



Mixing time of random walk on dynamical random cluster

Andrea Lelli¹ · Alexandre Stauffer^{1,2}

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Abstract

We study the mixing time of a random walker who moves inside a dynamical random cluster model on the d -dimensional torus of side-length n . In this model, edges switch at rate μ between *open* and *closed*, following a Glauber dynamics for the random cluster model with parameters p, q . At the same time, the walker jumps at rate 1 as a simple random walk on the torus, but is only allowed to traverse open edges. We show that for small enough p the mixing time of the random walker is of order n^2/μ . In our proof we construct a non-Markovian coupling through a multi-scale analysis of the environment, which we believe could be more widely applicable.

Keywords Mixing time · Random walk · Time inhomogeneous Markov chains · Random cluster

Mathematics Subject Classification Primary 60K35; Secondary 60K37

1 Introduction

We study the mixing time of a random walk on a dynamical random cluster model in \mathbb{T}_n^d , the d -dimensional torus of side-length n . In this model, each edge of \mathbb{T}_n^d can be in either of two states: *open* or *closed*. At time 0, we take the state of the edges to be distributed according to the random cluster measure with parameters $p \in (0, 1)$ and $q > 0$. That is, for any subset of edges $\omega \subset E(\mathbb{T}_n^d)$, with $E(\mathbb{T}_n^d)$ denoting the set of

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✉ Alexandre Stauffer
a.stauffer@bath.ac.uk

¹ Department of Mathematical Sciences, University of Bath, Bath, UK

² Department of Mathematics, King's College, London, UK

edges of the torus, the probability that the set of open edges at time 0 is equal to ω is

$$\nu(\omega) = \frac{1}{Z} p^{|\omega|} (1-p)^{|E(\mathbb{T}_n^d) \setminus \omega|} q^{\kappa(\omega)}, \quad (1.1)$$

where $\kappa(\omega)$ is the number of connected components obtained in the graph with vertex set \mathbb{T}_n^d and edge set ω , and $Z = Z(d, p, q) > 0$ is just a normalizing constant so that the above is a probability measure. Instead of representing the state of the edges by the set ω of open edges, we will often represent it by an element $\eta \in \{0, 1\}^{E(\mathbb{T}_n^d)}$, with $\eta(e) = 0$ meaning that the edge $e \in E(\mathbb{T}_n^d)$ is closed and $\eta(e) = 1$ meaning that e is open. Thus, given η , we have $\omega = \{e \in E(\mathbb{T}_n^d) : \eta(e) = 1\}$.

From time 0, edges change their state following a continuous-time Glauber dynamics. Thus, given a parameter $\mu > 0$, each edge $e \in E(\mathbb{T}_n^d)$ has a Poisson clock of rate μ , and when the clock of e rings, the state of e is resampled (open or closed) according to the conditional probability obtained from ν in (1.1) conditioned on the states of all the other edges. This resampling can be easily described: if the clock of e rings at time t , then the probability that e becomes open at time t is equal to

$$\begin{aligned} & p, \text{ if } e \text{ is not a cut-edge at time } t-, \\ & \frac{p}{p+(1-p)q}, \text{ if } e \text{ is a cut-edge at time } t-, \end{aligned} \quad (1.2)$$

where an edge e is called a *cut-edge* if modifying the state of e (while keeping the state of the other edges unaltered) causes a change in the number of connected components in the configuration. Note that whether an edge e is a cut-edge for a configuration η is, in fact, independent of $\eta(e)$. We let $\eta_t \in \{0, 1\}^{E(\mathbb{T}_n^d)}$ denote the configuration that gives the state of the edges at time t .

On top of this dynamic environment we place a random walker which starts from the origin of \mathbb{T}_n^d and has a Poisson clock of rate 1. When the clock of the walker rings, the walker chooses an edge uniformly at random from the set of edges that are adjacent to its current location, regardless of their states. If the chosen edge is open at that time, then the walker traverses the edge, otherwise the walker stays put. We denote by $X_t \in \mathbb{T}_n^d$ the position of the walker at time t , and let

$$\{M_t\}_{t \geq 0} = \{X_t, \eta_t\}_{t \geq 0},$$

denote the *full system* composed of the walker $\{X_t\}_{t \geq 0}$ and the environment $\{\eta_t\}_{t \geq 0}$. We note that $\{M_t\}_{t \geq 0}$ and $\{\eta_t\}_{t \geq 0}$ are Markov chains, while $\{X_t\}_{t \geq 0}$ is not.

One can check (for example, by reversibility) that if π denotes the uniform probability measure on \mathbb{T}_n^d then $\pi \times \nu$ is the unique stationary distribution of $\{M_t\}_t$.

Let T_{mix} denote the mixing time of the full system, starting from the worst-case initial state. In other words, given $x \in \mathbb{T}_n^d$ and $\xi \in \{0, 1\}^{E(\mathbb{T}_n^d)}$, let $T_{\text{mix}}^{x, \xi}$ be the smallest t such that, starting from $M_0 = (x, \xi)$, the total variation distance between the distribution of M_t and $\pi \times \nu$ is smaller than a given constant, which for concreteness we take to be $1/4$. Then $T_{\text{mix}} = \max_{x, \xi} T_{\text{mix}}^{x, \xi}$.

Our main result establishes that the mixing time is of order n^2/μ for all small enough p . We remark that p and q are considered to be constants independent of

n , while μ may depend on n ; in particular, a natural case in the context of dynamic networks is that $\mu \rightarrow 0$ as $n \rightarrow \infty$.

Theorem 1.1 *Given any $q > 0$ and any dimension $d \geq 1$, there exists a positive $p_0 > 0$ so that for all $p \in (0, p_0)$ there exists $C_1 = C_1(d, p, q) > 0$ for which*

$$T_{\text{mix}} \leq C_1 \frac{n^2}{\mu}, \quad \text{for all } \mu = \mu(n) > 0 \text{ and all } n \geq 1.$$

Previous works were restricted to the case $q = 1$, which is known as *dynamical percolation*. This is a much simpler case: both expressions in (1.2) are equal, so edges update independently of one another. This allowed the construction of very clean regeneration arguments to bound the mixing time and other quantities, especially in the subcritical regime (see related works below in Sect. 1.2). To the best of our knowledge, our work is the first to analyze the mixing time when edge updates are not independent.

The regeneration arguments developed for dynamical percolation cannot be applied when $q \neq 1$. We go around this by constructing a delicate *non-Markovian* coupling using a multi-scale analysis of the environment. We believe this to be a novel ideal and regard it as one of our main contributions.

In a nutshell, the idea is to develop a coupling between two random walkers moving on dynamical random clusters. To do this, we employ a multi-scale analysis of the environment alone, so as to control the evolution of the state of the edges. We then use this multi-scale analysis to identify *good* and *bad* regions of the environment. Depending on whether the walkers are passing through a space-time region that is good or bad, we employ a different coupling strategy for the walkers. It is crucial that the multi-scale analysis does not reveal all the information regarding the environment so that, conditioned on which regions are good and which regions are bad, there is still enough randomness in the environment for the coupling argument to be carried out. This leads to a quite delicate coupling, which is non-Markovian. In particular, in order to decide which coupling strategy to apply for the walkers at a time t , we will use *future* information about the environment (i.e., information about the state of the edges after time t). The reason is that, if the walkers are passing through a space-time region that is approaching (in time) a bad region of the environment, we will already need to start employing a coupling that prepares the walkers before they enter into the bad regions.

We believe that the idea of using a multi-scale analysis to develop a coupling argument can be more widely applicable to analyze the mixing time of random walks on particle systems. Our proof actually relies very little on specific properties of the random cluster dynamics, and can be readily adapted to other settings. We decided to carry out the proof for the random cluster dynamics for the sake of clarity, since the general conditions under which this argument can be made to work are rather cumbersome to state.

We will give a more thorough description of the main ideas of the proof in Sect. 1.4, since first, in Sect. 1.3, we will need to introduce an auxiliary process.

1.1 Lower bounds on the mixing time

We also derive matching lower bounds on the mixing time. We start by stating a straightforward generalization of the lower bound from [11].

We consider a larger class of models, which we refer to as *continuous-time random walks on general dynamical percolation*, where the word general is to mean that the percolation process may not be independent. Let $\{X_t, \eta_t\}_{t \geq 0}$ be a continuous-time Markov chain where the walker X_t jumps at rate 1 and can only traverse open edges of \mathbb{T}_n^d , and the environment $\{\eta_t\}_t$ is a Markov chain on $\{0, 1\}^{E(\mathbb{T}_n^d)}$ where edges refresh their state at rate μ independently of the walker. As usual μ may depend on n . Let π be the uniform distribution on \mathbb{T}_n^d and let ν be the stationary distribution of the Markov chain $\{\eta_t\}_t$.

We recall some fundamental definitions. The *spectral gap* γ of a reversible Markov chain is defined as

$$\gamma = \inf_f \frac{\mathcal{E}(f, f)}{\text{Var}(f)},$$

where the infimum is over all functions f from the state space to \mathbb{R} with $\text{Var}(f) \neq 0$, the variance $\text{Var}(f)$ being with respect to the stationary distribution of the chain, and $\mathcal{E}(f, f)$ is the so-called Dirichlet form. The *relaxation time* of the said Markov chain is defined as

$$T_{\text{rel}} = \gamma^{-1}.$$

Given a time interval $I \subset \mathbb{R}_+$, we say that an edge is I -open, if it is open at some time during I . Then, for any vertex $x \in \mathbb{T}_n^d$ and any time interval $I \subset \mathbb{R}_+$, we let $\mathcal{C}_x(I)$ denote the connected component of I -open edges from x . Finally, given a subset $S \subset \mathbb{T}_n^d$, let $\text{diam}(S) = \max_{x, y \in S} \|x - y\|_1$ be the diameter of S , where $\|x - y\|_1$ is the L_1 distance (or, equivalently, the length of the shortest path) between x and y in \mathbb{T}_n^d . We require the following two assumptions from the process $\{X_t, \eta_t\}_{t \geq 0}$:

$$\pi \times \nu \text{ is the stationary distribution of } \{X_t, \eta_t\}_{t \geq 0}, \quad (1.3)$$

and

$$\begin{aligned} \exists \delta > 0 \text{ and } C_2 > 0 \text{ such that for any } x \in \mathbb{T}_n^d, \text{ any } \mu = \mu(n) > 0 \\ \text{and any } n \geq 1 \text{ we have } \mathbb{E}_\nu \left(D_{x, \delta}^2 \right) \leq C_2, \end{aligned} \quad (1.4)$$

where $D_{x, \delta} = \text{diam}(\mathcal{C}_x([0, \delta]))$ and \mathbb{E}_ν denotes the expectation with respect to the stationary measure of the environment. The assumption in (1.3) just says that the stationary measure of the walker is uniform, while (1.4) gives that the environment is strictly subcritical.

Theorem 1.2 *Let $\{X_t, \eta_t\}_{t \geq 0}$ be a random walk in a general dynamical percolation satisfying (1.3) and (1.4) above. Then, there exist a constant $C_3 > 0$ depending only on d such that*

$$T_{\text{rel}} \geq \frac{C_3 \delta n^2}{C_2}.$$

A natural setting is when the environment starts from its stationary distribution. For this, let ν_t stand for the distribution of (X_t, η_t) where the walker starts from the origin and the environment starts from stationarity (that is, η_0 is distributed as ν). Then, $\|\nu_t - \pi \times \nu\|_{\text{TV}}$ is the total variation distance between ν_t and the stationary measure of $\{X_t, \eta_t\}_t$.

Theorem 1.3 *Let $\{X_t, \eta_t\}_{t \geq 0}$ be a random walk in a general dynamical percolation satisfying (1.3) and (1.4) above. Then, there exists a constant $C_4 > 0$ depending only on d such that, for any $\epsilon > 0$,*

$$\text{if } t \leq \frac{C_4}{C_2} \epsilon^{\frac{d+2}{d}} \delta n^2 \text{ then } \|\nu_t - \pi \times \nu\|_{\text{TV}} \geq 1 - \epsilon.$$

Moreover, there exists a constant $C_5 > 0$ depending only on d such that, for any $t \geq \delta$,

$$\mathbb{E}_{\nu_t} \left(\|X_t - X_0\|_1^2 \right) \leq C_5 C_2 \frac{t}{\delta},$$

where \mathbb{E}_{ν_t} stands for the expectation with respect to ν_t .

The proofs of Theorems 1.2 and 1.3 are identical to the ones for random walk on dynamical percolation from [11]. For the sake of completeness, we add the proofs in Sect. 8.

We want to apply the above theorems to derive lower bounds on the mixing time of a random walk in dynamical random cluster. It is clear that (1.3) holds in this case. We will show in Sect. 9 that (1.4) also holds, obtaining the corollary below. For any $q \geq 1$, let p_c^q be the critical probability for the appearance of an infinite cluster in the random cluster model on \mathbb{Z}^d .

Corollary 1.4 *If $\{X_t, \eta_t\}_{t \geq 0}$ is a random walker in the dynamic random cluster model, then for any $q \geq 1$ and any $p < p_c^q$, there exists a constant $c = c(d, q, p) > 0$ such that the relaxation time of the full system and the mixing time starting from a stationary environment is at least cn^2/μ . If $q < 1$, then for all small enough p the same conclusion holds.*

The proof of the lower bound is much simpler than that of the upper bound, allowing us to derive it in the whole subcritical regime when $q \geq 1$. In fact, when $q \geq 1$, the proof follows by using a sprinkling lemma to compare two random cluster configurations with densities $p < p'$ (Lemma 9.1), and the exponential decay of cluster sizes in the subcritical regime. When $q < 1$, exponential decay of cluster sizes is only known for small enough p . In fact, for $q < 1$, even the existence of a single critical value p_c^q separating a subcritical phase and a supercritical phase, has not yet been proved.

We expect the upper bound of order n^2/μ to hold in the whole subcritical regime as well, however our proof technique requires the percolation process to be a small perturbation of subcritical independent percolation, in a sense that we better explain in Remark 1.5, after introducing the \star -process.

1.2 Related works

We will restrict our discussion to works dealing with the mixing time of random walks on *dynamic* environments, as otherwise there is simply a plethora of works. We also remark that, if the environment is allowed to evolve in an arbitrary fashion (for example, by taking any sequence of graphs on a fixed vertex set), then several problems may arise. For example, there may not be a stationary distribution for the walker. Moreover, even if there is a stationary distribution, the distribution of the walker may not converge to stationarity, or the total variation distance to stationarity may not be monotone in time.

Random walk on dynamical percolation on \mathbb{T}_n^d . This model is equivalent to the model we described restricted to $q = 1$. This special case is already quite challenging but some results have been obtained recently. First note that, when $q = 1$, the two probabilities in (1.2) become equal, and when an edge updates, it does so *independently* of the other edges, becoming open with probability p or closed with probability $1 - p$. Though in this case edges evolve independently of one another, there are strong dependences between the location of the walker and the state of the edges (especially if $\mu \rightarrow 0$ as $n \rightarrow \infty$, since edges update very slowly in comparison to the rate of jump of the walker).

The random walk on dynamical percolation model was introduced by Peres, Stauffer and Steif [11], where it is shown that, in the whole subcritical regime,¹ the mixing time is of order n^2/μ . We remark that in [11] both upper and lower bounds of order n^2/μ were derived for T_{mix} . Recall that n^2 is the order of the mixing time of a simple random walk on the *static* torus (that is, where all edges are open at all times). So, in a subcritical dynamical percolation, the walker is delayed by a factor of $1/\mu$, which is the expected time that a single edge takes to refresh.

Later, Peres, Sousi and Steif [10] analyzed the supercritical regime and showed that, for p large enough, the mixing time is at most $(\log n)^a \left(n^2 + \frac{1}{\mu}\right)$ for some constant $a > 0$. Their upper bound is not believed to be tight: one expects that, in the whole supercritical regime, the mixing time is of order $n^2 + \frac{1}{\mu}$. This remains an interesting open problem. Their proof makes strong use of isoperimetric properties of the infinite cluster of supercritical percolation, which are only known for $q = 1$. With regard to the *critical* regime, the only known result is that the mixing time is of order at most $\frac{n^2}{\mu}$, which is the mixing time in the subcritical regime [7]. It is not inconceivable that the mixing time in the critical case is in fact of smaller order than $\frac{n^2}{\mu}$.

Random walk on dynamical percolation on other graphs. Sousi and Thomas [14] studied the case where the torus is replaced by the *complete graph*. This is a simpler

¹ That is, for any $p < p_c$ with $p_c = p_c(d)$ being the critical probability for the existence of an infinite cluster in percolation in \mathbb{Z}^d

case due to the lack of an underlying geometry, but for which a more detailed analysis can be carried out. They established the order of the mixing time in that case, and also the occurrence of a cut-off phenomenon. We remark that if the walker is at some vertex v and we know that an edge incident to v is updating to open, but we refrain from observing which of the edges incident to v is updating, then the other endpoint of this edge is uniform among all vertices (but v). So, by traversing this edge (call it e), after one additional step, the walker can find itself in a location that is essentially uniform, so very close to stationarity. Though suggestive, this is not enough to establish the mixing time, as one still needs to control that the walker “forgets” that e is now open (that is, the walker may be close to stationarity, but the full system is not). Still, this illustrates the kind of simplification that the lack of an underlying geometry brings.

The last work we mention for the random walk on dynamical percolation model is a recent result by Hermon and Sousi [7]. They developed a comparison principle and showed that, for *any* graph G , the so-called *spectral profile mixing time* for the random walk on dynamical percolation on G is at most $\frac{1}{\mu}$ times the spectral profile mixing time of simple random walk on (the static graph) \bar{G} .

In all the above results, it was crucial that when $q = 1$ edges update independently of one another. The main objective of our work is to develop a technique that can go beyond the dynamical percolation case and which can deal with environments whose edge updates may depend on one another, including the case of unbounded dependences such as in the dynamical random cluster.

Other models. We end this section by mentioning two lines of work. In the first one, Avena et al. [1, 2] studied a different dynamic on the environment, where instead of dynamical percolation one has a *dynamic configuration model*. This model has some intuitive similarities with the dynamical percolation on the *complete graph*, in the sense that it also lacks an underlying geometry. They studied the mixing time and the occurrence of a cut-off phenomenon in this setting, but restricted to a random walker that is non-backtracking. This helps the walker to move away from its current location, strongly reducing dependences between the walker and the environment.

Finally, the second line of work we mention is that of [12, 13]. They considered the case of a discrete-time random walk on a graph with a fixed set of vertices, but which evolves over time by means of an arbitrary sequence of graphs on that vertex set. The goal of their work is very different to ours; for example, they want to understand which conditions on the sequence of graphs one can impose to guarantee that the mixing time is polynomial. They also derive results for the hitting time and cover time. We refer to [12, 13] and references therein for a list of known results on dynamic graphs that go beyond the mixing time. We also refer the reader to [4] for results on a model similar to random walks on dynamical percolation on the complete graph.

1.3 The \star -process: retaining some randomness

Before giving the ideas of our proof, we need to describe a different representation of the full system, which is inspired by [11]. Recall that each edge has a Poisson clock of rate μ associated to it, which gives the times at which the edge is updated. To each update of an edge, we can decide whether the edge becomes open or closed by

sampling an independent random variable U with a uniform distribution in $(0, 1)$, and then making the edge open if and only if the edge is not a cut-edge and $U < p$ or the edge is a cut-edge and $U < \frac{p}{p+(1-p)q}$. Now, let

$$p_{\min} = \min \left\{ p, \frac{p}{p + (1-p)q} \right\} \quad \text{and} \quad p_{\max} = \max \left\{ p, \frac{p}{p + (1-p)q} \right\};$$

thus, $p_{\min} = \frac{p}{p+(1-p)q}$ if $q > 1$ and $p_{\min} = p$ if $q < 1$. Note that if U turns out to be in the interval $(0, p_{\min}) \cup (p_{\max}, 1)$ the outcome of the update (i.e., whether the edge becomes open or closed) is determined regardless of whether or not the edge is a cut-edge. In other words, the update is *oblivious* to the current configuration, and we will refer to those updates as \star -updates. We then let

$$p_{\star} = p_{\min} + 1 - p_{\max} \in (0, 1)$$

be the probability that a given update is a \star -update.

We now define the update of an edge e in two stages. First, we sample an independent random variable U_{\star} , which is uniform in $(0, 1)$, so that if $U_{\star} < p_{\star}$, then the update is a \star -update, otherwise it is not a \star -update. Next, we use the random variable U to determine whether e updates to open or closed. More precisely, in the case of a \star -update, we make e open if $U < \frac{p_{\min}}{p_{\star}}$, otherwise e becomes closed. In the case of a non- \star -update, we need to inspect the current configuration to see whether e is a cut-edge or not. In particular, we need to perform what we call an *exploration of e* , which means that we perform a local search from the endpoints of e that traverses only open edges in order to determine what are the open clusters of the endpoints of e . Hence, each update of e will be represented by a tuple (s, U_{\star}, U) , where $s > 0$ is the time at which the update occurs, $U_{\star} \in (0, 1)$ is the variable used to decide whether the update is a \star -update, and $U \in (0, 1)$ is the random variable governing whether the edge is to be updated open or closed.

We use this to introduce another Markov process which we denote by $\{M_t^{\star}\}_{t \geq 0} = \{X_t, \eta_t^{\star}\}_{t \geq 0}$, and which we refer to as the \star -process. This process will retain more randomness than $\{M_t\}_{t \geq 0}$ and its state space will be

$$\Omega^{\star} = \left\{ (v, \eta^{\star}) \in \mathbb{T}_n^d \times \{0, 1, \star\}^{E(\mathbb{T}_n^d)} : \eta^{\star}(e) \in \{0, 1\} \text{ for each edge } e \text{ adjacent to } v \right\}.$$

So an edge will be allowed to be in an additional state, called \star , which means that in its last update the edge underwent a \star -update. However, we do not allow that edges adjacent to the walker are in state \star .

The \star -process evolves as follows. If the Poisson clock of an edge e rings, we look at the variable U_{\star} associated with this update and determine whether the update is a \star -update. If the update is a \star -update and if e is not currently adjacent to the walker, then we make the state of e equal to \star . If e is adjacent to the walker, then we look at the variable U associated with this update and determine whether e is open or closed. Finally, if the update is not a \star -update, then we perform an exploration of e as mentioned above. The difference is that, in such an exploration, we may run into

edges that are in state \star . For each such edge, we immediately sample whether that edge is open or closed by using the random variable U associated with its last update. We proceed in this way until the exploration ends and we have fully determined the cluster of each endpoint of e . At this moment, we know whether or not e is a cut-edge, and we can use the random variable U associated to the update of e to determine whether e is to be made open or closed. There is still one final case to be described: when it is the clock of the walker that rings. Suppose this happens and the walker jumps from a vertex v to a vertex w . Then, if there are edges adjacent to w at state \star we sample the state of such edges (using the random variables U associated to their last update) and switch them to open or closed, appropriately.

Note that, conditioned on the position of the walker and on the state 0, 1 or \star of each edge, we gain no information concerning whether the edges in state \star are open or closed. In particular, each such edge is open with probability $\frac{p_{\min}}{p_{\star}}$ (which is the probability that the random variable U associated to their last update is at most $\frac{p_{\min}}{p_{\star}}$). Therefore, we do not need to keep track of the variables U related to the last \star -update of each edge, since we can sample U whenever needed independently of the whole trajectory of the process. The \star -process is thus a Markov process.

Remark 1.5 When $q = 1$, we have $p_{\min} = p_{\max}$, and so $p_{\star} = 1$: all updates are \star -updates, as in this case the random cluster model reduces to dynamical percolation. If $q \neq 1$, then as $p \rightarrow 0$ we have that $p_{\max} - p_{\min} \rightarrow 0$ and so $p_{\star} \rightarrow 1$. Therefore, for any fixed q and all small enough p , the dynamic random cluster model can be viewed as a small perturbation of dynamical percolation. We also obtain that edges of state \star are open with probability $\frac{p_{\min}}{p_{\star}} < p_c$, so they form a subcritical percolation process as well. Those are the properties that play an essential role in the constructions of the multi-scale analysis and the coupling used to establish the upper bound on the mixing time (Theorem 1.1).

1.4 Proof overview

We will only give an overview of the upper bound, which is our main result and by far the most involved proof. We start recalling the proof in [11] for the subcritical regime when $q = 1$. There they also define the \star -process (which they denote by \tilde{M}_t). Recall that, when $q = 1$, we have $p_{\star} = 1$, so all updates are \star -updates. With this, they define a stopping time τ_0 as the first time at which

all edges adjacent to the walker are closed, and all remaining edges are in state \star .
(1.5)

Then, one can define a sequence of times τ_1, τ_2, \dots so that τ_i is the first time after $\tau_{i-1} + C/\mu$, for some fixed constant $C > 0$, at which the event in (1.5) happens. These are regeneration times in the sense that the evolution of the full system from τ_i does not depend on what happened before τ_i . Once the full system is at a regeneration time τ_i , with positive probability the following sequence of events happen within time $\tau_i + C/\mu$:

- (i) an edge e adjacent to the walker opens
- (ii) when the walker jumps to the other endpoint of e , all the adjacent edges (which are in state \star) are sampled closed
- (iii) e remains open for some time of order $1/\mu$
- (iv) e closes before any of the other edges adjacent to e open, thereby locking the walker in one of e 's endpoints, and
- (v) the edges adjacent to the other endpoint of e (i.e., opposite to the location of the walker) refresh before the edges adjacent to the walker refresh.

When these events occur, the walker does nothing more than a jump to a uniformly random neighbor, and immediately gets back to a regeneration time (so $\tau_{i+1} = \tau_i + C/\mu$); such a regeneration time is then called a *simple random walk regeneration* since, at the end, what the walker did was just one step of a simple random walk in \mathbb{T}_n^d .

The proof in [11] then goes by showing that the $\tau_{i+1} - \tau_i$ are of order $\frac{1}{\mu}$. Therefore, after time $\frac{n^2}{\mu}$, the walker underwent an order of n^2 regeneration times, a positive fraction of which being simple random walk regeneration. So it is possible to couple the full system with another copy of the full system so that, whenever the walker does a simple random walk regeneration, we employ one of the standard couplings of simple random walks on the torus. On the other hand, if the regeneration time is not a simple random walk regeneration, we couple the motion of the two walkers from one regeneration time to the next identically, so that the distance between the walkers does not change. Since an order of n^2 steps is necessary to couple two simple random walks on \mathbb{T}_n^d , we get that performing an order of n^2 simple random walk regenerations is enough to couple the two processes, which translates to a mixing time of order n^2/μ .

If we try to mimic the steps above for the case $q \neq 1$, we immediately run into the issue that the event (1.5) now occurs very rarely. In fact, since non- \star -updates occur with positive probability, we will typically have a positive density of non- \star -edges. Therefore, it will take an exponential amount of time to reach a regeneration time as in (1.5), rendering this strategy useless.

We will devise a different strategy. We will, as before, construct a coupling between two copies of the full-system, where we see the edges “from the point of view of the walker” in the sense that whenever the edge $X_t + e$ updates at time t , where X_t is the position of the walker in the first copy, then in the second copy we will do the same update to the edge $\bar{X}_t + e$, where \bar{X}_t is the location of the walker in the second copy. Note that to establish the mixing time of the full system we need to couple the environments and the walkers. For simplicity, we concentrate our discussion here in the coupling of the walkers (which is the most delicate bit), and assume for now that somehow we managed to couple the two environments: that is, the two copies are coupled modulo a translation of the walkers. Note that, from this moment, if we were to employ the *identity coupling* (that is, the second copy mimics all the edge updates and jumps of the walker from the first copy) we would get that the environments will remain coupled (from the point of view of the walkers) but the distance between the walkers will not change, thereby not allowing the walkers to couple.

Our idea is to observe “a bit” the environment and, whenever the environment looks “favorable enough”, we attempt to do a coupling that could bring the walkers

closer together, which will be a standard coupling of simple random walks. We will refer to such moments as *simple random walk moments*, as an allusion to the simple random walk regenerations described above, but with the fundamental difference that they will not be regeneration times. On the other hand, when the environment is not favorable enough, then doing a simple random walk moment is a bit too risky, so instead we resort to identity coupling as a means to keeping the distance between the walkers unchanged and not spoiling the work done during the favorable regions of the environments.

But what does it mean for the environment to look favorable enough? In short terms, it will mean that the event (1.5) occurs *locally*. That is, at such times, all edges adjacent to the walkers will be closed and all edges in a small region around the walkers will be \star (for example, all edges inside a ball of radius 3 around the walkers, excluding the edges adjacent to the walkers). At such a time, with positive probability, the sequence of events described above for the simple random walk regeneration occurs, and therefore we could attempt to perform one of the standard couplings of simple random walks. However, there are two important caveats.

The first caveat is that if we succeed in doing a simple random walk moment with a coupling of simple random walks, then the distance between the walkers will change. This means that the translation that maps the location of one walker to the location of the second walker will change, and this map is what we use to match the edges of the first copy to the edges of the second copy, when we view the edges from the point of view of the walkers. As a consequence, the environments will immediately *decouple*. Of course, if we only had \star -edges (besides the ones adjacent to the walkers, as in the case $q = 1$), then the environments would not decouple since despite the change in the translation map, we would still match \star -edges in the first copy to \star -edges in the second copy, so we can easily maintain the environments coupled. But, since $q \neq 1$ implies a density of non- \star edges, the environments will necessarily decouple. Moreover, if we decide to just wait for the environments to recouple completely, this would take a time of order $\frac{\log n}{\mu}$, which is just too long: it will lead to an upper bound on the mixing time of order $\frac{n^2 \log n}{\mu}$. So we will not recouple the environments completely, but will work with partially coupled environments.

The second caveat is that a simple random walk moment occurs with *positive* probability, so it is also possible that it turns out that a simple random walk moment does not take place. Then, what could happen in this case? If the environments were completely coupled, then we are guaranteed that we can perform identity coupling and keep the distance between the walkers unchanged. But we have just seen that the environments will typically not be fully coupled. Yet, if we knew that the environments are coupled in a neighborhood around the walkers and that the walkers will not exit this neighborhood, then identity coupling is still doable. That will be our strategy, but to implement it we will require a more delicate definition of what a favorable enough environment means.

We will use a multi-scale analysis to control the environment. This will reveal *future information* regarding the environment; that is, we will observe some information about the environment from time 0 to some time t , and then decide how to couple

the walkers from time 0. Therefore, this construction will lead to a non-Markovian coupling.

A good picture to have in mind is that the environment is a process in space-time, where some regions are classified as *good* and others as *bad*. We observe these regions from time 0 to time t , and then start observing the walkers which are paths in space-time that start from time 0. Whenever we see that the walkers are passing through a good part of the environment, where good will also imply that the walkers will not move outside some neighborhood around their current locations, we will try to do a simple random walk moment. If successful, the distance between the walkers may change and the environments may decouple, but we will be able to recouple the environments within a neighborhood around the walkers using again that the walkers are passing through a good region in space-time. If, instead, the simple random walk moment is not successful, then the walkers may move more than just one step of a simple random walk, but again using the fact that the space-time region is good we will obtain that the walkers will not move too far away, in particular they will remain within a region where we know the environments were coupled. This will translate to a successful application of identity coupling.

On the other hand, if we see that the walkers are approaching a bad region of the environment, then we will want to do identity coupling but we will need to start preparing ourselves beforehand. The problem is that such a bad region could be of an arbitrarily large scale, and the larger its size is, the earlier we need to start preparing for it. So when we see that in space-time the path of the walker is getting dangerously near some bad region Q , we stop doing simple random walk moments even if in a smaller scale around the walkers the environment looks good. By switching off the simple random walk moments, we only apply identity coupling until the walkers reach Q or are again far enough from any bad region. We can show that such identity couplings will succeed and, since the translation map from one walker to the next will not change during this period, it will give enough time for the environments to couple in a region around the walkers that is as large as needed to contain Q . Then, with the environments properly coupled, if the walkers do enter Q , they can move as wildly as the environment there allows because we can perform identity coupling throughout Q . So the walkers survive the traversal of Q without changing their distance.

The above strategy is quite delicate, since in order to define which coupling (simple random walk moment or identity coupling) to use we need to observe which regions are good or bad from time 0 to time t . In particular, we need to compute the probability that a simple random walk moment is successful in a region Q given any possible assignment of good and bad to space-time regions so that Q is good. But such events will depend on one another. Thus, we will carefully split the definition of a good region into two parts, one that has a larger degree of independence, and another that is much more likely than the probability that a simple random walk moment is successful. In particular, we will need to estimate how such probabilities depend on p so that the coupling argument can be carried out. It becomes even more delicate to do this in more generality, which motivates our choice for carrying out the proof only for the random cluster dynamics.

Then one can imagine that the proof ends by showing that n^2 instances of a simple random walk moment are enough to guarantee that we can couple the walkers. This

is partially true. The fact is that, as mentioned above, we need to observe future information to carry out this coupling strategy. But in order to establish that the mixing time is at most t , we need to show that with a large enough probability the two copies of the full system are coupled at time t without revealing any information that goes beyond time t . So our strategy to finalize the proof is to choose an appropriate time $t' \in (0, t)$, reveal the information up to time t and do the coupling described above up to time t' , showing that within t' we have carried out an order of n^2 simple random walk moments, and that we coupled the walkers at time t' (the environments may, and typically will, be uncoupled except for a small region around the walkers). The whole analysis will be split into three phases, and the above will be carried out in the first two phases. We will be able to show that these first two phase succeed with positive probability.

Next, the goal is to try to do identity coupling from time t' to t in a similar manner as we were doing when approaching a bad region. In this second phase, identity coupling can only fail due to information that we have not observed because we are limited to observe the environment up to time t . We will show that, with positive probability, identity coupling will indeed succeed from t' to t , leading to a coupling of the full system at time t . This is the content of the third phase. If any of the three phases fail, then we just restart from scratch. We only need to repeat the phases a constant number of times to guarantee that the whole coupling succeeds with probability at least $3/4$.

1.5 Organization of the paper

In Sect. 2 we will introduce the multi-scale analysis that will allow to control the good regions of the environment. Then in Sect. 3 we will give a more thorough overview of the three phases of the proof of the upper bound, which will better explain the constructions from the tessellation of Sect. 2. Then in Sects. 4, 5 and 6 we will give the three phases of the coupling, with the second phase in Sect. 5 being the most delicate part where the non-Markovian coupling is developed. Then in Sect. 7 we put all phases together to complete the proof of the upper bound (Theorem 1.1). In Sect. 8 we establish the general lower bounds from Theorems 1.2 and 1.3, but which are essentially the same proofs as in [11]; this section is added for the sake of completeness. Finally, in Sect. 9 we apply these theorems to derive the lower bounds on the mixing time and relaxation time of random walks on the dynamical random cluster model (Corollary 1.4).

2 Multi-scale setup

We start defining a multi-scale tessellation of \mathbb{T}_n^d , which will consist of partitioning \mathbb{T}_n^d into boxes and defining the event that boxes are *good* or *bad*. Those events will be then used to define the favorable parts of the environment.

2.1 Tessellation

Let

$$\ell = p^{-\frac{1}{3d}}, \quad (2.1)$$

and m be a sufficiently large integer. For each scale $k \geq 1$ we tessellate \mathbb{T}_n^d into cubes of length ℓ_k where

$$\ell_1 = \ell \quad \text{and} \quad \ell_{k+1} = mk^2\ell_k. \quad (2.2)$$

The cubes will be indexed by integer vectors $i \in \mathbb{Z}^d$, and denoted $S_k^{\text{core}}(i) \subset \mathbb{T}_n^d$ with

$$S_k^{\text{core}}(i) = i\ell_k + [0, \ell_k)^d.$$

We will consider a tiling of \mathbb{T}_n^d with a hierarchy as each cube of scale k is contained inside a unique cube of scale $k+1$. For simplicity we will assume ℓ_k divides n for all k we will consider.² Moreover for any subset V of the vertices of \mathbb{T}_n^d , we denote by

$$E(V) = \{(v_1, v_2) \in E(\mathbb{T}_n^d) : v_1, v_2 \in V\}$$

the set of all edges incident only to vertices in V .

Now we define a multi-scale tessellation of time. At scale 1, we tessellate \mathbb{R} into intervals of length $t_1 = \frac{\sqrt{\ell}}{\mu}$ and then, for higher scales, we define

$$t_{k+1} = mk^2t_k, \quad k \geq 1.$$

We index the time intervals by $\tau \in \mathbb{Z}$ and denote them by $T_k^{\text{core}}(\tau)$, where

$$T_k^{\text{core}}(\tau) = [\tau t_k, (\tau + 1)t_k).$$

We will use the tessellations of space and time at scale k to define a space-time k -box later in (2.5). Before that, for any $i \in \mathbb{Z}^d$, $k \geq 1$, and $\tau \in \mathbb{Z}$, we define the *core* of the space-time k -box by

$$R_k^{\text{core}}(i, \tau) = S_k^{\text{core}}(i) \times T_k^{\text{core}}(\tau).$$

For any subset of $A \subset \mathbb{Z}^d$, we let ∂A denote its inner boundary. Then, in space-time, we define the spatial boundary of $R_k^{\text{core}}(i, \tau)$ by

$$\partial_s R_k^{\text{core}}(i, \tau) = \partial S_k^{\text{core}}(i) \times T_k^{\text{core}}(\tau). \quad (2.3)$$

² If that were not the case, one could consider for each k some cubes to have length between ℓ_k and $2\ell_k$ to fully tessellate the torus.

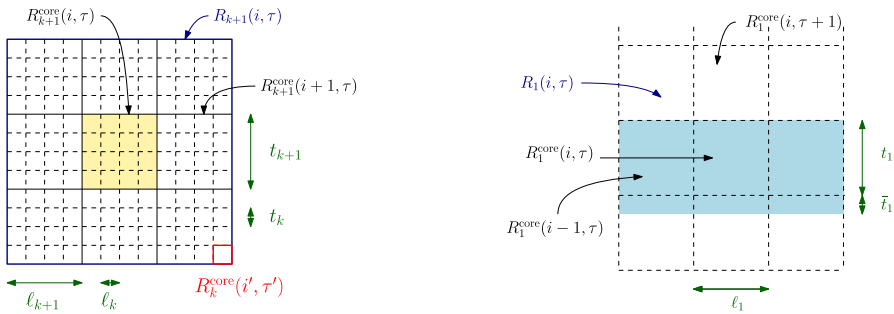


Fig. 1 On the left a space-time box $R_{k+1}(i, \tau)$ represented by a blue square, its core highlighted in yellow, its partition into cores of scale $k+1$ represented by solid black lines, and its partition into cores of scale k in dashed lines. The horizontal axis represents space and the vertical axis represents time. On the right a space-time box of scale 1 (highlighted in blue), and space-time cores of level 1 in dashed lines

For the time dimension, we define two time boundaries, the boundary ∂_t^+ corresponding to the largest unit of time in the box and the boundary ∂_t^- corresponding to the smallest unit of time in the box:

$$\begin{aligned} \partial_t^+ R_k^{\text{core}}(i, \tau) &= S_k^{\text{core}}(i) \times \sup \{T_k^{\text{core}}(\tau)\} = S_k^{\text{core}}(i) \times \{(\tau + 1)t_k\}, \text{ and} \\ \partial_t^- R_k^{\text{core}}(i, \tau) &= S_k^{\text{core}}(i) \times \inf \{T_k^{\text{core}}(\tau)\} = S_k^{\text{core}}(i) \times \{\tau t_k\}. \end{aligned} \quad (2.4)$$

For $k \geq 2$, each box $R_k^{\text{core}}(i, \tau)$ will be the central part of a larger box

$$R_k(i, \tau) = \bigcup_{(j, \tau') \in \{-1, 0, +1\}^{d+1}} R_k^{\text{core}}(i + j, \tau + \tau') = S_k(i) \times T_k(\tau), \quad (2.5)$$

where we let

$$S_k(i) = \bigcup_{j \in \{-1, 0, +1\}^d} S_k^{\text{core}}(i + j), \text{ and } T_k(\tau) = \bigcup_{\tau' \in \{-1, 0, +1\}} T_k^{\text{core}}(\tau + \tau'). \quad (2.6)$$

In words $R_k(i, \tau)$ is composed of a cube in space of side length $3\ell_k$ and a time interval of length $3t_k$, and it has $R_k^{\text{core}}(i, \tau)$ as its central part (see Fig. 1).

For scale $k = 1$, we will define boxes differently with respect to the time dimension. For this, let

$$\bar{t}_1 = \frac{\log^2 \ell}{\mu}, \quad (2.7)$$

and set for each $\tau \in \mathbb{Z}$

$$T_1(\tau) = [\tau t_1 - \bar{t}_1, (\tau + 1)t_1].$$

With S_1 defined as in (2.6) we can define $R_1(i, \tau) = S_1(i) \times T_1(\tau)$. We then define the space and time boundaries of $R_k(i, \tau)$ for each k, i, τ analogously to (2.3) and (2.4).

Finally we denote by $S_1^{\text{inn}}(i)$ the *inner part* of $S_1(i)$ which is obtained by removing all the vertices within distance $\frac{\gamma}{6} \log^2 \ell$ from the boundary of $S_1(i)$ (γ is a constant that will be clarified later in the definition); in symbols,

$$S_1^{\text{inn}}(i) = \left\{ v \in S_1(i) : \|v - w\|_1 > \frac{\gamma}{6} \log^2 \ell \text{ for all } w \in \partial S_1(i) \right\}. \quad (2.8)$$

2.2 Good boxes at scale 1

Definition 2.1 We say that an event A is restricted to a region $R \subset V(\mathbb{T}_n^d)$ and a time interval $[s_0, s_1]$ if A is measurable with respect to the σ -algebra generated by the updates of the edges from $E(R)$ from time s_0 to s_1 , together with the random variables U, U' associated to such updates.

Denote with $\mathcal{C}_x(t)$ the connected component of open edges containing vertex $x \in V$ at time t . Given a time interval $[s, s']$, we say that an edge is $[s, s']$ -open if there is at least one time during $[s, s']$ at which this edge is open. Then, we denote with $\mathcal{C}_x(s, s')$ the connected component of $[s, s']$ -open edges that contains x . If we denote $I = [s, s']$, then we employ the shorter notation $\mathcal{C}_x(I)$. Below we split the time interval of a box into two sets of sub-intervals, and then introduce the definition of *good boxes*.

Definition 2.2 Recall that $\bar{t}_1 = \frac{\log^2 \ell}{\mu}$. We define two other tessellations of disjoint time intervals. The first one has length $\frac{\log^2 \ell}{\mu}$:

$$\bar{T}_1(j) = [j\bar{t}_1, (j+1)\bar{t}_1), \quad \text{for } j \in \mathbb{Z}_+.$$

Moreover, we fix a constant $\gamma = \gamma(p, q, d) > 0$ such that $\frac{p_{\min}}{p_*} + \gamma < p_c$, where p_c is the critical probability for independent bond percolation on \mathbb{Z}^d , and introduce a tessellation of time of length $\frac{\gamma}{\mu}$:

$$\bar{\bar{T}}_1(j) = \left[j\frac{\gamma}{\mu}, (j+1)\frac{\gamma}{\mu} \right), \quad \text{for } j \in \mathbb{Z}_+.$$

We assume throughout this paper that \bar{t}_1 divides t_1 and that γ/μ divides \bar{t}_1 , so that $\bar{\bar{T}}_1$ is a finer tessellation than \bar{T}_1 , which in turn is a finer tessellation than T_1^{core} .

Remark 2.3 Note that the larger p is (that is, the closer p is to p_c) the smaller we need to take γ . However, as we will need to take p small enough in several places in the proof, we will set γ first (for example, it is enough to take $\gamma = \frac{1}{100}$). Then we make p small enough so that the condition on γ is satisfied.

The definition of good boxes will be done in steps. First we define some fundamental events that we will require from good boxes.

- (\mathcal{G}'_1) Given a box $R_1(i, \tau)$, let $\mathcal{G}'_1(i, \tau)$ be the event that, for any $e \in E(S_1(i))$, there are no non- \star updates on e during $T_1(\tau) \cap [0, \infty)$.
- (\mathcal{G}'_2) For each $j \in \mathbb{Z}_+$ and each spatial box $S_1(i)$, define $\mathcal{G}'_2(i, j)$ the event that, for any $e \in E(S_1(i))$, during $\bar{T}_1(j)$ edge e never gets a \star -update to become open.
- (\mathcal{G}'_3) For each $j \in \mathbb{Z}_+$ and each spatial box $S_1(i)$, let $\mathcal{G}'_3(i, j)$ be the event that, for each $e \in E(S_1^{\text{core}}(i))$, the number of \star -updates on edge e during $\bar{T}_1(j)$ is at least $\frac{1}{2}p_\star \log^2 \ell$ (for the values of ℓ and p we will consider this will always be at least 1).
- (\mathcal{G}'_4) For any $j \in \mathbb{Z}_+$ take the unique τ such that $\bar{\bar{T}}_1(j) \subset T_1^{\text{core}}(\tau)$. For any site x on the torus, if we regard all edges closed at time τt_1 and we only consider \star -updates disregarding all non- \star -updates, then let $\mathcal{G}'_4(x, j)$ be the event that $|\mathcal{C}_x(\bar{\bar{T}}_1(j))| < \log^2(\ell)$.

Now for a box $R_1(i, \tau)$, define

$$j(\tau) \text{ the value } j \text{ such that } \min \bar{T}_1(j) = \tau t_1 - \bar{t}_1; \quad (2.9)$$

in other words, it is the value such that $\bar{T}_1(j(\tau))$ starts at the initial time of $R_1(i, \tau)$. Note that both $\bar{T}_1(j(\tau))$ and $\bar{T}_1(j(\tau + 1))$ are contained in $T_1(\tau)$.

The event that a box $R_1(i, \tau)$ is good will be composed of four events, which we denote by $\mathcal{G}_1(i, \tau)$, $\mathcal{G}_2(i, \tau)$, $\mathcal{G}_3(i, \tau)$ and $\mathcal{G}_4(i, \tau)$. The first event regards only non- \star updates and is simply

$$\mathcal{G}_1(i, \tau) = \mathcal{G}'_1(i, \tau).$$

The second event regards the time intervals $\bar{T}_1(j(\tau))$ and $\bar{T}_1(j(\tau + 1))$, and is defined as

$$\mathcal{G}_2(i, \tau) = \mathcal{G}'_2(i, j(\tau)) \cap \mathcal{G}'_2(i, j(\tau + 1)) \bigcap_{i': S_1^{\text{core}}(i') \subset S_1(i)} \mathcal{G}'_3(i', j(\tau)) \cap \mathcal{G}'_3(i', j(\tau + 1)).$$

The next two events are confined to the time interval $T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1)$:

$$\mathcal{G}_3(i, \tau) = \bigcap_{\substack{i': S_1^{\text{core}}(i') \subset S_1(i) \\ j: \bar{T}_1(j) \subset (T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1))}} \mathcal{G}'_3(i', j)$$

and

$$\mathcal{G}_4(i, \tau) = \bigcap_{\substack{x \in S_1^{\text{inn}}(i) \\ j: \bar{\bar{T}}_1^j(\tau) \subset (T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1))}} \mathcal{G}'_4(x, j).$$

For convenience, we write

$$\mathcal{G}_{12}(i, \tau) = \mathcal{G}_1(i, \tau) \cap \mathcal{G}_2(i, \tau) \quad \text{and} \quad \mathcal{G}_{34}(i, \tau) = \mathcal{G}_3(i, \tau) \cap \mathcal{G}_4(i, \tau)$$

Lemma 2.4 *The family of events $\{\mathcal{G}_1(i, \tau)\}_{(i, \tau)}$, $\{\mathcal{G}_2(i, \tau)\}_{(i, \tau)}$ and $\{\mathcal{G}_{34}(i, \tau)\}_{(i, \tau)}$ are independent of one another.*

Proof The events $\mathcal{G}_1(\cdot, \cdot)$ depend only on non- \star updates, so it is independent of $\mathcal{G}_2(\cdot, \cdot)$, $\mathcal{G}_3(\cdot, \cdot)$ and $\mathcal{G}_4(\cdot, \cdot)$, which only regard \star -updates. Then, note that $\mathcal{G}_2(\cdot, \cdot)$ are events about \star -updates during $\bigcup_{\tau} \bar{T}_1(j(\tau))$, while $\mathcal{G}_3(\cdot, \cdot)$ and $\mathcal{G}_4(\cdot, \cdot)$ regard \star -updates during $\bigcup_{\tau} T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1)$. Since these two time intervals are disjoint, independence is obtained from standard properties of Poisson processes. \square

Lemma 2.5 *Let $R_1(i, \tau)$ be any box of scale 1. There exist constants $c, c' > 0$ so that for all small enough p we obtain*

$$\mathbb{P}(\mathcal{G}_{12}^c(i, \tau)) \leq c\ell^{d+\frac{1}{2}} p_{\max}.$$

and

$$\mathbb{P}(\mathcal{G}_{34}^c(i, \tau)) \leq \exp\left(-c' \log^2 \ell\right).$$

Proof We start with event $\mathcal{G}_{34}(i, \tau)$. For a given edge, an update that is not \star occurs at rate $(1 - p_{\star})\mu$, a \star -update occurs at rate $p_{\star}\mu$, and a \star -update that opens an edge occurs at rate $p_{\star} \frac{p_{\min}}{p_{\star}} \mu = p_{\min}\mu$. Moreover, there are at most $2d3^d \ell^d$ edges in $E(S_1(i))$.

For \mathcal{G}_3 , note that for a given edge $e \in S_1(i)$ and a given j , the number of \star -updates on e during $\bar{T}_1(j)$ is a Poisson random variable of mean $p_{\star} \log^2 \ell$. Therefore, using a standard Chernoff bound for Poisson random variables and the union bound over the edges in $S_1(i)$ and over the values of j , we obtain a constant $c_1 > 0$ such that

$$\mathbb{P}(\mathcal{G}_3^c(i, \tau)) \leq 2d3^d \ell^d \frac{(t_1 - \bar{t}_1)\mu}{\log^2 \ell} \exp\left(-c_1 p_{\star} \log^2 \ell\right).$$

Note that as p decreases to 0 we have that p_{\star} increases to 1 and ℓ increases to ∞ . So for all small enough p we obtain

$$\mathbb{P}(\mathcal{G}_3^c(i, \tau)) \leq \exp\left(-c_2 \log^2 \ell\right),$$

for some constant c_2 . Regarding $\mathcal{G}_4(i, \tau)$, for any $j \geq 0$ with $\bar{\bar{T}}_1(j) \subset T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1)$ and any edge e , note that the probability that e is open at the beginning of the interval $\bar{\bar{T}}_1(j)$, given that we only consider \star -updates and consider that all edges are closed at τt_1 , is at most $\frac{p_{\min}}{p_{\star}}$, since this is the probability that the last \star -update of e (if there was any) made e open. Thus, using that $1 - \exp(-p_{\star}\gamma)$ is the probability that e has a \star -update during $\bar{\bar{T}}_1^j$, we obtain that the probability that e is $\bar{\bar{T}}_1^j$ -open under the above assumptions is at most

$$\frac{p_{\min}}{p_{\star}} + 1 - \exp(-p_{\star}\gamma) \leq \frac{p_{\min}}{p_{\star}} + p_{\star}\gamma < p_c.$$

In other words, the set of \overline{T}_1^j -open edges (under the above assumptions) forms a subcritical percolation cluster. Therefore, using the exponential decay of cluster size for subcritical percolation, together with the union bound over all sites and values of j , we obtain a constant c'' such that

$$\mathbb{P}(\mathcal{G}_4^c(i, \tau)) \leq \ell^d \frac{t_1 \mu}{\gamma} \exp(-c_3 \log^2 \ell).$$

We note that c_3 increases as p decreases. So the bound on $\mathbb{P}(\mathcal{G}_{34}^c(i, \tau))$ follows by taking p small enough.

Now we turn to $\mathcal{G}_{12}(i, \tau)$. From the above considerations we obtain

$$\begin{aligned} \mathbb{P}(\mathcal{G}_{12}^c(i, \tau)) &\leq 1 - \exp\left(-2d3^d \ell^d \left((1 - p_\star)\mu(t_1 + \bar{t}_1) + p_{\min} \log^2 \ell\right)\right) \\ &\quad + 2d3^d \ell^d 2 \exp\left(-c_1 p_\star \log^2 \ell\right), \end{aligned}$$

where the term $1 - \exp(\cdot)$ comes from \mathcal{G}'_1 and \mathcal{G}'_2 , while the last term comes from \mathcal{G}'_3 as in the case of \mathcal{G}_3 above. Using that $t_1 + \bar{t}_1 \leq 2t_1$ and that $e^{-x} \geq 1 - x$ for all $x \in \mathbb{R}$, we obtain

$$\begin{aligned} \mathbb{P}(\mathcal{G}_{12}^c(i, \tau)) &\leq 2d3^d \ell^d \left((1 - p_\star)2\sqrt{\ell} + p_{\min} \log^2 \ell + 2 \exp\left(-c_1 p_\star \log^2 \ell\right)\right) \\ &\leq 4d3^d \ell^{d+\frac{1}{2}} (1 - p_\star + p_{\min}) \\ &= 4d3^d \ell^{d+\frac{1}{2}} (1 - p_\star + p_{\min}) = 4d3^d \ell^{d+\frac{1}{2}} p_{\max}, \end{aligned}$$

where we use that $\sqrt{\ell}$ is the term that dominates inside the parenthesis as p is made small enough (hence, ℓ is made large enough). So we can take p small enough so that $p_{\min} \log^2 \ell + 2 \exp(-c_1 p_\star \log^2 \ell) \leq 2p_{\min} \log^2 \ell \leq 2p_{\min} \sqrt{\ell}$. \square

We need one more step to define good boxes of scale 1. Using a standard result for percolation with bounded dependences [8], we will replace $\mathcal{G}_{34}(i, \tau)$ by a collection of i.i.d. Bernoulli random variables.

Lemma 2.6 *There exists a constant $C_6 = C_6(d) > 0$ such that letting $\{\widehat{\mathcal{G}}_{34}(i, \tau)\}_{(i, \tau)}$ be a collection of i.i.d. Bernoulli random variables of parameter $1 - \exp(-C_6 \log^2 \ell)$, Then for any p small enough we obtain that $\{\mathcal{G}_{34}(i, \tau)\}_{(i, \tau)}$ stochastically dominates $\{\widehat{\mathcal{G}}_{34}(i, \tau)\}_{(i, \tau)}$.*

Proof First note that if we fix i , then $\mathcal{G}_{34}(i, \tau)$ forms a collection of independent random variables as τ varies. So $\mathcal{G}_{34}(i, \tau)$ depends only on events $\mathcal{G}_{34}(i', \tau)$ such that $S_1(i) \cap S_1(i') \neq \emptyset$. Note that this is true even for the events \mathcal{G}'_4 since for $x \in S_1^{\text{inn}}(i)$, the event that the component of x is of length $\log^2 \ell$ is measurable with respect to the edges in $S_1(i)$. Given i , the number of i' such that $S_1(i) \cap S_1(i') \neq \emptyset$ is strictly smaller than $\Delta = 5^d$. Then, from Lemma 2.5 we obtain that by taking p small enough the marginal probability $w = \mathbb{P}(\mathcal{G}_{34}^c(i, \tau))$ can be made smaller than $\frac{(\Delta-1)\Delta^{-1}}{\Delta^\Delta}$, and

so we can apply [8, Theorem 1.3] to deduce that the family $\mathcal{G}_{34}(\cdot, \cdot)$ stochastically dominates a set of i.i.d. Bernoulli random variables with parameter

$$\begin{aligned}\rho &= \left(1 - \frac{w^{1/\Delta}}{(\Delta - 1)^{1-1/\Delta}}\right) \left(1 - (w(\Delta - 1))^{1/\Delta}\right) \\ &\geq 1 - \frac{w^{1/\Delta}}{(\Delta - 1)^{1-1/\Delta}} - (w(\Delta - 1))^{1/\Delta} \\ &= 1 - w^{1/\Delta} \left(\frac{\Delta}{(\Delta - 1)^{1-1/\Delta}}\right).\end{aligned}$$

Applying Lemma 2.5 for the value of w completes the proof since $w^{1/\Delta} \left(\frac{\Delta}{(\Delta - 1)^{1-1/\Delta}}\right) \leq \exp(-C_6 \log^2 \ell)$ for some constant $C_6 > 0$. \square

Now we are ready to define good boxes at scale 1.

Definition 2.7 (Good boxes of scale 1) Let $i \in \mathbb{Z}^d$ and $\tau \geq 0$. A box $R_1(i, \tau)$ is said to be *good* if the following event happens

$$\mathcal{G}(i, \tau) = \mathcal{G}_{12}(i, \tau) \cap \widehat{\mathcal{G}}_{34}(i, \tau).$$

For convenience, we assume that for $\tau < 0$ then $\mathcal{G}(i, \tau)$ holds for all $i \in \mathbb{Z}^d$. We also couple $\{\widehat{\mathcal{G}}_{34}(i, \tau)\}_{(i, \tau)}$ with $\{\mathcal{G}_{34}(i, \tau)\}_{(i, \tau)}$ so that,

$$\text{for each } (i, \tau), \text{ whenever } \widehat{\mathcal{G}}_{34}(i, \tau) \text{ holds, so does } \mathcal{G}_{34}(i, \tau). \quad (2.10)$$

The following lemma bounds the probability that a box is bad, and follows directly from the previous lemmas.

Lemma 2.8 Let $R_1(i, \tau)$ be any box of scale 1. There exists a constant $C_7 > 0$ so that for all small enough p we obtain

$$\mathbb{P}(\mathcal{G}^c(i, \tau)) \leq C_7 \ell^{d+\frac{1}{2}} p_{\max}.$$

Proof This follows from Lemmas 2.5 and 2.6 \square

Remark 2.9 Note that the event $\{R_1(i, \tau) \text{ is good}\}$ is restricted to the cube $S_1(i)$ and the interval $T_1(\tau)$. In fact, that is why in \mathcal{G}'_4 we assume that all edges are closed at time τt_1 , the initial time of the core $T_1^{\text{core}}(\tau)$; note that the fact that all edges are closed at time τt_1 is implied by \mathcal{G}_2 , but by explicitly adding the assumption in \mathcal{G}'_4 we make $\mathcal{G}_2(i, \tau)$ and $\mathcal{G}_4(i, \tau)$ independent of each other. Note also that the decision of whether a box is good is completely independent of the walker, it only depends on the updates of the dynamical random cluster process.

Recall that X_t denotes the position of the random walker at time t . In the lemma below, we will show that if the walker happens to be inside a good box, then it cannot move very quickly. This will allow us to have a better control on where the walker can be.

Lemma 2.10 *Let $t \geq 0$ be any given time and suppose $(X_t, t) \in R_1^{\text{core}}(i, \tau)$, where $R_1(i, \tau)$ is a good box. Then,*

$$\sup_{\substack{s \geq t \\ s \in T_1(\tau)}} \|X_t - X_s\|_1 \leq |\mathcal{C}_{X_t}(t)| + \lceil \frac{(\tau + 1)t_1 - t}{\gamma/\mu} \rceil \log^2 \ell,$$

In particular, if $\tau > 0$ then

$$\sup_{\substack{s \geq t \\ s \in T_1(\tau)}} \|X_t - X_s\|_1 \leq \log^2 \ell + \lceil \frac{(\tau + 1)t_1 - t}{\gamma/\mu} \rceil \log^2 \ell \leq \log^2 \ell + \frac{\sqrt{\ell}}{\gamma} \log^2 \ell \leq \frac{\ell}{10},$$

where the last inequality holds for all p small enough (thus, ℓ large enough) and where $\|x - y\|_1$ denotes the L^1 distance in the torus between the positions $x, y \in \mathbb{T}_n^d$; in particular, $\|x - y\|_1$ does not depend on whether edges are open or closed.

Proof This is a direct consequence of the event \mathcal{G}'_4 from the definition of good boxes, and the fact that good boxes do not have non- \star -updates. There is just one caveat. The event \mathcal{G}'_4 is not enforced in time intervals $\bar{T}_1(j)$ where $j = j(\tau)$ for some τ ; recall the definition of $j(\tau)$ from (2.9). So for example, $\mathcal{G}_4(i, \tau)$ does not include the event $\mathcal{G}'_4(j)$ for j such that $\bar{T}_1(j) = T_1(\tau) \cap T_1(\tau + 1)$, which is the only time interval of the type $\bar{T}_1(\cdot)$ inside $T_1(\tau)$. However, for such a j , we know that the connected components inside $\bar{T}_1(j - 1)$ are of size at most $\log^2 \ell$, and $\mathcal{G}_2(i, \tau)$ implies that during $\bar{T}_1(j)$ no edge of $S_1(i)$ gets an update to open. Therefore, the size of the connected components can only decrease during $\bar{T}_1(j)$ and the result follows. \square

2.3 Good boxes at larger scales

In this section we define the concept of good and bad boxes of scale larger than 1, but first we define a slightly relaxed version of *intersection* of boxes.

Definition 2.11 Since boxes are defined by semi-open intervals, we will consider boxes that are barely non-intersecting as intersecting. That is, we consider two boxes $R_k(i, \tau)$ and $R_k(i', \tau')$ as *non-intersecting* if and only if

$$\inf_{\substack{(j,s) \in R_k(i,\tau) \\ (j',s') \in R_k(i',\tau')}} \|(j,s) - (j',s')\|_1 \geq 2.$$

Definition 2.12 A k -box $R_k(i, \tau)$ with $k \geq 2$ is said to be *bad* if it contains at least two non-intersecting bad boxes of scale $k - 1$. Otherwise, $R_k(i, \tau)$ is called *good*.

Remark 2.13 The event $\{R_k(i, \tau) \text{ is bad}\}$ is strictly restricted to the cube $S_k(i)$ and the time interval $T_k(\tau)$. Moreover, by translation invariance, for any pair (i, τ) , (i', τ') and any scale k we have $\mathbb{P}(R_k(i, \tau) \text{ is bad}) = \mathbb{P}(R_k(i', \tau') \text{ is bad})$. Therefore if $R_k(i, \tau)$ and $R_k(i', \tau')$ are two non-intersecting boxes then

$$\mathbb{P}(R_k(i, \tau) \text{ and } R_k(i', \tau') \text{ are bad}) = \mathbb{P}(R_k(i, \tau) \text{ is bad})^2.$$

Definition 2.14 Define ρ_k as the probability that a k -box $R_k(i, \tau)$ is bad, that is,

$$\rho_k = \mathbb{P}(R_k(i, \tau) \text{ is bad}).$$

As noted in Remark 2.13, this probability does not depend on (i, τ) .

Recall that m is the variable that appears in the definition of ℓ_k from (2.2).

Lemma 2.15 For any $m > 0$, by setting p small enough we obtain

$$\rho_k \leq \rho_1^{2^{k-2}}$$

Proof We prove the lemma in a slightly stronger version: we prove that we can set values c_k , satisfying $c_k \geq \frac{1}{2}$ for all k , so that

$$\rho_k \leq \rho_1^{c_k 2^{k-1}}.$$

We prove this by induction. For $k = 1$ the statement is trivially satisfied by setting $c_1 = 1$. Assume the statement is true up to k . Now, by the definition of bad box we have

$$\begin{aligned} \rho_{k+1} &\leq \left(\left(\frac{3\ell_{k+1}}{\ell_k} \right)^d \frac{3t_{k+1}}{t_k} \right)^2 \rho_k^2 \\ &= 3^{2d+2} (mk^2)^{2d+2} \rho_k^2 \\ &\leq 3^{2d+2} (mk^2)^{2d+2} \left(\rho_1^{c_k 2^{k-1}} \right)^2 \\ &= 3^{2d+2} (mk^2)^{2d+2} \rho_1^{(c_k - c_{k+1}) 2^k} \rho_1^{c_{k+1} 2^k}. \end{aligned}$$

Setting $c_{k+1} = c_k - \frac{1}{10k^2}$ gives that

$$3^{2d+2} (mk^2)^{2d+2} \rho_1^{(c_k - c_{k+1}) 2^k} = 3^{2d+2} (mk^2)^{2d+2} \rho_1^{\frac{2^k}{10k^2}} \leq 1,$$

for all $k \geq 1$, provided ρ_1 is small enough with respect to m . Given m , ρ_1 can be made small enough by setting p small enough, as in Lemma 2.8. Notice that $c_k > c_1 - \sum_{i=1}^{\infty} \frac{1}{10i^2} \geq \frac{1}{2}$, which proves the lemma. \square

2.4 Enlargement of boxes

As we discussed in the proof overview, whenever the walkers are in a favorable region of the environment, we will try to use a simple random walk coupling to bring the walkers together. However, when the walkers are in an unfavorable region of the environment, which essentially means that the walkers are approaching a bad box (at some scale), then we will have to refrain from doing this simple random walk

coupling, and will just do a naïve identity coupling in order to let the environments couple around the walkers before they can reach the bad box. Here we will define two types of enlargements of boxes so that when the walker enters the enlargement of a bad box we will need to stop doing the simple random walk coupling.

Definition 2.16 (1-enlargement) The *1-enlargement* of a box $R_k(i, \tau)$ of scale k , is the set of boxes

$$R_k^{\text{enl1}}(i, \tau) = \bigcup_{(j, \beta) \in \{-3, -2, \dots, 3\}^{d+1}} R_k(i + j, \tau + \beta).$$

We also denote

$$S_k^{\text{enl1}}(i) = \bigcup_{j \in \{-3, -2, \dots, 3\}^d} S_k(i + j), \quad \text{and} \quad T_k^{\text{enl1}}(\tau) = \bigcup_{\beta \in \{-3, -2, \dots, 3\}} T_k(\tau + \beta).$$

Note that $R_k^{\text{enl1}}(i, \tau)$ is a parallelogram of spatial length $9\ell_k$ and time length $9t_k$ for $d \geq 2$ and $7t_1 + \bar{t}_1$ for $d = 1$.

Remark 2.17 The 1-enlargement is a $(d + 1)$ -dimensional parallelogram centered in $R_k(i, \tau)$ defined to obtain the following property. Let $R_k(i, \tau)$ be a bad box, whose whole 1-enlargement $R_k^{\text{enl1}}(i, \tau)$ is contained inside a good box of scale $k + 1$. Then, we know that the only boxes of scale k inside the $(k + 1)$ -box that can be bad are those intersecting $R_k(i, \tau)$. Let I be the set of tuples (i', τ') such that $R_k(i, \tau) \cap R_k(i', \tau') \neq \emptyset$. Note that $R_k(i', \tau') \subset R_k^{\text{enl1}}(i, \tau)$ for all $(i', \tau') \in I$. Moreover, the property that we get is that $\bigcup_{(i', \tau') \in I} R_k^{\text{core}}(i', \tau')$ does not exhaust $R_k^{\text{enl1}}(i, \tau)$ in the sense that $\bigcup_{(i', \tau') \in I} R_k^{\text{core}}(i', \tau')$ is separated from the outside of $R_k^{\text{enl1}}(i, \tau)$ by at least one layer of cores. We define this layer of cores as

$$R_k^{\partial \text{enl1}}(i, \tau) = R_k^{\text{enl1}}(i, \tau) \setminus \left(\bigcup_{(i', \tau') \in I} R_k^{\text{core}}(i', \tau') \right).$$

Definition 2.18 (2-enlargement) The *2-enlargement* of a box $R_k(i, \tau)$ of scale k is the set of boxes

$$R_k^{\text{enl2}}(i, \tau) = \bigcup_{\substack{j \in \{-20, -19, \dots, 20\}^d, \\ \beta \in \{-18, -17, \dots, 3\}}} R_k(i + j, \tau + \beta),$$

and we also denote

$$S_k^{\text{enl2}}(i) = \bigcup_{j \in \{-20, -19, \dots, 20\}^d} S_k(i + j), \quad T_k^{\text{enl2}}(\tau) = \bigcup_{\beta \in \{-18, -17, \dots, 3\}} T_k(\tau + \beta).$$

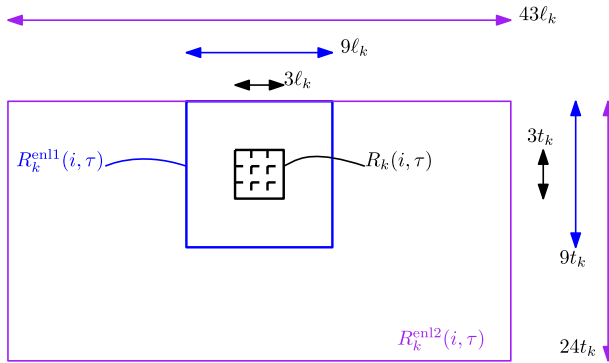


Fig. 2 In black a box $R_k(i, \tau)$, with its 1-enlargement in blue and its 2-enlargement in purple. The figure is not to scale and illustrates the case $k \geq 2$; recall that the time intervals are defined differently at scale 1

Note that the 2-enlargement is a larger $(d+1)$ -dimensional parallelogram centered in $R_k(i, \tau)$ so that $\sup \{T_k^{\text{enl1}}(\tau)\} = \sup \{T_k^{\text{enl2}}(\tau)\}$; See Fig. 2.

We will require a different type of boundary for the second enlargement.

Definition 2.19 (2-enlargement boundary) We define $\partial_s^{\text{enl2}} R_k^{\text{enl2}}(i, \tau)$, the 2-enlargement boundary of $R_k^{\text{enl2}}(i, \tau)$, as the set of space-time points $(x, s) \in R_k^{\text{enl2}}(i, \tau)$ such that $(x, s) \in R_k^{\text{core}}(i', \tau')$ for some (i', τ') with $R_k(i', \tau') \subset R_k^{\text{enl2}}(i, \tau)$ and $R_k(i', \tau') \cap \partial_s R_k^{\text{enl2}}(i, \tau) \neq \emptyset$.

2.5 Feasible paths

In this subsection we introduce the concept of feasible paths. For any graph $G = (V, E)$, we denote the neighbors of a vertex $v \in V$ by $\mathcal{N}_G(v) = \{w \in V : (v, w) \in E\}$. A path $\mathcal{P} : \mathbb{R}^+ \rightarrow V$ on a graph $G = (V, E)$ is a càdlàg function of time such that for any $s \in \mathbb{R}^+$, if we take s' to be the smallest value that is larger than s and such that $\mathcal{P}(s') \neq \mathcal{P}(s)$ then $\mathcal{P}(s') \in \mathcal{N}_G(\mathcal{P}(s))$. Note that a path, as defined above, does not consider whether edges are open or closed and is thus allowed to jump across closed edges. The same is true in the definition below. Recall the definition of the time intervals \overline{T}_1^j from Definition 2.2 and the inner part $S_1^{\text{inn}}(i)$ of box i from (2.8).

Definition 2.20 (Feasible path) A path \mathcal{P} is said to be *feasible* if for any times s, s' with $s, s' \in \overline{T}_1^j(\tau) \subset T_1^{\text{core}}(\tau)$ for some j and $\tau > 0$, and such that $\mathcal{P}(s) \in S_1^{\text{inn}}(i)$ for some $i \in \mathbb{Z}^d$ for which $R_1(i, \tau)$ is a good box, then

$$\|\mathcal{P}(s') - \mathcal{P}(s)\|_1 < \log^2 \ell.$$

Intuitively, a feasible path can move at most $\log^2 \ell$ during any interval \overline{T}_1^j in which it is inside good boxes. Even though the definition of feasible paths does not consider whether edges are open or closed, this is aligned with the fact that in good boxes the clusters are of size at most $\log^2 \ell$.

We will refer to a path that *leaves the box* $R_k(i, \tau)$ from the time boundary as a path \mathcal{P} such that $(\mathcal{P}(s), s) \in R_k^{\text{core}}(i, \tau)$ for some $s \geq 0$ and if $s' > s$ is the smallest value such that $(\mathcal{P}(s'), s') \notin R_k(i, \tau)$ then $\mathcal{P}(s') \in S_k(i)$. In other words, it is a path that exits $R_k(i, \tau)$ through $\partial_1^+ R_k(i, \tau)$. In the following two lemmas we will prove that a feasible path always leaves good boxes from the time boundary. Recall the definition of $\mathcal{C}_x(I)$, the connected component of x during a time interval I , from the paragraph preceding Definition 2.2.

We define the *spatial core* of a box as follows:

$$R_k^{\text{s-core}}(i, \tau) = S_k^{\text{core}}(i) \times T_k(\tau). \quad (2.11)$$

Lemma 2.21 *For all p small enough (hence, ℓ large enough) the following holds. Let (i, τ) be such that $R_1(i, \tau)$ is a good box. Let \mathcal{P} be a feasible path such that $(\mathcal{P}(s), s) \in R_1^{\text{s-core}}(i, \tau)$ for some s . Assume that either $s \geq \bar{t}_1$ or $|\mathcal{C}_{\mathcal{P}(s)}(s)| \leq \sqrt{\ell} \log^2 \ell$. Then, \mathcal{P} leaves $R_1(i, \tau)$ from the time boundary and*

$$\sup_{s' \in [s, \max T_1(\tau)]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 \leq \sqrt{\ell} \log^3 \ell.$$

Proof For any $(v, s) \in R_1^{\text{s-core}}(i, \tau)$ and any $(v', s') \in \partial_s R_1(i, \tau)$, $\|v - v'\|_1 \geq \ell$ as well as $|s' - s| \leq 3t_1$. Recall that $T_1(\tau)$ is split into intervals $\bar{T}_1^j(\tau)$ of length $\frac{\gamma}{\mu}$. Assume first that $s \geq \bar{t}_1$. Then the subinterval $\bar{T}_1^j(\tau)$ containing s consists only of positive times; hence, $\mathcal{C}_{\mathcal{P}(s)}(s)$ has cardinality at most $\log^2 \ell$. From the definition of feasible paths we obtain

$$\sup_{s' \in [s, \max T_1(\tau)]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 \leq \frac{3t_1\mu}{\gamma} \log^2 \ell.$$

For all small enough p we have $\frac{3t_1\mu}{\gamma} \log^2 \ell = \frac{3\sqrt{\ell}}{\gamma} \log^2 \ell \leq \ell \log^3 \ell$, recalling that γ is set before we take p small enough as in Remark 2.3.

If $s \in [0, \bar{t}_1)$ then $s \in \bar{T}_1^1(0) \subset T_1(0)$ and we use that during $T_1^1(0)$ no edge in $S_1(i)$ opens. Therefore,

$$\sup_{s' \in [s, \max T_1(\tau)]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 \leq |\mathcal{C}_{\mathcal{P}(s)}(s)| + \frac{3t_1\mu}{\gamma} \log^2 \ell \leq \sqrt{\ell} \log^3 \ell.$$

□

The next lemma is the analogue of Lemma 2.21 for higher scales.

Lemma 2.22 *For all m large enough and all p small enough with respect to m , the following holds. Let $k \geq 2$ and let (i, τ) be such that $R_k(i, \tau)$ is a good box. Let \mathcal{P} be a feasible path such that $(\mathcal{P}(s), s) \in R_k^{\text{s-core}}(i, \tau)$ for some s . Assume that either $s \geq \bar{t}_1$ or $|\mathcal{C}_{\mathcal{P}(s)}(s)| \leq \sqrt{\ell} \log^2 \ell$. Then \mathcal{P} leaves $R_k(i, \tau)$ from the time boundary. Moreover, while the path is inside the box, from time s up to time $(\tau + 2)t_k$, the path must be within distance $\ell_k/9$ from $\mathcal{P}(s)$.*

Proof For any $(v, s) \in R_k^{\text{s-core}}(i, \tau)$ and any $(v', s') \in \partial_s R_k(i, \tau)$, $\|v - v'\|_1 \geq \ell_k$ as well as $|s' - s| \leq 3t_k$. We do a proof by induction on k . The case $k = 1$ is a direct consequence of Lemma 2.21. We will actually assume a slightly stronger induction hypothesis. Take $c_1 = \frac{1}{100}$, $c_2 = \frac{1}{10}$ and, for $j \geq 2$, set $c_{j+1} = c_j + \frac{11}{mj^2}$. We take m large enough so that $c_j \leq \frac{1}{9}$ for all $j \geq 1$. Now, for a scale k , assume that if $\mathcal{P}(s) \in R_k^{\text{s-core}}(i, \tau)$ for some (i, τ) and some s as in the statement of the lemma, then $\sup_{s' \in [s, (\tau+2)t_k]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 \leq c_k \ell_k$. We want to prove the above for scale $k+1$.

We split into two cases, starting with $k \geq 2$ (thus, $k+1 \geq 3$). Let now \mathcal{P} be a feasible path such that $(\mathcal{P}(s), s) \in R_{k+1}^{\text{s-core}}(i, \tau)$, and $R_{k+1}(i, \tau)$ is a good box. Thus there are no pairs of non intersecting bad boxes of scale k inside $R_{k+1}(i, \tau)$. By Remark 2.17, if $R_{k+1}(i, \tau)$ contains at least one bad box $R_k(i', \tau')$, then all bad boxes contained in $R_{k+1}(i, \tau)$ are contained in $R_k^{\text{enl}}(i', \tau')$. Inside $R_k^{\text{enl}}(i', \tau')$ a feasible path has no restriction on how quickly it can move and it could potentially traverse $S_k^{\text{enl}}(i')$ instantaneously. The remaining boxes of scale k that are in $R_{k+1}(i, \tau)$ are good and by the inductive hypothesis we can use that in these ones the maximum displacement of the path is bounded above by $c_k \ell_k$, so for $k \geq 2$ it follows that

$$\begin{aligned} \sup_{s' \in [s, (\tau+2)t_{k+1}]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 &\leq 9\ell_k + \left(12 + \frac{2t_{k+1}}{2t_k}\right) c_k \ell_k \\ &\leq 9\ell_k + \left(12 + k^2 m\right) c_k \ell_k \\ &= \left(\frac{9}{mk^2} + \frac{12c_k}{mk^2} + c_k\right) \ell_{k+1} \\ &\leq c_{k+1} \ell_{k+1}. \end{aligned}$$

The term $\left(12 + \frac{2t_{k+1}}{2t_k}\right)$ accounts for the following boxes. Each time the path \mathcal{P} finds itself at the starting time of a k -core, it spends at least time $2t_k$ inside the corresponding k -box, and after that amount of time it finds itself at the starting time of another k -core. This gives at most $\frac{2t_{k+1}}{2t_k}$ k -cores for which we can apply the induction hypothesis. There are situations, however, that we cannot guarantee that the path \mathcal{P} is at the *starting* time of a k -core. One such situation is the very first box. We can still apply the induction hypothesis in such cases, since the hypothesis requires only that the path is inside the spatial core, regardless of it being the starting time of a core or not, but can give rise to at most 12 additional boxes to the counting: the first box, the boxes right before and after $R_k^{\text{enl}}(i', \tau')$, and the 9 time intervals contained in $R_k^{\text{enl}}(i', \tau')$.

Now we turn to the last case, which is $k = 1$ (that is, $k+1 = 2$). We proceed in the same way as before, but taking care of the fact that boxes at scale 1 have a different length in the time dimension. We have

$$\begin{aligned} \sup_{s' \in [s, (\tau+2)t_2]} \|\mathcal{P}(s) - \mathcal{P}(s')\|_1 &\leq 9\ell_1 + \left(12 + \frac{2t_2}{t_1}\right) c_1 \ell_1 \\ &\leq 9\ell_1 + (12 + 2m) c_1 \ell_1 \end{aligned}$$

$$= \left(\frac{9}{m} + \frac{12c_1}{m} + 2c_1 \right) \ell_2 \leq c_2 \ell_2.$$

The proof is concluded by taking m large enough to guarantee that $c_{k+1} = \frac{1}{100} + \frac{1}{10} + \frac{20}{m} \sum_{i=2}^k \frac{1}{i^2} < \frac{1}{9}$ for all k . \square

Next we will prove that if a feasible path enters the 2-enlargement of a box $R_k(i, \tau)$ from its spatial boundary, and all the k -boxes inside $R_k^{\text{enl}2}(i, \tau) \setminus R_k^{\text{enl}1}(i, \tau)$ are good, then the path remains far from the box $R_k(i, \tau)$. For this lemma, recall the definition of $\partial_s^{\text{enl}2} R_k^{\text{enl}2}(i, \tau)$ the 2-enlargement boundary of $R_k(i, \tau)$ from Definition 2.19.

Lemma 2.23 *Let \mathcal{P} be a feasible path such that $(\mathcal{P}(s), s) \in \partial_s^{\text{enl}2} R_k^{\text{enl}2}(i, \tau)$ for some (i, τ) and $s \geq 0$. Assume that either $s \geq \bar{t}_1$ or $|\mathcal{C}_{\mathcal{P}(s)}(s)| \leq \sqrt{\ell} \log^2 \ell$. Assume also that $R_k(i', \tau')$ is good for all $R_k(i', \tau') \subset R_k^{\text{enl}2}(i, \tau) \setminus R_k^{\text{enl}1}(i, \tau)$. Then,*

$$\inf_{\substack{v \in S_k^{\text{enl}1}(i) \\ s' \geq s, s' \in T_k^{\text{enl}2}(\tau)}} \|\mathcal{P}(s') - v\|_1 \geq 12\ell_k.$$

Proof By hypothesis every box $R_k(\cdot, \cdot) \subset R_k^{\text{enl}2}(i, \tau) \setminus R_k^{\text{enl}1}(i, \tau)$ is good. For any $s' \in T_k^{\text{enl}2}(\tau)$, $|s - s'| \leq 24t_k$. Assume without loss of generality that during $[s, s']$ the path never visits a box $R_k(\cdot, \cdot)$ which is not contained in $R_k^{\text{enl}2}(i, \tau)$; otherwise we can carry out the proof separately to each portion of the path that only traverses boxes $R_k(\cdot, \cdot)$ contained in $R_k^{\text{enl}2}(i, \tau)$. Let (i', τ') be such that $(\mathcal{P}(s), s) \in R_k^{\text{core}}(i', \tau')$. Since $R_k(i', \tau')$ is a good box, letting $s'' = \sup \{T_k(\tau')\}$ we have that $\|\mathcal{P}(s) - \mathcal{P}(s'')\|_1 \leq \frac{\ell_k}{9}$ by Lemma 2.22. If $s'' < s'$, we can iterate the above argument obtaining that $\|\mathcal{P}(s) - \mathcal{P}(s')\|_1 \leq 24\frac{\ell_k}{9}$, where 24 amounts for the largest number of iterations. Since boxes have length $2t_k$ in the time dimension, it would be enough to replace 24 by 12 for $k \geq 2$, but we just use the larger bound 24 to accommodate also the $k = 1$ case, for which the length of a box in the time dimension is smaller.

Now since $\frac{24\ell_k}{9}$ is smaller than $13\ell_k$, which is the distance between $\mathcal{P}(s)$ and the spatial boundary of $S_k^{\text{enl}1}(i)$ enlarged by all boxes $R_k(\cdot, \cdot)$ that intersects it, the path can only traverse good k -boxes while inside $R_k^{\text{enl}2}(i, \tau)$. In addition, for any $v \in S_k^{\text{enl}1}(i)$ one has

$$\|v - \mathcal{P}(s')\|_1 \geq \|v - \mathcal{P}(s)\|_1 - \frac{24\ell_k}{9} \geq 15\ell_k - \frac{24\ell_k}{9} \geq 12\ell_k.$$

\square

2.6 Great boxes

We will need a stroger notion for boxes of scale 1, which we will call *great* boxes.

Definition 2.24 A box $R_1(i, \tau)$ is said to be k -great if for all $k' \leq k$, for all $R_{k'}(i', \tau')$ such that $R_{k'}^{\text{enl}2}(i', \tau')$ intersects $R_1(i, \tau)$ then $R_{k'}(i', \tau')$ is good. Moreover, we define

$$G_k = \{(i, \tau) : R_1(i, \tau) \text{ is } k\text{-great}\}$$

to be the set of k -great boxes.

Later we will see that the walker has to traverse a feasible path. The next lemma will be used to say that if the walker traverses a k -box that is good with a large neighborhood of good k -boxes, then it is necessarily the case that the walker has to traverse enough k -great boxes. Such great boxes will be the places where we will attempt a simple random walk coupling later.

Lemma 2.25 *Let \mathcal{P} be a feasible path such that $(\mathcal{P}(\tau t_k), \tau t_k) \in \partial_t^- R_k^{\text{core}}(i, \tau)$ for some (i, τ) . Assume that either $\tau > 0$ or $|\mathcal{C}_{\mathcal{P}(\tau t_k)}(\tau t_k)| \leq \sqrt{\ell} \log^2 \ell$. Then there exists $C_8 > 0$ such that letting $r = C_8 \frac{t_k}{t_1}$ we can find times $s_1 < \dots < s_r$ and distinct space-time indices $(i_1, \tau_1), \dots, (i_r, \tau_r)$ such that the following all hold:*

- *For all j , $(\mathcal{P}(s_j), s_j) \in \partial_t^- R_1^{\text{core}}(i_j, \tau_j)$ and \mathcal{P} exits $R_1(i_j, \tau_j)$ from the time boundary.*
- *$R_1(i_j, \tau_j) \subset R_k(i, \tau)$ are k -great for all j .*
- *$\tau_1 t_1 \geq \tau t_k$ and $\tau_j \geq \tau_{j-1} + 2$ for all $j \in \{2, 3, \dots, r\}$.*

Proof First, note that if a box $R_1(\tilde{i}, \tilde{\tau})$ is $(k-1)$ -great and is contained in $R_k(i, \tau)$ then it is also k -great. We will prove the statement of the lemma replacing C_8 with c_k , some function of k . Then the lemma follows by showing that there is a universal value C_8 such that $0 < C_8 \leq c_k$ for all k . We will do a proof by induction on k . Case $k = 1$ is trivially verified by choosing $c_1 = 1$ because in this case $r = c_1 = 1$ and we take $(i_1, \tau_1) = (i, \tau)$.

Now, for $k \geq 2$, assume the lemma is true up to scale $k-1$ and consider a feasible path that at time τt_k is inside $\partial_t^- R_k^{\text{core}}(i, \tau)$ such that every box of scale k whose 2-enlargement intersects $R_k(i, \tau)$ is good. By Remark 2.17, the bad boxes of scale $k-1$ inside $R_k(i, \tau)$ (if there are any) are all contained in $R_{k-1}^{\text{enl}1}(i', \tau')$ for some i', τ' . We then regard all boxes of scale 1 which are in at least one of the 2-enlargement of the boxes contained in $R_{k-1}^{\text{enl}1}(i', \tau')$ as potentially not $(k-1)$ -great. By Lemma 2.22 we know that \mathcal{P} crosses $\partial_t^+ R_k(i, \tau)$ before $\partial_s R_k(i, \tau)$. In words the path stays for time at least $2t_k$ in the box $S_k(i)$.

Since the path starts from $\partial_t^- R_k^{\text{core}}(i, \tau)$, it starts on $\partial_t^- R_{k-1}^{\text{core}}(i'', \tau'')$ for some i'' and τ'' . In this $(k-1)$ -box we can apply the inductive hypothesis, so after time $2t_{k-1}$ the path has gone through at least $c_{k-1} \frac{t_{k-1}}{t_1}$ distinct $(k-1)$ -great boxes. Since this path remains inside $R_k(i, \tau)$, we immediately obtain that such boxes are all k -great boxes. When the path reaches $\partial_t^+ R_{k-1}(i'', \tau'')$ we have that the path is now on $\partial_t^- R_{k-1}^{\text{core}}(i'' + j, \tau'' + 2) \subset R_k(i, \tau)$ for some $j \in \mathbb{Z}^d$ and from here we can reapply the inductive hypothesis. So it remains to count how many times we can iterate this procedure before $2t_k$ amount of time has passed.

To do this, we first count how much time the path can spend inside the 2-enlargement of a bad $(k-1)$ -box. It suffices to count how much time is spanned by the boxes whose

2-enlargements intersect $R_{k-1}^{\text{enl}}(i', \tau')$, which is $|T_{k-1}^{\text{enl1}}(\cdot)| + 2|T_{k-1}^{\text{enl2}}(\cdot)| = 57t_{k-1}$. Hence, the number of times the above procedure can be iterated is at least

$$\frac{2t_k - 57t_{k-1}}{2t_{k-1}} = \frac{t_k}{t_{k-1}} \left(1 - \frac{57}{2m(k-1)^2} \right).$$

From the inductive hypothesis, the path will traverse at least

$$\frac{t_k}{t_{k-1}} \left(1 - \frac{57}{2m(k-1)^2} \right) c_{k-1} \frac{t_{k-1}}{t_1} = c_k \frac{t_k}{t_1}$$

$(k-1)$ -great boxes, by setting $c_k = \left(1 - \frac{57}{2m(k-1)^2} \right) c_{k-1}$. These boxes are k -great by the properties of $R_k(i, \tau)$. The lemma is then concluded by setting

$$C_8 = \lim_{k \rightarrow \infty} c_k = \prod_{i=1}^{\infty} \left(1 - \frac{57}{2mi^2} \right) > 0.$$

□

3 Overview of the proof

In this section we give a high-level description of the proof. Consider two processes $\{M_t^*\}_{t \geq 0} = \{X_t, \eta_t^*\}_{t \geq 0}$ and $\{\bar{M}_t^*\}_{t \geq 0} = \{\bar{X}_t, \bar{\eta}_t^*\}_{t \geq 0}$ with starting states $M_0^*, \bar{M}_0^* \in \Omega^*$. We will construct a coupling of the two processes so that for some time $T = \Delta_1 + \Delta_2 + \Delta_3$ of order $\frac{n^2}{\mu}$ the two configurations agree with positive probability. Since $\{M_t^*\}_{t \geq 0}$ can be recovered from $\{M_t^*\}_{t \geq 0}$, by sampling independently the edges with status \star , we will obtain our result.

The coupling will consist of three different phases which we will describe in a high level way below. The coupling of each phase will have a small, albeit positive, probability of failing. If the coupling of a phase fails, we declare the whole three-phase procedure to have failed, let the two processes evolve arbitrarily until time T and restart everything again from phase 1. The detailed analysis of each phase will be given in sections 4, 5 and 6. Then in Sect. 7, we will put all phases together and complete the proof of Theorem 1.1.

3.1 First phase: the local coupling

During the first phase we let the two processes evolve *independently*, and wait for the first time the graphs of the two processes agree on a ball of radius 2ℓ around the walkers, that is, we wait for a time t such that

$$\eta_t^*(X_t + e) = \bar{\eta}_t^*(\bar{X}_t + e),$$

for all edges $e \in E(B_{2\ell}^\infty(0))$, where $B_r^\infty(x)$ is the vertices inside the L_∞ ball of radius r around x . We will show in Lemma 4.1 that this will happen within time Δ_1 with large enough probability, where Δ_1 has order $\frac{\log^2 n}{\mu}$. This is the shortest of the three phases.

If the first phase does not end within time Δ_1 , we declare the whole three-phase procedure to have failed. This phase will be handled in Sect. 4.

3.2 Second phase: the non-Markovian coupling of the walkers

This is the most involved phase. After the first phase has been completed successfully, the graphs of the two processes are the same on a ball of radius 2ℓ around the walkers. Then, in the second phase we wish to couple the motion of the walkers. We use the tessellation to decide when to couple the walkers identically (so that they jump in the same way) and when to perform a better coupling aiming to decrease the distance between the walkers.

Intuitively, whenever the walkers are passing through a “bad” region of the environment (which in our case will be the 2-enlargement of a bad box) we will just do identity coupling to make sure the distance between the walkers does not increase. In fact, we will only be able to do identity coupling because we will use the annulus between the 2-enlargement of the bad box and the bad box itself (which is composed of good boxes) to give time for the graphs around the walkers to get coupled in both configurations, allowing identity coupling to be carried out. If instead the two walkers are in a great box, then we try to do a better coupling, which we shall refer to as a *simple random walk moment*.

More precisely, translate time so that the second phase starts from time 0. Then, we create the multi-scale tessellation described in Sect. 2 up to time $\Delta_2 + \Delta_3$ where Δ_2 and Δ_3 are of order $\frac{n^2}{\mu}$. We will fix a largest scale k_{\max} and will look at how many times the walkers enter k_{\max} -great boxes.

When the walkers are in great boxes, Lemma 5.14 will give that the environment is favourable enough so that with positive probability the displacement of the walkers will have the same distribution as that of a simple random walk on \mathbb{T}_n^d (i.e., where all edges are open). Phase 2 ends at time Δ_2 where we check whether the walkers are coupled and the graphs are coupled on a ball of radius 2ℓ around the walkers.

Lemma 2.25 says that the walkers will cross an order of n^2 great boxes during $[0, \Delta_2]$ and, therefore, by time Δ_2 the walkers are expected to have done an order of n^2 simple random walk steps. Since two simple random walkers on \mathbb{T}_n^d can be coupled in a way that they coalesce after a time of order n^2 , we can ensure that with high probability phase 2 ends successfully. The details are carried out in Sect. 5.

3.3 Third phase: the coupling of the graphs

The third phase starts at time Δ_2 ; as before we translate time so that the second phase starts at time 0. At the beginning of the third phase the walkers are coupled and the graphs are coupled as well on a ball of radius 2ℓ around them. The idea of this phase

is to keep performing identity couplings until the graphs couple together everywhere. We will show that this simple idea works.

There is one tricky issue. During the second phase, we needed to construct the tessellation all the way to time $\Delta_2 + \Delta_3$, while the second phase ends already at time Δ_2 . The reason for this is that, in order to know whether we can perform a simple random walk moment, we need to observe a little bit of future information about the environment. Therefore, as we performed the second phase, we observed some information from the updates after the end of phase two.

So the goal of the third phase is simply to let time pass until we get to a point where no information regarding future times has been observed, meanwhile doing identity coupling. With this, during the third phase we aim to keep the walkers coupled at all times, while we finish to couple the graphs before time $\Delta_2 + \Delta_3$.

We do not use any further information from the tessellation than what we already observed for phase 2. The delicate point is that in order to apply identity coupling of the walkers, as we explained in the second phase, we have to ensure that the graphs around the walkers are coupled. How large a region we require to be coupled depends on the environment of good and bad boxes that is ahead of the walker, but now we cannot observe anything beyond what we have already observed in phase two; otherwise we would keep observing future information.

As hinted above, we just proceed with identity coupling “blindly”. That is, we perform identity couplings up to time $\Delta_2 + \Delta_3$ assuming that any information that we have not yet observed is “good”, and simply “hope for the best”. It will turn out that this procedure succeeds with large probability leaving the two processes completely coupled (both the graphs and the walkers) by time $\Delta_2 + \Delta_3$. The details of this phase are given in Sect. 6.

3.4 What if a phase fails?

If any of the three phases does not successfully end, we let the two processes run independently (modulo what has already been observed) until the end of the third phase. This is needed as we might have observed some information about the environment up to that time. After that, we repeat the procedure from phase 1. Since the three phases succeed with positive probability, we only need to repeat the whole procedure a constant number of times. The end of the proof of the upper bound is given in Sect. 7.

4 First phase

During the first phase we let the processes $M_t^\star = (X_t, \eta_t^\star)$ and $\overline{M}_t^\star = (\overline{X}_t, \overline{\eta}_t^\star)$ evolve independently. Let $\Psi_t : V \rightarrow V$ be the translation that maps X_t into \overline{X}_t ; we will abuse notation and use the same Ψ_t to denote the corresponding translation map $E \rightarrow E$ of the edges. For any $i \in \mathbb{Z}_+ \cup \{\infty\}$, we define

$$E(B_r^i(v)) = \{(v_1, v_2) \in E : v_1, v_2 \in B_r^i(v)\} \quad \text{and} \quad B_r^i(v) = \{v_1 \in V : \|v - v_1\|_i \leq r\};$$

thus $E(B_r^i(v))$ is the set of edges in the ball of radius r around v according to the norm L_i . We omit i from the superscript whenever $i = 1$. Define the event

$$\mathcal{B}_t = \left\{ \forall e \in E(B_1(X_t)), \eta_t^\star(e) = \bar{\eta}_t^\star(\Psi_t(e)) = 0 \right\} \\ \cap \left\{ \forall e \in E(B_{2\ell}^\infty(X_t)) \setminus E(B_1(X_t)), \eta_t^\star(e) = \bar{\eta}_t^\star(\Psi_t(e)) = \star \right\} \quad (4.1)$$

that the edges in an L_∞ ball of radius 2ℓ around the walkers are all \star at time t , except for the ones adjacent to the walkers which are closed; recall that ℓ is the size of the core of boxes of scale 1, whose value is given in (2.1). Let

$$\tau_B = \inf \{ t \geq 0 : \mathcal{B}_t \text{ holds} \}. \quad (4.2)$$

Note that τ_B is a stopping time. Define $\Delta_1 = \frac{C_9 \log^2 n}{\mu}$, for some constant $C_9(p) > 0$, and define the event

$$F_1 = \{ \tau_B < \Delta_1 \}, \quad (4.3)$$

which we shall take as the event that phase 1 succeeds. This event is a bit more restricted than the one announced in the previous section, but this will be convenient for us in the next phase.

We then run phase 1 until τ_B or Δ_1 , whichever occurs first. If it turns out that F_1 does not occur, phase 1 is then stopped at time Δ_1 and we declare the whole procedure to have failed at time Δ_1 . In this case, we do not proceed to the second phase, and define Δ_1 as the *failing time* of the procedure and, as we will explain more thoroughly in Sect. 7, we will restart from phase 1 from Δ_1 .

The following lemma establishes the probability that the first phase is successful.

Lemma 4.1 (*Phase 1 success probability*) *For any $\delta > 0$, there exists $p_0 = p_0(d, \delta) > 0$ such that for any $p < p_0$, there exists $C_9(p, d, \delta) > 0$ in the definition of Δ_1 so that for any initial configurations $\eta_0^\star, \bar{\eta}_0^\star \in \Omega^\star$ we obtain*

$$\mathbb{P}(F_1^c) \leq \delta$$

for all large enough n .

Before showing that phase 1 succeeds with good probability, we need to establish a simple result on percolation. We then prove Lemma 4.1 in Sect. 4.2.

4.1 Percolation on cylinders and open upwards paths

Let $G = (V, E)$ be a finite graph whose maximum degree is d_{\max} ; in our case, it would be enough to take G to be the d -dimensional torus \mathbb{T}_n^d of side length n , where nearest-neighbors are defined according to the ℓ_∞ norm. We consider the discrete cylinder $V_{\text{cyl}} = V \times \mathbb{Z}_+$ and define a site percolation process on V_{cyl} with parameter $q > 0$. In other words, we declare each site of V_{cyl} to be *open* with probability q , independently of one another; a vertex that is not open is said to be *closed*. For two vertices of $x, y \in V$, we write $x \sim y$ to denote that the graph distance between x and y is at most 1 in G ; thus, for example, $x \sim x$ for all $x \in V$.

Definition 4.2 (open upwards path) An *open upwards path* in V_{cyl} is defined as a sequence of sites $(i_0, \tau_0), (i_1, \tau_1), (i_2, \tau_2), \dots, (i_r, \tau_r)$ such that $i_j \in V$, $\tau_j \in \mathbb{Z}_+$, $i_j \sim i_{j+1}$ and the following holds for all j . If (i_j, τ_j) is open, then $\tau_{j+1} = \tau_j + 1$; otherwise, $\tau_{j+1} - \tau_j \in \{0, 1\}$. In other words, the path is compelled to move “upwards” in the cylinder when it visits open sites.

Note that an open upwards path is allowed to visit a vertex more than once. We say that an open upwards path $(i_0, \tau_0), (i_1, \tau_1), (i_2, \tau_2), \dots, (i_r, \tau_r)$ *traverses m levels* if $\tau_r - \tau_0 = m$. When ϱ is close to 1, an open upwards path cannot visit too many closed sites. This is quantified in the next lemma.

Lemma 4.3 (Open upwards path) Let $(i_0, 0) \in V_{\text{cyl}}$ be fixed. For any $\alpha > 0$, there exists $\varrho' = \varrho'(\alpha, d_{\max}) \in (0, 1)$ such that if $\varrho > \varrho'$ then the probability that there exists an open upwards path from $(i_0, 0)$ that traverses m levels and visits at least αm distinct closed sites is at most e^{-cm} for some constant $c = c(\alpha, d_{\max}) > 0$.

Before proving the above result, we need the following estimate on the number of subgraphs of G that contain a given vertex.

Lemma 4.4 Given a vertex $v \in V$, let A_r be the number of induced connected subgraphs of V containing v and having r vertices. There exists a constant $c = c(d_{\max}) > 0$ such that $A_r \leq e^{cr}$ for all r .

Proof This proof is quite standard and a version for the lattice can be found in [6, Proof of Theorem 4.20]; we include a proof here for the sake of completeness. Let $A_{r,s}$ be the number of induced connected subgraphs of V containing v and having r vertices and s boundary vertices, where a boundary vertex is a vertex that does not belong to the subgraph but has a neighbor who does. Hence, $A_r = \sum_s A_{r,s}$. Note that for any $\varrho \in (0, 1)$, if we perform percolation on V , we obtain

$$\sum_{r,s} A_{r,s} \varrho^r (1 - \varrho)^s = 1. \quad (4.4)$$

For any vertex $u \in V$ denote $d(u)$ its degree in G . Let $S \in V$ be one subgraph counted in $A_{r,s}$, denote $m(S)$ the number of edges between vertices of S and ∂S the number of edges between vertices in S and vertices in $V \setminus S$. Then,

$$d_{\max} r \geq \sum_{u \in S} d(u) = 2m(S) + \partial S \geq 2(r - 1) + s.$$

Thus, $s \leq (d_{\max} - 2)r + 2$. Plugging this result into (4.4) and taking $\varrho = \frac{1}{2}$ we obtain

$$1 \geq \sum_{r,s} A_{r,s} \varrho^r (1 - \varrho)^{(d_{\max}-2)r+2} = \sum_{r,s} A_{r,s} 2^{-(d_{\max}-1)r-2} = \sum_r A_r 2^{-(d_{\max}-1)r-2}.$$

Thus, $A_r \leq 2^{(d_{\max}-1)r+2}$ for each r . □

Proof of Lemma 4.3 Let m be an integer and, for convenience, set $\tau_0 = 0$. Consider an open upwards path $(i_0, \tau_0), (i_1, \tau_1), (i_2, \tau_2), \dots, (i_r, \tau_r)$ such that $\tau_r - \tau_0 = m$; that is, the path traverses m levels. Let s_1 be the number of distinct closed sites visited by the path before it traverses 1 level, and for $j \geq 2$ let s_j be the number of distinct closed sites visited by the path after having traversed $j - 1$ levels and before traversing j levels. So $\sum_{j=1}^m s_j$ is the total number of distinct closed sites visited by the path. Note that, the sites counted in each s_j must be of the form $(\cdot, j - 1)$ and must form a connected set with respect to the relation \sim over G . Using Lemma 4.4, given s_1, s_2, \dots, s_m , the number of possible ways to pick the set of distinct sites within $(i_0, \tau_0), (i_1, \tau_1), \dots$ is

$$(d_{\max} + 1)^m \prod_{j=1}^m s_j e^{c' s_j},$$

where c' is the constant from Lemma 4.4, and the term s_j in the product counts the number of sites at level j that can be selected to be the last vertex visited by the path before going to level $j + 1$. Then, $(d_{\max} + 1)^m$ accounts for the number of ways to choose the first site at level j given the last site at level $j - 1$; this amounts to at most $d_{\max} + 1$ choices per level. If we fix $\sum_{j=1}^m s_j = S$, the number of ways to select the s_j is $\binom{S+m-1}{S}$. Finally, given all sites in the path with s_j as defined above, the probability that this path is an open upwards path is at most $\prod_{j=1}^m (1 - \varrho)^{s_j}$ since each site counted in the s_j must be closed. Therefore, the expected number of open upwards paths that traverse m levels and visit at least αm closed sites is at most

$$\begin{aligned} & \sum_{S \geq \alpha m} \binom{S+m-1}{S} (d_{\max} + 1)^m \prod_{j=1}^m e^{c' s_j} s_j (1 - \varrho)^{s_j} \\ &= (d_{\max} + 1)^m \sum_{S \geq \alpha m} \binom{S+m-1}{S} \left(c'' e^{c''} (1 - \varrho) \right)^S, \end{aligned}$$

where we used that given c there exists a constant c'' such that $z e^{c' z} \leq c'' e^{c'' z}$ for all z . It is enough to use the trivial bound $\binom{S+m-1}{S} \leq 2^{S+m-1}$ in the above expression to obtain the upper bound

$$(2(d_{\max} + 1))^m \sum_{S \geq \alpha m} \left(2c'' e^{c''} (1 - \varrho) \right)^S \leq 2 (2(d_{\max} + 1))^m \left(2c'' e^{c''} (1 - \varrho) \right)^{\alpha m},$$

with the inequality hold whenever ϱ is close enough to 1 so that $2c'' e^{c''} (1 - \varrho) \leq \frac{1}{2}$. Then the lemma holds by setting ϱ further closer to 1 so that $4(d_{\max} + 1) \left(2c'' e^{c''} (1 - \varrho) \right)^\alpha \leq e^{-c}$. \square

By using a result by Liggett, Schonmann and Stacey [8], the above result can be extended to percolation on $\mathbb{T}_d \times \mathbb{Z}_+$ with bounded dependences.

Lemma 4.5 *Let $C \geq 1$ be a constant. Consider a site percolation process on $\mathbb{T}_d \times \mathbb{Z}_+$ where the probability that a given site is open depends on at most C other sites. Then Lemma 4.3 holds with the lower bound on q depending on C .*

Proof For any \tilde{q} , provided q is large enough we can apply Liggett, Schonmann and Stacey [8, Theorem 0.0] to obtain that the dependent site percolation process stochastically dominates an independent site percolation process of parameter \tilde{q} . The lemma then follows by applying Lemma 4.3 to this independent site percolation process. \square

4.2 Proof of Lemma 4.1

Now we are in a position to establish the occurrence of the first phase.

Proof of Lemma 4.1 Let τ'_B be the first time $t \geq t_1$ such that X_t and \bar{X}_t are both isolated, meaning that all edges adjacent to them are closed. We will show that τ'_B occurs before time $\Delta_1 - t_1 + \bar{t}_1$.

For each process M_t^\star and \bar{M}_t^\star we create a tessellation of $\mathbb{T}_n^d \times [0, \Delta_1]$ into boxes of scale 1 using the values for ℓ and t_1 from Sect. 2. The event that a given box is good is defined as in Definition 2.7. We let M_t^\star and \bar{M}_t^\star evolve independently of one another until a stopping time st_1 where X_{st_1} and \bar{X}_{st_1} are both in good boxes and $s \geq 1$. Note that good 1-boxes form a dependent site percolation process on $\mathbb{T}_n^d \times \mathbb{Z}_+$ so that we can apply Lemma 4.5. Let $(i_0, 1)$ and $(\bar{i}_0, 1)$ be the boxes visited by X_t and \bar{X}_t at time $t = t_1$. Now, since random walks must traverse a feasible path, and since feasible paths leave good boxes from the time boundary (cf. Lemma 2.21), we obtain that from $(i_0, 1)$ and $(\bar{i}_0, 1)$ the random walks X_t and \bar{X}_t must traverse an open upwards path. Therefore, the probability that up to level $m = \Delta_1/t_1 - 1$ we have that X_t and \bar{X}_t each visited more than $\frac{m}{3}$ bad 1-boxes is at most $2e^{-cm}$ provided p is small enough (which makes the probability that a 1-box being good large enough). Under this event, there must exist $\frac{m}{3}$ instances of time $s \leq m$ at which X_{st_1} and \bar{X}_{st_1} are both in good 1-boxes. when this happens, at time $st_1 + \bar{t}_1$ both $X_{st_1 + \bar{t}_1}$ and $\bar{X}_{st_1 + \bar{t}_1}$ are isolated in a vertex (i.e., all edges adjacent to them are closed). Therefore,

$$\mathbb{P}(\tau'_B \geq \Delta_1 - t_1 + \bar{t}_1) \leq n^{2d} \mathbb{P}(s > m) \leq 2ne^{-cm},$$

where the term n^{2d} accounts for the number of choices for i_0 and \bar{i}_0 .

Now let \mathcal{F} be the σ -algebra generated by M_t^\star and \bar{M}_t^\star during $t \in [0, \tau'_B]$. We want to establish a lower bound on the probability that $\mathcal{B}_{\tau'_B + \bar{t}_1}$ given \mathcal{F} . Since X_t and \bar{X}_t are isolated in $t = \tau'_B$, it is enough to compute the probability that all edges inside a L_∞ ball of radius 2ℓ around the walkers do a \star -update but no non- \star update, and the edges adjacent to the walkers do not open or do a non \star -update during $[\tau'_B, \tau'_B + \bar{t}_1]$. This probability is

$$\exp(-\bar{t}_1 \mu (1 - p_\star) 2(4\ell)^d) (1 - \exp(-\bar{t}_1 \mu p_\star))^{2(4\ell)^d} \exp(-\bar{t}_1 \mu p_{\min} 4d),$$

where the first term is the probability that no edge in the L_∞ ball of radius 2ℓ around the walkers does a non \star -update, the second term is the probability that those edges

do a \star -update and the last term is the probability that the edges adjacent to the walkers do not open. Therefore,

$$\begin{aligned}\mathbb{P}(F_1^c) &\leq 2ne^{-cm} + 1 - \exp\left(-\bar{t}_1\mu(1-p_\star)2(4\ell)^d - \bar{t}_1\mu p_{\min}4d\right) \\ &\quad (1 - \exp(-\bar{t}_1\mu p_\star))^{2(4\ell)^d} \\ &\leq 2ne^{-cm} + 1 - \exp\left(-\bar{t}_1\mu(1-p_\star)2(4\ell)^d - \bar{t}_1\mu p_{\min}4d\right) \\ &\quad - 2(4\ell)^d \exp(-\bar{t}_1\mu p_\star).\end{aligned}$$

Recall that $\Delta_1 = \frac{C_9 \log^2 n}{\mu}$, $\bar{t}_1 = \frac{\log^2 \ell}{\mu}$ and $t_1 = \frac{\sqrt{\ell}}{\mu}$, where ℓ is just a large enough constant that is set before letting p be small enough. Now we show that we can make the above smaller than δ . We start with the term $2(4\ell)^d \exp(-\bar{t}_1\mu p_\star)$, which can be made, say, smaller than $\frac{\delta}{3}$. We will do this by adjusting ℓ only, but this term involves also p though p_\star . However, note that p_\star goes to 1 as p goes to 0. So, since $\bar{t}_1\mu$ is of order $\log^2 \ell$, we can choose ℓ large enough so that $2(4\ell)^d \exp(-\bar{t}_1\mu p_\star) \leq \frac{\delta}{3}$ for all p so that $p_\star \geq \frac{1}{2}$. After fixing ℓ , we can take p close enough to 0, which makes p_{\min} goes to 0 and p_\star goes to 1, so that $\exp(-\bar{t}_1\mu(1-p_\star)2(4\ell)^d - \bar{t}_1\mu p_{\min}4d) \geq 1 - \frac{\delta}{3}$. Finally, after fixing ℓ and p , we can take n large enough so that $2ne^{-cm} \leq \frac{\delta}{3}$ since $m = \frac{\Delta_1}{t_1} - 1$ is of order $\log^2 n$ as a function of n . This concludes the first phase. \square

5 The second phase: non Markovian coupling

To describe the coupling during the second phase we will use the full multi-scale space-time tessellation described in Sect. 2. For simplicity, we translate time so that this phase starts at time 0 and that X_0 is at the origin. Hence, \bar{X}_0 can be arbitrary, and η_0^\star and $\bar{\eta}_0^\star$ can be any configuration for which the event \mathcal{B}_0 from (4.1) holds.

5.1 Largest scale

We begin by creating the multi-scale space-time tessellation of $\mathbb{T}_n^d \times [0, \Delta_2]$ and with largest scale

$$k_{\max} = \log_2 \log n. \quad (5.1)$$

We consider a positive constant $C_1 0(p) > 0$ to be chosen later so that $t_{k_{\max}}$ divides $C_{10} \frac{n^2}{\mu}$, and define

$$\Delta_3 = \Delta_2 + \frac{n^2}{\mu} \quad \text{and} \quad \Delta_2 = C_{10} \frac{n^2}{\mu}. \quad (5.2)$$

The following lemma shows that with large probability there are no bad boxes of scale k_{\max} or larger. This will allow us to restrict our analysis to boxes of scale at

most k_{\max} . We will consider all the boxes contained into the tessellation $\mathbb{T}_n^d \times [0, \Delta_3]$, which in particular are all the boxes intersecting the tessellation of $\mathbb{T}_n^d \times [0, \Delta_2]$.

Lemma 5.1 *For any $\delta > 0$, there exists $p_0 = p_0(\delta, d) > 0$ such that for all $p < p_0$ and n large enough*

$$\mathbb{P}\left(R_k(i, \tau) \text{ is bad for some } R_k(i, \tau) \subset \mathbb{T}_n^d \times [0, \Delta_3], \text{ with } k \geq k_{\max}\right) \leq \rho_1^{2^{k_{\max}-3}}.$$

Proof The number ζ_k of boxes of scale k in $\mathbb{T}_n^d \times [0, \Delta_3]$ is trivially bounded as

$$\zeta_k \leq \left(\frac{n}{\ell_k}\right)^d \frac{\Delta_3}{t_k} \leq (C_{10} + 1)n^{d+2}.$$

Using Lemma 2.15 the probability that there exists a box of scale k_{\max} or bigger that is bad is bounded above by

$$\sum_{k \geq k_{\max}} \zeta_k \rho_k \leq (C_{10} + 1)n^{d+2} \sum_{k \geq k_{\max}} \rho_1^{2^{k-2}} \leq 2(C_{10} + 1)n^{d+2} \rho_1^{2^{k_{\max}-2}}.$$

Using the value of k_{\max} and the fact that ρ_1 can be made arbitrarily small by taking p small concludes the proof. \square

5.2 The coupling

Recall the map Ψ_t introduced in Sect. 4 which maps X_t into \overline{X}_t . In order to define the coupling of the two processes, we will use a different map Φ_t . The idea is that our new map will be equal to Ψ_t in good parts of the environment, but when the walker enters the enlargement of a bad box, we will stop changing Φ_t and will keep it “frozen” until the walkers exit the enlargements of all bad boxes. The idea is that in the enlargement of bad boxes we want to couple the graphs in a large region around the walkers so that if the walkers enter a bad box, then they do so with their graphs coupled within the box. We stop updating Φ_t because when Φ_t changes many edges uncouple.

More precisely, given a time t , denote with

$$\bar{s}_t = \sup\{s \leq t : (X_s, s) \text{ is inside } ak_{\max}\text{-great box}\}$$

the last time before t the walker is in a k_{\max} great box. We will consider the new map Φ_t defined as

$$\Phi_t = \Psi_{\bar{s}_t}.$$

We will show that this change of map actually will not create any problems; in fact, we will show that $\Phi_t \equiv \Psi_t$ for all t because in the way we construct the coupling, when the walkers are in the enlargement of a bad box, we will succeed in applying identity coupling, hence the translation map remains constant. So, the introduction

of Φ_t here is a formalism so that the coupling procedure is well defined. This will imply that our application of identity coupling later on will be successful, which in turn implies that $\Phi_t \equiv \Psi_t$.

As soon as the second phase begins we check whether the box $R_1(i, 0)$, such that $(X_0, 0) \in R_1^{\text{core}}(i, 0)$, is k_{\max} -great (the reason we do this will be clarified later, see Remark 5.9). If that is the case then we can begin the coupling procedure relative to the second phase. The coupling is composed of two parts: the coupling of the graphs (that is, the coupling of η_t^* and $\bar{\eta}_t^*$) and the coupling of the walkers.

5.2.1 Coupling of the graphs

We let the process $\{\eta_t^*\}_{t \geq 0}$ evolve. Denote with $\mathcal{C}_v(t)$ (resp., $\bar{\mathcal{C}}_v(t)$) the cluster that contains vertex v at time t in the process η_t^* (resp., $\bar{\eta}_t^*$). When an update (s, U', U) occurs at an edge e in η_s^* we update the process $\bar{\eta}_s^*$ as follows.

- If the update is a \star -update we refrain from looking at U and instead simply set $\eta_s^*(e) = \star$ and $\bar{\eta}_s^*(\Phi_s(e)) = \star$.
- If the update is not a \star -update we must check in both configurations η_s^* and $\bar{\eta}_s^*$ whether e is a cut-edge or not. We do this by looking at the connected components of the endpoints v_1, v_2 of the edge e . If an edge e' is such that $\eta_s^*(e') = \star$ and e' is incident to a vertex in $\mathcal{C}_{v_1}(s) \cup \mathcal{C}_{v_2}(s)$, we sample its current status, open or closed, according to its last update. Note that this last update is itself a tuple $(\bar{s}, \bar{U}', \bar{U})$, so this step boils down to checking the value of \bar{U} . If $\bar{\eta}_s^*(\Phi_s(e')) = \star$ we set $\bar{\eta}_s^*(\Phi_s(e')) = \eta_s^*(e')$ as well. We continue this procedure until the components of v_1 and v_2 have been fully explored in η_s^* and proceed analogously for the process $\bar{\eta}_s^*$ until the components of $\Phi_s(v_1)$ and $\Phi_s(v_2)$ have been fully explored. A potential disagreement $\eta_s^*(e) \neq \bar{\eta}_s^*(\Phi_s(e))$ can happen only if, by revealing the components of $v_1, v_2, \Phi_s(v_1)$ and $\Phi_s(v_2)$, we find that e is a cut-edge in η_s^* but $\Phi_s(e)$ is not a cut-edge in $\bar{\eta}_s^*$, or vice-versa.

In this way edges whose status is \star can always be coupled equivalently whereas non \star -updates cause the reveal of the status of other edges, potentially creating disagreements between the two configurations.

Remark 5.2 (Momentaneous change of coupling) At some times we will carry out a different coupling of the environment. This will be done by simply introducing another map $\tilde{\Phi}$ of the environments, and the coupling of the graphs will go as described above with Φ_t replaced with $\tilde{\Phi}$ until we specify that Φ_t is again the map to be used.

5.2.2 Coupling of the walkers

During this discussion the reader should refer to Fig. 3.

Our goal is to define a coupling that can bring the walkers together. For this we will use the multi-scale tessellation. The coupling of the walkers will be composed of two different couplings. When the walker X_s enters the core of a great box $R_1^{\text{core}}(i, \tau)$, we will try to take advantage of the nice environment that a great box provides to perform a coupling that we refer to as a *simple random walk moment*. This coupling

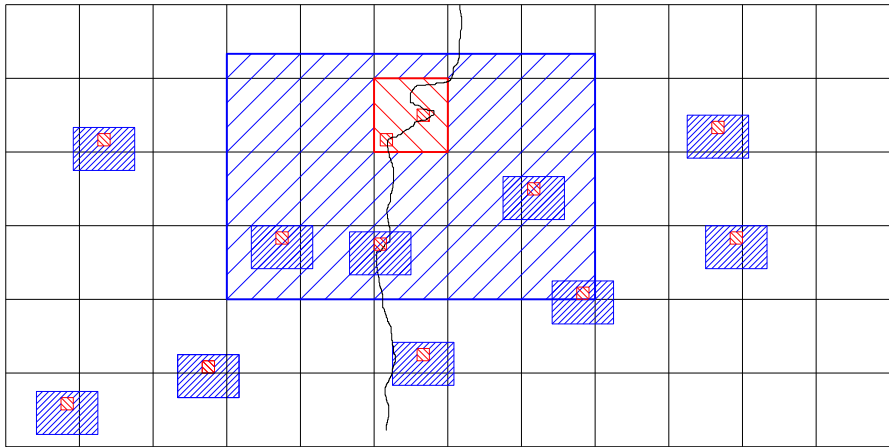


Fig. 3 In red the bad boxes, in blue their enlargement, in black the tessellation and the walker's trajectory. In bad boxes there is no control over the displacement of the walker, whereas in good boxes, the walker always leaves the box from its time boundary. Whenever the walker enters the enlargement of a bad box, we start doing identity coupling. Otherwise, the walker is in great boxes, and we attempt to check whether a SRWM occurs, in which case a coupling of simple random walks is performed

aims to change the distance between the walkers, so that eventually the walkers may find themselves at the same site.

On the other hand, whenever X_s is not in a great box, then we do not have a good enough control on the environment around the walker to do a simple random walk moment. In such cases, we will just resort to a simple identity coupling that keeps the distance between the walkers unchanged. An identity coupling will only be able to be performed if the environment around the walkers are the same. For this, we define the following event:

$$B'_t = \left\{ \forall e \in E(B_{\ell/2}^\infty(X_t)), \eta_t^*(e) = \overline{\eta}_t^*(\Phi_t(e)) \right\}. \quad (5.3)$$

If B'_s holds for all $s \in (s_1, s_2)$, then in this time interval the walkers can perform the same jumps and not change their relative distance. In other words, identity coupling is successful. In fact if the environment around the walkers is the same (as a matter of fact we only need the environments to agree on a ball of radius 1 around the walkers), by doing identity coupling the walkers are able to perform the same jumps.

So the proof is now split into three steps. Since Φ_t does not change when the walker enters the 2-enlargement of a bad box, we will show in Sect. 5.3 that when Φ_t does not change the graph couples. Next, we deal with showing that identity coupling can be successfully implemented as the walker enters the 2-enlargement of a bad box (i.e., when the walker is not in a great box). This is carried out in Sect. 5.4. Then in Sect. 5.5 we deal with the simple random walk moments.

5.3 Coupling of the graphs with Φ_t unchanged

Given $I \subset \mathbb{Z}^d$ and $k \geq 1$, let

$$\mathcal{S}_k(I) = \bigcup_{i \in I} S_k(i) \quad \text{and} \quad \mathcal{S}_k^\diamond(I) = \bigcup_{i \in I} S_k^\diamond(i),$$

for any $\diamond \in \{\text{core}, \text{enl1}, \text{enl2}\}$. Recall the value m in the definition of ℓ_k in (2.2). Recall also \bar{t}_k from (2.7). Then, for $k \geq 2$, we define

$$\bar{t}_k = \frac{6t_k}{(k-1)^2 m}. \quad (5.4)$$

We start this section showing that the graph gets coupled in regions of good boxes if Φ_t does not change.

Lemma 5.3 (*Graphs couple in good boxes*) *Let m be large enough, and then let ℓ be large enough with respect to m . Let $R_k(i, \tau)$ be a good box, and let s_1 be any time instance so that $[s_1, s_1 + 2\bar{t}_k] \subset T_k(\tau)$. If Φ_t does not change during $[s_1, s_1 + 2\bar{t}_k]$, then*

$$\text{there exists } t \in [s_1, s_1 + 2\bar{t}_k] \text{ such that } \eta_t^*(e) = \bar{\eta}_t^*(\Phi_t(e)) \text{ for all } e \in S_k(i). \quad (5.5)$$

Proof If $k = 1$ then the proof follows since each edge of $S_k(i)$ receives only \star -updates and gets updated at least once during $[s_1, s_1 + 2\bar{t}_k]$. For $k \geq 2$, we assume that the statement of the lemma holds up to scale $k - 1$. Let $s_2 = \max T_k(\tau)$. Let τ' be the first time index such that $\tau' t_{k-1} \in [s_1, s_2]$ and all boxes $R_{k-1}(\cdot, \tau') \subset R_k(i, \tau)$ are good. Let I be the set of indices containing all $(k - 1)$ -boxes that are inside $S_k(i)$; more precisely,

$$I = \{i' : S_{k-1}(i') \subset S_k(i)\}.$$

Then, by induction, by time $\tau' t_{k-1} + 2\bar{t}_{k-1}$ we obtain that $S_{k-1}(I) = S_k(i)$ has been coupled.

Now it remains to show that $\tau' t_{k-1} + 2\bar{t}_{k-1} \leq s_1 + 2\bar{t}_k$. Note that since $R_k(i, \tau)$ is a good box, there exist $\hat{\tau}, \hat{\tau}'$ such that all $(k - 1)$ -bad boxes contained in $R_k(i, \tau)$ are contained in $R_{k-1}^{\text{enl1}}(\hat{\tau}, \hat{\tau}')$. Since the amount of time spanned by the enlargement at scale $k - 1$ is $9t_{k-1}$, we obtain that $\tau' t_{k-1} \leq s_1 + 9t_{k-1} + t_{k-1}$, where the last t_{k-1} is to account for the possibility that s_1 is not a multiple of t_{k-1} . Hence, using the notation $a_+ = \max\{a, 1\}$ for consistency with the case $k = 2$, and noting that $\bar{t}_1 \leq \frac{6t_1}{(k-2)_+^2 m}$ provided ℓ is made large enough once m has been fixed, we have

$$\begin{aligned} \tau' t_{k-1} + 2\bar{t}_{k-1} &\leq s_1 + 10t_{k-1} + 2\bar{t}_{k-1} \\ &\leq s_1 + t_{k-1} \left(10 + 2 \frac{6}{(k-2)_+^2 m} \right) \end{aligned}$$

$$\begin{aligned}
 &= s_1 + \frac{t_k}{m(k-1)^2} \left(10 + \frac{12}{(k-2)_+^2 m} \right) \\
 &= s_1 + \frac{\bar{t}_k}{6} \left(10 + \frac{12}{(k-2)_+^2 m} \right) \\
 &\leq s_1 + 2\bar{t}_k.
 \end{aligned}$$

□

Recall the definition of $S_1^{\text{inn}}(i)$ from (2.8). For $k \geq 2$, define

$$S_k^{\text{inn}}(i) = \bigcup_{j: S_{k-1}^{\text{enl}2}(i) \subset S_k(j)} S_{k-1}^{\text{core}}(j). \quad (5.6)$$

For a set of indices I , we write

$$S_k^{\text{inn}}(I) = \bigcup_{j \in I} S_k^{\text{inn}}(j).$$

Note that by taking m large enough, then $S_k^{\text{core}}(I) \subset S_k^{\text{inn}}(I) \subset S_k(I)$. We start with a simple result about the connected component of a vertex.

Lemma 5.4 *Let $m \geq 2$ and let ℓ be large enough with respect to m . Let $I \subset \mathbb{Z}^d$ be a set of indices, $k \geq 1$ a scale and $\tau \geq 1$ a time index such that $R_k(i, \tau)$ is a good box for all $i \in I$. Then, for any $v \in S_k^{\text{inn}}(I)$ and any $t \in T_k(\tau)$, the connected component of v is contained in $B_{5\ell_k/m}^\infty(v)$, where we recall that $B_r^\infty(v)$ is the L_∞ ball of radius r around v .*

Proof For $k = 1$, the result follows by the fact that components have size at most $\log^2 \ell$ in good 1-boxes when $\tau \geq 1$, and ℓ is large enough so $\log^2 \ell \leq 5\ell/m$. For $k \geq 2$, let (i', τ') be such that $v \in S_{k-1}^{\text{core}}(i')$ and $t \in T_{k-1}(\tau') \subset T_k(\tau)$; there could be more than one choice for τ' , it is irrelevant which one we pick. Note that since $v \in S_k^{\text{inn}}(I)$

$$S_{k-1}(i') \subset S_{k-1}^{\text{enl}2}(i') \subset S_k(i) \text{ for some } i \in I.$$

If $R_{k-1}(i', \tau')$ is good, then the connected component of v is contained in $B_{5\ell_{k-1}/m}(v) \subset B_{5\ell_k/m}(v)$ by applying the induction hypothesis at scale $k-1$ and set of indices $\{i'\}$. Otherwise, note that by Remark 2.17 we have that $R_{k-1}^{\text{enl}1}(i', \tau')$ contains all bad boxes in $R_k(i, \tau)$. If the connected component of v is contained in $S_{k-1}^{\text{enl}1}(i')$ then it is contained in $B_{5\ell_{k-1}}(v) \supset S_{k-1}^{\text{enl}1}(i')$. Since $5\ell_{k-1} = 5\frac{\ell_k}{m(k-1)^2} \leq \frac{5\ell_k}{m}$ the lemma holds on this case as well. In the final case, when the connected component of v is not contained in $S_{k-1}^{\text{enl}1}(i')$, it may sound contradictory but we can get an even smaller bound for the component of v . The reason is that there must exist i'' such that v is at the same component of a vertex u with $u \in S_{k-1}^{\text{core}}(i'')$ and

$S_{k-1}^{\text{core}}(i'') \cap S_{k-1}^{\text{enl1}}(i') = \emptyset$ but $S_{k-1}(i'') \cap S_{k-1}^{\text{enl1}}(i') \neq \emptyset$. But since (u, t) is in the box $R_{k-1}(i'', \tau')$, and $S_{k-1}(i'') \subset S_{k-1}^{\text{enl2}}(i') \subset S_k(i)$, we have that $R_{k-1}(i'', \tau')$ is a good box. Thus, by induction we obtain that the connected component of v is inside $B_{5\ell_{k-1}/m}(u) \subset B_{10\ell_{k-1}/m}(v)$. Since $10 \frac{\ell_{k-1}}{m} = 10 \frac{\ell_k}{m^2(k-1)^2} \leq \frac{5\ell_k}{m}$ for all k as long as $m \geq 2$, the proof is completed. \square

With the help of the above lemma, we can show that the graph cannot uncouple in regions surrounded by good boxes.

Lemma 5.5 (Graphs remain coupled if Φ_t does not change) *Let m be large, and let ℓ be large enough with respect to m . Let $R_k(i, \tau)$ be a good box, and let $s_1 < s_2$ with $s_1, s_2 \in T_k(\tau)$ and $s_1 \leq (\tau + 1)t_k$. If Φ_t does not change during $t \in [s_1, s_2]$ and*

$$\eta_{s_1}^*(e) = \bar{\eta}_{s_1}^*(\Phi_{s_1}(e)) \text{ for all } e \in E(S_k(i)),$$

then $\eta_t^(e) = \bar{\eta}_t^*(\Phi_t(e))$ for all $e \in E(S_k^{\text{inn}}(i))$ and all $t \in [s_1, s_2]$.*

Proof For $k = 1$ the lemma is obvious, since for any $e \in E(S_k(i))$, e only receives \star -updates during $T_k(\tau)$. Therefore, $\eta_t^*(e) = \bar{\eta}_t^*(\Phi_t(e))$ for all $t \in [s_1, s_2]$. For $k \geq 2$, assume the lemma holds up to scale $k - 1$. Let

$$\mathcal{T} = \{\tau' : T_{k-1}(\tau') \cap [s_1, s_2] \neq \emptyset \text{ and } T_{k-1}(\tau') \subset T_k(\tau)\}.$$

Let $\tau_1 = \min \mathcal{T}$. Note that either

$$s_1 \in T_{k-1}^{\text{core}}(\tau_1) \quad \text{or} \quad s_1 \in T_{k-1}^{\text{core}}(\tau_1 - 1), \quad (5.7)$$

where the latter happens when s_1 is near the starting time of $T_k(\tau)$. Because s_1 cannot be near the ending time of $T_k(\tau)$ due to the condition $s_1 \leq (\tau + 1)t_k$, we obtain that \mathcal{T} is not empty. We will first show that

$$\begin{aligned} S_{k-1}^{\text{inn}}(j) \text{ is coupled during } [s_1, s_2] \text{ for all } j \text{ such that } S_{k-1}(j) \subset S_k(i) \\ \text{and for which } R_{k-1}(j, \tau') \text{ is good for all } \tau' \in \mathcal{T}. \end{aligned} \quad (5.8)$$

To see this, let $r_1 = \sup T_{k-1}(\tau_1)$ and note that $r_1 \geq s_1 + t_{k-1}$ because of (5.7). Now, induction gives that $S_{k-1}^{\text{inn}}(j)$ remains coupled up to time r_1 . We would like to reapply the induction hypothesis on the box $S_{k-1}(j)$ in the next time step, but for this we need $S_{k-1}(j)$ to be coupled, not only $S_{k-1}^{\text{inn}}(j)$. Thus, we first apply Lemma 5.3 from time $r_1 - 2\bar{t}_{k-1}$ to obtain that there exists a time $r'_1 \in [r_1 - 2\bar{t}_{k-1}, r_1]$ for which the whole of $S_{k-1}(j)$ is coupled. Let τ_2 be such that $r'_1 \in T_{k-1}^{\text{core}}(\tau_2)$ and note that $\tau_2 \geq \tau_1 + 1$. Thus, we repeat the induction hypothesis and the application of Lemma 5.3 to obtain a sequence of τ_i, r_i and r'_i until a certain value $r'_i \in [s_2 - 2\bar{t}_{k-1}, s_2]$. At that time, the induction hypothesis gives that $S_{k-1}^{\text{inn}}(j)$ is coupled at time s_2 , establishing (5.8).

Now we turn to establish the lemma. If $R_k(i, \tau)$ has no bad $(k - 1)$ -box intersecting the time interval $[s_1, s_2]$, then (5.8) and the fact that $(k - 1)$ -boxes overlap give that $\bigcup_{j: S_{k-1}(j) \subset S_k(i)} S_{k-1}^{\text{inn}}(j) \supset S_k^{\text{inn}}(i)$ is coupled during $[s_1, s_2]$.

Now assume that $R_k(i, \tau)$ contains bad $(k-1)$ -boxes that intersect $[s_1, s_2]$. From Remark 2.17, there exists $R_{k-1}(i', \tau')$ so that all $(k-1)$ -bad boxes contained in $R_k(i, \tau)$ are contained in $R_{k-1}^{\text{enl}}(i', \tau')$. Let

$$\mathcal{J} = \left\{ j : S_{k-1}(j) \subset S_k(i) \text{ and } S_{k-1}(j) \not\subset S_{k-1}^{\text{enl}}(i') \right\},$$

and note that $R_{k-1}(j, \tau'')$ is good for all $j \in \mathcal{J}$ and $\tau'' \in \mathcal{T}$. Therefore, (5.8) gives that $S_{k-1}^{\text{inn}}(j)$ is coupled during $[s_1, s_2]$ for all $j \in \mathcal{J}$. The remaining of the proof is split into two cases. First assume that $S_{k-1}^{\text{enl}}(i')$ is separated from infinity by \mathcal{J} , which means that any path from $S_{k-1}^{\text{enl}}(i')$ to the outside of $S_k(i)$ must enter $S_{k-1}^{\text{core}}(j)$ for some $j \in \mathcal{J}$. In fact, letting

$$\mathcal{J}' = \left\{ j \in \mathcal{J} : S_{k-1}^{\text{core}}(j) \subset S_{k-1}^{\text{enl}}(i') \right\},$$

we get that the path must enter $S_{k-1}^{\text{core}}(j)$ for some $j \in \mathcal{J}'$. Besides, Lemma 5.4 gives that for all $v \in S_{k-1}^{\text{inn}}(\mathcal{J})$ and all $s \in T_{k-1}^{\text{enl}}(\tau')$ we have that the connected component of v is contained in $B_{5\ell_{k-1}/m}^\infty(v)$. Therefore, all connected components intersecting $S_{k-1}^{\text{enl}}(i')$ must be contained in $\bigcup_{v \in S_{k-1}^{\text{enl}}(i')} B_{5\ell_{k-1}/m}^\infty(v)$, which is a spatial region contained in the interior of $S_k^{\text{inn}}(J')$. Therefore, since $S_{k-1}^{\text{inn}}(J) \supset S_{k-1}^{\text{inn}}(J')$ remains coupled throughout $[s_1, s_2]$ by (5.8), non- \star updates inside $S_{k-1}^{\text{enl}}(i')$ cannot uncouple the graph.

Turning to the second case, we assume that $S_{k-1}^{\text{enl}}(i')$ is not separated from infinity by \mathcal{J} . This means that $S_{k-1}^{\text{enl}}(i')$ is so close to the boundary of $S_k(i)$ that it does not intersect $S_k^{\text{inn}}(i)$. More formally, for any $v \in S_{k-1}^{\text{enl}}(i)$ we have that $B_{10\ell_{k-1}}^\infty(v)$ cannot be contained in $S_k(i)$. But this implies that any i'' with $S_{k-1}^{\text{core}}(i'') \subset S_{k-1}^{\text{enl}}(i')$ we have that $S_{k-1}^{\text{enl}2}(i'') \not\subset S_k^{\text{inn}}(i)$. Therefore, applying (5.8) to the boxes in \mathcal{J} already gives that $S_k^{\text{inn}}(i)$ is coupled during $[s_1, s_2]$. \square

5.4 Identity coupling

We prove that, by doing identity coupling, as long as the particle X_t is in a point $(v, t) \in \mathbb{T}_n^d \times \mathbb{R}^+$ in space-time that is part of a 1-box $R_1(\cdot, \cdot)$ that is good, it is always possible to keep the distance between X_t and \bar{X}_t constant. Recall the event \mathcal{B}'_t from (5.3), the event \mathcal{B}_t from (4.1), and the definition of the spatial core of a box in (2.11). We will need a weaker version of \mathcal{B}'_t which we define as

$$\mathcal{B}''_t = \left\{ \forall e \in E \left(B_{\ell/3}^\infty(X_t) \right), \eta_t^*(e) = \bar{\eta}_t^*(\Phi_t(e)) \right\}. \quad (5.9)$$

Recall that in the second phase we assume that $R_0(0, 0)$ is a k_{\max} -great box and \mathcal{B}_0 holds; we do not restate these conditions on the lemmas.

Lemma 5.6 (Identity coupling succeeds in good boxes) *Let s_1 be a time so that \mathcal{B}_{s_1}'' holds and $(X_{s_1}, s_1) \in R_1^{\text{S-core}}(i, \tau)$ with $R_1(i, \tau)$ being a good box. Let $s_2 \in T_1(\tau)$,*

$s_2 \geq s_1$. If we attempt to do identity coupling for the entire time interval $[s_1, s_2]$, then the coupling succeeds and $\Psi_{s_1} \equiv \Psi_t$ for all $t \in [s_1, s_2]$.

Proof Let B be the L_∞ ball of radius $\ell/3$ around X_{s_1} ; B is a fixed region in space, not changing in time. The edges in $E(B)$ are coupled at time s_1 since \mathcal{B}'_{s_1} holds. By Lemma 2.22 the walker never leaves $S_1(i) \subset B$ during the time interval $[s_1, s_2]$; if $s_1 \leq \bar{t}_1$, then we know that the component of the walker is at most $\log^2 \ell$ since \mathcal{B}_0 holds and the box $R_1(0, 0)$ is k_{\max} -great by the properties of the second phase. Since there is no non- \star update in $E(B)$ during $[s_1, s_2]$, B remains coupled with its translate throughout and identity coupling is successful. \square

The lemma below is a composition of the previous lemma when the walker traverses a sequence of good 1-boxes. We assume that the stronger event \mathcal{B}'_t holds at the start time to be able to guarantee that \mathcal{B}''_t holds during the entire time interval covered by the lemma.

Lemma 5.7 (Identity coupling succeeds in sequences of good boxes) *Let s_1 be a time so that \mathcal{B}'_{s_1} holds. Let $s_2 > s_1$ be such that during $[s_1, s_2]$ the walker only traverses 1-boxes that are good. Then, if we attempt to do identity coupling for the entire time interval $[s_1, s_2]$, the coupling succeeds, \mathcal{B}''_t holds and $\Psi_{s_1} \equiv \Psi_t$ for all $t \in [s_1, s_2]$. Moreover, \mathcal{B}'_t holds for all $t \in [s_1 + 2\bar{t}_1, s_2]$.*

Proof Let $R_1(i, \tau)$ be the box the walker is in its core at time s_1 . Since $R_1(i, \tau)$ is a good box, Lemma 5.6 gives that identity coupling works up to the end of $T_1(\tau)$ and Lemma 5.3 gives that $S_1(i)$ couples at some time during $[s_1, s_1 + 2\bar{t}_1]$. Moreover, for any $t \in [s_1, s_1 + 2\bar{t}_1]$, \mathcal{B}''_t holds since \mathcal{B}'_{s_1} holds. For $t \in T_1(\tau) \cap [s_1 + 2\bar{t}_1, \infty)$ we have that $\mathcal{B}'_t \subset \mathcal{B}''_t$ holds by Lemma 5.5. Hence, if $R_1(i', \tau')$ is the box whose core the walker is in when exiting $R_1(i, \tau)$, we can apply Lemma 5.6 again to show that identity coupling succeeds. Repeating this argument over and over again establishes the lemma. \square

Now we analyze what happens in the neighborhood around a bad 1-box, supposing that the walker enters the 2-enlargement of that box. Two things can happen, either the walker enters the 2-enlargement of the box from the space boundary $\partial_s R_1^{\text{enl}2}(\cdot, \cdot)$ or it enters from the time boundary $\partial_t^- R_1^{\text{enl}2}(\cdot, \cdot)$. If it is from the space boundary, then the walker does not get too close to the bad box and \mathcal{B}'_t would still be verified for all t . Moreover, as long as the walker is in the 2-enlargement, identity coupling can be applied successfully. In the other case, if the walker enters from the time boundary, then it could eventually reach the bad box but the environment in the 2-enlargement of the bad box will be coupled before that. In particular, the environments will be coupled at all times in $T_1^{\text{enl}1}(\cdot, \cdot)$ thanks to the abundance of \star -updates in $T_1^{\text{enl}2}(\cdot, \cdot) \setminus T_1^{\text{enl}1}(\cdot, \cdot)$. This reasoning gives that the relative distance between the walkers does not change and the graphs remain coupled in $E(S_1^{\text{enl}2}(i))$ when the walker cross a bad box of scale 1.

In the lemma below we will require m to be large enough so that the following holds:

For any (k, i, τ) and any (i', τ') so that $R_{k+1}^{\text{core}}(i', \tau')$ intersects $R_k(i, \tau)$

we obtain that

$$\begin{aligned} R_{k+1}(i', \tau') \text{ contains all } k\text{-boxes that intersects } R_k^{\text{enl2}}(i, \tau), \\ \text{and } S_{k+1}^{\text{inn}}(i') \text{ contains } S_k^{\text{enl2}}(i). \end{aligned} \quad (5.10)$$

Lemma 5.8 (*Identity coupling in enlargement of bad boxes*) *Let m be large enough so that (5.10) holds. Let $R_k(i, \tau)$ be a bad box of scale k such that $R_{k+1}(i', \tau')$ is good for some $(k+1)$ -box for which $R_{k+1}^{\text{core}}(i', \tau') \cap R_k(i, \tau) \neq \emptyset$. Denote with $\tau_k^+ = \max T_k^{\text{enl2}}(\tau)$, and with $\tau_k^- = \min T_k^{\text{enl2}}(\tau)$. Let s_c be a time at which the walker enters $R_k^{\text{enl2}}(i, \tau)$ so $X_{s_c} \in S_k^{\text{enl2}}(i)$ but $X_{s_c-} \notin S_k^{\text{enl2}}(i)$ or $s_c = \tau_k^-$. Let $s_e = \inf\{t \in (s_c, \tau_k^+]: X_t \notin S_k^{\text{enl2}}(i)\}$ the first time the walker exits $R_k^{\text{enl2}}(i, \tau)$ after s_c ; we take the convention that $s_e = \tau_k^+$ if $X_t \in S_k^{\text{enl2}}(i)$ for all $t \in [s_c, \tau_k^+]$. Thus*

$$\text{if } \mathcal{B}'_{s_c} \text{ holds, then for all } t \in [s_c, s_e] \Psi_t \text{ remains unchanged and } \mathcal{B}'_t \text{ holds;} \quad (5.11)$$

consequently, identity coupling succeeds during $[s_c, s_e]$. Moreover,

$$\text{if } s_c > \tau_k^-, \text{ the walker does not enter } R_k^{\text{enl1}}(i, \tau). \quad (5.12)$$

Ultimately, letting $\mathcal{J} = \{j: S_k(j) \cap S_k^{\text{enl2}}(i) \neq \emptyset\}$,

$$\begin{aligned} \text{if } s_c = \tau_k^-, \text{ then } \eta_t^*(e) = \bar{\eta}_t^*(\Phi_t(e)) \text{ for all } e \in E(S_k^{\text{inn}}(\mathcal{J})) \\ \text{and all } t \in [s_c, s_e] \text{ with } t \geq \min T_k^{\text{enl1}}(\tau). \end{aligned} \quad (5.13)$$

The proof uses induction on k , so we treat the case $k = 1$ separately.

Proof of Lemma 5.8 for $k = 1$ We start with the case $s_c > \tau_1^-$, meaning that the walker entered the 2-enlargement of the bad box from $\partial_s R_1^{\text{enl2}}(i, \tau)$. We need to establish (5.11) and (5.12) in this case. We establish (5.12) by showing that the walker never gets closer than 12ℓ from $S_1^{\text{enl1}}(i)$. To see this, from (5.10) we have that $R_2(i', \tau')$ contains the 2-enlargement of $R_1(i, \tau)$, and $R_2(i', \tau')$ is a good box. Moreover, Remark 2.17 gives that the 1-enlargement of $R_1(i, \tau)$ contains all bad 1-boxes inside $R_2(i', \tau')$, and Lemma 2.23 gives that the distance between the walker and $S_1^{\text{enl1}}(i)$ is at least 12ℓ , establishing (5.12). To establish (5.11), note that the walker only traverses good boxes during $[s_c, s_e]$, so (5.11) follows from 5.7.

Now we consider the case $s_c = \tau_1^-$, and need to establish (5.11) and (5.13). The idea in this case is to use the time interval between s_c and $\min T_1^{\text{enl1}}(\tau)$, which is large enough for the graphs to couple. In fact, applying Lemma 5.3 to the box $R_2(i', \tau')$ from time s_c , we obtain a time $s \in [s_c, s_c + 2\bar{t}_2]$ so that $S_2(i')$ is coupled. From this time onwards Lemma 5.5 gives that $S_2^{\text{inn}}(i') \supset S_1^{\text{enl2}}(i)$ remains coupled up to time s_e . From Lemma 2.22 we know that the walker does not leave $S_2^{\text{inn}}(i')$ during $[s_c, s_e]$. So if identity coupling succeeds up to time $s_c + 2\bar{t}_2$, then it succeeds up to time s_e . Moreover, note that $2\bar{t}_2 = 12t_2/m = 12t_1$ is smaller than the distance between s_c and $\min T_1^{\text{enl1}}(\tau)$, which is $15t_1$. So $S_2^{\text{inn}}(i')$ couples before the walker can enter $R_1^{\text{enl1}}(i, \tau)$ and (5.13) is established.

It remains to show that the coupling succeeds and \mathcal{B}_t'' holds for all $t \in [s_c, s_c + 2\bar{t}_2]$, completing the proof of (5.11). For this, we only need to note that during this time interval the walker only traverses good 1-boxes, so (5.11) follows from Lemma 5.7. \square

Proof of Lemma 5.8 for $k \geq 2$ We have already established the case $k = 1$. Now we proceed via induction. Assume all claims of the lemma are proved up to scale $k - 1$. Let $R_k(i, \tau)$ be a bad box and $R_{k+1}(i', \tau')$ as in the statement of the lemma be a good box. All bad boxes in $R_{k+1}(i', \tau')$ are contained in $R_k^{\text{enl}}(i, \tau)$.

We first prove the case $s_c > \tau_k^-$, which requires establishing (5.11) and (5.12). In this case we use the same argument as in the case $k = 1$; that is, (5.12) follows from Lemma 2.23. To show that identity coupling can be performed and \mathcal{B}_t' holds, notice that if at time s_c the walker is inside a bad box $R_{k''}(i'', \tau'')$ for some $k'' < k$, then since $R_1(0, 0)$ is k_{\max} -great, we have that in a previous time the walker was in the boundary of $R_{k''}^{\text{enl}2}(i'', \tau'')$. If there are more than one tuple (k'', i'', τ'') satisfying the property above, we take the one with the largest k'' (breaking ties arbitrarily if there still are more than one such tuples). Since the walker must have entered the 2-enlargement $R_{k''}^{\text{enl}2}(i'', \tau'')$ at some time s_c'' , we obtain by induction that while traversing the bad box $R_{k''}(i'', \tau'')$ identity coupling is successful and \mathcal{B}_t' holds up to the end of $T_{k''}(\tau'')$, since (5.13) implies \mathcal{B}_t' . Therefore, when the walker leaves $R_{k''}(i'', \tau'')$, we can apply the induction hypothesis again if the walker is inside another bad box. It remains to check that identity coupling can be performed while the walker passes through space-time locations that belong to good boxes at all scale, in particular, while the walker passes through good 1-boxes. But since \mathcal{B}_t' holds at that time, identity coupling succeeds by Lemma 5.7, concluding the proof of (5.11).

We now prove the case $s_c = \tau_k^-$, which requires establishing (5.11) and (5.13). Assume that $s_e \geq \min T_k^{\text{enl}}(\tau)$, otherwise (5.11) follows from the same argument above and (5.13) is irrelevant. We can do the same argument as for $k = 1$; i.e., we show that the time interval between s_c and $\min T_k^{\text{enl}}(\tau)$ is large enough for the graphs to couple. By Lemma 5.3 we obtain a time $s \in [s_c, s_c + 2\bar{t}_{k+1}]$ so that $S_{k+1}(i')$ is coupled and, by Lemma 5.5, $S_{k+1}^{\text{inn}}(i') \supset S_k^{\text{enl}2}(i)$ remains coupled until s_e . Since Lemma 2.22 gives that the walker does not leave $S_{k+1}^{\text{inn}}(i')$ during $[s_c, s_e]$, if identity coupling succeeds up to time $s_c + 2\bar{t}_{k+1}$, then it succeeds up to time s_e . Besides, $2\bar{t}_{k+1} = 12 \frac{t_{k+1}}{k^2 m} = 12t_k$ is smaller than the distance between s_c and $\min T_k^{\text{enl}}(\tau)$, which is $15t_k$. So $S_{k+1}^{\text{inn}}(i')$ couples before the walker can enter $R_k^{\text{enl}}(i, \tau)$ and (5.13) is established. To establish (5.11), we need to show that \mathcal{B}_t'' holds for all $t \in [s_c, s_c + 2\bar{t}_{k+1}]$, but during this time the walker only traverses good 1-boxes, so (5.11) follows from Lemma 5.7. \square

Remark 5.9 The 2-enlargement of a bad box is chosen so that whenever the walker crosses it, by doing identity coupling the two processes have time to couple the environment before the walker crosses the bad box. For this exact reason we want the first box whose core the walker is at, at the beginning of the second phase, to be k_{\max} -great, so we know that the walker does not start inside the enlargement of a bad box, meaning that if the walker encounters a bad box during the second phase, it must first traverse its enlargement.

5.5 Simple random walk moment

Now we handle the case when the walker traverses great boxes, during which we do not perform identity coupling but try a different coupling. This coupling will be based on what we call a *simple random walk moment* (SRWM), which is a given condition of the evolution of the environment that makes the walker performs a simple random walk step.

Definition 5.10 (Simple random walk moment) Let $R_1(i, \tau)$ be a great box such that $(X_{\tau t_1}, \tau t_1) \in R_1^{\text{core}}(i, \tau)$. We consider three consecutive intervals I_1, I_2, I_3 of lengths

$$|I_1| = \frac{t_1}{2} \quad \text{and} \quad |I_2| = |I_3| = \frac{1}{\mu},$$

such that I_1 begins at time $\tau t_1 = \min T_1^{\text{core}}(\tau)$; note that $\tau t_1 + \sum_{j=1}^3 |I_j| < (\tau+2)t_1 = \max T_1^{\text{core}}(\tau)$. Let $v \in S_1(i)$ be the position of the walker $X_{\tau t_1}$; note that since $R_1(i, \tau)$ is a good box then all edges adjacent to v at time τt_1 are closed. All the events below consider only \star -updates during $I_1 \cup I_2 \cup I_3$, ignoring all non- \star updates. Then, a *simple random walk moment* (SRWM) is said to occur in $R_1(i, \tau)$ if the following events happen consecutively:

- (E_1) During I_1 , one of the edges adjacent to v , say $e = (v, u)$, receives an update to become open, and the edges adjacent to u with status \star are sampled closed. Moreover, the other edges adjacent to v or u do not open during I_1 , and after e opens, e does not close for at least time $\frac{C_1 1}{\mu}$.
- (E_2) During I_2 , edge e closes and does not open, while the edges adjacent to e , that were closed, do not open; note that at the end of I_2 , the walker is in either u or v .
- (E_3) During I_3 , the edges adjacent to u or v do a \star -update, and the edges adjacent to the walker do not open.

See Fig. 4 for an illustrative realization of a simple random walk moment. Define

$$\mathbb{I}_{(i, \tau)}^{\text{SRWM}} \text{ to be the indicator for the event that SRWM occurs in } R_1(i, \tau). \quad (5.14)$$

Remark 5.11 Given v , the position of the walker at time τt_1 , the event SRWM depends only on the updates in $E(S_1(i))$ during the time interval $I_1 \cup I_2 \cup I_3$. In particular, it does not depend on the jumps of the walkers during $I_1 \cup I_2 \cup I_3$, and does not depend on non- \star updates that could occur during $I_1 \cup I_2 \cup I_3$.

Note that from τt_1 to time $\tau t_1 + |I_1 \cup I_2 \cup I_3|$ the walker essentially performed a simple random walk step since the edge e adjacent to v that is chosen to open during I_1 is a uniformly random edge. Now assume that the walker enters $R_1^{\text{core}}(i, \tau)$ with $R_1(i, \tau)$ being a k_{\max} -great box; i.e., $(X_{\tau t_1}, \tau t_1) \in R_1^{\text{core}}(i, \tau)$. We define the coupling we employ in this situation.

Definition 5.12 (Coupling on great boxes) At time τt_1 both walkers are trapped at some vertices $v = X_{\tau t_1} \in S_1^{\text{core}}(i)$ and $\bar{v} = \Phi_{\tau t_1}(v)$. Then we perform the following steps.

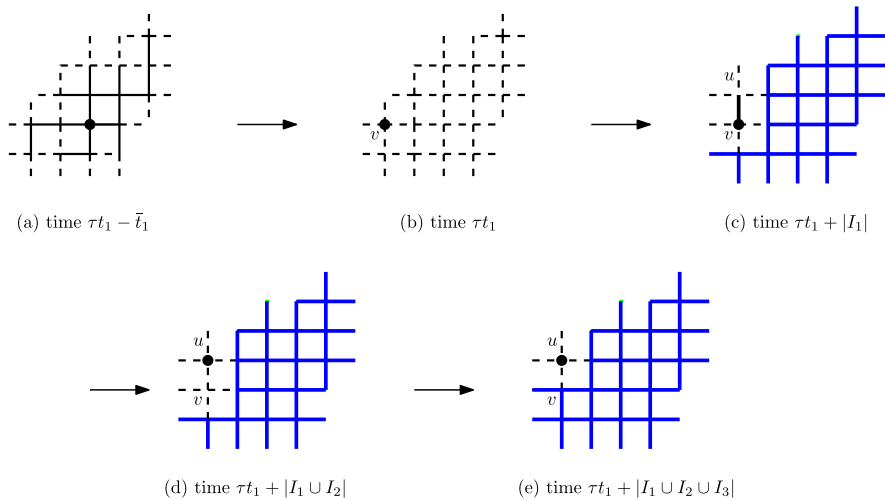


Fig. 4 A possible realization of SRWM in a k_{\max} -great box. **a** configuration at time $\tau t_1 - \bar{t}_1$, with dashed lines representing closed edges, solid lines representing open edges, and the black ball representing the walker. **b** During $[\tau t_1 - \bar{t}_1, \tau t_1]$ all edges close, trapping the walker in a vertex v . **c** During I_1 , edge (u, v) adjacent to the walker opens, edges adjacent to u are closed, and the other edges receive a \star -update. Blue lines represent edges that are updated \star . **d** Edge (u, v) closes at some time during I_2 , trapping the walker in one of its endpoints, in this case endpoint u . **e** During I_3 all edges adjacent to v receive a \star -update, while the edges adjacent to the walker do not open

1. Sample whether a simple random walk moment occurs in $R_1(i, \tau)$. If not, sample the updates of the edges in $S_1(i)$ during $I_1 \cup I_2 \cup I_3$ from the distribution conditioned on $\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 0$, apply the coupling of the graphs from Sect. 5.2.1 and apply identity coupling for the walkers. Identity coupling succeeds since the graphs are coupled inside $S_1(i)$ and we obtain that Φ_t does not change during $I_1 \cup I_2 \cup I_3$. This concludes the coupling when $\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 0$.
2. If $\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1$, choose a coordinate $j \in \{1, 2, \dots, d\}$ and a sign $s \in \{-1, +1\}$ uniformly at random. If v and \bar{v} agree in that coordinate, let $e = (v, v + se_j)$ and $\bar{e} = (\bar{v}, \bar{v} + se_j)$ be the edges chosen to open during I_1 in the configurations η^\star and $\bar{\eta}^\star$, respectively, where e_1, e_2, \dots, e_d stands for the standard basis of \mathbb{Z}^d . In this case, during I_2 , we let the walkers perform the same jumps across e and \bar{e} (i.e., we perform identity coupling), and note that Φ_t maps e into \bar{e} during this time. Then we couple the graphs using Φ_t , as described in Sect. 5.2.1, until the end of I_3 . In this case, the map Φ_t does not change during $I_1 \cup I_2 \cup I_3$.
3. If $\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1$, and v and \bar{v} do not agree in the j th coordinate, we set $e = (v, v + se_j)$ and $\bar{e} = (\bar{v}, \bar{v} - se_j)$. This is the most delicate case as we will need to change the coupling of the graphs from the time e opens to the end of $I_1 \cup I_2 \cup I_3$. For this, we will use the map $\tilde{\Phi}$ which maps v to \bar{v} and is a translation map in all coordinates but the j th one, where it is a reflection map around e . In particular, $\tilde{\Phi}$ maps e onto \bar{e} . Then the graphs will be coupled as in Remark 5.2; that is, the graphs are coupled as in Sect. 5.2.1 but using the map $\tilde{\Phi}$ instead of Φ_t . Note that any update to e translates to an update of \bar{e} , so they open at the same time and close at the same time. Let ζ be the time that e and \bar{e} open for the first time during I_1 . Then, they remain open

during $[\zeta, \zeta + C_{11}/\mu]$ since $\mathbb{I}_{(i,\tau)}^{\text{SRWM}} = 1$. We couple the position of the walkers at time $\zeta + C_{11}/\mu$ as follows. Let δ be the probability that $X_{\zeta+C_{11}/\mu} = v$ and $1 - \delta$ be the probability that $X_{\zeta+C_{11}/\mu} = u$. Then we make $X_{\zeta+C_{11}/\mu} = \bar{X}_{\zeta+C_{11}/\mu}$ with probability $\min\{\delta, 1 - \delta\}$; otherwise, we sample them accordingly. Then, we let the graph and the walkers evolve up to the end of the interval $I_1 \cup I_2 \cup I_3$, coupling the jumps of the walkers so that they jump at the same times after time $\zeta + C_{11}/\mu$; note that the walkers do not move after e and \bar{e} close for the first time after $\zeta + C_{11}/\mu$.

Now, let $s = \tau t_1 + |I_1 \cup I_2 \cup I_3|$ be the end time of the simple random walk moment. Note that if SRWM occurs then $\|X_s - \bar{X}_s\|_1$ may differ from $\|X_{\tau t_1}^1 - X_{\tau t_1}^2\|_1$, and as a result the translation map Φ_s may be different from $\Phi_{\tau t_1}$ as well. So it could be the case that an edge that was coupled before the simple random walk moment (in the sense that $\eta_{\tau t_1}(e') = \bar{\eta}_{\tau t_1}(\Phi_{\tau t_1}(e'))$) may get uncoupled because the map Φ changes. On the other hand, after I_1 all the edges in the box receive a \star update. So at the end of the SRWM, all edges in $S_1(i)$ are \star with the only exception being the edges adjacent to the walker which are closed. So the configurations are coupled locally, in particular, \mathcal{B}'_s holds. Moreover, as $R_1(i, \tau)$ is great (so it is also good) the particles will stay in $S_1(i)$ for the whole time interval $T_1(\tau)$. In other words we obtain that the edges in $S_1(i)$, where the random walk moment is occurring, are coupled after the simple random walk moment ends.

More formally, we will implement this by assigning a “hidden” random variable to each 1-box, which tells whether the box will undergo a SRWM should the walker pass there. We will not try the above coupling at each great box the walker enters, since we do need a bit of time separation between two simple random walk moments because of the overlapping of the boxes. But whenever we decide to attempt a simple random walk moment inside a great box the walker is in, the hidden random variable will tell whether SRWM occurs. The main point is that we can obtain a lower bound on $\mathbb{P}\left(\mathbb{I}_{(i,\tau)}^{\text{SRWM}} = 1\right)$ that is uniform on the location of the walker at time τt_1 . Because of this uniform bound, we can couple the outcome of the hidden variable with the evolution of the processes M_t^\star and \bar{M}_t^\star so that the simple random walk moment takes place, regardless of the location of the walker within the box. The content of the hidden variable is just a Bernoulli random variable of parameter $C_{12} p^{\frac{6d-1}{6d}}$, which is the bound we derive in Lemma 5.14 below, so the event of successfully performing a SRWM stochastically dominates the hidden variable. Whenever we decide to look at the hidden variable of a box, we perform the coupling described above. Otherwise, we just do identity coupling.

Before establishing a bound on $\mathbb{P}\left(\mathbb{I}_{(i,\tau)}^{\text{SRWM}} = 1\right)$ we need to show that the environments recouple locally after a SRWM.

Lemma 5.13 (Recoupling the graphs after SRWM) *Let $R_1(i, \tau)$ be a k_{\max} -great box such that $X_{\tau t_1} \in S_1^{\text{core}}(i)$ and $\mathcal{B}''_{\tau t_1}$ holds. Suppose the walkers perform successfully a simple random walk moment. Then*

$$\eta_t^\star(e) = \bar{\eta}_t^\star(\Phi_t(e)) \text{ for alle } e \in E(S_1^{\text{enl2}}(i)) \text{ and all } t \in [s, (\tau + 1)t_1],$$

where $s = \tau t_1 + |I_1 \cup I_2 \cup I_3| + \bar{t}_1$.

Proof $R_1(i, \tau)$ is k_{\max} -great, and in particular 1-great. Thus, every 1-box $R(j, \tau)$ such that $S_1(j) \cap S_1^{\text{enl}2}(i) \neq \emptyset$ is good. After $s - \bar{t}_1 = \tau t_1 + |I_1 \cup I_2 \cup I_3|$, we have that the edges in $S_1(i)$ are coupled and we start performing identity coupling of the walkers. The coupling is successful so Φ_t does not change from that moment onwards and all edges in $S_1(j)$ with $S_1(j) \cap S_1^{\text{enl}2}(i) \neq \emptyset$ receives a \star -update and couples. \square

Now we bound the probability of a SRWM. Recall from Definition 2.7 that the event that a box $R_1(i, \tau)$ is good is based on the events $\mathcal{G}_{12}(i, \tau)$ and $\widehat{\mathcal{G}}_{34}(i, \tau)$. Define \mathcal{J} to be the set of all tuples (i, τ) such that $R_1(i, \tau)$ is a box of the tessellation of the second phase. Let $\Sigma = \{0, 1\}^{2\mathcal{J}}$ be the set of all possible assignments of occurrence or non occurrence to the events $\mathcal{G}_{12}(i, \tau)$ and $\widehat{\mathcal{G}}_{34}(i, \tau)$. Then for each $\sigma \in \Sigma$ and each (i, τ) , the values $\sigma_{12}(i, \tau)$ and $\sigma_{34}(i, \tau)$ will be used to specify whether the events $\mathcal{G}_{12}(i, \tau)$ and $\widehat{\mathcal{G}}_{34}(i, \tau)$ occur, respectively. In this way, given $\sigma \in \Sigma$, we abuse notation and denote by σ the event that the realizations of $\mathcal{G}_{12}(i, \tau)$ and $\widehat{\mathcal{G}}_{34}(i, \tau)$ match the values of $\sigma_{12}(i, \tau)$ and $\sigma_{34}(i, \tau)$ for each $(i, \tau) \in \mathcal{J}$, and write $\mathbb{P}(\cdot \mid \sigma)$ for the corresponding conditional probability. More precisely,

$$\mathbb{P}(\cdot \mid \sigma) = \mathbb{P}\left(\cdot \mid \bigcap_{(i, \tau) \in \mathcal{J}} \{\sigma_{12}(i, \tau) = \mathbb{1}(\mathcal{G}_{12}(i, \tau))\} \cap \{\sigma_{34}(i, \tau) = \mathbb{1}(\widehat{\mathcal{G}}_{34}(i, \tau))\}\right).$$

Note that once we condition on some $\sigma \in \Sigma$, then which boxes of all scales are good or bad is a deterministic function of σ . Let \mathcal{F}_t be the σ -algebra generated by the trajectory of the walker X_s and the value of the map Ψ_s , $s \in [0, t]$, and all the updates of the graph up to time t . Let $\Sigma_{i, \tau} \subset \Sigma$ be the set of assignments σ for which $R_1(i, \tau)$ is a k_{\max} -great box.

Lemma 5.14 *Let (i, τ) be such that $R(i, \tau)$ is a k_{\max} -great box. There exists $p_0 > 0$ and $C_{12} > 0$ such that for all $p < p_0$, for all $\sigma \in \Sigma_{i, \tau}$, and all $F \in \mathcal{F}_{\tau t_1}$ for which $\mathbb{P}(\sigma \cap F) > 0$, then the probability of performing a simple random walk moment in $R_1(i, \tau)$ is*

$$\mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \sigma\right) \geq C_{12} p^{\frac{6d-1}{6d}}. \quad (5.15)$$

Proof Start with the following simplification of σ . Recall the definition of $j(\tau)$ from (2.9). So $j(\tau)$ and $j(\tau + 1)$ are the first and last interval of the type $\bar{T}_1(\cdot)$ inside $T_1(\tau)$. Recall that $\sigma_{34}(\cdot, \cdot)$ correspond to the events $\widehat{\mathcal{G}}_{34}(\cdot, \cdot)$, which are i.i.d. events coupled with the events $\mathcal{G}_{34}(\cdot, \cdot)$. Since $\mathcal{G}_{34}(\cdot, \cdot)$ are independent of $\mathcal{G}_{12}(\cdot, \cdot)$ by Lemma 2.4, we have that also $\widehat{\mathcal{G}}_{34}(\cdot, \cdot)$ are independent of $\mathcal{G}_{12}(\cdot, \cdot)$. Moreover, for any x , we have that $\mathcal{G}_{34}(i', \tau)$ is independent of $\mathcal{F}_{\tau t_1}$ since $\mathcal{G}_{34}(i', \tau)$ only considers updates on the edges during the interval $T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1)$. Since for any fixed x we have that $\mathcal{G}_{34}(i', \tau)$ are independent for different s , we have that $\widehat{\mathcal{G}}_{34}(i', \tau)$ is independent of $\mathcal{F}_{\tau t_1}$. So now we collect in \mathcal{J}_{34} all tuples from \mathcal{J} for which $\mathbb{I}_{(i, \tau)}^{\text{SRWM}}$ depend on $\sigma_{34}(\cdot, \cdot)$:

$$\mathcal{J}_{34} = \{(i', \tau) : S_1(i') \cap S_1(i) \neq \emptyset\} \quad \text{and} \quad \mathcal{J}' = \mathcal{J} \setminus \mathcal{J}_{34}.$$

We will not need to split $\sigma_{12}(\cdot, \cdot)$ into two groups since those events are already independent of $\mathbb{I}_{(i, \tau)}^{\text{SRWM}}$.

For any $\sigma \in \Sigma$ denote

$$\begin{aligned} \mathcal{S} &= \mathcal{S}(\sigma) = \bigcap_{(i', \tau') \in \mathcal{J}} \sigma_{12}(i', \tau') \bigcap_{(i', \tau') \in \mathcal{J}'} \sigma_{34}(i', \tau') \\ \mathcal{S}_{34} &= \mathcal{S}_{34}(\sigma) = \bigcap_{(i', \tau') \in \mathcal{J}_{34}} \sigma_{34}(i', \tau'). \end{aligned}$$

Then,

$$\begin{aligned} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \sigma\right) &= \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \mathcal{S} \cap \mathcal{S}_{34}\right) \\ &\geq \frac{\mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \mathcal{S}\right) - \mathbb{P}\left(\mathcal{S}_{34}^c \mid F \cap \mathcal{S}\right)}{\mathbb{P}\left(\mathcal{S}_{34} \mid F \cap \mathcal{S}\right)} \\ &\geq \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \mathcal{S}\right) - \mathbb{P}\left(\mathcal{S}_{34}^c \mid F \cap \mathcal{S}\right). \end{aligned}$$

Note that $T_1^{\text{core}}(\tau) \setminus T_1(\tau + 1) \supset I_1 \cup I_2 \cup I_3$, so $\mathbb{I}_{(i, \tau)}^{\text{SRWM}}$ does not depend on $F \cap \mathcal{S}$ given the position of the walker at time τt_1 . Letting $S'_1(i) = \bigcup_{u \in S_1^{\text{core}}(i)} B_{\log^2 \ell}^\infty(u)$, which are the places where the walker can be at time τt_1 , we write

$$\begin{aligned} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \mathcal{S}\right) &= \sum_{v \in S'_1(i)} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \mathcal{S} \cap \{X_{\tau t_1} = v\}\right) \\ &\quad \mathbb{P}\left(X_{\tau t_1} = v \mid F \cap \mathcal{S}\right) \\ &= \sum_{v \in S'_1(i)} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid X_{\tau t_1} = v\right) \mathbb{P}\left(X_{\tau t_1} = v \mid F \cap \mathcal{S}\right) \end{aligned}$$

We are left with the following lower bound on $\mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \sigma\right)$:

$$\begin{aligned} &\sum_{v \in S'_1(i)} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid X_{\tau t_1} = v\right) \mathbb{P}\left(X_{\tau t_1} = v \mid F \cap \mathcal{S}\right) - \mathbb{P}\left(\mathcal{S}_{34}^c \mid F \cap \mathcal{S}\right) \\ &\geq \inf_{v \in S'_1(i)} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid X_{\tau t_1} = v\right) - \mathbb{P}\left(\mathcal{S}_{34}^c \mid F \cap \mathcal{S}\right). \end{aligned} \quad (5.16)$$

We start with the first term in (5.16) to derive a lower bound on $\mathbb{P}\left(\mathbb{I}_{(i,\tau)}^{\text{SRWM}} = 1 \mid X_{\tau t_1} = v\right)$ that is uniform in v . Since $\mathbb{I}_{(i,\tau)}^{\text{SRWM}}$ is composed of the events E_1, E_2 and E_3 , which are independent of one another since they involve disjoint time intervals, we will derive a lower bound for each of them. For the event E_1 , we will require that an edge adjacent to v (call it e) opens during the first half of I_1 , so that e has time to remain open for time C_{11}/μ during I_1 . Recall that I_1 has length $t_1/2$, so its first half has length $t_1/4$, and the rate at which an edge opens due to a \star -update is $\mu p_{\star} \frac{p_{\min}}{p_{\star}} = \mu p_{\min}$, and the rate at which an edge close due to a \star -update is $1 - p_{\max}$. We obtain

$$\mathbb{P}\left(E_1 \mid X_{\tau t_1} = v\right) = \left(1 - e^{-2d\mu p_{\min} \frac{t_1}{4}}\right) \left(\frac{1 - p_{\max}}{p_{\star}}\right)^{2d-1} e^{-(4d-2)\mu p_{\min} \frac{t_1}{2}} e^{-\mu(1-p_{\max}) \frac{C_{11}}{\mu}}.$$

In the product above, the first term corresponds to an edge adjacent to the walker (call it e) opening during the first half of I_1 , the second term is the probability that all $2d - 1$ edges adjacent to e are closed at that time, the third term is the probability that none of the $4d - 2$ edges adjacent to e open until the end of I_1 , and the fourth term is the probability that e remains open for at least time C_{11}/μ . Recalling that $t_1 = \frac{\sqrt{\ell}}{\mu}$ and that $\ell = p^{-\frac{1}{3d}}$ we obtain

$$\mathbb{P}\left(E_1 \mid X_{\tau t_1} = v\right) = \left(1 - e^{-dp_{\min} \frac{\sqrt{\ell}}{2}}\right) \left(\frac{1 - p_{\max}}{p_{\star}}\right)^{2d-1} e^{-(2d-1)p_{\min}\sqrt{\ell}} e^{-C_{11}(1-p_{\max})}.$$

Using that $p_{\star} \leq 1$ in the second term, $p_{\min} \in \left[\frac{p}{1+q}, p\right]$ in the first and third terms, and $p_{\max} \geq 0$ in the fourth term, and then making p small enough so that $p_{\min} \leq p_{\max} \leq \frac{1}{2}$ and $e^{-d\frac{p}{1+q}\frac{\sqrt{\ell}}{2}} \leq 1 - \frac{dp\sqrt{\ell}}{4(1+q)}$ we obtain

$$\begin{aligned} \mathbb{P}\left(E_1 \mid X_{\tau t_1} = v\right) &\geq \left(1 - e^{-d\frac{p}{1+q}\frac{\sqrt{\ell}}{2}}\right) (1 - p_{\max})^{2d-1} e^{-(2d-1)p\sqrt{\ell}} e^{-C_{11}} \\ &\geq \frac{dp\sqrt{\ell}}{4(1+q)} 2^{-2d+1} e^{-(2d-1)p\sqrt{\ell}} e^{-C_{11}} \\ &= \frac{e^{-C_{11}}d}{2^{2d+1}(1+q)} p\sqrt{\ell} e^{-(2d-1)p\sqrt{\ell}}. \end{aligned}$$

Now note that $p\sqrt{\ell} = p^{1-\frac{1}{6d}}$ goes to 0 as $p \rightarrow 0$. Thus, we can take p small enough so that $p\sqrt{\ell} e^{-(2d-1)p\sqrt{\ell}} \geq \frac{p\sqrt{\ell}}{2}$ to obtain

$$\mathbb{P}\left(E_1 \mid X_{\tau t_1} = v\right) \geq \frac{e^{-C_{11}}d}{2^{2d+2}(1+q)} p\sqrt{\ell} = \frac{e^{-C_{11}}d}{2^{2d+2}(1+q)} p^{1-\frac{1}{6d}}. \quad (5.17)$$

The event E_1 is the main one governing the probability that SRWM occurs, since it involves the opening of an edge, which has small probability. For E_2 and E_3 we will just derive simple bounds that will not go to 0 as $p \rightarrow 0$. Recall that I_2 and I_3 are time intervals of length $1/\mu$, so

$$\mathbb{P}\left(E_2 \mid X_{\tau_{t_1}} = v\right) = \left(1 - e^{-\mu(1-p_{\max})\frac{1}{\mu}}\right) e^{-\mu p_{\min}\frac{1}{\mu}} e^{-(4d-2)\mu p_{\min}\frac{1}{\mu}},$$

where the first term is the probability that e has a \star -update to close, the second term is the probability that e does not get a \star -update to open, and the final term is the probability that all $4d-2$ edges adjacent to e do not receive a \star -update to open. Recall that p_{\min} and p_{\max} both go to 0 as $p \rightarrow 0$, so we obtain that

$$\mathbb{P}\left(E_2 \mid X_{\tau_{t_1}} = v\right) \geq 1 - \frac{1}{2e}. \quad (5.18)$$

Regarding E_3 , we obtain

$$\mathbb{P}\left(E_3 \mid X_{\tau_{t_1}} = v\right) = \left(1 - e^{-(2d-1)\mu p_{\star}\frac{1}{\mu}}\right) e^{-2d\mu p_{\min}\frac{1}{\mu}} \geq 1 - \frac{1}{2e^{2d-1}}, \quad (5.19)$$

where the inequality follows for all small enough p since $p_{\star} \rightarrow 1$ and $p_{\min} \rightarrow 0$ as $p \rightarrow 0$. Putting (5.17), (5.18) and (5.19) together we have a constant $c = c(d, q)$ so that for all small enough p we obtain

$$\mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid X_{\tau_{t_1}} = v\right) \geq cp^{1-\frac{1}{6d}}.$$

Plugging the bound above into (5.16), we obtain

$$\mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \sigma\right) \geq cp^{1-\frac{1}{6d}} - \mathbb{P}\left(\mathcal{S}_{34}^c \mid F \cap \mathcal{S}\right). \quad (5.20)$$

Now as we explained in the beginning of the proof, \mathcal{S}_{34} is independent of $\mathcal{F}_{\tau_{t_1}}$ and of \mathcal{S} . Moreover, \mathcal{S}_{34} is composed of an intersection of independent events $\widehat{\mathcal{G}}_{34}(\cdot, \tau)$ since $\sigma \in \Sigma_{i, \tau}$ so that $R_1(i, \tau)$ is k_{\max} -great. Therefore,

$$\begin{aligned} \mathbb{P}\left(\mathbb{I}_{(i, \tau)}^{\text{SRWM}} = 1 \mid F \cap \sigma\right) &\geq cp^{1-\frac{1}{6d}} - \mathbb{P}\left(\bigcup_{(i', \tau) \in \mathcal{J}_{34}} \widehat{\mathcal{G}}_{34}^c(i', \tau)\right) \\ &\geq cp^{1-\frac{1}{6d}} - \sum_{(i', \tau) \in \mathcal{J}_{34}} \mathbb{P}\left(\widehat{\mathcal{G}}_{34}^c(i', \tau)\right) \\ &\geq cp^{1-\frac{1}{6d}} - 5^d \exp\left(-C_6 \log^2 \ell\right), \end{aligned}$$

where the last inequality follows from Lemma 2.6. Since as $p \rightarrow 0$ the second term is much smaller than the first one, the lemma follows. \square

5.6 Concluding the second phase

Recall that for simplicity we are assuming that $(X_0, 0) = (0, 0)$, and recall the value of Δ_2 from (5.2). Denote with $I_d : V \rightarrow V$ the identity map, then we define

$$F_2 = \{(0, 0) \text{ isk}_{\max}\text{-great}\} \cap \{\Phi_{\Delta_2} = I_d\} \cap \mathcal{B}'_{\Delta_2}. \quad (5.21)$$

If F_2 is verified, the second phase is successful and the third phase can start, otherwise we let the two processes evolve independently until the end of *phase 3*, and only then restart the coupling from phase 1.

Lemma 5.15 *Assume F_1 is verified at time 0. For any $\delta > 0$ and for all p small enough, there exists $C_{10} = C_{10}(d, p, \delta) > 0$ in the definition of Δ_2 and $n_0 < \infty$ such that for all $n > n_0$*

$$\mathbb{P}(F_2) \geq 1 - \delta.$$

Proof Let \mathcal{P} be any feasible path and consider

$$\begin{aligned} \Upsilon_1^{\mathcal{P}} &= \inf \left\{ \tau > 0 : (\mathcal{P}(\tau t_1), \tau t_1) \in R_1^{\text{core}}(i, \tau) \text{ where } R_1(i, \tau) \text{ is } \text{ak}_{\max}\text{-great box} \right\}, \\ \Upsilon_j^{\mathcal{P}} &= \inf \left\{ \tau > \Upsilon_{j-1}^{\mathcal{P}} : (\mathcal{P}(\tau t_1), \tau t_1) \in R_1^{\text{core}}(i, \tau) \text{ where } R_1(i, \tau) \text{ is } \text{ak}_{\max}\text{-great box} \right\}, \end{aligned}$$

for $j \geq 2$. For any feasible path \mathcal{P} we let $\kappa_{\mathcal{P}}$ be the largest value such that $\Upsilon_{\kappa_{\mathcal{P}}}^{\mathcal{P}} \leq \frac{\Delta_2}{t_1} - 2$. Recall that Σ represents the set of all possible realizations of occurrences and non occurrences for the events $\mathcal{G}_{12}(\cdot, \cdot)$ and $\widehat{\mathcal{G}}_{34}(\cdot, \cdot)$, so the good and bad boxes at all scales are deterministic functions of σ . Let $\mathfrak{F}(\sigma)$ be the set of all feasible paths for a given σ . Given the uniform bound from Lemma 5.14, we let Y_1, Y_2, \dots be a sequence of i.i.d. Bernoulli random variables of parameter $C_{12}p^{\frac{6d-1}{6d}}$ where Y_j gives whether the j th SRWM will succeed when we try to perform it during the coupling. Let

$$\zeta = \frac{C_8 C_{10}}{4\sqrt{\ell}} n^2 = \frac{C_8}{4t_1} \Delta_2, \quad (5.22)$$

where C_8 is from Lemma 2.25 and C_{10} from the definition of Δ_2 in (5.2). Define the following events

$$\begin{aligned} E_1 &= \left\{ \sum_{j=1}^{\zeta} Y_j \geq c_0 n^2 \right\}, \text{ with } c_0 \text{ to be chosen later, and} \\ E_2 &= \{\kappa_{\mathcal{P}} \geq \zeta \text{ for all feasible paths } \mathcal{P} \in \mathfrak{F}(\sigma)\}. \end{aligned}$$

In this stage we want to couple the position of the walkers. From Lemma 5.8, by doing identity coupling whenever the walkers are not in a great box, their relative distance does not change. Their relative distance changes only when they are in a great box and a simple random walk moment is successfully performed. Let $E_{\text{coup}} = \{\Phi_{\Delta_2} = I_d\} \cap \mathcal{B}'_{\Delta_2}$. Hence

$$\mathbb{P}(F_2^c) \leq \mathbb{P}((0, 0) \text{ is not } k_{\max}\text{-great}) + \mathbb{P}(E_1 \cap E_2 \cap E_{\text{coup}}^c) + \mathbb{P}(E_1^c) + \mathbb{P}(E_2^c).$$

We start by bounding the first term. Notice that $\{(0, 0) \text{ is not } k_{\max}\text{-great}\}$ does not depend on the configuration at time 0. Moreover, at time 0, the walkers are stuck in a vertex, so X_s has to leave $R_1(0, 0)$ from the time boundary if $R_1(0, 0)$ is a good box. Using Lemmas 2.8 and 2.15 to bound ρ_j , we obtain

$$\mathbb{P}((0, 0) \text{ is not } k_{\max}\text{-great}) \leq c_d \sum_{j=1}^{k_{\max}} \rho_j \leq c_d \left(\rho_1 + \sum_{j=2}^{k_{\max}} \rho_1^{2^{j-2}} \right) \leq 3c_d \rho_1 \leq \frac{\delta}{4},$$

where c_d is a constant that counts the number of boxes whose 2-enlargement intersects $R_1(0, 0)$, and the last inequality follows for all p small enough. Next we bound

$$\mathbb{P}(E_1 \cap E_2 \cap E_{\text{coup}}^c) \leq \frac{\delta}{4}.$$

Under $E_1 \cap E_2$, we know we performed at least $c_0 n^2$ simple random walk moments. So, $\mathbb{P}(E_1 \cap E_{\text{coup}}^c)$ can be bounded by the probability that two random walkers performing SRW on \mathbb{T}_n^d are not coupled after $c_0 n^2$ steps. Taking $c_0 = c_0(d, \delta)$ large enough we obtain that they have coupled with probability at least $1 - \frac{\delta}{4}$.

Next, we bound $\mathbb{P}(E_2^c)$. From Lemma 5.1, with probability at least $1 - \rho_1^{2^{k_{\max}-3}}$, all k_{\max} -boxes in the tessellation are good. Thus, Lemma 2.25 gives that while traversing the first good k_{\max} -great box any feasible paths will traverse at least

$$C_8 \frac{t_{k_{\max}}}{t_1}$$

k_{\max} -great 1-boxes. After the feasible path exits the first k_{\max} -great box, it enters into another one and we obtain again another set of k_{\max} -great 1-boxes. The total number of steps we can iterate this procedure up to reaching time Δ_2 is $\frac{\Delta_2}{2t_{k_{\max}}} - 1 \geq \frac{\Delta_2}{4t_{k_{\max}}}$. Therefore, any feasible path must traverse at least

$$C_8 \frac{t_{k_{\max}}}{t_1} \frac{\Delta_2}{4t_{k_{\max}}} = C_8 \frac{\Delta_2}{4t_1} = \zeta$$

k_{\max} -great boxes. Hence,

$$\mathbb{P}(E_2^c) \leq \rho_1^{2^{k_{\max}-3}} \leq \frac{\delta}{4},$$

by simply having n large enough.

Finally we bound $\mathbb{P}(E_1^c)$. This is a simple Chernoff bound for the sum of independent Bernoulli random variables, where $\mathbb{P}(Y_j = 1) = C_{12} p^{\frac{6d-1}{6d}}$. Since

$$\begin{aligned}\mathbb{E}\left(\sum_{j=1}^{\zeta} Y_j\right) &= C_{12} p^{\frac{6d-1}{6d}} \zeta = \frac{C_8 C_{10} C_{12}}{4} \frac{p^{\frac{6d-1}{6d}}}{\sqrt{\ell}} n^2 = \frac{C_8 C_{10} C_{12}}{4} p^{\frac{1}{6d} + \frac{6d-1}{6d}} n^2 \\ &= \frac{C_8 C_{10} C_{12}}{4} p n^2.\end{aligned}$$

Now we take C_{10} large enough so that the above is larger than $2c_0 n^2$, which gives a constant c so that

$$\mathbb{P}(E_1^c) \leq \exp\left(-c \mathbb{E}\left(\sum_{j=1}^{\zeta} Y_j\right)\right) \leq \exp(-c 2c_0 n^2) \leq \frac{\delta}{4},$$

where the last inequality follows by taking n large. \square

To conclude the second phase, once the walkers are coupled after one SRWM, we just perform identity coupling up to time Δ_2 . If t is a time where a SRWM ended, notice that Lemma 5.13 gives that \mathcal{B}'_t holds. So we succeed performing identity coupling up to Δ_2 by Lemmas 5.6, 5.7 and 5.8.

6 Third phase

The third phase starts at time Δ_2 , at which time the walkers are coupled and \mathcal{B}'_{Δ_2} holds. During the third phase we let \overline{M}_t^\star mimic the evolution of M_t^\star by doing identity coupling on both the motion of the walkers and the updates of the edges. We now check whether the processes are fully coupled by time $\Delta_3 = \Delta_2 + \frac{n^2}{\mu}$.

Define

$$F_3 = \left\{ X_{\Delta_3} = \overline{X}_{\Delta_3} \text{ and } \eta_{\Delta_3}^\star(e) = \overline{\eta}_{\Delta_3}^\star(e) \forall e \in E(\mathbb{T})_n^d \right\}. \quad (6.1)$$

If F_3 is not verified, we restart the coupling at time Δ_3 from phase 1.

Lemma 6.1 *For any $\delta > 0$, if p is small enough and n large enough, we obtain*

$$\mathbb{P}(F_3) \geq 1 - \delta. \quad (6.2)$$

Proof Recall that boxes contained in $[0, \Delta_3]$ have been sampled as good or bad during the second phase. By Lemmas 5.6, 5.7 and 5.8, identity coupling is successful provided we cannot enter a bad box without first entering its 2-enlargement. Therefore, for the walkers to get uncoupled during $[\Delta_2, \Delta_3]$, it must so happen that the walkers entered a bad box of some scale k whose 2-enlargement intersects $[0, \Delta_2]$ and which was not observed during the second phase because it is not contained in $[0, \Delta_3]$. We now count the number of such boxes.

We start by deriving bounds on ℓ_k and t_k , the size of the boxes of scale k , for which the above can happen. When $k \geq k_{\max}$, we can choose n large enough so that for any m, ℓ fixed

$$2\ell k^{2k} \leq \ell_k = m^k (k!)^2 \ell \leq \ell k^{3k},$$

$$2\sqrt{\ell} k^{2k} \leq \mu t_k = m^k (k!)^2 \sqrt{\ell} \leq \sqrt{\ell} k^{3k}.$$

Recall that $k_{\max} = \log_2 \log n$ from (5.1). Then, $\mu t_{k_{\max}} \leq \sqrt{\ell} k_{\max}^{3 \log_2 \log n}$ is much smaller than a polynomial in n . Therefore, any box whose enlargement intersects $[0, \Delta_2]$ and is not contained in $[0, \Delta_3]$ must be of scale larger than k_{\max} . So

$$\mathbb{P}(F_3) \geq 1 - \mathbb{P}\left(\exists k > k_{\max} : R_k(i, \tau) \text{ is bad and } \Delta_3 \in T_k^{\text{enl}2}(\tau)\right).$$

Next, using the bounds we derived above for ℓ_k and t_k , the number ζ_k of boxes of scale k that intersect $\mathbb{T}_n^d \times \Delta_3$ is bounded above and below by

$$\zeta_k \geq \left(\frac{n}{3\ell_k}\right)^d 24 \geq \frac{24n^d}{3^d \ell^d k^{3dk}},$$

$$\zeta_k \leq 1 + 24 \left(\frac{n}{\ell_k}\right)^d \leq 1 + \frac{24n^d}{\ell^d k^{2dk}}.$$

In the upper bound of ζ_k we add a 1 to the fraction to consider the case when k is so large that we cannot find a box all contained in the tessellation. Using Lemma 2.15 the probability that there exists a box of scale k_{\max} or bigger that is bad is bounded above by

$$\sum_{k \geq k_{\max}} \zeta_k \rho_k \leq \sum_{k \geq k_{\max}} \zeta_k \rho_1^{2^{k-2}},$$

moreover using the inequalities above for ζ_k it is easy to see that, for any $k \geq k_{\max}$,

$$\sum_{k \geq k_{\max}} \zeta_k \rho_k \leq 2\zeta_{k_{\max}} \rho_1^{2^{k_{\max}-2}}.$$

Since $2^{k_{\max}} = \log n$, by taking p small enough we make ρ_1 small enough, which gives that

$$\mathbb{P}(F_3) \geq 1 - \delta.$$

□

7 Completing the proof of Theorem 1.1

Proof of Theorem 1.1 Let $\{M_t^*\}_{t \geq 0}$ and $\{\overline{M}_t^*\}_{t \geq 0}$ denote two copies of the process, each starting from an arbitrary configuration in $\mathbb{T}_n^d \times \{0, 1\}^{E(\mathbb{T}_n^d)}$. Recall the events F_1 , F_2 and F_3 from (4.3), (5.21) and (6.1). If the three events hold then $X_{\Delta_3} = \overline{X}_{\Delta_3}$ and $\eta_{\Delta_3}^* \equiv \overline{\eta}_{\Delta_3}$. So from Δ_3 onwards we can keep the processes coupled. We can now set

$\delta = \frac{1}{12}$ so that $F_1 \cap F_2 \cap F_3$ all hold with probability at least $\frac{3}{4}$. If any of the above fails, we just let the processes evolve independently up to time Δ_3 and restart from scratch. Since Δ_3 is of order n^2/μ from (5.2), we obtain that the mixing time is of order n^2 concluding the proof. \square

8 Proof of the lower bound (Theorems 1.2 and 1.3)

The proof of the lower bounds are identical to the ones in [11]. We add them here for completion.

Proof of Theorem 1.2 First we introduce a discrete time Markov chain $\tilde{M}_k = (\tilde{X}_k, \tilde{\eta}_k)$ which is defined by sampling the continuous time chain $M_t = (X_t, \eta_t)$ on intervals of length δ ; that is,

$$\tilde{X}_k = X_{k\delta} \quad \text{and} \quad \tilde{\eta}_k = \eta_{k\delta},$$

where δ is given from (1.4). Let $\tilde{\gamma} = \tilde{\gamma}(\tilde{M})$ and $\gamma = \gamma(M)$ be the spectral gaps of the discrete time and continuous time chain, respectively. We obtain

$$1 - \tilde{\gamma} = \exp(-\delta\gamma).$$

For all $\tilde{\gamma} \leq 1/2$ we simply use the bound $\gamma \leq \frac{2\tilde{\gamma}}{\delta}$. The lower bound on the relaxation time follows by taking the function $f(x, \xi) = d(x, 0)$, so $f(X_t, \eta_t)$ is the distance between the walker and the origin of \mathbb{T}_n^d . Since the stationary distribution of the walker is uniform by (1.3), it follows that $\text{Var}(f) \geq cn^2$ for some constant $c > 0$. Moreover, from (1.4), we have

$$\begin{aligned} \mathcal{E}(f, f) &= \frac{1}{2} \sum_{x, \xi} \pi(x) \nu(\xi) \sum_{x', \xi'} P((x, \xi), (x', \xi')) (d(x, 0) - d(x', 0))^2 \\ &\leq \frac{1}{2} \mathbb{E}_{x \sim \pi} (D_{x, \delta}^2) \leq \frac{1}{2} C_2, \end{aligned}$$

where $\mathbb{E}_{x \sim \pi}$ denotes the expectation where x is a random variable sampled according to π , the uniform measure on \mathbb{T}_n^d . From the above we obtain

$$\tilde{\gamma} \leq \frac{C_2}{2cn^2}.$$

If the above is at most $1/2$ we obtain

$$\gamma \leq \frac{C_2}{cn^2\delta}.$$

Otherwise, if $\tilde{\gamma} \geq 1/2$ we obtain that γ is of order $1/\delta$. The above establishes the relaxation time of the chain. \square

Proof of Theorem 1.3 We use the following nice result from [9], which appeared implicitly already in [3].

Lemma 8.1 *Let $\{Y_k\}_{k \in \mathbb{Z}}$ be a discrete-time, stationary, reversible Markov chain with finite state space \mathcal{S} , and let $h: \mathcal{S} \rightarrow \mathbb{R}^m$ for some $m \in \mathbb{Z}_+$. Then, for each $k \geq 0$*

$$\mathbb{E} \left(\|h(Y_k) - h(Y_0)\|_{L_2}^2 \right) \leq k \mathbb{E} \left(\|h(Y_1) - h(Y_0)\|_{L_2}^2 \right),$$

where $\|\cdot\|_{L_2}$ denotes the Euclidean norm on \mathbb{R}^m .

Letting $g_n: \mathbb{T}_n^d \rightarrow \mathbb{R}^{2d}$ the function

$$g_n(x_1, x_2, \dots, x_d) = (n \cos(2\pi x_1/n), n \sin(2\pi x_1/n), \dots, n \cos(2\pi x_d/n), n \sin(2\pi x_d/n)).$$

For $x \in \mathbb{T}_n^d$ and $\xi \in \{0, 1\}^{E(\mathbb{T}_n^d)}$ we let $h(x, \xi) = g_n(x)$. Then, noting that g_n is bi-Lipschitz with some constant c we have

$$\begin{aligned} \mathbb{E}_{\pi \times \nu} \left(\|\tilde{X}_k - \tilde{X}_0\|_1^2 \right) &\leq c^2 \mathbb{E}_{\pi \times \nu} \left(\left(g_n(\tilde{X}_k) - g_n(\tilde{X}_0) \right)^2 \right) \\ &\leq c^2 k \mathbb{E}_{\pi \times \nu} \left(\left(g_n(\tilde{X}_1) - g_n(\tilde{X}_0) \right)^2 \right) \\ &\leq c^4 k \mathbb{E}_{\pi \times \nu} \left(\|\tilde{X}_1 - \tilde{X}_0\|_2^2 \right) \leq c^4 k \mathbb{E}_{\pi \times \nu} \left(D_{\tilde{X}_0, \delta}^2 \right) \leq c^4 C_2 k. \end{aligned}$$

Hence for any $t \geq \delta$ we have

$$\mathbb{E}_{\pi \times \nu} \left(\|X_t - X_0\|_1^2 \right) \leq c^4 C_2 \lceil \frac{t}{\delta} \rceil \leq 2c^4 C_2 \frac{t}{\delta}.$$

Now for the total variation starting from a stationary environment, we simply make

$$\begin{aligned} \|v_t - \pi \times \nu\|_{\text{TV}} &\geq \mathbb{P} \left(\|X_t - X_0\|_1 \leq \epsilon^{1/d} n \right) \left(1 - \frac{2\epsilon}{3} \right) \\ &= \left(1 - \mathbb{P} \left(\|X_t - X_0\|_1 \leq \epsilon^{1/d} n \right) \right) \left(1 - \frac{2\epsilon}{3} \right) \\ &\geq \left(1 - \frac{\mathbb{E}(\|X_t - X_0\|_1^2)}{\epsilon^{2/d} n^2} \right) \left(1 - \frac{2\epsilon}{3} \right) \\ &\leq \left(1 - \frac{2c^4 C_2 t}{\delta \epsilon^{2/d} n^2} \right) \left(1 - \frac{2\epsilon}{3} \right). \end{aligned}$$

Therefore, if $t \leq \frac{\epsilon^{2/d} \delta n^2}{6c^4 C_2}$ we have that $\|v_t - \pi \times \nu\|_{\text{TV}} \geq 1 - \epsilon$. \square

9 Proof of Corollary 1.4

In order to apply the above to the random walk on dynamical random cluster model, we first need a certain sprinkling lemma for the random cluster model. Given $q \geq 1$ and $p > 0$, let $\nu_{p,q}$ be the measure of a random cluster model with parameters p, q . Let η be a configuration sampled from $\nu_{p,q}$. We construct a sprinkling by associating to each edge e an independent Bernoulli random variable $Z(e)$ of parameter ϵ . Define the configurations

$$(\eta + Z)(e) = \mathbb{1}(\eta(e) + Z(e) \geq 1) \text{ for } e \in E(\mathbb{T}_n^d)$$

and

$$(\eta - Z)(e) = \mathbb{1}(\eta(e)(1 - Z(e)) = 1) \text{ for } e \in E(\mathbb{T}_n^d).$$

So $\eta + Z$ (resp., $\eta - Z$) is the configuration obtained from η by opening (resp., closing) all edges e with $Z(e) = 1$. Given two elements ξ, ξ' of $\{0, 1\}^{E(\mathbb{T}_n^d)}$ we say that $\xi \leq \xi'$ if $\xi(e) \leq \xi'(e)$ for all $e \in E(\mathbb{T}_n^d)$.

Lemma 9.1 (Sprinkling lemma) *Let $q \geq 1$, $0 < p < p' < 1$ and $\epsilon > 0$ be fixed. Let $\{Z(e) : e \in E(\mathbb{T}_n^d)\}$ be a collection of i.i.d. Bernoulli random variables of parameter ϵ . Let η and η' be random configurations with distributions $\nu_{p,q}$ and $\nu_{p',q}$, respectively. If*

$$\epsilon + (1 - \epsilon)p \leq p' \quad \text{and} \quad \epsilon + (1 - \epsilon)\frac{p}{p + (1 - p)q} \leq \frac{p'}{p' + (1 - p')q} \quad (9.1)$$

then there exists a coupling between ν, ν', Z such that $(\eta + Z) \leq \eta'$. Similarly, if

$$(1 - \epsilon)p' \geq p \quad \text{and} \quad (1 - \epsilon)\frac{p'}{p' + (1 - p')q} \geq \frac{p}{p + (1 - p)q} \quad (9.2)$$

then there exists a coupling between ν, ν', Z such that $(\eta' - Z) \geq \eta$.

Proof Let $\{\eta_t\}_t$ and $\{\eta'_t\}_t$ be the single-site Glauber dynamics Markov chains on the random cluster model with parameters (p, q) and (p', q) , respectively. Let $\{Z_t\}_t$ be a Glauber dynamics Markov chain on the state space $\{0, 1\}^{E(\mathbb{T}_n^d)}$ with stationary distribution given by a product of Bernoulli measures with parameter ϵ . Start with arbitrary configurations such that $\eta_0 \equiv \eta'_0$ and $Z_0(e) = 0$ for all $e \in E(\mathbb{T}_n^d)$. Assume that $\eta_t + Z_t \leq \eta'_t$ at some time t . We will show that we can couple the next transition of the chains so that $\eta_{t+1} + Z_{t+1} \leq \eta'_{t+1}$. This establishes the lemma. For any edge e and configuration η , let

$$\alpha(e, \eta) = \mathbb{1}(e \text{ is a cut-edge in } \eta).$$

In the coupling we will choose the same edge to be updated in all chains. Let e be such an edge. Then, note that

$$\begin{aligned} \mathbb{P}((\eta_{t+1} + Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) \\ = \epsilon + (1 - \epsilon) \left(\alpha(e, \eta_t) \frac{p}{p + (1 - p)q} + (1 - \alpha(e, \eta_t)) p \right). \end{aligned}$$

Since $\eta_t \leq \eta'_t$ we have that $\alpha(e, \eta_t) \geq \alpha(e, \eta'_t)$. So if $\alpha(e, \eta_t) = 0$ we have that $\alpha(e, \eta'_t) = 0$, which gives

$$\begin{aligned} \mathbb{P}((\eta_{t+1} + Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) &= \epsilon + (1 - \epsilon)p \leq p' \\ &= \mathbb{P}(\eta'_{t+1}(e) = 1 \mid \eta_t, \eta'_t, Z_t), \end{aligned}$$

where in the first inequality we used (9.1). If $\alpha(e, \eta_t) = 1$, then we use the second part of (9.1) and that $\frac{p'}{p' + (1 - p')q} \leq p'$ to write

$$\begin{aligned} \mathbb{P}((\eta_{t+1} + Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) &= \epsilon + (1 - \epsilon) \frac{p}{p + (1 - p)q} \\ &\leq \frac{p'}{p' + (1 - p')q} \\ &\leq \alpha(e, \eta'_t) \frac{p'}{p' + (1 - p')q} + (1 - \alpha(e, \eta'_t)) p' \\ &= \mathbb{P}(\eta'_{t+1}(e) = 1 \mid \eta_t, \eta'_t, Z_t). \end{aligned}$$

Therefore, it follows that we can couple the next transition of the Markov chains so that $\eta_{t+1} \leq (\eta_{t+1} + Z_{t+1}) \leq \eta'_{t+1}$. Consequently, we can couple the stationary measures of such chains to obtain that $(\eta + Z) \leq \eta'$.

For the second part of the lemma, we use the same strategy and analyze the transition probabilities for $\eta'_t - Z_t$. We have

$$\begin{aligned} \mathbb{P}((\eta'_{t+1} - Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) \\ = (1 - \epsilon) \left(\alpha(e, \eta'_t) \frac{p'}{p' + (1 - p')q} + (1 - \alpha(e, \eta'_t)) p' \right). \end{aligned}$$

If $\alpha(e, \eta'_t) = 1$ then $\alpha(e, \eta_t) = 1$, yielding

$$\begin{aligned} \mathbb{P}((\eta'_{t+1} - Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) &= (1 - \epsilon) \frac{p'}{p' + (1 - p')q} \geq \frac{p}{p + (1 - p)q} \\ &= \mathbb{P}(\eta_{t+1} = 1 \mid \eta_t, \eta'_t, Z_t). \end{aligned}$$

If $\alpha(e, \eta'_t) = 0$ then

$$\begin{aligned} \mathbb{P}((\eta'_{t+1} - Z_{t+1})(e) = 1 \mid \eta_t, \eta'_t, Z_t) &= (1 - \epsilon)p' \geq p \\ &\geq \alpha(e, \eta_t) \frac{p}{p + (1 - p)q} + (1 - \alpha(e, \eta_t)) p \\ &= \mathbb{P}(\eta_{t+1} = 1 \mid \eta_t, \eta'_t, Z_t). \end{aligned}$$

Therefore, there exists a coupling such that $\eta_{t+1} \leq (\eta'_{t+1} - Z_{t+1})$ and we obtain $(\eta' - Z) \geq \eta$. \square

Proof of Corollary 1.4 We only need to check that assumptions (1.3) and (1.4) hold for the random walk on dynamical random cluster model. For any q, p we have that (1.3) holds. For $q \geq 1$, (1.4) holds for all $p < p_c^q$ using the following argument. Take $p' = \frac{p+p_c^q}{2} \in (p, p_c^q)$. Take $\epsilon > 0$ small enough so that (9.1) is satisfied. We choose $\delta = \epsilon/\mu$ and take η to be a random cluster configuration of parameters p, q . Note that the probability that a given edge gets refreshed during $[0, \delta]$ is

$$1 - e^{-\mu\delta} = 1 - e^{-\epsilon} \leq \epsilon.$$

Therefore, if $Z(e)$ is a Bernoulli random variable of parameter ϵ , we can couple $Z(e)$ with the refresh clocks of the dynamical random cluster so that if e gets refreshed during $[0, \delta]$ then $Z(e) = 1$. Therefore, this coupling gives that $C_x([0, \delta])$ is contained in the cluster of x inside the configuration $\eta + Z$, which by Lemma 9.1 is contained inside η' , a random cluster configuration with parameters p', q . Then it follows by the sharpness of the phase transition [5] that the cluster of x in η' has an exponential decay, establishing (1.4) and allowing us to obtain the conclusions of Theorems 1.2 and 1.3 for the random cluster model with $q \geq 1$.

Regarding the case $q < 1$, one can deduce the exponential decay of the cluster $\eta + Z$ only when p is small enough. This becomes rather trivial as regardless of the state of the other edges, we obtain that an edge e is open during $[0, \delta]$ with probability at most

$$\max \left\{ \frac{p}{p + (1-p)q}, p \right\} + \epsilon = \frac{p}{p + (1-p)q} + \epsilon$$

which for small enough p can be made smaller than p_c , the critical probability for independent percolation. \square

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Declarations

Competing interests The authors declare no competing interests.

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