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Number variance for hierarchical random walks and related fluctuations*

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Abstract

We study an infinite system of independent symmetric random walks on a hierarchical group, in particular, the c -random walks. Such walks are used, e.g., in mathematical physics and population biology. The number variance problem consists in investigating if the variance of the number of “particles” $N_n(L)$ lying in the ball of radius L at a given step n remains bounded, or even better, converges to a finite limit, as $L \rightarrow \infty$. We give a necessary and sufficient condition and discuss its relationship to transience/recurrence property of the walk. Next we consider normalized fluctuations of $N_n(L)$ around the mean as $n \rightarrow \infty$ and L is increased in an appropriate way. We prove convergence of finite dimensional distributions to a Gaussian process whose properties are discussed. As the c -random walks mimic symmetric stable processes on \mathbb{R} , we compare our results with those obtained by Hambly and Jones (2007, 2009), who studied the number variance problem for an infinite system of independent symmetric stable processes on \mathbb{R} . Since the hierarchical group is an ultrametric space, corresponding results for symmetric stable processes and hierarchical random walks may be analogous or quite different, as has been observed in other contexts. An example of a difference in the present context is that for the stable processes a fluctuation limit process is a Gaussian process which is not Markovian and has long range dependent stationary increments, but the counterpart for hierarchical random walks is Markovian, and in a special case it has independent increments.

Key words: hierarchical random walk, hierarchical group, ultrametric, number variance, fluctuation, limit theorem.

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1 Introduction

Stochastic models on the d -dimensional Euclidean lattice \mathbb{Z}^d play a fundamental role in mathematical physics and population biology. Many of the basic questions of interest in these fields lead to mathematical challenges, many of which remain open. The class of hierarchical models first introduced by Dyson [16] in 1969 have played an important role for gaining insight into these problems. One reason for this is that the models can be parametrized in the hierarchical lattice in such a way as to provide a “caricature” for the Euclidean lattice of dimension d of interest (e.g. $d = 4$ and dimensions “infinitesimally close” to 4 in the study of ferromagnetic models, see e.g. [5]). The other reason is that the structure of these models is such that it has been possible to carry out the renormalization group analysis in a rigorous way in the hierarchical lattice. For example, this has been achieved in [5] for the study of ferromagnetism. It has also been used in the study of self-avoiding random walks (e.g. [4]), and Anderson localization in disordered media (e.g. [2, 28]). In population models it has been used to study mutually catalytic branching (e.g. [6, 15]), and interacting diffusions (e.g. [13, 14]). Random walks on hierarchical lattices, called *hierarchical random walks*, play a key role in the analysis of these systems. For example, they arise in the analysis of the moment structures, and the associated hierarchical Laplacian plays the role of the Euclidean Laplacian in the study of the Anderson model. For this reason a deeper study of hierarchical random walks is of interest per se, as is the case for random walks on the Euclidean lattice. Recurrence and transience of hierarchical random walks have been studied in [10] (and references therein). Lévy processes on totally disconnected groups have been studied from an abstract point of view in [17], and scaling limits of hierarchical random walks are examples of such Lévy processes, which play a key role in [15].

Some other references where stochastic models based on hierarchical structures have been studied are [20, 21, 26] (interacting diffusions and population models), [9, 12] (branching systems), [1] (contact processes), [7, 8, 27] (percolation), [24, 25] (search algorithms). Other references are found in [11].

The hierarchical random walks we will consider in this paper are of a type first introduced by Spitzer [30] (p.93) in a special case, and more generally by Sawyer and Felsenstein [29] in the context of population genetics. We will focus on “number variance” properties and related fluctuations for these random walks. The number variance problem, which is explained below, has been investigated by Hambly and Jones [22, 23] for α -stable processes on \mathbb{R} , and we will compare our results with theirs, in particular regarding properties that are qualitatively different.

The state space for our random walks is a hierarchical group. For an integer $M \geq 2$, the *hierarchical group of order M* (called *hierarchical lattice* in Physics) is defined by

$$\Omega_M = \{x = (x_1, x_2, \dots) : x_i \in \{0, 1, \dots, M-1\}, \sum_i x_i < \infty\},$$

with addition componentwise mod M . It is a countable Abelian group which is also described as the direct sum of a countable number of copies of the cyclic group of order M . The *hierarchical distance* $|\cdot|$ on Ω_M is defined by

$$|x - y| = \begin{cases} 0 & \text{if } x = y, \\ \max\{i : x_i \neq y_i\} & \text{if } x \neq y. \end{cases}$$

It is translation-invariant, and satisfies the strong (non-Archimedean) triangle inequality

$$|x - y| \leq \max\{|x - z|, |z - y|\} \quad \text{for any } x, y, z.$$

This means that $(\Omega_M, |\cdot|)$ is an ultrametric space. In such a space two balls are either disjoint or one is contained in the other, and this is the cause of some significant differences from mathematical models based on \mathbb{R}^d or \mathbb{Z}^d . In particular, as opposed to the Euclidean case, it is not possible to go far by a sequence of small steps, and the only way to go far is to make jumps of ever bigger sizes. A picture of $(\Omega_M, |\cdot|)$ is the set of leaves at the top of an infinite tree where each inner node at each level $j \geq 1$ has one predecessor at level $j + 1$ and M successors at level $j - 1$. The *distance* between two individuals (leaves) at level $j = 0$ is the depth in the tree to their most recent common ancestor, and it measures the degree of relatedness between the two individuals.

A random walk $\xi = (\xi_n)_{n \geq 0}$ on Ω_M starting at 0 is defined by $\xi_0 = 0, \xi_n = \rho_1 + \dots + \rho_n, n \geq 1$, where ρ_1, ρ_2, \dots are independent copies of ρ , which is a random element of Ω_M with distribution of the form

$$P(\rho = y) = \frac{r_{|y|}}{M^{|y|-1}(M-1)}, \quad y \neq 0, \quad P(\rho = 0) = 0, \quad (1.1)$$

where $(r_j)_{j=1,2,\dots}$ is a probability law on $\{1, 2, \dots\}$. That is, the jumps of ξ_n are taken by first choosing distance j with probability r_j , and then choosing a point with uniform probability among those at distance j from the previous position of the walk (note that $M^{j-1}(M-1)$ is the number of points at distance j from a given point of Ω_M). These random walks are the most general symmetric random walks on Ω_M . We assume that (r_j) is not restricted to a bounded set, so that the walk can reach arbitrarily large distances (since it is not possible to go far with small steps). We refer to a hierarchical random walk determined by (r_j) as r_j -rw. A particular r_j -rw is the c -rw, where c is a constant such that $0 < c < M$, and

$$r_j = \left(1 - \frac{c}{M}\right) \left(\frac{c}{M}\right)^{j-1}, \quad j = 1, 2, \dots, \quad (1.2)$$

A c -rw mimics a standard α -stable process on \mathbb{R}^d in the sense that both have analogous behaviors. The analogies, which refer to time and space behaviors, are mentioned in the Appendix. They are given in terms of degrees γ , in particular, $\gamma < 0$ corresponds to recurrence and $\gamma > 0$ to transience. The α -stable process on \mathbb{R}^d has degree

$$\gamma = \frac{d}{\alpha} - 1, \quad (1.3)$$

and the c -rw on Ω_M has degree

$$\gamma = \frac{\log c}{\log(M/c)}. \quad (1.4)$$

Correspondences are:

$$\text{recurrence: } d < \alpha \leftrightarrow c < 1, d = \alpha \leftrightarrow c = 1,$$

$$\text{transience: } d > \alpha \leftrightarrow c > 1.$$

See also Remark 2.13(b) concerning the order of the fluctuations in (1.5) for α -stable process and in (2.12) for c -rw.

Other r_j -rw's we will consider are defined in the Appendix and used for examples.

The number variance problem studied in [22, 23] for α -stable processes on \mathbb{R} has a motivation from physics (see the Introduction of [22]). The model is as follows in simplified form. An independent copy of the standard α -stable process, representing the motion of a particle, starts at time 0 from

each point $u_j = j - \varepsilon, j \in \mathbb{Z}$, where ε is a uniformly distributed random variable on $[0, 1]$ (the ε -displacement provides spatial homogeneity). Let $N_t[0, L]$ denote the number of particles that at time $t > 0$ lie in the interval $[0, L]$, and consider its variance, $\text{Var}N_t[0, L]$. The number variance problem refers to the behavior of $\text{Var}N_t[0, L]$ for fixed t as $L \rightarrow \infty$. It is shown that for $\alpha < 1$ (transient case) the variance tends to ∞ , and for $1 < \alpha \leq 2$ (recurrent case) it has a finite limit up to an additive fluctuation bounded by 1. The latter behavior is called *saturation*. For $\alpha = 1$ the process is recurrent and the variance tends to infinity. The rescaled fluctuation process defined by

$$Z_t(s) = \frac{N_t[0, st^{1/\alpha}] - EN_t[0, st^{1/\alpha}]}{t^{1/2\alpha}}, \quad s \geq 0, \quad (1.5)$$

is discussed, and a continuous interpolation of this process is shown to converge weakly on a space of continuous functions, as $t \rightarrow \infty$, to a centered Gaussian process with stationary increments, which is not Markovian and has long range dependence. Hence this process has a resemblance to fractional Brownian motion.

In this paper we will prove counterparts to the results of [22, 23] for a system of independent hierarchical random walks starting from each point of Ω_M . We will do this for a general class of random walks, and we will also consider a model where at the initial time there is a Poisson number of particles at each site. As with other models involving systems of hierarchical random walks, one of our objectives is to exhibit analogies and differences with the results for the Euclidean model, which can be compared due to the abovementioned correspondences.

The system of independent stable processes in [22, 23] is a very simplified model motivated by determinantal processes arising from random matrix theory, and although the results do not necessarily have a significance in connection with the motivation from physics, the formulation and the results are interesting as a new class of stochastic models. Our work should be regarded from that point of view as well.

Now we give a summary of our results.

We consider a system of independent r_j -rw's (particles) starting from each point of Ω_M , and denote by $N_n(L)$ the number of particles that at step n lie in the ball $\{x \in \Omega_M : |x| \leq L\}$, $L \in \mathbb{Z}_+$ (in [22, 23] the interval $[0, L]$ was considered, but essentially nothing changes if it is replaced by $[-L, L]$). For general r_j we give a necessary and sufficient condition for boundedness of $\text{Var}N_n(L)$ and for existence of its limit as $L \rightarrow \infty$. It turns out that

$$\lim_{L \rightarrow \infty} \text{Var}N_n(L) = 2n \frac{1}{M-1} \lim_{L \rightarrow \infty} M^L r_L$$

(Theorem 2.2). In particular, for the c -rw this limit exists, and it is finite if and only if $c \leq 1$, i.e., if the walk is recurrent. For the non-critical cases, $c < 1$ and $c > 1$, this result corresponds to the properties of the number variance for α -stable processes mentioned above, but the correspondence breaks down for the critical case, $c = 1$: the process is recurrent and the variance has a finite limit. For general r_j , boundedness of $\text{Var}N_n(L)$ and recurrence are not equivalent (see Remark 2.4(c)). For the initial Poisson system the situation is simpler, $\text{Var}N_n(L)$ always tends to infinity. It can be shown that the same thing happens for the α -stable process.

Next we investigate the fluctuations of $N_n(L)$ around the mean as $n \rightarrow \infty$, when $L = L(n)$ is increased in an appropriate way. Analogously as in (1.5), where a "time" parameter s was introduced, we introduce a "time" parameter t , taking $L(n) + R(t)$, where $R(t)$ is a suitable non-decreasing function of t . Thus, $L(n)$ and $R(t)$ here correspond to t and s in [22, 23], respectively. The additive form,

rather than the multiplicative one, is explained below. We assume that $\lim_{j \rightarrow \infty} r_{j+1}/r_j = a > 0$. It turns out that the cases $a < 1$ and $a = 1$ are significantly different (see Theorem 2.8). For $a < 1$, the most appropriate $L(n)$ has logarithmic growth and the norming is of the order $n^{\theta/2}$, where $\theta = \log M / \log \frac{1}{a}$. In general the fluctuation processes may not converge, but convergent subsequences can be chosen (in the sense of finite dimensional distributions), and a family of processes indexed by a parameter $\kappa \in [1, \frac{1}{a}]$ is obtained. These processes have stepwise trajectories, they are Gaussian with covariance of the form $f(s \vee t)g_\kappa(s \wedge t)$, in particular they are Markovian. If $a = 1$, the fluctuation process converges to a Gaussian process with independent increments. In contrast, for the α -stable process on \mathbb{R} the fluctuation limit is a non-Markov process with stationary increments [22, 23]. The Markov property in our case is a consequence of the ultrametric structure of Ω_M . In Remark 2.9(c) we give an intuitive explanation for this.

We study some further properties of our limit processes and discuss a probabilistic interpretation of the parameter θ , which is closely related to the degree γ .

Finally, for the Poisson model we show that with general r_j the fluctuation limit exists and it is, up to a constant, the same process as that obtained in the previous case for $a = 1$.

Due to the ultrametric structure of Ω_M , the calculations involved in the proofs are quite different from those in [22, 23].

2 Results

Fix an integer $M \geq 2$ and let Ω_M be the corresponding hierarchical group as defined in the Introduction. We consider a system of independent r_j -rw's on Ω_M starting from each point $u \in \Omega_M$. Recall that $N_n(L)$ is the number of particles lying in the ball $B_L = \{x \in \Omega_M : |x| \leq L\}$ at step n , i.e.,

$$N_n(L) = \sum_{u \in \Omega_M} \mathbf{1}_{B_L}(u + \xi_n^u), \quad (2.1)$$

where $\{\xi_n^u\}_{u \in \Omega_M}$ are independent copies of r_j -rw's starting at 0.

Proposition 2.1.

$$(a) \quad EN_n(L) = M^L, \quad n, L = 0, 1, 2, \dots, \quad (2.2)$$

$$(b) \quad \text{Var}N_n(L) \sim 2nM^L P(|\rho| > L) \text{ as } L \rightarrow \infty, \quad (2.3)$$

$$(c) \quad P(|\rho| \leq L)^{2n-1} n P(|\rho| > L) M^L \leq \text{Var}N_n(L) \leq 2M^L, \quad n, L = 0, 1, 2, \dots, \quad (2.4)$$

(\sim means that the quotient of both sides tends to 1).

The number variance problem is solved in the following theorem.

Theorem 2.2. *For any $n = 1, 2, \dots$, $\limsup_{L \rightarrow \infty} \text{Var}N_n(L) < \infty$ if and only if $\limsup_{L \rightarrow \infty} M^L r_L < \infty$. Moreover, $\lim_{L \rightarrow \infty} \text{Var}N_n(L)$ exists (finite or not) if and only if $\lim_{L \rightarrow \infty} M^L r_L$ exists, and in this case*

$$\lim_{L \rightarrow \infty} \text{Var}N_n(L) = 2n \frac{1}{M-1} \lim_{L \rightarrow \infty} M^L r_L. \quad (2.5)$$

This theorem applied to the examples in the Appendix gives the following results.

Example 2.3. (a) c -rw:

$$\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \begin{cases} 0 & \text{if } c < 1, \\ 2n & \text{if } c = 1, \\ \infty & \text{if } c > 1. \end{cases} \quad (2.6)$$

(b) j^β -rw:

$$\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \begin{cases} \infty & \text{if } \beta > 0, \\ 2n & \text{if } \beta = 0. \end{cases}$$

(for $\beta = 0$ it is obviously a c -rw with $c = 1$)

(c) $r_j = Dj^{-\beta}$, $\beta > 1$:

$$\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \infty.$$

Remark 2.4. (a) Let us compare the result (2.6) for the c -rw with the solution of the number variance problem for the α -stable process in [22, 23]. With $c \leq 1$ (recurrent case) we obtain “true” limits, whereas for the α -stable process with $\alpha > 1$ the variances are bounded but do not converge due to a fluctuating term (see [23]). On the other hand, with $c < 1$ the limit is trivial (zero), whereas for the α -stable process with $\alpha > 1$ the variances are bounded away from 0. The only non-trivial limit for the c -rw is obtained in the critical case $c = 1$, and it has no counterpart for the α -stable process, since for $\alpha = 1$ the variances tend to infinity.

(b) For the c -rw, finiteness of the variance limit is equivalent to recurrence, whereas for the α -stable process this equivalence breaks down in the critical case $\alpha = 1$.

(c) Example 2.3(b) shows that in general finiteness of the variance limit is not equivalent to recurrence (see the Appendix).

We now give a result for the Poisson case.

Proposition 2.5. Assume that initially at each site there is a Poisson number of particles, and these numbers are i.i.d. Then for any system of r_j -rw's, $\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \infty$.

Remark 2.6. (a) It can be shown that an analogous result holds for α -stable processes.

(b) We will come back to the Poisson system later on. Poisson systems seem to be the most natural as random initial configurations. However, it can be shown that for each initial configuration determined by i.i.d. random variables $\{v_x\}_{x \in \Omega_M}$ (v_x particles at site x) which are truly random, i.e., with $\text{Var } v_x > 0$, we have $\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \infty$.

So far we have considered $N_n(L)$ for n fixed. Now we will vary both n and L , more precisely, we want to investigate the normalized fluctuations of $N_n(L)$ as $n \rightarrow \infty$, simultaneously increasing L in an appropriate way.

We make the following assumption on the random walk:

$$r_{j+1} \leq r_j \text{ for } j \geq \text{some } j_0, \text{ and } \lim_{j \rightarrow \infty} \frac{r_{j+1}}{r_j} = a > 0. \quad (2.7)$$

This assumption is satisfied for all our examples (see Example 2.3 and the Appendix).

By Proposition 2.1(c), it is natural to take

$$L(n) = \sup \left\{ L \in \mathbb{Z}_+ : h(L) \geq \frac{1}{n} \right\}, \quad (2.8)$$

where

$$h(L) = P(|\rho| > L) = \sum_{j=L+1}^{\infty} r_j, \quad (2.9)$$

so that, by (2.4), $\text{Var}N_n(L(n))$ has the same rate of increase as $M^{L(n)}$. Hence $\sqrt{M^{L(n)}}$ is the natural normalization for the fluctuations of $N_n(L(n))$.

$L(n)$ has the following properties.

Lemma 2.7. (a) if $a < 1$, then

$$\lim_{n \rightarrow \infty} \frac{L(n)}{\log_a \frac{1}{n}} = 1. \quad (2.10)$$

(b)

$$1 \leq nh(L(n)) \quad \text{and} \quad \limsup_{n \rightarrow \infty} nh(L(n)) \leq \frac{1}{a}. \quad (2.11)$$

As in [22, 23], we want to investigate a fluctuation process introducing a new “time” parameter. In [22, 23] this was done with a multiplicative parameter s , but in our case an additive parameter $R(t)$ is appropriate. This is caused by the hierarchical structure of the state space, which for $a < 1$ implies the logarithmic growth of the radius of the balls (Lemma 2.7(a)). This is because the volume of a ball of radius L in \mathbb{R} grows like L , and the volume of a ball of radius L in Ω_M grows like M^L (see the beginning of the proofs). Consider any non-decreasing function $R : \mathbb{R}_+ \rightarrow \mathbb{Z}$ such that $\lim_{t \rightarrow \infty} R(t) = \infty$. For $n = 1, 2, \dots$ we define the fluctuation process as

$$X_n(t) = \frac{N_n((L(n) + R(t))^+) - EN_n((L(n) + R(t))^+)}{\sqrt{M^{L(n)}}}, \quad (2.12)$$

(see (2.2)).

In what follows \Rightarrow_f denotes weak convergence of finite-dimensional distributions, and $\lfloor x \rfloor$ is the integer part of $x \in \mathbb{R}$.

For a given by (2.7), we denote

$$b = \frac{M - a}{M - 1}. \quad (2.13)$$

Theorem 2.8. Assume (2.7).

(a) Suppose $a < 1$. Let $(n_i)_i$ be any subsequence such that

$$\lim_{i \rightarrow \infty} n_i h(L(n_i)) = \kappa \quad (2.14)$$

for some $\kappa \in [1, \frac{1}{a}]$. Then $X_{n_i} \Rightarrow_f X^{(\kappa, R)}$, where $X^{(\kappa, R)}$ is a centered Gaussian process with covariance

$$EX^{(\kappa, R)}(s)X^{(\kappa, R)}(t) = M^{R(s \wedge t)} g_\kappa(s \vee t), \quad (2.15)$$

where

$$g_\kappa(t) = 1 - \left(\frac{M-1}{M}\right)^2 \left(\sum_{j=0}^{\infty} \frac{e^{-\kappa b a^{R(t)} a^j}}{M^j}\right)^2 - \frac{(M-1)^3}{M^4} \sum_{j=0}^{\infty} \frac{1}{M^j} \left(\sum_{k=0}^{\infty} \frac{e^{-\kappa b a^{R(t)} a^{j+k+1}} - e^{-\kappa b a^{R(t)} a^j}}{M^k}\right)^2. \quad (2.16)$$

(b) If $a = 1$, then $\lim_{n \rightarrow \infty} nh(L(n)) = 1$ and $X_n \Rightarrow_f X^{(R)}$, where $X^{(R)}$ is a centered Gaussian process with covariance

$$EX^{(R)}(s)X^{(R)}(t) = (1 - e^{-2})M^{R(s \wedge t)}. \quad (2.17)$$

Remark 2.9. (a) Existence of a subsequence $(n_i)_i$ in part (a) follows immediately from Lemma 2.7(b). For example, for c -rw we have $L(n) = \lfloor \log_{c/M} \frac{1}{n} \rfloor$ and the condition (2.14) is satisfied for any subsequence $(n_i)_i$ such that $\log_{c/M} \frac{1}{n_i} - \lfloor \log_{c/M} \frac{1}{n_i} \rfloor$ converges (to $\log_{c/M} \frac{1}{\kappa}$), and any $\kappa \in [1, \frac{M}{c}]$ can be obtained in this way.

(b) As it will be seen in the proof, part (b) of the theorem holds under the weaker (than $r_{j+1}/r_j \rightarrow 1$) assumption

$$\lim_{j \rightarrow \infty} \frac{r_{j+1}}{h(j)} = 0. \quad (2.18)$$

(c) It is well known that Gaussian processes with covariances of the form (2.15) have the Markov property. This is in sharp contrast with the corresponding result in the Euclidean case [22, 23]. The Markov property in our case may be given a “particle picture” interpretation as follows. Due to ultrametricity (recall that it is not possible to go far with small steps), the random walks $u + \xi_n^u$ in (2.1) tend to stay at the same distance from the origin when they are far away from their starting positions (this follows from [10], Proposition 3.5.2), and since the probability of making a jump out of or into a large ball is small, the number of particles in a ball of radius $L(n) + R(t)$ with large n (see (2.12)) will not depend much on the past history of the numbers of particles in the ball, conditionally on the present number.

(d) The limit process obtained in part (b) has independent increments.

(e) The function $R(t)$ accounts for the time scaling of the limit process only. From the form of the covariances it is seen that the most natural forms of $R(t)$ are

$$R(t) = \left\lfloor \log_a \frac{1}{t} \right\rfloor \text{ if } a < 1, \quad (2.19)$$

$$R(t) = \lfloor \log_M t \rfloor \text{ if } a = 1. \quad (2.20)$$

We then have, for $a < 1$,

$$t^\theta \leq M^{R(t)} \leq M t^\theta, \quad (2.21)$$

where

$$\theta = \frac{\log M}{\log \frac{1}{a}}, \quad (2.22)$$

and

$$t \leq M^{R(t)} \leq M t \text{ if } a = 1. \quad (2.23)$$

It is also seen that the exponents in (2.16) are “close” to $\frac{1}{t}$, namely

$$\frac{1}{t} \leq a^{R(t)} < \frac{1}{at}. \quad (2.24)$$

In (2.21), (2.23) and (2.24), the left-hand inequalities are equalities for $t = a^{-n}$, $n \in \mathbb{Z}$.

Let us denote by $X^{(\kappa)}$ and X the limit processes corresponding to R given by (2.19) and (2.20), respectively. Observe that in spite of the fact that such R 's are not defined at $t = 0$, the processes $X^{(\kappa)}$ and X themselves are right continuous in L^2 at 0 with $X^{(\kappa)}(0) = 0$ and $X(0) = 0$. This follows from (2.21), (2.23), and since the function g_κ is bounded.

Remark 2.10. The process X has the representation

$$X_t = \sqrt{1 - e^{-2}} W_{M^j} \text{ for } M^j \leq t < M^{j+1}, \quad j \in \mathbb{Z},$$

where W is a standard Brownian motion.

Properties of the process $X^{(\kappa)}$ are summarized in the next proposition.

Proposition 2.11. (a) $X^{(\kappa)}$ is determined by a Gaussian sequence of random variables $(\zeta_j)_{j \in \mathbb{Z}}$,

$$X_t^{(\kappa)} = \zeta_j \text{ for } a^{-j} \leq t < a^{-(j+1)},$$

where the ζ_j have the representation

$$\zeta_j = g_\kappa(a^{-j}) \sum_{i=-\infty}^j v_i \sqrt{\frac{M^i}{g_\kappa(a^{-i})} - \frac{M^{i-1}}{g_\kappa(a^{-(i-1)})}}, \quad (2.25)$$

with $(v_i)_{i \in \mathbb{Z}}$ i.i.d. standard normal.

(b) $X^{(\kappa)}$ has a long range dependence property with exponent 1 in the sense that

$$0 < \limsup_{\tau \rightarrow \infty} \tau |E(X^{(\kappa)}(t) - X^{(\kappa)}(s))(X^{(\kappa)}(t + \tau) - X^{(\kappa)}(s + \tau))| < \infty. \quad (2.26)$$

(c) For $\theta > 1$ (i.e., $1 > a > \frac{1}{M}$), if we define

$$Y_m^{(\kappa)}(t) = a^{m(\theta-1)/2} X^{(\kappa)}(a^{-m}t), \quad t \geq 0,$$

then

$$Y_m^{(\kappa)} \Rightarrow_f C(\kappa) \sum_{j \in \mathbb{Z}} Y(a^{-j}) \mathbf{1}_{[a^{-j}, a^{-(j+1)})} \text{ as } m \rightarrow \infty,$$

where

$$Y(t) = \frac{\sqrt{\theta + 1}}{t} \int_0^t u^{\theta/2} dW_u, \quad t \geq 0.$$

Remark 2.12. (a) Part (c) corresponds to Proposition 4.16 in [22]. In that paper the limit process is a fractional Brownian motion, whereas in our case Y is a diffusion process satisfying the equation

$$dY(t) = -\frac{1}{t}Y(t)dt + (\sqrt{\theta + 1})t^{\theta/2-1}dW_t.$$

Its covariance has the particularly simple form

$$EY(s)Y(t) = (s \wedge t)^\theta (s \vee t)^{-1}. \quad (2.27)$$

(b) The condition $\theta > 1$ in part (c) is required only to ensure that the process Y is well defined at 0. For $\theta \leq 1$ the assertion of part (c) remains true on $[\varepsilon, \infty)$, $\varepsilon > 0$.

(c) We think that part (a) can be given a nicer form, analogous to that of the limit of $Y_m^{(\kappa)}$ in part (c), i.e., that the process $X^{(\kappa)}$ can be “interpolated” between the time points a^{-j} by a Gaussian diffusion process. To derive such a representation it would suffice to prove that the function g_κ defined by (2.16), with $a^{R(t)}$ replaced by $\frac{1}{t}$, is decreasing in t . We have not been able to prove it.

Remark 2.13. (a) Part (c) of Proposition 2.11 and Remark 2.12(b) show that the behavior of $X^{(\kappa)}(t)$ as $t \rightarrow \infty$ changes drastically depending on whether $\theta > 1$ or $\theta < 1$ (for $\theta > 1$, $\lim_{t \rightarrow \infty} \text{Var}X^{(\kappa)}(t) = \infty$, and for $\theta < 1$, $\lim_{t \rightarrow \infty} \text{Var}X^{(\kappa)}(t) = 0$). In fact, the parameter θ has a deeper probabilistic interpretation which is related to the recurrence/transience properties of the underlying random walk. These properties are in a sense characterized by the degree γ of a random walk (see the Appendix). From (A.2) it follows that $\theta = \gamma + 1$. In particular, this implies that for $\theta > 1$ the random walk is transient, and for $\theta < 1$ it is recurrent. In the critical case $\theta = 1$ both situations can occur. Moreover, by Theorem 2.2 it is not hard to check that if $\theta > 1$, then $\lim_{L \rightarrow \infty} \text{Var}N_n(L) = \infty$, and if $\theta < 1$, then $\lim_{L \rightarrow \infty} \text{Var}N_n(L) = 0$, while, as it was seen in Example 2.3, if $\theta = 1$ the number variance can be finite or infinite. We recall, however, that in general the condition $\limsup_{L \rightarrow \infty} \text{Var}N_n(L) < \infty$ is not equivalent to recurrence of the random walk.

(b) The equality $\theta = \gamma + 1$ implies that for c -rw's the parameter θ corresponds to $\frac{1}{\alpha}$ for the symmetric α -stable process in \mathbb{R} (see (1.3), and (A.2) in the Appendix). Hence for c -rw's the rate of the norming in (2.12), $\sqrt{M^{L(n)}} (\sim n^{\theta/2})$, corresponds exactly to $t^{1/2\alpha}$, which is the norming in the fluctuation theorem in [22] (see (1.5)).

Finally, we give the result for the Poisson system.

Theorem 2.14. *Under the assumptions of Proposition 2.5, for any r_j -rw, any functions $L : \mathbb{Z}_+ \rightarrow \mathbb{Z}_+$ and $R : \mathbb{R}_+ \rightarrow \mathbb{Z}$, non-decreasing with $\lim_{n \rightarrow \infty} L(n) = \infty$ and $\lim_{t \rightarrow \infty} R(t) = \infty$, if X_n is defined by (2.12), then $X_n \Rightarrow_f CX^{(R)}$, where $X^{(R)}$ is given in Theorem 2.8(b) and C is a constant.*

Remark 2.15. From the proof of this theorem it will be seen that the same result is true if the initial number of particles at site $u \in \Omega_M$ is η^u , where $(\eta^u)_{u \in \Omega_M}$ are i.i.d. random variables with $E(\eta^u)^{2+\delta} < \infty$ for some $\delta > 0$, and $E\eta^u = \text{Var} \eta^u$.

3 Proofs

Let us denote

$$\begin{aligned} B_L(u) &= \{x \in \Omega_M : |x - u| \leq L\}, \quad B_L = B_L(0), \\ S_L(u) &= \{x \in \Omega_M : |x - u| = L\}, \quad S_L = S_L(0), \\ p_n^u(L) &= P(u + \xi_n \in B_L). \end{aligned}$$

We write $|A|$ for the number of points in a bounded subset A of Ω_M (no confusion arises with the hierarchical distance). Note that

$$|B_L(u)| = M^L, \quad |S_L(u)| = (M-1)M^{L-1}.$$

By ultrametricity of Ω_M , if $u \in B_L$, then $u + \xi_n \in B_L$ if and only if $\xi_n \in B_L$. On the other hand, if $|u| = L+k, k=1, 2, \dots$, then $B_L \subset S_{L+k}(u)$, hence

$$p_n^u(L) = P(u + \xi_n \in B_L \cap S_{L+k}(u)) = P(|\xi_n| = L+k) \frac{|B_L|}{|S_{L+k}|}.$$

Thus we have

$$p_n^u(L) = \begin{cases} P(|\xi_n| \leq L) & \text{if } u \in B_L, \\ P(|\xi_n| = L+k) \frac{1}{(M-1)M^{k-1}} & \text{if } |u| = L+k, k=1, 2, \dots \end{cases} \quad (3.1)$$

Proof of Proposition 2.1 By (2.1) and (3.1),

$$EN_n(L) = \sum_{u \in B_L} p_n^u(L) + \sum_{u \notin B_L} p_n^u(L) = |B_L|.$$

This proves (a) (this formula is surely known, but we have not found a reference).

In order to investigate $\text{Var}N_n(L)$ we study the tail of ξ_n . We have

$$\begin{aligned} P(|\xi_n| > L) &\leq \sum_{k=1}^n P(|\xi_1| \leq L, \dots, |\xi_{k-1}| \leq L, |\xi_k| > L) \\ &= \sum_{k=1}^n P(|\rho| \leq L)^{k-1} P(|\rho| > L), \end{aligned} \quad (3.2)$$

by ultrametricity. Again by ultrametricity, the event that in the first n steps there is just one farthest jump whose length is j and it occurs at step k has the form

$$A_{k,j} = \{|\xi_1| < j, \dots, |\xi_{k-1}| < j, |\xi_k| = j, \xi_{k+1} \in B_{j-1}(\xi_k), \dots, \xi_n \in B_{j-1}(\xi_k)\},$$

hence

$$\begin{aligned} P(|\xi_n| > L) &\geq \sum_{k=1}^n \sum_{j=L+1}^{\infty} P(A_{k,j}) = \sum_{k=1}^n \sum_{j=L+1}^{\infty} P(|\rho| < j)^{n-1} P(|\rho| = j) \\ &\geq nP(|\rho| \leq L)^{n-1} P(|\rho| > L). \end{aligned} \quad (3.3)$$

Formulas (3.2) and (3.3) imply that

$$P(|\xi_n| > L) \sim nP(|\rho| > L) \quad \text{as } L \rightarrow \infty. \quad (3.4)$$

Using (2.1) and (3.1) we have

$$\text{Var}N_n(L) = I(L) + II(L),$$

where

$$\begin{aligned} I(L) &= M^L P(|\xi_n| \leq L) P(|\xi_n| > L), \\ II(L) &= M^L \sum_{k=1}^{\infty} P(|\xi_n| = L+k) \left(1 - \frac{1}{(M-1)M^{k-1}} P(|\xi_n| = L+k) \right). \end{aligned}$$

Hence (b) follows immediately from (3.4).

The upper estimate in (c) is clear from the previous calculations. The lower bound is an easy consequence of

$$\text{Var}N_n(L) \geq II(L) \geq M^L P(|\xi_n| > L)(1 - P(|\xi_n| > L))$$

and (3.2), (3.3). □

Proof of Theorem 2.2 All the statements follow from (2.3) and the trivial formulas

$$M^L P(|\rho| > L) = M^L r_{L+1} + M^L P(|\rho| > L+1), \quad M^L P(|\rho| > L) = \sum_{j=1}^{\infty} \frac{r_{j+L} M^{j+L}}{M^j}.$$

□

Proof of Proposition 2.5 If the number of particles at each site is Poisson with parameter λ , then

$$\text{Var}N_n(L) = \lambda \sum_{u \in \Omega_M} p_n^u(L) = \lambda EN_n(L) = \lambda M^L,$$

by (2.2), hence the assertion follows. □

Proof of Lemma 2.7 (a) By (2.7) and (2.9) it is easy to see that for any $0 < \varepsilon < a \wedge (1-a)$ there exist positive constants C_1, C_2 such that for sufficiently large n ,

$$C_1(a - \varepsilon)^n \leq h(n) \leq C_2(a + \varepsilon)^n. \tag{3.5}$$

Then, for large n ,

$$h \left(\left\lceil \log_{a+\varepsilon} \frac{1}{n} + \log_{a+\varepsilon} \frac{1}{C_2} \right\rceil + 1 \right) \leq C_2(a + \varepsilon)^{\log_{(a+\varepsilon)} 1/n C_2} = \frac{1}{n},$$

which implies that

$$L(n) \leq \log_{a+\varepsilon} \frac{1}{n} + \log_{a+\varepsilon} \frac{1}{C_2} + 1,$$

hence

$$\limsup_{n \rightarrow \infty} \frac{L(n)}{\log_a \frac{1}{n}} \leq \frac{\log a}{\log(a + \varepsilon)}.$$

Analogously, using the left-hand side of (3.5) we obtain

$$\liminf_{n \rightarrow \infty} \frac{L(n)}{\log_a \frac{1}{n}} \geq \frac{\log a}{\log(a - \varepsilon)}.$$

Hence (2.10) follows

(b) The lower bound in (2.11) is obvious by (2.8). The assumption (2.7) clearly implies

$$\lim_{n \rightarrow \infty} \frac{r_{n+k}}{r_n} = a^k, \quad k = 1, 2, \dots, \quad (3.6)$$

hence for $a < 1$ we have

$$\left(\sum_{j=0}^{\infty} (a + \varepsilon)^j \right)^{-1} \leq \liminf_{n \rightarrow \infty} \frac{r_{n+1}}{h(n)} \leq \limsup_{n \rightarrow \infty} \frac{r_{n+1}}{h(n)} \leq \left(\sum_{j=0}^k a^j \right)^{-1}$$

for any $0 < \varepsilon < 1 - a$ and $k = 1, 2, \dots$. The right-hand side inequality also holds for $a = 1$. Therefore, for all $0 < a \leq 1$,

$$\lim_{n \rightarrow \infty} \frac{r_{n+1}}{h(n)} = 1 - a. \quad (3.7)$$

By (2.8) and (2.9),

$$nh(L(n)) < 1 + nr_{L(n)+1},$$

so (2.11) follows easily from (3.7). \square

Proof of Theorem 2.8 (a) We will prove

$$\lim_{i \rightarrow \infty} \text{Cov}(X_{n_i}(s), X_{n_i}(t)) = M^{R(s \wedge t)} g_{\kappa}(s \vee t) \quad (3.8)$$

and

$$\lim_{n \rightarrow \infty} \frac{1}{(M^{L(n)})^{1+\delta/2}} \sum_{u \in \Omega_M} E \left| \mathbf{1}_{B_{(L(n)+R(t))^+}(u + \xi_n^u)} - P(u + \xi_n^u \leq (L(n) + R(t))^+) \right|^{2+\delta} = 0, \quad \delta > 0. \quad (3.9)$$

Then by (2.1), (2.12) and independence of random walks the result follows from the central limit theorem in the Lyapunov version (see e.g. [3]).

Denote

$$L(i, t) = (L(n_i) + R(t))^+. \quad (3.10)$$

Fix $s \leq t$. By (2.1) and independence of random walks,

$$\begin{aligned} \text{Cov}(X_{n_i}(s), X_{n_i}(t)) &= M^{-L(n_i)} \sum_{u \in \Omega_M} p_{n_i}^u(L(i, s)) \left(1 - p_{n_i}^u(L(i, t)) \right) \\ &= M^{-L(n_i)} (I + II), \end{aligned} \quad (3.11)$$

where $I = \sum_{u \in B(L(i, t))} \dots$ and $II = \sum_{u \notin B(L(i, t))} \dots$. We have

$$\begin{aligned} I &= \sum_{u \in B(L(i, s))} \dots + \sum_{u \in B(L(i, t)) \setminus B(L(i, s))} \dots \\ &= M^{L(i, s)} P(|\xi_{n_i}| \leq L(i, s)) P(|\xi_{n_i}| > L(i, t)) \\ &\quad + \sum_{j=L(i, s)+1}^{L(i, t)} \sum_{u \in S_j} P(|\xi_{n_i}| = j) \frac{|B_{L(i, s)}|}{|S_j|} \left(1 - P(|\xi_{n_i}| \leq L(i, t)) \right) \\ &= M^{L(i, s)} P(|\xi_{n_i}| \leq L(i, t)) P(|\xi_{n_i}| > L(i, t)), \end{aligned} \quad (3.12)$$

by (3.1). Similarly

$$\begin{aligned} II &= \sum_{j=1}^{\infty} \sum_{u \in S_{L(i,t)+j}} \dots \\ &= M^{L(i,s)} \sum_{j=1}^{\infty} P(|\xi_{n_i}| = L(i,t) + j) \left(1 - P(|\xi_{n_i}| = L(i,t) + j) \frac{1}{(M-1)M^{j-1}} \right). \end{aligned} \quad (3.13)$$

By (3.10)-(3.13), for sufficiently large i , we obtain

$$\begin{aligned} \text{Cov}(X_{n_i}(s), X_{n_i}(t)) &= M^{R(s)} \left[P(|\xi_{n_i}| > L(i,t))(1 + P(|\xi_{n_i}| \leq L(i,t))) \right. \\ &\quad \left. - \frac{1}{M-1} \sum_{j=1}^{\infty} \frac{P^2(|\xi_{n_i}| = L(i,t) + j)}{M^{j-1}} \right]. \end{aligned} \quad (3.14)$$

It is known ([29], see also [10]) that

$$P(\xi_n = u) = -\frac{f_k^n}{M^k} + (M-1) \sum_{j=k+1}^{\infty} \frac{f_j^n}{M^j}, \quad \text{if } |u| = k > 0,$$

where

$$f_k = 1 - h(k-1) - \frac{r_k}{M-1}. \quad (3.15)$$

Hence

$$\begin{aligned} P(|\xi_n| = k) &= \frac{M-1}{M} \left(-f_k^n + (M-1) \sum_{j=1}^{\infty} \frac{f_{j+k}^n}{M^j} \right) \\ &= \frac{M-1}{M} \sum_{j=0}^{\infty} \frac{f_{j+k+1}^n - f_{j+k}^n}{M^j}. \end{aligned} \quad (3.16)$$

This implies that

$$P(|\xi_n| > L) = \frac{M-1}{M} \sum_{j=0}^{\infty} \frac{1}{M^j} (1 - f_{L+j+1}^n). \quad (3.17)$$

By (2.14), (3.6), (3.7) and (3.10) we have

$$\lim_{i \rightarrow \infty} n_i \left(h(L(i,t) + j - 1) + \frac{r_{L(i,t)+j}}{M-1} \right) = \kappa b a^{R(t)+j-1}, \quad (3.18)$$

where b is defined by (2.13). Hence

$$\lim_{i \rightarrow \infty} f_{L(i,t)+j}^{n_i} = e^{-\kappa b a^{R(t)+j-1}}, \quad (3.19)$$

by (3.15).

Combining (3.16)-(3.19) and (3.14) (it is clear that we can pass to the limits under the sums) we arrive at

$$\begin{aligned}
& \lim_{i \rightarrow \infty} \text{Cov}(X_{n_i}(s), X_{n_i}(t)) \\
&= M^{R(s)} \left[\left(\frac{M-1}{M} \sum_{j=0}^{\infty} \frac{1 - e^{-\kappa b a^{R(t)} a^j}}{M^j} \right) \left(1 + \frac{M-1}{M} \sum_{j=0}^{\infty} \frac{e^{-\kappa b a^{R(t)} a^j}}{M^j} \right) \right. \\
&\quad \left. - \frac{(M-1)^3}{M^4} \sum_{j=0}^{\infty} \frac{1}{M^j} \left(\sum_{k=0}^{\infty} \frac{e^{-\kappa b a^{R(t)} a^{j+1+k}} - e^{-\kappa b a^{R(t)} a^j}}{M^k} \right)^2 \right] \\
&= M^{R(s)} g_{\kappa}(t) \tag{3.20}
\end{aligned}$$

(see (2.16)). This proves (3.8).

It is easy to see that the expression under the limit on the left hand side of (3.9) is estimated from above by

$$\frac{C}{M^{L(n)(1+\delta/2)}} \sum_{u \in \Omega_M} p_n^u((L(n) + R(t))^+) = \frac{C}{M^{L(n)(1+\delta/2)}} M^{(L(n)+R(t))^+},$$

by (2.2), hence (3.9) follows.

The proof of part (a) is complete.

(b) The fact that

$$\lim_{n \rightarrow \infty} n(h(L(n)) = 1 \tag{3.21}$$

follows immediately from (2.11). This and (3.7) imply that

$$\lim_{n \rightarrow \infty} n \left(h((L(n) + R(t))^+ + j - 1) + \frac{r_{(L(n)+R(t))^+ + j}}{M-1} \right) = 1 \tag{3.22}$$

(cf. (3.18)), hence

$$\lim_{n \rightarrow \infty} f_{(L(n)+R(t))^+ + j}^n = e^{-1}$$

for $j = 0, 1, 2, \dots$. Therefore, a counterpart of (3.20) is

$$\lim_{n \rightarrow \infty} \text{Cov}(X_n(s), X_n(t)) = (1 - e^{-2}) M^{R(s \wedge t)}. \tag{3.23}$$

Now, the \Rightarrow_f convergence (the Lyapunov condition (3.9)) is obtained in the same way as before.

Observe that (3.21), (3.22), and hence (3.23) as well, follow from (3.7) in this case. (Recall that (3.7) with $a = 1$ implies (2.11), see the end of the proof of Lemma 2.7(b).) Therefore, as stated in Remark 2.9 (b), assumption (2.7) can be replaced by (2.18). \square

Proof of Proposition 2.11 (a) From positive-definiteness of the covariance function (2.15) it is easy to see that $M^i/g_{\kappa}(a^{-i})$ is increasing in i , moreover, by (2.16), $\lim_{i \rightarrow -\infty} M^i/g_{\kappa}(a^{-i}) = 0$. Hence the result is obtained by a direct computation.

(b) By (2.15), for $s < t < s + \tau < t + \tau$ we have

$$\begin{aligned}
& E(X^{(\kappa)}(t) - X^{(\kappa)}(s))(X^{(\kappa)}(t + \tau) - X^{(\kappa)}(s + \tau)) \\
&= (M^{R(t)} - M^{R(s)})(g_{\kappa}(t + \tau) - g_{\kappa}(s + \tau)),
\end{aligned}$$

hence it suffices to investigate the second factor. Obviously, for large τ , the time points $s + \tau, t + \tau$ belong either to the same interval of the form $[a^{-k}, a^{-(k+1)})$, or to two neighboring such intervals. Since, by (2.19) and (2.16), g_κ is constant on such intervals, it is enough to consider a sequence of the form

$$\tau_m = a^{-m} - d_m, \quad s < d_m \leq t, \quad m = 1, 2, \dots$$

Then, for large m we have

$$\tau_m(g_\kappa(\tau_m + t) - g_\kappa(\tau_m + s)) = \tau_m a^m a^{-m} (\tilde{g}_\kappa(a^m) - \tilde{g}_\kappa(a^{m-1})),$$

where $\tilde{g}_\kappa(r)$ is obtained from (2.16) by putting r instead of $a^{R(t)}$. The mean value theorem implies that $\frac{1}{r}(\tilde{g}_\kappa(r) - \tilde{g}_\kappa(\frac{r}{a}))$ has a finite positive limit as $r \rightarrow 0$. This proves (2.26).

(c) Since the processes $Y_m^{(\kappa)}$ are centered Gaussian, it suffices to prove convergence of covariances. By (2.19) and (2.22),

$$\begin{aligned} R(a^{-m}t) &= m + R(t), \quad m = 1, 2, \dots, \\ M^m &= a^{-m\theta}. \end{aligned} \tag{3.24}$$

This, (2.15) and (2.19) imply that for $s \leq t$,

$$\text{Cov}(Y_m^{(\kappa)}(s), Y_m^{(\kappa)}(t)) = M^{R(s)} a^{-m} g_\kappa(a^{-m}t).$$

By (3.24) we also have

$$\lim_{m \rightarrow \infty} \frac{1 - e^{-\kappa b a^{R(a^{-m}t)} a^j}}{a^m} = \kappa b a^{R(t)} a^j.$$

Hence, using the form of g_κ given by (3.20), and by (2.13), it is easy to see that

$$\lim_{m \rightarrow \infty} \text{Cov}(Y_m^{(\kappa)}(s), Y_m^{(\kappa)}(t)) = 2\kappa M^{R(s)} a^{R(t)}, \quad s \leq t.$$

Hence (c) follows (cf (2.27)). □

Proof of Theorem 2.14 Let η^u be the number of particles at site u at time 0. The r.v.'s $(\eta^u)_{u \in \Omega_M}$ are i.i.d. Let $(\xi^{u,k})_{u \in \Omega_M, k \in \mathbb{N}}$ be independent copies of r_j -rw. Then

$$N_n(L) = \sum_{u \in \Omega_M} \sum_{k=1}^{\eta^u} \mathbf{1}_{\{u + \xi_n^{u,k} \in B_L\}},$$

and for $s \leq t$, denoting $L(n, t) = (L(n) + R(t))^+$,

$$\begin{aligned} \text{Cov}(X_n(s), X_n(t)) &= \frac{1}{M^{L(n)}} \sum_{u \in \Omega_M} \left[E\eta^u P(u + \xi_n \in B_{L(n,s)}) \right. \\ &\quad \left. + (\text{Var } \eta^u - E\eta^u) P(u + \xi_n \in B_{L(n,s)}) P(u + \xi_n \in B_{L(n,t)}) \right] \\ &= \frac{E\eta^0}{M^{L(n)}} M^{L(n,s)}, \end{aligned}$$

by (2.2). Hence

$$\lim_{n \rightarrow \infty} \text{Cov}(X_n(s), X_n(t)) = E\eta^0 M^{R(s)}.$$

Next, using the inequality $(a_1 + \dots + a_m)^{2+\delta} \leq m^{1+\delta}(a_1^{2+\delta} + \dots + a_m^{2+\delta})$, $a_i \geq 0$, we easily obtain

$$\lim_{n \rightarrow \infty} \frac{1}{M^{(1+\delta/2)L(n)}} \sum_{u \in \Omega_M} E \left(\sum_{k=1}^{\eta^u} \mathbf{1}_{B_{L(n,t)}}(u + \xi_n^{u,k}) \right)^{2+\delta} = 0.$$

Convergence of finite-dimensional distributions now follows from the central limit theorem (Lyapunov criterion). \square

Appendix: The degree γ

The development and applications of hierarchical random walks are outlined in [11], and their recurrence/transience and some other properties are studied in [10] (and references therein).

In order to deal with the standard α -stable process on \mathbb{R}^d and discrete time hierarchical random walks in a unified way, the continuous time version of a walk with unit rate holding time was taken in [10]. Its transition probability from 0 to y in time $t > 0$ is given by

$$p_t(0, y) = e^{-t} \sum_{n=0}^{\infty} \frac{t^n}{n!} p^{(n)}(0, y),$$

where $p^{(n)}(0, y)$ is the transition probability from 0 to y in n steps. Generally, for a Lévy process on a Polish space S (with additive group structure in the cases considered in [10]) with semigroup \mathcal{T}_t , the *degree* of the process is defined by

$$\gamma = \sup\{\zeta > -1 : G^{\zeta+1}\varphi < \infty \text{ for all } \varphi \in \mathcal{B}_b^+(S)\},$$

where $\mathcal{B}_b^+(S)$ is the space of bounded non-negative measurable functions with bounded support on S , and

$$G^\zeta \varphi = \frac{1}{\Gamma(\zeta)} \int_0^\infty t^{\zeta-1} \mathcal{T}_t \varphi dt, \quad \zeta > 0,$$

is the fractional power of the potential (or Green) operator of the process. For $\zeta = \gamma$ it may happen that $G^{\gamma+1}\varphi < \infty$ or $G^{\gamma+1}\varphi = \infty, \varphi \neq 0$. Recurrence corresponds to $\gamma \in (-1, 0)$ and transience corresponds to $\gamma \in (0, \infty]$. The value $\gamma = 0$ is special since both cases $G\varphi < \infty$ and $G\varphi = \infty, \varphi \neq 0$, can happen. The value $\gamma = \infty$ occurs, for example, for a simple asymmetric random walk on \mathbb{Z} . For $\gamma \geq 0$, γ is also given by

$$\gamma = \sup\{\zeta \geq 0 : EL_{B_R}^\zeta < \infty \text{ for all } R > 0\},$$

where L_{B_R} is the last exit time of the process (starting at 0) from an open ball B_R of radius R centered at 0. Thus, $\gamma > 0$, called *degree of transience* in this case [10], gives information on how fast the process escapes to infinity. For a discrete time random walk on a discrete space, instead of $G^\zeta \varphi$ it is natural to consider

$$\frac{1}{\Gamma(\zeta)} \sum_{n=0}^{\infty} n^{\zeta-1} p^{(n)}(0, 0), \tag{A.1}$$

but since

$$\int_0^\infty t^\zeta p_t(0,0)dy = \sum_{n=0}^\infty \frac{\Gamma(n+\zeta+1)}{n!} p^{(n)}(0,0),$$

and

$$\frac{\Gamma(n+\zeta+1)}{n!n^\zeta} \rightarrow 1 \quad \text{as } n \rightarrow \infty,$$

the degree of the random walk is the same in discrete and continuous time.

For the standard α -stable process on \mathbb{R}^d , γ is given by (1.3), and the range of possible values of γ for those processes is restricted to the interval $[-\frac{1}{2}, \infty)$. For the c -rw defined by (1.1), (1.2), γ is given by (1.4), and the range of possible values of γ is $(-1, \infty)$. Hence this class of hierarchical random walks is richer than the class of standard α -stable processes and corresponding symmetric random walks on \mathbb{Z}^d . The c -rw's mimic the behavior of α -stable processes because they have the same recurrence/transience properties for equal values of their degrees (in particular, for $\gamma = 0$, $G\varphi = \infty$ holds for both of them), and they have the same spatial asymptotic decay of powers of their potential operators in terms of the "Euclidean radial distance" (see [10], Remark 3.2.2(b)). The correspondence of degrees allows to choose c in order to study "caricatures" of α -stable processes by means of c -rw's, including non-integer values of the dimension d (which is one of the reasons for using hierarchical random walks in statistical physics, see [10, 11] and references therein). But there are also differences, for example, the distance from 0 of the c -rw has a different behavior from the α -stable Bessel process (see [10], Remark 3.5.7). Definitions of k -weak transience and k -strong transience for each integer $k \geq 1$ were given in [9], and they were related to occupation time fluctuation limits of branching systems with k levels of branching for $k = 0, 1, 2$. The analogies between the results for α -stable processes on \mathbb{R}^d and c -rw's on Ω_M with equal values of γ for the two processes occur because they have the same k -weak/strong transience behaviors. The correspondence between the degrees of the α -stable process and the c -rw is also used in this paper to compare our results with those of the α -stable process in [22, 23]. See in particular Remark 2.13(b).

The assumption $r_{j+1}/r_j \rightarrow a$ as $j \rightarrow \infty$, $0 < a \leq 1$, implies that the degree of the r_j -rw is given by

$$\gamma = \frac{\log M}{\log \frac{1}{a}} - 1 \quad \text{for } a < 1, \quad \gamma = \infty \quad \text{for } a = 1. \quad (\text{A.2})$$

This is obtained from [10] (Proposition 3.2.7). Hence

$$\gamma \begin{cases} < 0, \\ = 0, \\ > 0, \end{cases} \quad \text{if } a \begin{cases} < \frac{1}{M}, \text{ recurrent} \\ = \frac{1}{M}, \\ > \frac{1}{M}, \text{ transient.} \end{cases}$$

The parameter $\theta = \gamma + 1$ (see (2.22)) therefore has a meaning in terms of recurrence/transience.

Examples:

1. c -rw: $0 < c < M$,

$$r_j = \left(1 - \frac{c}{M}\right) \left(\frac{c}{M}\right)^{j-1}, \quad j = 1, 2, \dots, \quad a = \frac{c}{M}, \quad \gamma = \frac{\log c}{\log(M/c)},$$

recurrent for $c \leq 1$, transient for $c > 1$.

2. j^β -rw: $\beta \geq 0$,

$$r_j = D \frac{j^\beta}{M^j}, \quad j = 1, 2, \dots,$$

where D is a normalizing constant,

$$a = \frac{1}{M}, \quad \gamma = 0,$$

recurrent for $\beta \leq 1$, transient for $\beta > 1$ (follows from the recurrence criterion in [19], see also Example 3.2.6 in [10], which is slightly different) .

3.

$$r_j = D j^{-\beta}, \quad j = 1, 2, \dots,$$

where $\beta > 1$ and D is a normalizing constant,

$$a = 1, \quad \gamma = \infty, \quad \text{transient.}$$

Remark. In [10] r_j is written in the form

$$r_j = D \frac{c_j}{N^{j/\mu}}, \quad j = 1, 2, \dots,$$

where μ is a positive constant, (c_j) is a sequence of positive numbers, and D is a normalizing constant. The parameter μ is useful for some applications (see [11, 12]).

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