| $\epsilon/T$ | $\lambda$  | $\zeta$  |
|-------------|-----------|---------|
| 0.5         | 762.2     | 2.36    |
| 1.0         | 154.1     | 3.15    |
| 1.5         | 64.5      | 4.11    |
| 2.0         | 36.4      | 5.13    |

where

$$D = \frac{L_0}{t_0^{1/\zeta}}. \tag{43}$$

Comparing with Eqs. (30) and (31) and identifying $\psi$ with $\varphi$, we find that the observed growth is of the HH-type, except for the replacement of the pure growth exponent $z$ with the disorder-dependent quantity $\zeta$ (which is distinct from $\tau$). The top line of Eq. (42) accounts for the prior observations \cite{19, 21, 22} of disorder-dependent power-law growth, without need of invoking the argument from Eq. (33) up to Eq. (35). This is the pre-asymptotic regime followed by the asymptotically logarithmic one in the second line of Eq. (42).

In addition, comparing Eq. (29) with Eq. (40) we can make the identification

$$\kappa = \frac{Tb}{\varphi}. \tag{44}$$

The first important conclusion, from the analysis of the numerical data presented above, is the validation, at least at the qualitative level, of the HH scenario. Clearly, it would be important to check also on the quantitative HH prediction $\varphi = 1/4$. However, it turns out to be difficult to obtain the numerical value of $\varphi$ from fitting the data sets in Fig. 3. We have considered three trial values $\varphi = 1/4, 1/2, 1$, and for each one of them we have fitted $\zeta, b$. In Fig. 4 we have superposed these fits on the data sets for $z_{\text{eff}}$ vs. $L$ for $\epsilon/T = 0.5, 1, 1.5$. In the following we use $\varphi = 1$, since this provides the best fit using the least-squares criterion. However, differences are too small to commit to one particular value and we leave the issue open. A precise determination of $\varphi$, as it is evident from Fig. 4, requires to push the accuracy of the numerical computation far beyond what we have done. The corresponding values of $\zeta$ and $\lambda = 1/b$ are listed in Table 3 and plotted in the left and right panels of Fig. 5. Notice that since $L_0 \sim 1$, as it is evident
Figure 4: Comparison of $z_{\text{eff}}$ computed from Eq. (37) with $\psi = 1/4, 1/2, 1$ (continuous lines) and data from the simulations.
from Fig. 1 the assumed inequality $L_0 \ll \lambda$ is satisfied. Furthermore, $\lambda$ is found to decrease with increasing disorder as in Eq. (22) with

$$\lambda \sim (T/\epsilon)^2. \quad (45)$$

Hence, comparing with Eq. (29), we obtain $\kappa \sim \epsilon^2 / T$.

Summarizing, our data for $z_{\text{eff}}$ are evidence for the existence of a crossover of which, however, we can access only the pre-asymptotic region, since the values of the crossover length $\lambda$ are at the edge of the larger values of $L(t)$ reached in the simulations for all disorder values (see Fig. 6). Hence, numerically we can see only the onset of the crossover. Nonetheless, a number of conclusions can be drawn. As Eq. (45) shows, disorder is relevant implying the crossover to disorder-dominated behavior. The pre-asymptotic algebraic growth with a disorder-dependent exponent $\zeta$ signals that disorder, although being globally relevant, acts marginally in the neighborhood of the pure (un-
Figure 6: Fully developed crossover obtained extrapolating the growth law from Eq. (38) (continuous lines) with parameters $\zeta$ and $\lambda$ from Table 3 and $\psi = 1$. Symbols depict the numerical data from the simulation. Asterisks mark the values of the crossover length $\lambda$. 

\[ L(t) \]
4.2 Autocorrelation Function

Having established the existence of a crossover in the growth law, and that disorder is a relevant perturbation, we can anticipate that SU will not hold (see Sec. 2). If the autocorrelation function did obey SU as in Eq. (10), keeping in mind that $\alpha_C = 0$ for pure and disordered systems, the plot of $C(t, t_w)$ against $L(t)/L(t_w)$ (for different values of $\epsilon/T$ and $t_w$) should collapse all the data sets on a single master curve. Figure 7 shows that this is not the case. Instead, there is collapse only when $t_w$ is varied keeping $\epsilon/T$ fixed, but the master curves of these partial collapses, corresponding to the different values of $\epsilon/T$, are distinct. What we have called partial collapse has also been reported in the context of the diluted Ising ferromagnet [17, 37], without reaching, however, a definite conclusion on the SU validity.

The behavior in Fig. 7 can be accounted for on the basis of the two-variable scaling $C(t, t_w, \epsilon) = \tilde{F}_C(x, y)$ in Eq. (12) with $\alpha_C = 0$. The pure case $\tilde{F}_C(x, \infty) = F_C(x)$ corresponds to the lower black curve, where the collapse is very good since all the dependence on $t_w$ takes place through $x$. However, in the presence of disorder, $y$ is finite. The remaining curves in the plot are explained considering that with the $t_w$-values used in the simulations, the variation of $y$ is small while keeping $\epsilon/T$ fixed. This produces the quasi-collapse on the coloured curves, each one corresponding to a distinct value of $\epsilon/T$. But the variation of $y$ becomes sizable when $\epsilon/T$ (and thereby $\lambda$) is changed, producing the separation of the curves with different colors. The trend in Fig. 7 shows that $\tilde{F}_C(x, y)$ increases as $y$ decreases, for fixed $x$. Hence, the bundle of curves generated by letting $y$ vary is bounded below

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The statements that the perturbation is marginal and that the pure fixed point is unstable are compatible when nonlinear terms in the expansion around the fixed point are taken into account [36].
Figure 7: Autocorrelation function for different disorder values. The $t_w$ values have been taken according to the rule $t_w(k) = t_{\text{min}}(t_m/2t_{\text{min}})^{(k-1)/10}$ where $t_{\text{min}} = 200$, $k$ is an integer running from 1 to 11 and $t_m$ is the maximum time in Table 1. Accordingly, the waiting times are logarithmically spaced between $t_w = 200$ and $t_w = t_m/2$. 
by the pure scaling function $F_C(x)$, and must be bounded above by the numerically inaccessible scaling function $\hat{F}_C(x)$ of the disorder-dominated fixed point. Exactly the same behavior is observed in the autocorrelation function of the $d = 1$ RFIM [10].

Within the framework of our findings, we conclude that previously reported observations of SU in the autocorrelation function for the $d = 3$ RFIM [14, 15] and the $d = 2$ RBIM [22] are due to time-ranges and disorder values deep inside the pre-asymptotic regime, where it is not possible to detect the onset of the crossover.

### 4.3 Autoresponse Function

The qualitative difference between the autoresponse and autocorrelation functions is that the exponent $\alpha_\chi$ in the scaling relation

$$\chi(t, t_w, \epsilon) = L_w^{-\alpha_\chi} \tilde{F}_\chi(x, y),$$

is not constrained to vanish, like $\alpha_C$, by geometrical requirements. Hence, in principle, the crossover could take the most general form of Eq. (14) with $\hat{\alpha}_\chi \neq \alpha_\chi$.

Before analysing the data it is useful to summarize the structure of scaling in the pure case [28], where Eq. (16) reads

$$\chi(t, t_w) = L_w^{-\alpha_\chi} F_\chi(x)$$

and for large $x$ the scaling function decays with the same exponent appearing in the prefactor

$$F_\chi(x) \sim x^{-\alpha_\chi}. \quad (48)$$

As a matter of fact, the measurement of the decaying slope of $\chi(x, t_w)$ for large $x$ is the fast numerical method to extract the scaling exponent $\alpha_\chi$. The physical origin of this feature can be understood introducing the effective response associated with a single interface\(^5\), which is defined by [27]

$$\chi(t, t_w) = \rho(t) \chi_{\text{eff}}(t, t_w) \quad (49)$$

where $\rho(t) = L^{-1}(t)$ is the interface density. Then, from the scaling form in Eq. (47) follows

$$\chi_{\text{eff}}(t, t_w) \sim L_w^{(1-\alpha_\chi)} x F_\chi(x). \quad (50)$$

\(^5\)For instance, a configuration with a single interface has been introduced in Eq. (16).
On the other hand, if there is only one interface in the system the response is not expected to age, yielding a time-translation invariant $\chi_{\text{eff}}(t - t_w)$. Therefore, regarding the left hand side as dependent on time through $L(t - t_w)$ and considering the large $x$ regime where it is possible to replace $L(t - t_w)$ by $L(t)$, the dependence on $L_w$ must drop out of the right hand side implying $xF_\chi(x) \sim x^{(1 - \alpha_x)}$. This, in turn, implies the validity of Eq. (48) and

$$\chi_{\text{eff}}(t, t_w) \sim \chi_{\text{eff}}(t, t_w) \sim L^{1 - \alpha_x}. \quad (51)$$

Let us now go to the data in presence of disorder. We have computed $\chi(t, t_w, \epsilon)$ with the algorithm derived in Ref. [38], which does not require the switching on of the external perturbation, thus improving the precision and the efficiency of the numerical computation. The use of algorithms of this type is indispensable for the study of linear response when the system is quenched to $T = 0$. The quantity $L^{\alpha_x} \chi(t, t_w, \epsilon)$ has been plotted in Fig. 8 against $L/L_w$ for different values of $\epsilon/T$ and $t_w$, using for all disorder values $\alpha_x = 0.6$, which is the value producing the best data collapse in the pure case (top black curves). The structure of the plot follows the same pattern as in Fig. 7 with good partial collapse when $\epsilon/T$ is kept fixed and distinct master curves when $\epsilon/T$ is varied, for the same reasons explained in the discussion of the autocorrelation function. The collapse of the curves with the pure exponent $\alpha_x$ shows clearly that the numerical data belong to the preasymptotic region controlled by the pure fixed point up to the onset of the crossover, since the $y$ dependence of the scaling function $\tilde{F}(x, y)$ is manifest in Fig. 8.

Furthermore, since the different curves in the plot appear to be obtained one from the other by a vertical shift, in this region of parameters the dependence on $x$ and $y$ can be taken to be of the multiplicative type

$$\tilde{F}(x, y) = f(x)g(y) \quad (52)$$

with $f(x)$ decaying algebraically for large $x$, exactly as in Eq. (48) for the pure case. In order to highlight this feature, in the inset of Fig. 8 we have plotted $\tilde{F}(x, y)/\tilde{F}(x = 2, y)$ obtaining the superposition of the different curves, that is the elimination of the $y$-dependence, which is a nice check on the validity of the assumption made in Eq. (52). Therefore, the curve in the inset is the normalized plot of $f(x)$, whose decay in agreement with Eq. (48) is a confirmation i) that scaling is controlled by the pure fixed point and ii) of the validity of the single interface structure of the response.
Figure 8: Zero-field-cooled susceptibility for different disorder values and $t_w$ values chosen according to the rule in the caption of Fig. 7.
As a further check of this latter point, we have measured the ZFCS \( \chi_{\text{single}}(t, t_w, \epsilon) \) after preparing an initial state with a single interface\(^6\) and for different disorder values. The plot in Fig. 9, which also includes the pure case, does indeed show that \( \chi_{\text{single}}(t, t_w, \epsilon) \) grows with a power law in agreement with Eq. (51). The exponent does not seem to depend on \( \epsilon/T \) and it is consistent with the expected value \( 1 - \alpha_{\chi} \simeq 0.4 \). Conversely, in front of the power there is a prefactor \( A \) dependent on \( \epsilon/T \) (see Table 4), which enters into the factor \( g(y) \), in Eq. (52), and is responsible for the separation of the master curves in Fig. 8.

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\(^6\)The initial single interface is prepared along a diagonal in the lattice, in such a way that spins on the interface experience zero Weiss field. If the interface was prepared along the horizontal or vertical axis, the Weiss field would not be zero and no flip could take place with the accelerated algorithm presented at the beginning of this section.
Table 4: Amplitude $A$ of $\chi\text{single}(t, t_w)$ extracted from the data of Fig. 9 for different disorder values.

| $\epsilon/T$ | $A$   |
|------------|------|
| 0.0        | 0.00235 |
| 0.5        | 0.00184 |
| 1.0        | 0.00125 |
| 1.5        | 0.00093 |

In summary, since our data do not allow to access the full crossover region, we cannot make statements on the interesting question whether the exponent $\hat{\alpha}_\chi$ will be different or not from $\alpha_\chi$ when the disorder-dominated asymptotics are reached. However, precisely because what we observe is the pre-asymptotic regime controlled by the pure fixed point, this result is important in itself, as explained in detail in Ref. [11], since it brings new and independent evidence on the value of $\alpha_\chi$ in the pure system. In fact, together with the $d = 1$ result in Eq. (23), this result substantiates the conjecture put forward [27] for the pure system that $\alpha_\chi$ depends on dimensionality according to

$$\alpha_\chi = (d - 1)/2,$$

when $d \leq 3$. The numerical finding $\alpha_\chi \simeq 0.6$ does not exactly coincide with the value $\alpha_\chi = 0.5$, predicted by Eq. (53) for $d = 2$. This means that either the phenomenological formula or the precision of the numerical computation still need improvement. However, it seems clear that the alternative conjecture [39, 40, 41] predicting

$$\alpha_\chi = 1,$$

with no dependence on dimensionality [40], is ruled out by the data from the $d = 1$ and $d = 2$ RBIM. Preliminary results on this point have been presented in Ref. [11].

5 Summary and Discussion

Let us conclude this paper with a summary and discussion of the results presented here. We have reported results from the most comprehensive sim-
ulations to date of domain growth in the $d = 2$ ferromagnetic RBIM with non-conserved Glauber kinetics. We undertook this computationally expensive study with a specific purpose in mind, viz., to obtain a clear understanding of possible crossovers in the domain growth law and other statistical properties of the evolution morphology. As regards the domain growth law, our results are incompatible with algebraic growth scenario of Ref. [19] and are compatible with a crossover from a pre-asymptotic regime having power-law growth with a disorder-dependent exponent to an asymptotic logarithmic regime $[L(t) \sim (\ln t)^{1/\varphi}]$. The pre-asymptotic regime is consistent with existing results for the RBIM [19, 21, 22]. To the best of our knowledge, the predictive power of our result for the late-stage result is new and allow us to give numerical support to the HH scenario. However, in spite of the huge numerical effort, we cannot conclusively determine the exponent $\varphi$ which characterizes the logarithmic regime. Clearly, more work is needed in this direction.

We use RG arguments to understand the crossover in the growth law in terms of the competition between a pure fixed point and a disordered fixed point. Within this framework the corresponding crossover in two-time quantities, like the autocorrelation function $C(t, t_w, \epsilon)$ and the autoresponse function $\chi(t, t_w, \epsilon)$, is expected. The scaling function for $C(t, t_w, \epsilon)$ shows the onset of the crossover from a pure form to a disordered one. Even though the numerical data are limited to the pre-asymptotic region, there is clear evidence that the SU hypothesis does not hold for two-time quantities. A similar behavior is observed in the data for $\chi(t, t_w, \epsilon)$, where the crossover region is not entered deep enough for $\tilde{F}_\chi(x, y)$ to develop the $y$-singularity which eventually should change the scaling exponent from the pure value $\alpha_\chi$ to the disorder value $\hat{\alpha}_\chi$. Nonetheless, the measurement of the pure exponent $\alpha_\chi \simeq 0.6$ is quite an interesting result in itself, adding new evidence in support of the phenomenological conjecture of Eq. (53).

At a more general level, the RG framework formulated in this paper, in the context of the $d = 2$ RBIM, proves to be comprehensive enough to give also a natural explanation for the known results for the $d = 1$ RFIM and $d = 1$ RBIM. As such, it is a good candidate for the understanding of coarsening in disordered systems, whose predictive power could be tested at higher dimensionality or in other systems, like the site diluted Ising model, where the crossover region perhaps is more easily accessible than in the RBIM [17].
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