

Hierarchical Cooperation Achieves Optimal Capacity Scaling in *Ad Hoc* Networks

Ayfer Özgür, Olivier Lévêque, *Member, IEEE*, and David N. C. Tse, *Member, IEEE*

Abstract— n source and destination pairs randomly located in an area want to communicate with each other. Signals transmitted from one user to another at distance r apart are subject to a power loss of $r^{-\alpha}$ as well as a random phase. We identify the scaling laws of the information-theoretic capacity of the network when nodes can relay information for each other. In the case of dense networks, where the area is fixed and the density of nodes increasing, we show that the total capacity of the network scales linearly with n . This improves on the best known achievability result of $n^{2/3}$ of Aeron and Saligrama. In the case of extended networks, where the density of nodes is fixed and the area increasing linearly with n , we show that this capacity scales as $n^{2-\alpha/2}$ for $2 \leq \alpha < 3$ and \sqrt{n} for $\alpha \geq 3$. The best known earlier result of Xie and Kumar identified the scaling law for $\alpha > 4$. Thus, much better scaling than multihop can be achieved in dense networks, as well as in extended networks with low attenuation. The performance gain is achieved by intelligent node cooperation and distributed multiple-input multiple-output (MIMO) communication. The key ingredient is a hierarchical and digital architecture for nodal exchange of information for realizing the cooperation.

Index Terms—*Ad hoc* wireless networks, capacity scaling, digital communications, distributed multiple-input multiple-output (MIMO), hierarchical cooperation.

I. INTRODUCTION

THE seminal paper by Gupta and Kumar [3] initiated the study of scaling laws in large *ad hoc* wireless networks. Their by-now-familiar model considers n nodes randomly located in the unit disk, each of which wants to communicate to a random destination node at a rate $R(n)$ bits per second.

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They ask what is the maximally achievable scaling of the total throughput $T(n) = nR(n)$ with the system size n . They showed that classical multihop architectures with conventional single-user decoding and forwarding of packets cannot achieve a scaling better than $O(\sqrt{n})$, and that a scheme that uses only nearest neighbor communication can achieve a throughput that scales as $\Theta(\sqrt{n/\log n})$. This gap was later closed by Franceschetti *et al.* [4], who showed using percolation theory that the $\Theta(\sqrt{n})$ scaling is indeed achievable.

The Gupta–Kumar model makes certain assumptions on the physical-layer communication technology. In particular, it assumes that the signals received from nodes other than one particular transmitter are interference to be regarded as noise degrading the communication link. Given this assumption, direct communication between source and destination pairs is not preferable, as the interference generated would preclude most of the other nodes from communicating. Instead, the optimal strategy is to confine to nearest neighbor communication and maximize the number of simultaneous transmissions (spatial reuse). However, this means that each packet has to be retransmitted many times before getting to the final destination, leading to a sublinear scaling of system throughput.

A natural question is whether the Gupta–Kumar scaling law is a consequence of the physical-layer technology or whether one can do better using more sophisticated physical-layer processing. More generally, what is the *information-theoretic* scaling law of *ad hoc* networks? This question was first addressed by Xie and Kumar [5]. They showed that whenever the power path loss exponent α of the environment is greater than 6 (i.e., the received power decays faster than r^{-6} with the distance r from the transmitter), then the nearest neighbor multihop scheme is in fact order-optimal. They also showed that the same conclusion holds if the power path loss is exponential in the distance r , a channel model proposed recently by Franceschetti *et al.* [6]. The work [5] was followed by several others [7]–[10], [2]. Successively, they improved the threshold on the path-loss exponent α for which multihop is order-optimal ($\alpha > 5$ in [7], $\alpha > 4.5$ in [10], and $\alpha > 4$ in [2]). However, the question is open for the important range of α between 2 and 4, $\alpha = 2$ corresponding to free-space attenuation.

It is important to realize that the achievability result in [3] and the upper bounds in [5] (and follow-up works) are actually based on different network models. The paper [3] deals with *dense* networks, where the total area is fixed and the density of nodes increases. The paper [5] and the subsequent works, on the other hand, focus on *extended* networks, which scale to cover an increasing area with the density of nodes fixed. A way to understand the difference between the engineering implications of

these two network scalings is by drawing a parallelism with the classical notions of *interference-limitedness* and *coverage-limitedness*, the two operating regimes of cellular networks. Cellular networks in urban areas tend to have dense deployments of base stations so that signals are received at the mobiles with sufficient signal-to-noise ratio (SNR) but performance is limited by *interference* between transmissions in adjacent cells. Cellular networks in rural areas, on the other hand, tend to have sparse deployments of base stations so that performance is mainly limited by the ability to transmit enough power to reach all the users with sufficient SNR. Analogously, in the dense network scaling, all nodes can communicate with each other with sufficient SNR; performance can only be limited by interference, if at all. The Gupta–Kumar scaling of \sqrt{n} comes precisely from such interference limitation. In the extended network scaling, the source and destination pairs are at increasing distance from each other, and so both interference limitation and power limitation can come into play. The network can be either coverage-limited or interference-limited. The information-theoretic limits on performance proved in [5], [7]–[10], [2], for extended networks are all based on bounding the maximum amount of power that can be transferred across the network and then showing that multihop achieves that bound. Hence, what was shown by these works is that for $\alpha > 4$, when signals attenuate fast enough, an extended network is fundamentally coverage-limited: even with optimal cooperative relaying, the amount of power transferred across the network cannot be larger than that achieved by multihop. For α between 2 and 4, when attenuation is lower and power transfer become easier, the question remains open whether the network is coverage-limited or interference-limited.

Viewing the earlier results in this light, a natural first step in completing the picture is to return to the simpler dense scaling as a vehicle to focus exclusively on the issue of interference. Can the interference limitation implied by the Gupta–Kumar result be overcome by more sophisticated physical-layer processing? In a recent work [1], Aeron and Saligrama have showed that the answer is indeed yes: they exhibited a scheme which yields a throughput scaling of $\Theta(n^{2/3})$ bits per second. However, it is not clear if one can do even better. The first main result in this paper is that, for any value of $\alpha \geq 2$, one can in fact achieve arbitrarily close to *linear* scaling: for any $\epsilon > 0$, we present a scheme that achieves an aggregate rate of $\Theta(n^{1-\epsilon})$. This is a surprising result: a linear scaling means that there is essentially *no* interference limitation; the rate for *each* source–destination pair does not degrade significantly even as one puts more and more nodes in the network. It is easy to show that one cannot get a better capacity scaling than $O(n \log n)$, so our scheme is close to optimal.

To achieve linear scaling, one must be able to perform *many* simultaneous long-range communications. A physical-layer technique which achieves this is multiple-input multiple-output (MIMO): the use of multiple transmit and receive antennas to multiplex several streams of data and transmit them simultaneously. MIMO was originally developed in the point-to-point setting, where the transmit antennas are collocated at a single transmit node, each transmitting one data stream, and the receive antennas are collocated at a single receive node, jointly processing the vector of received observations at the antennas.

A natural approach to apply this concept to the network setting is to have both source nodes and destination nodes cooperate in *clusters* to form distributed transmit and receive antenna arrays, respectively. In this way, mutually interfering signals can be turned into useful ones that can be jointly decoded at the receive cluster and spatial multiplexing gain can be realized. In fact, if *all* the nodes in the network could cooperate for free, then a classical MIMO result [11], [12] says that a sum-rate scaling proportional to n could be achieved. However, this may be overoptimistic: communication between nodes is required to set up the cooperation and this may drastically reduce the useful throughput. The Aeron–Saligrama scheme is MIMO based and its performance is precisely limited by the cooperation overhead between receive nodes. Our main contribution is to introduce a new *multiscale, hierarchical* cooperation architecture without significant overhead. Such cooperation first takes place between nodes within very small local clusters to facilitate MIMO communication over a larger spatial scale. This can then be used as a communication infrastructure for cooperation within larger clusters at the next level of the hierarchy. Continuing in this fashion, cooperation can be achieved at an almost global scale.

The result for dense networks builds the foundation for understanding extended networks in the low-attenuation regime of the path loss exponent α between 2 and 4. Cooperative MIMO communication provides not only a degree of freedom gain but also a power gain, obtained by combining signals received at the different nodes. This power gain is not very important in the dense setting, since there is already sufficient SNR in any direct communication between individual nodes and the capacity is only *logarithmic* in the SNR. In the extended setting, however, this power gain becomes very important, since the power transferred between an individual source and destination pair vanishes due to channel attenuation. The operation is in the low-SNR regime where the capacity is *linear* in the SNR. Cooperation between nodes can significantly boost the power transfer. In fact, it can be shown that the capacity of long range n by n cooperative MIMO transmission scales exactly like the total received power. This total received power scales like $n^{2-\alpha/2}$. We show that a simple modification of our hierarchical cooperation scheme to the extended setting can achieve a network total throughput arbitrarily close to this cooperative MIMO scaling. Thus, for $\alpha < 3$, our scheme performs strictly better than multihop.

Can we do better? Recall that earlier results in [5], [7]–[10], [2] are all based on upper-bounding the amount of power transferred across cut-sets of the network. It turns out that their upper bounds are tight when $\alpha > 4$ but not tight for α between 2 and 4. By evaluating exactly the scaling of the power transferred, we show that it matches the performance of the hierarchical scheme for α between 2 and 3 and that of the multihop scheme for $\alpha > 3$. More precisely, we obtain the following tight characterization for the scaling exponent for all α in the extended case:

$$e(\alpha) := \lim_{n \rightarrow \infty} \frac{\log C_n(\alpha)}{\log n} = \begin{cases} 2 - \frac{\alpha}{2}, & 2 \leq \alpha \leq 3 \\ \frac{1}{2}, & \alpha > 3 \end{cases}$$

where $C_n(\alpha)$ is the total capacity of the network. In particular, when $\alpha = 2$, linear capacity scaling can be achieved, even in the extended case. Note that the capacity is limited by the power

transferred for *all* $\alpha \geq 2$; hence, extended networks are fundamentally *coverage-limited*, even for α between 2 and 4. For $\alpha > 3$, multihop is sufficient in transferring the optimal amount of power; for $\alpha < 3$, when the attenuation is slower, cooperative MIMO is needed to provide the power gain and also enough degrees of freedom to operate in the power-efficient regime. Just like in the dense setting, interference limitation does not play a significant role, as far as capacity scaling is concerned. Cooperative MIMO takes care of that.

Our approach to the problem is to first look at the dense case to isolate the issue of interference and then to tackle the extended case. But the dense scaling is also of interest in its own right. It is relevant whenever one wants to design networks to serve many nodes, all within communication range of each other (within a campus, an urban block, etc.). This scaling is also a reasonable model to study problems such as *spectrum sharing*, where many users in a geographical area are sharing a wide band of spectrum. Consider the scenario where we segregate the total bandwidth into many orthogonal bands, one for each separate network supporting a *fixed* number of users. As we increase the number of users, the number of such segregated networks increases but the *spectral efficiency*, in bits per second per hertz (bits/s/Hz), does not scale with the *total* number of users. In contrast, if we build one large *ad hoc* network for all the users on the entire bandwidth, then our result says that the spectral efficiency actually increases *linearly* with the number of users. The gain is coming from a *network* effect via cooperation between the many nodes in the system.

The rest of the paper is structured as follows. In Section II, we present the model and discuss the various assumptions. Section III contains the main result for dense networks and an outline of the proposed architecture together with a back-of-the-envelope analysis of its performance. The details of its performance analysis are given in Section IV. Section V characterizes the scaling law for extended networks. Section VI discusses the limitations of the model and the results. Section VII contains our conclusions.

II. MODEL

There are n nodes uniformly and independently distributed in a square of unit area in the dense scaling (Sections III and IV) and a square of area n in the extended scaling (Section V). Every node is both a source and a destination. The sources and destinations are paired up one-to-one in an arbitrary fashion. Each source has the same traffic rate $R(n)$ to send to its destination node and a common average transmit power budget of P Joules per symbol. The total throughput of the system is $T(n) = nR(n)$.¹

We assume that communication takes place over a flat channel of bandwidth W hertz around a carrier frequency of f_c , $f_c \gg W$. The complex baseband-equivalent channel gain between node i and node k at time m is given by

$$H_{ik}[m] = \sqrt{G} r_{ik}^{-\alpha/2} \exp(j\theta_{ik}[m]) \quad (1)$$

¹In the sequel, whenever we say a total throughput $T(n)$ is achievable, we implicitly mean that a rate of $T(n)/n$ is achievable for every source–destination pair.

where r_{ik} is the distance between the nodes, $\theta_{ik}[m]$ is the random phase at time m , uniformly distributed in $[0, 2\pi]$ and $\{\theta_{ik}[m], 1 \leq i \leq n, 1 \leq k \leq n\}$ is a collection of independent and identically distributed (i.i.d.) random processes. The $\theta_{ik}[m]$'s and the r_{ik} 's are also assumed to be independent. The parameters G and $\alpha \geq 2$ are assumed to be constants; α is called the path-loss exponent. For example, under free-space line-of-sight propagation, Friis' formula applies and

$$|H_{ik}[m]|^2 = \frac{G_{Tx} \cdot G_{Rx}}{(4\pi r_{ik}/\lambda_c)^2} \quad (2)$$

so that

$$G = \frac{G_{Tx} \cdot G_{Rx} \cdot \lambda_c^2}{16\pi^2}, \quad \alpha = 2$$

where G_{Tx} and G_{Rx} are the transmitter and receiver antenna gains, respectively, and λ_c is the carrier wavelength.

Note that the channel is random, depending on the location of the users and the phases. The locations are assumed to be fixed over the duration of the communication. The phases are assumed to vary in a stationary ergodic manner (fast fading).² We assume that the channel gains are known at all the nodes. The signal received by node i at time m is given by

$$Y_i[m] = \sum_{k=1}^n H_{ik}[m] X_k[m] + Z_i[m]$$

where $X_k[m]$ is the signal sent by node k at time m and $Z_i[m]$ is white circularly symmetric Gaussian noise of variance N_0 per symbol.

Several comments about the model are in order.

- The path-loss model is based on a *far-field* assumption: the distance r_{ik} is assumed to be much larger than the carrier wavelength. When the distance is of the order or shorter than the carrier wavelength, the simple path-loss model obviously does not hold anymore as path loss can potentially become path “gain.” The reason is that near-field electro-magnetics now come into play.
- The phase $\theta_{ik}[m]$ depends on the distance between the nodes modulo the carrier wavelength. The random-phase model is thus also based on a far-field assumption: we are assuming that the nodes' separation is at a much larger spatial scale compared to the carrier wavelength, so that the phases can be modeled as completely random and independent of the actual positions.
- It is realistic to assume the variation of the phases since they vary significantly when users move a distance of the order of the carrier wavelength (fractions of a meter). The positions determine the path losses and they, on the other hand, vary over a much larger spatial scale. So the positions are assumed to be fixed.
- We essentially assume a line-of-sight type environment and ignore multipath effects. The randomness in phases is sufficient for the long range MIMO transmissions needed

²With more technical efforts, we believe our results can be extended to the slow-fading setting where the phases are fixed as well. See the remark at the end of Appendix I for further discussion on this point.

in our scheme. With multipaths, there is a further randomness due to random constructive and destructive interference of these paths. It can be seen that our results extend to the multipath case.

We will discuss further the limitations of this model in Section VI after we present our results.

III. MAIN RESULT FOR DENSE NETWORKS

We first give an information-theoretic upper bound on the achievable scaling law for the aggregate throughput in the network. Before starting to look for good communication strategies, Theorem 3.1 establishes the best we can hope for.

Theorem 3.1: The aggregate throughput in the network with n nodes is bounded above by

$$T(n) \leq K' n \log n$$

with high probability (i.e., probability going to 1 as n grows) for some constant $K' > 0$ independent of n .

Proof: Consider a source–destination pair (s, d) in the network. The transmission rate $R(n)$ from source node s to destination node d is upper-bounded by the capacity of the single-input multiple-output (SIMO) channel between source node s and the rest of the network. Using a standard formula for this channel (see, e.g., [13, eq. (5.26)]), we get

$$\begin{aligned} R(n) &\leq \log \left(1 + \frac{P}{N_0} \sum_{\substack{i=1 \\ i \neq s}}^n |H_{is}|^2 \right) \\ &= \log \left(1 + \frac{P}{N_0} \sum_{\substack{i=1 \\ i \neq s}}^n \frac{G}{r_{is}^\alpha} \right). \end{aligned}$$

It is easy to see that in a random network with n nodes uniformly distributed on a fixed two-dimensional area, the minimum distance between any two nodes in the network is larger than $\frac{1}{n^{1+\delta}}$ with high probability for large n , for any $\delta > 0$. To see this, consider one specific node in the network which is at distance larger than $\frac{1}{n^{1+\delta}}$ to all other nodes in the network. This is equivalent to saying that there are no other nodes inside a circle of area $\frac{\pi}{n^{2+2\delta}}$ around this node. The probability of such an event is $(1 - \frac{\pi}{n^{2+2\delta}})^{n-1}$. The minimum distance between any two nodes in the network is larger than $\frac{1}{n^{1+\delta}}$ only if this condition is satisfied for all nodes in the network. Thus, by the union bound we have

$$\begin{aligned} P \left(\text{minimum distance in the network is smaller than } \frac{1}{n^{1+\delta}} \right) \\ \leq n \left(1 - \left(1 - \frac{\pi}{n^{2+2\delta}} \right)^{n-1} \right) \end{aligned}$$

which decreases to zero as $1/n^{2\delta}$ with increasing n .

Using this fact on the minimum distance in the network, we obtain

$$R(n) \leq \log \left(1 + \frac{GP}{N_0} n^{\alpha(1+\delta)+1} \right) \leq K' \log n$$

for some constant $K' > 0$ independent of n for all-source destination pairs in the network with high probability. The theorem follows. \square

In the view of what is ultimately possible, established by Theorem 3.1, we are now ready to state the main result of this paper.

Theorem 3.2: Let $\alpha \geq 2$. For any $\epsilon > 0$, there exists a constant $K_\epsilon > 0$ independent of n such that with high probability, an aggregate throughput

$$T(n) \geq K_\epsilon n^{1-\epsilon}$$

is achievable in the network for all possible pairings between sources and destinations.

Theorem 3.2 states that it is actually possible to perform arbitrarily close to the bound given in Theorem 3.1. The two theorems together establish the capacity scaling for the network up to logarithmic terms. Note how dramatically different is this new linear capacity scaling law from the well-known throughput scaling of $\Theta(\sqrt{n})$ implied by [3], [4] for the same model. Note also that the upper bound in Theorem 3.1 assumes a genie-aided removal of interference between simultaneous transmissions from different sources. By proving Theorem 3.2, we will show that it is possible to mitigate such interference without a genie but with cooperation between the nodes.

The proof of Theorem 3.2 relies on the construction of an explicit scheme that realizes the promised scaling law. The construction is based on recursively using the following key lemma, which addresses the case when $\alpha > 2$ and whose proof is relegated to Section IV.

Lemma 3.1: Consider $\alpha > 2$ and a network with n nodes subject to interference from external sources. The signal received by node i is given by

$$Y_i = \sum_{k=1}^n H_{ik} X_k + Z_i + I_i$$

where I_i is the external interference signal received by node i . Assume that $\{I_i, 1 \leq i \leq n\}$ is a collection of uncorrelated zero-mean stationary and ergodic random processes with power P_{I_i} upper-bounded by

$$P_{I_i} \leq K_I, \quad 1 \leq i \leq n$$

for a constant $K_I > 0$ independent of n . Let us assume there exists a scheme such that for each n , with probability at least $1 - e^{-n^{c_1}}$ achieves an aggregate throughput

$$T(n) \geq K_1 n^b$$

for every possible source–destination pairing in this network of n nodes. K_1 and c_1 are positive constants independent of n and the source–destination pairing, and $0 \leq b < 1$. Let us also assume that the per-node average power budget required to realize this scheme is upper-bounded by P/n as opposed to P .

Then one can construct another scheme for this network that achieves a *higher* aggregate throughput

$$T(n) \geq K_2 n^{\frac{1}{2-b}}$$

for every source–destination pairing in the network, where $K_2 > 0$ is another constant independent of n and the pairing.

Moreover, the failure rate for the new scheme is upper-bounded by $e^{-n^{c_2}}$ for another positive constant c_2 while the per-node average power needed to realize the scheme is also upper-bounded by P/n .

Lemma 3.1 is the key step to build a hierarchical architecture. Since $\frac{1}{2-b} > b$ for $0 \leq b < 1$, the new scheme is always better than the old. We will now give a rough description of how the new scheme can be constructed given the old scheme, as well as a back-of-the-envelope analysis of the scaling law it achieves. The following section is devoted to its precise description and performance analysis.

The constructed scheme is based on clustering and long-range MIMO transmissions between clusters. We divide the network into clusters of M nodes. Let us focus for now on a particular source node s and its destination node d . s will send M bits to d in three steps.

- 1) Node s will distribute its M bits among the M nodes in its cluster, one for each node.
- 2) These nodes together can then form a distributed transmit antenna array, sending the M bits *simultaneously* to the destination cluster where d lies.
- 3) Each node in the destination cluster obtained one observation from the MIMO transmission, and it quantizes and ships the observation back to d , which can then do joint MIMO processing of all the observations and decode the M transmitted bits.

From the network point of view, all source–destination pairs have to eventually accomplish these three steps. Step 2 is long-range communication and only one source–destination pair can operate at a time. Steps 1 and 3 involve local communication and can be parallelized across source–destination pairs. Combining all this leads to three phases in the operation of the network:

Phase 1. Setting Up Transmit Cooperation: Clusters work in parallel. Within a cluster, each source node has to distribute M bits to the other nodes, 1 bit for each node, such that at the end of the phase, each node has 1 bit from each of the source nodes in the same cluster. Since there are M source nodes in each cluster, this gives a traffic demand of exchanging $M(M-1) \sim M^2$ bits. (Recall our assumption that each node is a source for some communication request and destination for another.) The key observation is that this is similar to the original problem of communicating between n source and destination pairs, but on a network of size M . More specifically, this traffic demand of exchanging M^2 bits is handled by setting up M subphases, and assigning M source–destination pairs for each subphase. Since our channel model is scale invariant, note that the scheme given in the hypothesis of Lemma 3.1 can be used in each subphase by simply scaling down the power with cluster area. Having aggregate throughput M^b , each subphase is completed in M^{1-b} time slots while the whole phase takes M^{2-b} time slots. See Fig. 1.

Phase 2. MIMO Transmissions: We perform successive long-distance MIMO transmissions between source–destination pairs, one at a time. In each of the MIMO transmissions, say one between s and d , the M bits of s are simultaneously transmitted by the M nodes in its cluster to the M nodes in the cluster of d . Each of the long-distance MIMO transmissions are

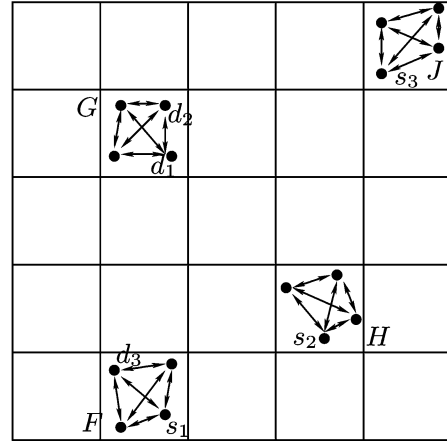


Fig. 1. Nodes inside clusters $F, G, H,$ and J are illustrated while exchanging bits in Phases 1 and 3. Note that in Phase 1 the exchanged bits are the source bits whereas in Phase 3 they are the quantized MIMO observations. Clusters work in parallel. In this and in Fig. 2, we highlight three source–destination pairs $s_1 - d_1, s_2 - d_2,$ and $s_3 - d_3,$ such that nodes s_1 and d_3 are located in $F,$ nodes s_2 and s_3 are located in H and J respectively, and nodes d_1 and d_2 are located in G .

repeated for each source–destination pair in the network, hence we need n time slots to complete the phase. See Fig. 2.

Phase 3. Cooperate to Decode: Clusters work in parallel. Since there are M destination nodes inside the clusters, each cluster received M MIMO transmissions in Phase 2, one intended for each of the destination nodes in the cluster. Thus, each node in the cluster has M received observations, one from each of the MIMO transmissions, and each observation is to be conveyed to a different node in its cluster. Nodes quantize each observation into fixed Q bits so there are now a total of at most QM^2 bits to exchange inside each cluster. Using exactly the same scheme as in Phase 1, we conclude the phase in QM^{2-b} time slots. See again Fig. 1.

Assuming that each destination node is able to decode the transmitted bits from its source node from the M quantized signals it gathers by the end of Phase 3, we can calculate the rate of the scheme as follows: Each source node is able to transmit M bits to its destination node, hence nM bits in total are delivered to their destinations in $M^{2-b} + n + QM^{2-b}$ time slots, yielding an aggregate throughput of

$$\frac{nM}{M^{2-b} + n + QM^{2-b}}$$

bits per time slot. Maximizing this throughput by choosing $M = n^{\frac{1}{2-b}}$ yields $T(n) = \frac{1}{2+Q} n^{\frac{1}{2-b}}$ for the aggregate throughput which is the result in Lemma 3.1.

Clusters can work in parallel in Phases 1 and 3 because for $\alpha > 2$, the aggregate interference at a particular cluster caused by other active nodes is bounded; moreover, the interference signals received by different nodes in the cluster are zero-mean and uncorrelated satisfying the assumptions of Lemma 3.1. For $\alpha = 2$, the aggregate interference scales like $\log n$, leading to a slightly different version of the lemma, whose proof is given towards the end of Section IV.

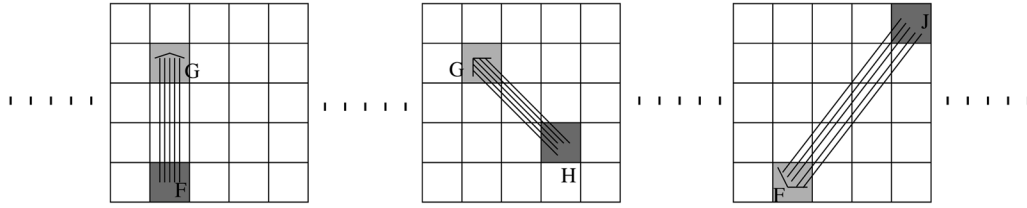


Fig. 2. Successive MIMO transmissions are performed between clusters. The first figure depicts MIMO transmission from cluster F to G , where bits originally belonging to s_1 are simultaneously transmitted by all nodes in F to all nodes in G . The second MIMO transmission is from H to G , while now bits of source node s_2 are transmitted by nodes in H to nodes in G . The third graph illustrates MIMO transmission from cluster J to F .

Lemma 3.2: Consider $\alpha = 2$ and a network with n nodes subject to interference from external sources. The signal received by node i is given by

$$Y_i = \sum_{k=1}^n H_{ik} X_k + Z_i + I_i$$

where I_i is the external interference signal received by node i . Assume that $\{I_i, 1 \leq i \leq n\}$ is a collection of uncorrelated zero-mean stationary and ergodic random processes with power P_{I_i} upper-bounded by

$$P_{I_i} \leq K_I \log n, \quad 1 \leq i \leq n$$

for a constant $K_I \geq 0$ and independent of n . Let us assume there exists a scheme such that for each n with failure probability at most $e^{-n^{c_1}}$, achieves an aggregate throughput

$$T(n) \geq K_1 \frac{n^b}{\log n}$$

for every source–destination pairing in this network. K_1 and c_1 are positive constants independent of n and the source–destination pairing, and $0 \leq b < 1$. Let us also assume that the average power budget required to realize this scheme is upper-bounded by P/n , as opposed to P .

Then one can construct another scheme for this network that achieves a *higher* aggregate throughput scaling

$$T(n) \geq K_2 \frac{n^{\frac{1}{2-b}}}{(\log n)^2}$$

for every source–destination pairing, where $K_2 > 0$ is another constant independent of n and the pairing. Moreover, the failure rate for the new scheme is upper-bounded by $e^{-n^{c_2}}$ for another positive constant c_2 while the per-node average power needed to realize the scheme is also upper-bounded by P/n .

We can now use Lemmas 3.1 and 3.2 to prove Theorem 3.2.

Proof of Theorem 3.2: We only focus on the case of $\alpha > 2$. The case of $\alpha = 2$ proceeds similarly.

We start by observing that the simple scheme of transmitting directly between the source–destination pairs one at a time (time-division multiple access (TDMA)) satisfies the requirements of Lemma 3.1. The aggregate throughput is $\Theta(1)$, so $b = 0$. The failure probability is 0. Since each source is only transmitting $\frac{1}{n}$ th of the time and the distance between the source and its destination is bounded, the average power consumed per node is of the order of $\frac{1}{n}$.

As soon as we have a scheme to start with, Lemma 3.1 can be applied recursively, yielding a scheme that achieves higher throughput at each step of the recursion. More precisely, starting with a TDMA scheme with $b = 0$ and applying Lemma 3.1 recursively h times, one gets a scheme achieving $\Theta(n^{\frac{h}{h+1}})$ aggregate throughput. Given any $\epsilon > 0$, we can now choose h such that $\frac{h}{h+1} \geq 1 - \epsilon$ and we get a scheme that achieves $\Theta(n^{1-\epsilon})$ aggregate throughput scaling with high probability. This concludes the proof of Theorem 3.2. \square

Gathering everything together, we have built a hierarchical scheme to achieve the desired throughput. At the lowest level of the hierarchy, we use the simple TDMA scheme to exchange bits for cooperation among small clusters. Combining this with longer range MIMO transmissions, we get a higher throughput scheme for cooperation among nodes in larger clusters at the next level of the hierarchy. Finally, at the top level of the hierarchy, the cooperation clusters are almost the size of the network and the MIMO transmissions are over the global scale to meet the desired traffic demands. Fig. 3 shows the resulting hierarchical scheme with a focus on the top two levels.

It is important to understand the aspects of the channel model which the scheme made use of in achieving the linear capacity scaling:

- the random channel phases enable the long-range MIMO transmissions;
- the path attenuation decay law $1/r^\alpha$ ($\alpha \geq 2$) ensures that the *aggregate* signals from far away nodes are much weaker than signals from close-by nodes; this enables spatial reuse.

Note that the second property is exactly the same one which allows multihop schemes to achieve the \sqrt{n} -scaling in the paper by Gupta–Kumar [3] and in many others after that. Although the gain between nearby nodes becomes unbounded as $n \rightarrow \infty$ in the model, the received signal-to-interference-plus-noise ratio (SINR) is always bounded in the scheme at all levels of the hierarchy. The scheme does not communicate with unbounded SINR, although it is possible in the model.

IV. DETAILED DESCRIPTION AND PERFORMANCE ANALYSIS

In this section, we concentrate in more detail on the scheme that proves Lemmas 3.1 and 3.2. We first focus on Lemma 3.1 and then extend the proof to Lemma 3.2. As we have already seen in the previous section, we start by dividing the unit square into smaller squares of area $A_c = \frac{M}{n}$. Since the node density is n , there will be on average M nodes inside each of these small

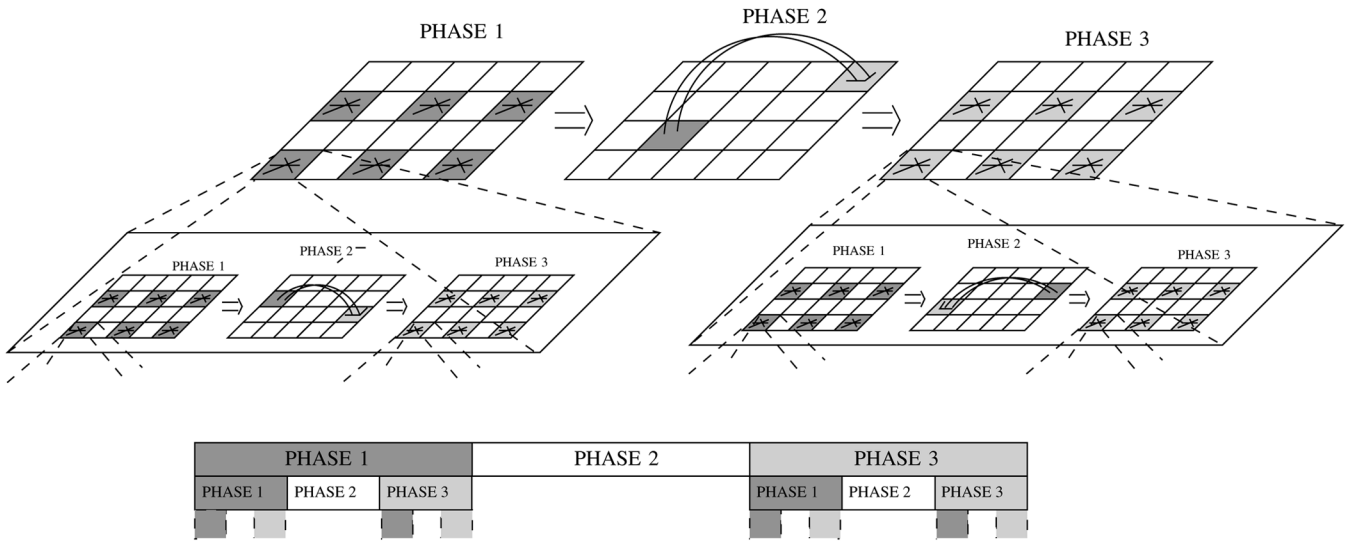


Fig. 3. The upper graph illustrates the salient features of the three-phase hierarchical scheme. The time division in this hierarchical scheme is explicitly given the graph below.

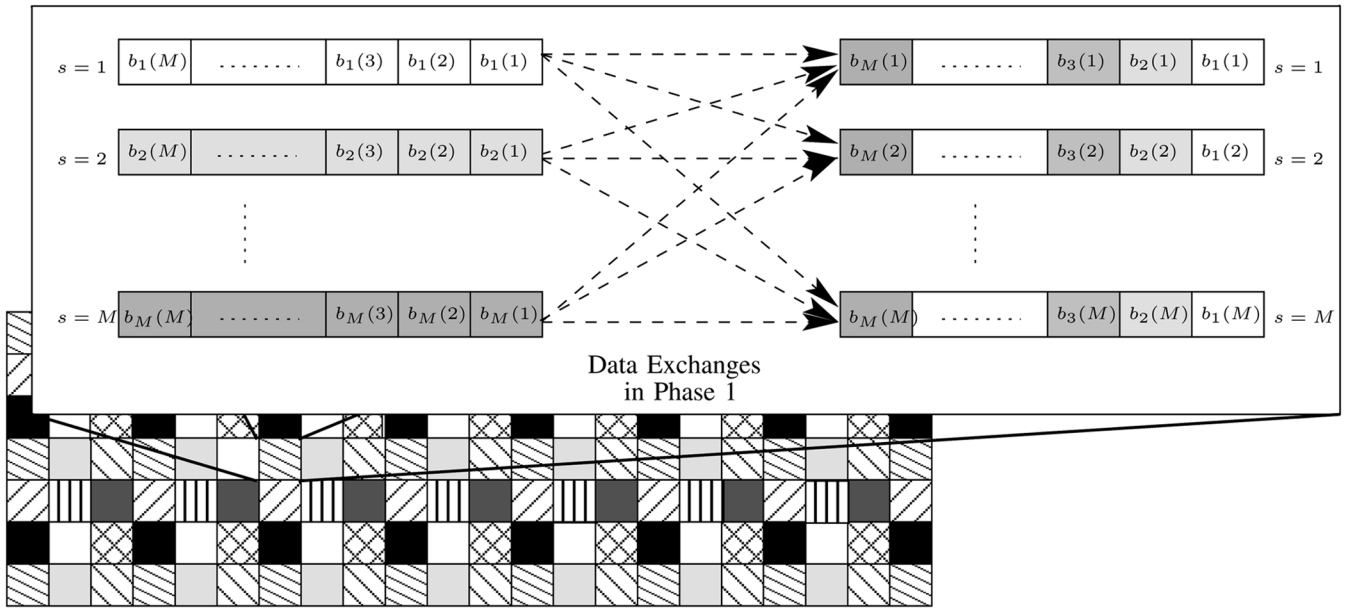


Fig. 4. Buffers of the nodes in a cluster are illustrated before and after the data exchanges in Phase 1. The data stream of the source nodes are distributed to the M nodes in the network as depicted. $b_s(j)$ denotes the j th subblock of the source node s . Note the 9-TDMA scheme that is employed over the network in this phase.

squares. The following lemma upper-bounds the probability of having large deviations from the average. Its proof is relegated to the end of the section.

Lemma 4.1: Let us partition a unit area network of size n into cells of area A_c , where A_c can be a function of n . The number of nodes inside each cell is between $((1 - \delta)A_c n, (1 + \delta)A_c n)$ with probability larger than $1 - \frac{1}{A_c} e^{-\Lambda(\delta)A_c n}$ where $\Lambda(\delta)$ is independent of n and satisfies $\Lambda(\delta) > 0$ when $\delta > 0$.

Applying Lemma 4.1 to the squares of area M/n , we see that all squares contain order M nodes with probability larger than $1 - \frac{n}{M} e^{-\Lambda(\delta)M}$. We assume $M = n^\gamma$, where $0 < \gamma \leq 1$, in which case this probability tends to 1 as n increases. This condition is sufficient for the following analysis on scaling laws to hold. However, in order to simplify the presentation, we assume that

there are exactly M nodes in each square. The clustering is used to realize a distributed MIMO system in three successive steps.

Phase 1. Setting Up Transmit Cooperation: In this phase, source nodes distribute their data streams over their clusters and set up the stage for the long-range MIMO transmissions that we want to perform in the next phase. Clusters work in parallel according to the 9-TDMA scheme depicted in Fig. 4, which divides the total time for this phase into nine time slots and assigns simultaneous operation to clusters that are sufficiently separated. The nine different patterns used to color the clusters in Fig. 4 correspond to these nine-time slots. The clusters with the same pattern are operating simultaneously in the same time slot while the other clusters stay inactive. Note that with this scheduling, in every time slot there are at least two inactive clusters between any two clusters that are active.

Let us focus on one specific source node s located in cluster S with destination node d in cluster D . Node s will divide a block of length LM bits of its data stream into M subblocks each of length L bits, where L can be arbitrarily large but bounded. The destination of each subblock in Phase 1 depends on the relative position of clusters S and D .

- 1) If S and D are either the same cluster or are not neighboring clusters: one subblock is to be kept in s and the remaining $M - 1$ subblocks are to be transmitted to the other $M - 1$ nodes located in S , one subblock for each node.
- 2) If S and D are neighboring clusters: Divide the cluster S into two halves, each of area $A_c/2$, one half located close to the border with D and the other half located farther from D . The M subblocks of source node s are to be distributed to the $M/2$ nodes located in the second half cluster (farther from D), each node gets two subblocks.³

Since the above traffic is required for every source node in cluster S , we end up with a highly uniform traffic demand of delivering $M \times LM$ bits in total to their destinations. A key observation is that the problem can be separated into subproblems, each similar to our original problem, but on a network size M and area A_c . More specifically, the traffic of transporting LM^2 bits can be handled by organizing M sessions and assigning M source–destination pairs for each session. (Note that due to the nonuniformity arising from point 2) above, one might be able to assign only $M/2$ source–destination pairs in a session and hence need to handle the traffic demand of transporting LM^2 bits by organizing up to $2M$ sessions in the extreme case instead of M .) The assigned source–destination pairs in each session can then communicate L bits. Since our channel model is scale invariant, the scheme in the hypothesis of Lemma 3.1 can be used to handle the traffic in each session, by simply scaling down the powers of the nodes by $(A_c)^{\alpha/2}$. Hence, the power used by each node will be bounded above by $\frac{P(A_c)^{\alpha/2}}{M}$. The scheme is to be operated simultaneously inside all the clusters in the 9-TDMA scheme, so we need to ensure that the resultant inter-cluster interference satisfies the properties in Lemma 3.1.

Lemma 4.2: Consider clusters of size M and area A_c operating according to 9-TDMA scheme in Fig. 4 in a network of size n . Let each node be constrained to an average power $\frac{P(A_c)^{\alpha/2}}{M}$. For $\alpha > 2$, the interference power received by a node from the simultaneously operating clusters is upper-bounded by a constant K_{I_1} independent of n . For $\alpha = 2$, the interference-power is bounded by $K_{I_2} \log n$ for K_{I_2} independent of n . Moreover, the interference signals received by different nodes in the cluster are zero-mean and uncorrelated.

The proof of this lemma is given at the end of the present section. Let us for now concentrate on the case $\alpha > 2$. By Lemma 4.2, the inter-cluster interference has bounded power and is uncorrelated across different nodes. Thus, the strategy in the hypothesis of Lemma 3.1 can achieve an aggregate rate $K_1 M^b$ in each session for some $K_1 > 0$, with probability larger than $1 - e^{-M^{c_1}}$. Using the union bound, with probability larger than

³Note that at this point, we need to show that there are $\Theta(M/2)$ nodes inside the half clusters, which is a stronger result than the earlier discussion on having $\Theta(M)$ nodes in each cluster. The result follows similarly from Lemma 4.1 together with the union bound.

$1 - 2ne^{-M^{c_1}}$, the aggregate rate $K_1 M^b$ is achieved inside all sessions in all clusters in the network. (Recall that the number of sessions in one cluster can be $2M$ in the extreme case and there are n/M clusters in total.) With this aggregate rate, each session can be completed in at most $(L/K_1)M^{1-b}$ channel uses and $2M$ successive sessions are completed in $(2L/K_1)M^{2-b}$ channel uses. Using the 9-TDMA scheme, the phase is completed in less than $(18L/K_1)M^{2-b}$ channel uses all over the network with probability larger than $1 - 2ne^{-M^{c_1}}$.

Phase 2. MIMO Transmissions: In this phase, we are performing the actual MIMO transmissions for all the source–destination pairs serially, i.e., one at a time. A MIMO transmission from source s to destination d involves the M (or $M/2$) nodes in the cluster S containing s (referred to as the source cluster for this MIMO transmission) to the M (or $M/2$) nodes of the cluster D containing d (referred to as the destination cluster of the MIMO transmission).

Let the distance between the midpoints of the two clusters be r_{SD} . If S and D are the same cluster, we skip the step for this source–destination pair $s-d$. Otherwise, we operate in two slightly different modes depending on the relative positions of S and D . Each mode is a continuation of the operations performed in the first phase. First consider the case where S and D are not neighboring clusters. In this case, the M nodes in cluster S independently encode the L bits-long subblocks they possess, originally belonging to node s , into C symbols by using a randomly generated Gaussian code \mathcal{C} that respects an average transmit power constraint $\frac{P(r_{SD})^\alpha}{M}$. The nodes then transmit their encoded sequences of length C symbols simultaneously to the M nodes in cluster D . The nodes in cluster D quantize the signals they observe during the C transmissions and store these quantized signals (that we will simply refer to as *observations* in the following text), without trying to decode the transmitted symbols. In the case where S and D are neighbors, the strategy is slightly modified so that the MIMO transmission is from the $M/2$ nodes in S , that possess the subblocks of s after Phase 1, to the $M/2$ nodes in D that are located in the farther half of the cluster to S . Each of these $M/2$ nodes in S possess two subblocks that come from s . They encode each subblock into C symbols by again using a Gaussian code of power $\frac{P(r_{SD})^\alpha}{M}$. The nodes then transmit the $2C$ symbols to the $M/2$ nodes in D that, in turn, sample their received signals and store the observations. The observations accumulated at various nodes in D at the end of this step are to be conveyed to node d during the third phase.

After concluding the step for the pair $s-d$, the phase continues by repeating the same step for the next source node $s + 1$ in S and its destination d' . Note that the destination cluster for this new MIMO transmission is, in general, a different cluster D' , which is the one that contains the destination node d' . The MIMO transmissions are repeated until the data originated from all source nodes in the network are transmitted to their respective destination clusters. Since the step for one source–destination pair takes either C or $2C$ channel uses, completing the operation for all n source nodes in the network requires at most $2C \times n = 2Cn$ channel uses.

Phase 3. Cooperate to Decode: In this phase, we aim to provide each destination node, the observations of the symbols that have been originally intended for it. With the MIMO transmis-

sions in the second phase, these observations have been accumulated at the nodes of its cluster. As before, let us focus on a specific destination node d located in cluster D . Note that depending on whether the source node of d is located in a neighboring cluster or not, either each of the M nodes in D have C observations intended for d , or $M/2$ of the nodes have $2C$ observations each. Note that these observations are some real numbers that need to be quantized and encoded into bits before being transmitted. Let us assume that we are encoding each block of C observations into CQ bits, by using fixed Q bits per observation on the average. The situation is symmetric for all M destination nodes in D , since the cluster received M MIMO transmissions in the previous phase, one for each destination node. (The destination nodes that have source nodes in D are an exception. Recall from Phase 1 and Phase 2 that in this case, each node in D possesses subblocks of the original data stream for the destination node, not MIMO observations. We will ignore this case by simply assuming $L \leq CQ$ in the below computation.) The arising traffic demand of transporting $M \times CQM$ bits in total is similar to Phase 1 and can be handled by using exactly the same scheme in less than $(2CQ/K_1)M^{2-b}$ channel uses. Recalling the discussion on the first phase, we conclude that the phase can be completed in less than $(18CQ/K_1)M^{2-b}$ channel uses all over the network with probability larger than $1 - 2ne^{-M^{c_1}}$.

Note that if it were possible to encode each observation into fixed Q bits without introducing any distortion, which is obviously not the case, the following lemma on MIMO capacity, would suggest that with the Gaussian code \mathcal{C} used in Phase 2 satisfying $L/C \geq \kappa$ for some constant $\kappa > 0$, the transmitted bits could be recovered by an arbitrarily small probability of error from the observations gathered by the destination nodes at the end of Phase 3. The lemma is proven in Appendix I.

Lemma 4.3: The mutual information achieved by the $M \times M$ MIMO transmission between any two clusters grows at least linearly with M .

The following lemma states that there is actually a way to encode the observations using a fixed number of bits per observation and at the same time, not to degrade the performance of the overall channel significantly, that is, to still get a linear capacity growth for the resulting *quantized MIMO* channel.

Lemma 4.4: There exists a strategy to encode the observations at a fixed rate Q bits per observation and get a linear growth of the mutual information for the resultant $M \times M$ quantized MIMO channel.

We leave the proof of the lemma to the end of Appendix II. However, the following small lemma, whose proof is given at the end of the present section, may provide motivation for the stated result. Lemma 4.5 points out a key observation on the way we choose our transmit powers in the MIMO phase. It is central to the proof of Lemma 4.4 and states that the observations have bounded power, that does not scale with M . This in turn suggests that one can use a fixed number of bits to encode them without degrading the scaling performance of the scheme.

Lemma 4.5: In Phase 2, the power received by each node in the destination cluster is bounded below and above by constants P_1 and P_2 , respectively, that are independent of M .

Putting it together, we have seen that the three phases described effectively realize virtual MIMO channels achieving spatial multiplexing gain M between the source and destination nodes in the network. Using these virtual MIMO channels, each source is able to transmit ML bits in

$$\begin{aligned} T_t &= T(\text{Phase 1}) + T(\text{Phase 2}) + T(\text{Phase 3}) \\ &= \frac{18L}{K_1}M^{2-b} + 2Cn + \frac{18CQ}{K_1}M^{2-b} \end{aligned}$$

total channel uses where $L/C \geq \kappa$ for some $\kappa > 0$ independent of M (or n). This gives an aggregate throughput of

$$\begin{aligned} T(n) &= \frac{nML}{(18L/K_1)M^{2-b} + 2Cn + (18CQ/K_1)M^{2-b}} \\ &\geq K_2 n^{\frac{1}{2-b}} \end{aligned} \tag{3}$$

for some $K_2 > 0$ independent of n , by choosing $M = n^{\frac{1}{2-b}}$ with $0 \leq b < 1$, which is the optimal choice for the cluster size as a function of b . A failure arises if there are not order M nodes in each cluster or the scheme used in Phases 1 and 3 fails to achieve the promised throughput. Combining the result of Lemma 4.1 with the computed failure probabilities for Phases 1 and 3 yields

$$P_f \leq 4ne^{-M^{c_1}} + \frac{n}{M}e^{-\Lambda(\delta)M} \leq e^{-n^{c_2}}$$

for some $c_2 > 0$.

Next, we show that per-node average power used by the new scheme is also bounded above by P/n : for Phases 1 and 3, we know that the scheme employed inside the clusters uses average per-node power bounded above by $PA_c^{\alpha/2}/M$. Indeed, $A_c = M/n$, and for $\alpha \geq 2$ we have

$$\frac{PA_c^{\alpha/2}}{M} = \frac{P}{M} \left(\frac{M}{n}\right)^{\alpha/2} = \frac{P}{n} \left(\frac{M}{n}\right)^{\alpha/2-1} \leq \frac{P}{n}.$$

In Phase 2, each node is transmitting with power $\frac{P(r_{SD})^\alpha}{M}$ in at most a fraction M/n of the total duration of the phase, while keeping silent during the rest of the time. This yields a per-node average power $\frac{P(r_{SD})^\alpha}{n}$. Recall that r_{SD} is the distance between the midpoints of the source and destination clusters and $r_{SD} < 1$, which yields the upper bound P/n on the per-node average power also for the second phase.

In order to conclude the proof of Lemma 3.1, we should note that the new scheme achieves the same aggregate throughput scaling when the network experiences interference from the exterior. In Phases 1 and 3, this external interference with bounded power will simply add to the inter-cluster interference experienced by the nodes. For the MIMO phase, this will result in uncorrelated background-noise-plus-interference at the receiving nodes which is not necessarily Gaussian. In Appendices I and II we prove the results stated in Lemmas 4.3 and 4.4 for this more general case. This concludes the proof of Lemma 3.1. \square

Proof of Lemma 3.2: The scheme that proves Lemma 3.2 is quite similar to the one described above. Lemma 4.2 states that when $\alpha = 2$, the inter-cluster interference power experienced during Phases 1 and 3 is upper-bounded by $K'_{I_2} \log n = K'_{I_2} \log M$. According to the assumptions of Lemma 3.2, there is furthermore the external interference with power bounded by

$K_I \log n$ that is adding to the inter-cluster interference. Under these conditions, the scheme in the hypothesis of Lemma 3.2 achieves an aggregate rate $K_1 \frac{M^b}{\log M}$ when used to handle the traffic in Phases 1 and 3. For the second phase, we have the following lemma which provides a lower bound on the spatial multiplexing gain of the quantized MIMO channel under the interference experienced. The proof of the lemma is relegated to the end of Appendix II.

Lemma 4.6: Let the MIMO signal received by the nodes in the destination cluster be corrupted by an interference of power $K_I \log M$, uncorrelated over different nodes, and independent of the transmitted signals. There exists a strategy to encode these corrupted observations at a fixed rate Q bits per observation and get an $M/\log M$ growth of the mutual information for the resulting $M \times M$ quantized MIMO channel.

A capacity of $M/\log M$ for the resulting MIMO channel implies that there exists a code \mathcal{C} that encodes L -bits-long subblocks into $C \log M$ symbols, where $L/C \geq \kappa'$ for a constant $\kappa' > 0$, so that the transmitted bits can be decoded at the destination nodes with arbitrarily small probability of error for L and C sufficiently large. Hence, starting again with a block of LM bits in each source node, the LM^2 bits in the first phase can be delivered in $(L/K_1)M^{2-b} \log M$ channel uses. In the second phase, the L -bits-long subblocks now need to be encoded into $C \log M$ symbols, hence the transmission for each source–destination pair takes $C \log M$ channel uses, the whole phase taking $Cn \log M$ channel uses. Note that there are now $CM^2 \log M$ observations encoded into $CQM^2 \log M$ bits that need to be transported in the third phase. With the scheme of aggregate rate $K_1 \frac{M^b}{\log M}$, we need $(CQ/K_1)M^{2-b}(\log M)^2$ channel uses to complete the phase. Choosing $M = n^{\frac{1}{2-b}}$ gives an aggregate throughput of $K_2 n^{\frac{1}{2-b}}/(\log n)^2$ for the new scheme. This concludes the proof of Lemma 3.2. \square

We continue with the proofs of the lemmas introduced in the section.

Proof of Lemma 4.1: The proof of the lemma is a standard application of Chebyshev's inequality. Note that the number of nodes in a given cell is a sum of n i.i.d. Bernoulli random variables B_i , such that $\mathbb{P}(B_i = 1) = A_c$. Hence

$$\begin{aligned} & \mathbb{P}\left(\sum_{i=1}^n B_i \geq (1+\delta)A_c n\right) \\ &= \mathbb{P}\left(e^s \sum_{i=1}^n B_i \geq e^{s(1+\delta)} A_c n\right) \\ &\leq (\mathbb{E}[e^{sB_1}])^n e^{-s(1+\delta)A_c n} \\ &= (e^s A_c + (1 - A_c))^n e^{-s(1+\delta)A_c n} \\ &\leq e^{-A_c n(s(1+\delta) - e^s + 1)} \\ &= e^{-A_c n \Lambda_+(\delta)} \end{aligned}$$

where $\Lambda_+(\delta) = (1+\delta)\log(1+\delta) - \delta$ by choosing $s = \log(1+\delta)$. The proof for the lower bound follows similarly by considering the random variables $-B_i$. The conclusion follows from the union bound. \square

Proof of Lemma 4.2: Consider a node v in cluster V operating under the 9-TDMA scheme in Fig. 5. The interfering

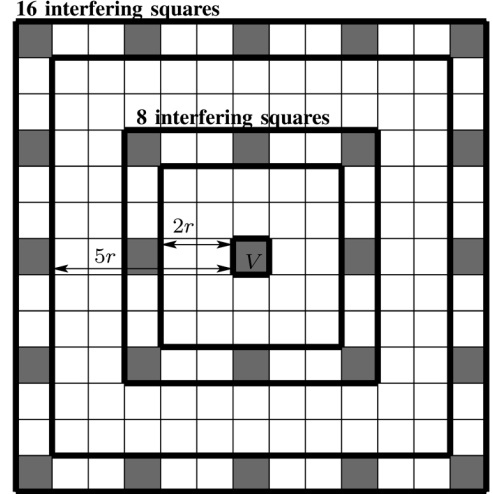


Fig. 5. Grouping of interfering clusters in the 9-TDMA Scheme.

signal received by this node from the simultaneously operating clusters \mathcal{U}_V is given by

$$I_v = \sum_{U \in \mathcal{U}_V} \sum_{j \in U} H_{vj} X_j$$

where H_{vj} are the channel coefficients given by (1) and X_j is the signal transmitted by node j which is located in a simultaneously operating cluster U . First note that the signals I_v and $I_{v'}$ received by two different nodes v and v' in V are uncorrelated since the channel coefficients H_{vj} and $H_{v'j}$ are independent for all j . The power of the interfering signal I_v is given by

$$P_I = \sum_{U \in \mathcal{U}_V} \sum_{j \in U} \frac{GP_j}{(r_{vj})^\alpha}$$

by using the fact that channel coefficients corresponding to different nodes j are independent. As illustrated by Fig. 5, the interfering clusters \mathcal{U}_V can be grouped based on their distance to V such that each group $\mathcal{U}_V(i)$ contains $8i$ clusters or less and all clusters in group $\mathcal{U}_V(i)$ are separated by a distance larger than $(3i-1)\sqrt{A_c}$ from V for $i = 1, 2, \dots$ where A_c is the cluster area. The number of such groups can be simply bounded by the number of clusters n/M in the network. Thus

$$\begin{aligned} P_I &< \sum_{i=1}^{n/M} \sum_{U \in \mathcal{U}_V(i)} \sum_{j \in U} \frac{GP_j}{((3i-1)\sqrt{A_c})^\alpha} \\ &\leq \sum_{i=1}^{n/M} 8i \frac{GP}{(3i-1)^\alpha} \end{aligned} \quad (4)$$

where we have used the fact that the powers of the signals are bounded by $P_j \leq PA_c^{\alpha/2}/M, \forall j$. The sum in (4) is convergent for $\alpha > 2$, thus is bounded by a constant K_{I_1} . For $\alpha = 2$, the sum can be bounded by $K_{I_2} \log n$ where K_{I_2} is a constant independent of n . \square

Proof of Lemma 4.5: We consider only the case where the source cluster S and the destination cluster D are not neighbors. The argument for the other case follows similarly. The signal

received by a destination node d located in cluster D during MIMO transmission from source cluster S is given by

$$Y_d = \sum_{s=1}^M H_{ds} X_s + Z_d$$

where X_s is the signal sent by a source node $s \in S$ constrained to power $\frac{P(r_{SD})^\alpha}{M}$ and Z_d is $\sim \mathcal{N}_C(0, N_0)$. The power of this signal is given by

$$\begin{aligned} \mathbb{E}[|Y_d|^2] &= \sum_{s=1}^M |H_{ds}|^2 \frac{P(r_{SD})^\alpha}{M} + N_0 \\ &= \sum_{s=1}^M \frac{GP}{M} \left(\frac{r_{SD}}{r_{sd}} \right)^\alpha + N_0 \end{aligned}$$

where we use the fact that all H_{ds} , X_s , and Z_d are independent. Observe that

$$r_{SD} - \sqrt{2A_c} \leq r_{sd} \leq r_{SD} + \sqrt{2A_c}$$

while $r_{SD} \geq 2\sqrt{A_c}$. These two relations yield

$$\left(\frac{\sqrt{2}}{\sqrt{2}+1} \right)^\alpha \leq \left(\frac{r_{SD}}{r_{sd}} \right)^\alpha \leq \left(\frac{\sqrt{2}}{\sqrt{2}-1} \right)^\alpha$$

which in turn yields the following lower and upper bounds for the received power at each destination node:

$$\begin{aligned} P_1 &\equiv \left(\frac{\sqrt{2}}{\sqrt{2}+1} \right)^\alpha GP + N_0 \\ &\leq \mathbb{E}[|Y_d|^2] \leq \left(\frac{\sqrt{2}}{\sqrt{2}-1} \right)^\alpha GP + N_0 \equiv P_2. \end{aligned} \quad (5)$$

□

V. EXTENDED NETWORKS

A. Bursty Hierarchical Scheme Does Better Than Multihop for $\alpha < 3$

So far, we have considered *dense* networks, where the total geographical area is fixed and the density of nodes increasing. Another natural scaling is the *extended* case, where the density of nodes is fixed and the area is increasing, a $\sqrt{n} \times \sqrt{n}$ square. This models the situation where we want to scale the network to cover an increasing geographical area.

As compared to dense networks, the distance between nodes is increased by a factor of \sqrt{n} , and hence for the same transmit powers, the received powers are all decreased by a factor of $n^{\alpha/2}$. Equivalently, by rescaling space, an extended network can just be considered as a dense network on a unit area but with the average power constraint per node reduced to $P/n^{\alpha/2}$ instead of P .

Lemmas 3.1 and 3.2 state that the average power per node required to run our hierarchical scheme in dense networks is not the full power P but P/n . In light of the observation above, this immediately implies that when $\alpha = 2$, we can directly apply our scheme to extended networks and achieve a *linear* scaling. For

extended networks with $\alpha > 2$, our scheme would not satisfy the equivalent power constraint $P/n^{\alpha/2}$ and we are now in the power-limited regime (as opposed to the degrees-of-freedom-limited regime). However, we can consider a simple “bursty” modification of the hierarchical scheme which runs the hierarchical scheme a fraction

$$\frac{1}{n^{\alpha/2-1}}$$

of the time with power P/n per node and remains silent for the rest of the time. This meets the given average power constraint of $P/n^{\alpha/2}$, and achieves an aggregate throughput of

$$\frac{1}{n^{\alpha/2-1}} \cdot n^{1-\epsilon} = n^{2-\alpha/2-\epsilon} \text{ bits/second.}$$

(The idea of using burstiness in improving the low-SNR performance of relay networks was introduced in [14] in the context of single-relay networks.)

Note that the quantity $n^{2-\alpha/2} = n^2 \cdot n^{-\alpha/2}$ can be interpreted as the total power transferred between a size n transmit cluster and a size n receive cluster, n^2 node pairs in all, with a power attenuation of $n^{-\alpha/2}$ for each node pair. This power transfer is taking place at the top level of the hierarchy (see Fig. 3). The fact that the achievable rate is proportional to the power transfer further emphasizes that our scheme is power-limited rather than degrees-of-freedom-limited in extended networks.

Let us compare our scheme to multihop. For $\alpha < 3$, it performs strictly better than multihop, while for $\alpha > 3$, it performs worse. Summarizing these observations, we have the following achievability theorem for extended networks, the counterpart to Theorem 3.2 for dense networks.

Theorem 5.1: Consider an extended network on a $\sqrt{n} \times \sqrt{n}$ square. There are two cases.

- $2 \leq \alpha < 3$: For every $\epsilon > 0$, with high probability, an aggregate throughput

$$T(n) \geq Kn^{2-\alpha/2-\epsilon}$$

is achievable in the network for all possible pairings between sources and destinations. $K > 0$ is a constant independent of n and the source–destination pairing.

- $\alpha \geq 3$: With high probability, an aggregate throughput

$$T(n) \geq K\sqrt{n}$$

is achievable in the network for all possible pairings between sources and destinations. $K > 0$ is a constant independent of n and the source–destination pairing.

Note that because of the bursty transmission strategy, the hierarchical scheme has a high peak-to-average power ratio. However, although we talk in terms of time in the above discussion, such burstiness can just as well be implemented over frequency with only a fraction of the total bandwidth W used. For example, this can be implemented in an orthogonal frequency-division multiplexing (OFDM) system, using a subset of the subcarriers at any one time, but putting more energy in the active subcarriers. This way, the peak power remains constant over time.

B. Cutset Upper Bound for Random Source–Destination Pairings

Can we do better than the scaling in Theorem 5.1? So far we have been considering arbitrary source–destination pairings but clearly there are some pairings for which a much better scaling can be achieved. For example, if the source–destinations are all nearest neighbor to each other, then a linear capacity scaling can be achieved for any α . Thus, for the extended network case, we need to narrow down the class of source–destination pairings to prove a sensible upper bound. In this subsection, we will therefore focus on *random* source–destination pairings, assuming that the pairs are chosen according to a random permutation of the set of nodes, without any consideration on node locations. We prove a high probability upper bound that matches the achievability result in Theorem 5.1, within a polynomial factor of arbitrarily small exponent. Theorem 5.2 together with Theorem 5.1 identify therefore the capacity scaling law in extended networks for *all* values of $\alpha \geq 2$. The rest of the section is devoted to the proof of the theorem.

Theorem 5.2: Consider an extended network of n nodes with random source–destination pairing. For any $\epsilon > 0$, the aggregate throughput is bounded above by

$$T(n) \leq \begin{cases} K'n^{2-\alpha/2+\epsilon}, & 2 \leq \alpha \leq 3 \\ K'n^{1/2+\epsilon}, & \alpha > 3 \end{cases}$$

with high probability for a constant $K' > 0$ independent of n .

Note that the hierarchical scheme is achieving near global cooperation. In the context of dense networks, this yields a near linear number of degrees of freedom for communication. In the context of extended networks, in addition to the degrees of freedom provided, this scheme allows almost all nodes in the network to cooperate in transferring energy between any source–destination pair. In fact, we saw that in extended networks with $\alpha > 2$, our scheme is *power-limited* rather than degrees-of-freedom-limited. A natural place to look for a matching upper bound is to consider a *cut-set* bound on how much power can flow across the network. Our proof of Theorem 5.2 is a careful evaluation of such a cut-set bound.

Proof of Theorem 5.2: We start by considering several properties that are satisfied with high probability in the random network. The following lemma is similar in spirit to Lemma 4.1 for dense networks and can be proved using a similar technique. In parallel to the dense case, it forms the groundwork for our following discussion.

Lemma 5.1: The random network with random source–destination pairing satisfies the following properties with high probability.

- a) Let the network area be divided into n squarelets of unit area. Then, there are less than $\log n$ nodes inside all squarelets.
- b) Let the network area be divided into $\frac{n}{2 \log n}$ squarelets each of area $2 \log n$. Then, there is at least one node inside all squarelets.

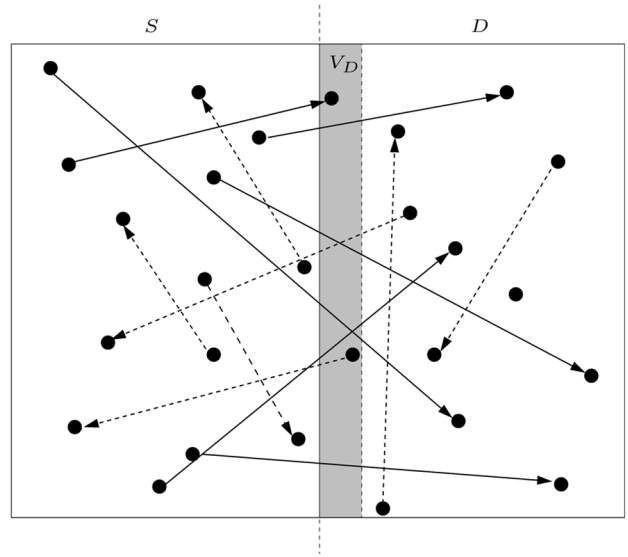


Fig. 6. The cut-set considered in the Proof of Theorem 5.2. The communication requests that pass across the cut from left to right are depicted in bold lines.

- c) Consider a cut dividing the network area into two equal halves. The number of communication requests with sources on the left half network and destinations on the right half network is between $((1 - \delta)n/4, (1 + \delta)n/4)$, for any $\delta > 0$.

We consider a cut dividing the $\sqrt{n} \times \sqrt{n}$ network area into two equal halves (see Fig. 6). We are interested in bounding above the sum of the rates of communications passing through the cut from left to right. By Part c) of the lemma, this sum-rate is equal to 1/4th of the total throughput $T(n)$ with high probability. The maximum achievable sum-rate between these source–destination pairs is bounded above by the capacity of the MIMO channel between nodes S located to the left of the cut and nodes D located to the right. Under the fast fading assumption, we have⁴

$$\sum_{k \in S, i \in D} R_{ik} \leq \max_{\substack{Q(H) \geq 0 \\ \mathbb{E}(Q_{kk}(H)) \leq P, \forall k \in S}} \mathbb{E}(\log \det(I + HQ(H)H^*)) \tag{6}$$

where

$$H_{ik} = \frac{\sqrt{G}e^{j\theta_{ik}}}{r_{ik}^{\alpha/2}}, \quad k \in S, i \in D.$$

$Q(\cdot)$ is a mapping from the set of possible channel realizations H to the set of positive semi-definite transmit covariance matrices. The diagonal element $Q_{kk}(H)$ corresponds to the power allocated to the k th node at channel state H . A natural way to upper-bound (6) is by relaxing the individual power constraint to a total transmit power constraint. In the present context, however, this is not convenient: some nodes in S are close to the cut and some are far apart, so the impact of these nodes on the system performance is quite different. A total transmit power constraint allows the transfer of power from the nodes far apart

⁴Here and in the following, we set the noise variance N_0 to be equal to 1, in order to ease the notation.

to those nodes that are close to the cut, resulting in a loose bound. Instead, we will relax the individual power constraints to a total *weighted* power constraint, where the weight assigned to a node is set to be the total *received* power on the other side of the cut per watt of transmit power from that node. However, before doing that, we need to isolate the contribution of some nodes in D that are located very close to the cut. Typically, there are few nodes on both sides of the cut that are located at a distance as small as order $\frac{1}{\sqrt{n}}$ from the cut. If included, the contribution of these few pairs to the total received power would be excessive, resulting in a loose bound in the discussion below.

Let V_D denote the set of nodes located on the $1 \times \sqrt{n}$ rectangular area immediately to the right of the cut. Note that there are no more than $\sqrt{n} \log n$ nodes in V_D by Part a) of Lemma 5.1. By generalized Hadamard's inequality, we have

$$\log \det(I + HQ(H)H^*) \leq \log \det(I + H^{(1)}Q(H)H^{(1)*}) + \log \det(I + H^{(2)}Q(H)H^{(2)*}) \quad (7)$$

where $H^{(1)}$ and $H^{(2)}$ are obtained by partitioning the original matrix H : $H^{(1)}$ is the rectangular matrix with entries $H_{ik}, k \in S, i \in V_D$ and $H^{(2)}$ is the rectangular matrix with entries $H_{ik}, k \in S, i \in D \setminus V_D$. In turn, (6) is bounded above by

$$\begin{aligned} & \sum_{k \in S, i \in D} R_{ik} \\ & \leq \max_{\substack{Q(H^{(1)}) \geq 0 \\ \mathbb{E}(Q_{kk}(H^{(1)})) \leq P, \forall k \in S}} \mathbb{E} \left(\log \det(I + H^{(1)}Q(H^{(1)})H^{(1)*}) \right) \\ & + \max_{\substack{Q(H^{(2)}) \geq 0 \\ \mathbb{E}(Q_{kk}(H^{(2)})) \leq P, \forall k \in S}} \mathbb{E} \left(\log \det(I + H^{(2)}Q(H^{(2)})H^{(2)*}) \right). \end{aligned} \quad (8)$$

The first term in (8) can be easily upper-bounded by applying Hadamard's inequality once more or, equivalently, by considering the sum of the capacities of the individual multiple-input single-output (MISO) channels between nodes in S and each node in V_D . A discussion similar to the proof of Theorem 3.1 that makes use of the fact that the minimum distance between any two nodes in the network is larger than $\frac{1}{n^{1/2+\delta}}$ with high probability for any $\delta > 0$, yields the following upper bound for the first term:

$$\begin{aligned} & \max_{\substack{Q(H^{(1)}) \geq 0 \\ \mathbb{E}(Q_{kk}(H^{(1)})) \leq P, \forall k \in S}} \mathbb{E} \left(\log \det(I + H^{(1)}Q(H^{(1)})H^{(1)*}) \right) \\ & \leq K' \sqrt{n} (\log n)^2 \end{aligned}$$

where $K' > 0$ is a constant independent of n .

The second term in (8) is the capacity of the MIMO channel between nodes in S and nodes in $D \setminus V_D$. This is the term that dominates in (8) and thus its scaling determines the scaling of (6). The result is given by the following lemma, which completes the proof of Theorem 5.2.

Lemma 5.2: Let $P_{\text{tot}}(n)$ be the total power received by all the nodes in $D \setminus V_D$, when nodes in S are transmitting *independent* signals at full power. Then for every $\epsilon > 0$

$$\begin{aligned} & \max_{\substack{Q(H^{(2)}) \geq 0 \\ \mathbb{E}(Q_{kk}(H^{(2)})) \leq P, \forall k \in S}} \mathbb{E} \left(\log \det(I + H^{(2)}Q(H^{(2)})H^{(2)*}) \right) \\ & \leq n^\epsilon P_{\text{tot}}(n). \end{aligned}$$

Moreover, the scaling of the total received power can be evaluated to be

$$P_{\text{tot}}(n) \leq \begin{cases} K' n (\log n)^3, & \alpha = 2 \\ K' n^{2-\alpha/2} (\log n)^2, & 2 < \alpha < 3 \\ K' \sqrt{n} (\log n)^3, & \alpha = 3 \\ K' \sqrt{n} (\log n)^2, & \alpha > 3 \end{cases}$$

with high probability for a constant $K' > 0$ independent of n . \square

Lemma 5.2 says two things of importance. First, it says that independent signaling at the transmit nodes is sufficient to achieve the cut-set upper bound, as far as scaling is concerned. There is therefore no need, in order for the transmit nodes to cooperate, to do any sort of transmit beamforming. This is fortuitous since our hierarchical MIMO performs only independent signaling across the transmit nodes in the long-range MIMO phase. Second, it identifies the total received power under independent transmissions as the fundamental quantity limiting performance. Depending on α , there is a dichotomy on how this quantity scales with the system size. This dichotomy can be interpreted as follows.

The total received power is dominated either by the power transferred between nodes near the cut (order 1 distance) or by the power transferred between nodes far away from the cut. There are relatively fewer node *pairs* near the cut than away from the cut (order \sqrt{n} versus order n^2), but the channels between the nodes near the cut are considerably stronger than between the nodes far away from the cut. When the attenuation parameter α is less than 3, the received power is dominated by transfer between nodes far away from the cut. The hierarchical scheme, which involves at the top level of the hierarchy MIMO transmissions between clusters of size $n^{1-\epsilon}$ at distance \sqrt{n} apart, achieves arbitrarily closely the required power transfer and is therefore optimal in this regime. When $\alpha \geq 3$, the received power in the cut-set bound is dominated by the power transfer by the nodes near the cut. This can be achieved by nearest neighbor multihop and multihop is therefore optimal in this regime.

It should be noted that earlier works identified thresholds on α above which nearest neighbor multihop is order-optimal ($\alpha > 6$ first established in [5], then subsequently refined to hold for $\alpha > 5$ in [7], $\alpha > 4.5$ in [10], and $\alpha > 4$ in [2]). All of them essentially use the same cut-set bound as we did. The fact that they did not identify the tightest threshold (which we are showing to be 3) is because their upper bounds on the cut-set bound are not tight.

Proof of Lemma 5.2: We are interested in the scaling of the MIMO capacity

$$\max_{\substack{Q(H^{(2)}) \geq 0 \\ \mathbb{E}(Q_{kk}(H^{(2)})) \leq P, \forall k \in S}} \mathbb{E} \left(\log \det \left(I + H^{(2)} Q \left(H^{(2)} \right) H^{(2)*} \right) \right). \quad (9)$$

Let us rescale each column k of the matrix by the (square root of the) *total received power* on the right from source node k on the left. Let indeed P_k denote the total received power in $D \setminus V_D$ of the signal sent by user $k \in S$

$$P_k = PG \sum_{i \in D \setminus V_D} r_{ik}^{-\alpha} := PG d_k.$$

The expression (9) is then equal to

$$\max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\tilde{Q}_{kk}(\tilde{H})) \leq P_k, \forall k \in S}} \mathbb{E}(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*))$$

where

$$\tilde{H}_{ik} = \frac{e^{j\theta_{ik}}}{r_{ik}^{\alpha/2}} \frac{1}{\sqrt{d_k}}, \quad i \in D \setminus V_D, k \in S.$$

The above expression is in turn bounded above by

$$\max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E}(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*))$$

where

$$P_{\text{tot}}(n) = \sum_{k \in S} P_k = PG \sum_{k \in S, i \in D \setminus V_D} r_{ik}^{-\alpha}.$$

Let us now define, for given $n \geq 1$ and $\varepsilon > 0$, the set

$$B_{n,\varepsilon} = \{ \|\tilde{H}\|^2 > n^\varepsilon \}$$

where $\|A\|$ denotes the largest singular value of the matrix A . Note that the matrix \tilde{H} is better conditioned than the original channel matrix $H^{(2)}$: all the diagonal elements of $\tilde{H} \tilde{H}^*$ are roughly of the same order (up to a factor $\log n$), and it can be shown that there exists $K'_1 > 0$ such that

$$\mathbb{E}(\|\tilde{H}\|^2) \leq K'_1 (\log n)^3$$

for all n . In Appendix III, we show the following more precise statement.

Lemma 5.3: For any $\varepsilon > 0$ and $p \geq 1$, there exists $K'_1 > 0$ such that for all n ,

$$\mathbb{P}(B_{n,\varepsilon}) \leq \frac{K'_1}{n^p}.$$

It follows that

$$\begin{aligned} & \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E}(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*)) \\ & \leq \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E} \left(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*) 1_{B_{n,\varepsilon}} \right) \\ & \quad + \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E} \left(\text{Tr}(\tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*) 1_{B_{n,\varepsilon}^c} \right). \quad (10) \end{aligned}$$

The first term in (10) refers to the event that the channel matrix \tilde{H} is accidentally ill-conditioned. Since the probability of such an event is polynomially small by Lemma 5.3, the contribution of this first term is actually negligible. In the second term in (10), the matrix \tilde{H} is well-conditioned, and this term is actually proportional to the maximum power transfer from left to right. Details follow below.

For the first term in (10), we use Hadamard's inequality and obtain (11a) at the bottom of the page, where \tilde{H}_i is the i th row of \tilde{H} . Equation (11b) follows by Jensen's inequality, which in turn is bounded above by (11d), since

$$\|\tilde{H}_i\|^2 = \sum_{k \in S} r_{ik}^{-\alpha} \frac{1}{d_k} \leq \sum_{k \in S} 1 \leq n.$$

The fact that the minimum distance between the nodes in S and $D \setminus V_D$ is at least 1 yields $P_{\text{tot}}(n) \leq PGn^2$. Noting that $x \mapsto x \log(1 + 1/x)$ is increasing on $[0, 1]$ and using Lemma 5.3, we obtain finally that for any $p \geq 1$, there exists $K'_1 > 0$ such that

$$\begin{aligned} & \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E} \left(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*) 1_{B_{n,\varepsilon}} \right) \\ & \leq K'_1 n^{1-p} \log \left(1 + \frac{n^{3+p}}{K'_1} \right), \end{aligned}$$

which decays polynomially to zero with arbitrary exponent as n tends to infinity.

$$\max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E} \left(\log \det(I + \tilde{H} \tilde{Q}(\tilde{H}) \tilde{H}^*) 1_{B_{n,\varepsilon}} \right) \leq \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \mathbb{E} \left(\sum_{i \in D \setminus V_D} \log(1 + \tilde{H}_i \tilde{Q}(\tilde{H}) \tilde{H}_i^*) \middle| B_{n,\varepsilon} \right) \mathbb{P}(B_{n,\varepsilon}) \quad (11a)$$

$$\leq \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \sum_{i \in D \setminus V_D} \log(1 + \mathbb{E}(\|\tilde{H}_i\|^2 \text{Tr} \tilde{Q}(\tilde{H}) \mid B_{n,\varepsilon})) \mathbb{P}(B_{n,\varepsilon}) \quad (11b)$$

$$\leq \max_{\substack{\tilde{Q}(\tilde{H}) \geq 0 \\ \mathbb{E}(\text{Tr} \tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n)}} \sum_{i \in D \setminus V_D} \log(1 + \mathbb{E}(\|\tilde{H}_i\|^2 \text{Tr} \tilde{Q}(\tilde{H})) / \mathbb{P}(B_{n,\varepsilon})) \mathbb{P}(B_{n,\varepsilon}) \quad (11c)$$

$$\leq n \log \left(1 + \frac{n P_{\text{tot}}(n)}{\mathbb{P}(B_{n,\varepsilon})} \right) \mathbb{P}(B_{n,\varepsilon}) \quad (11d)$$

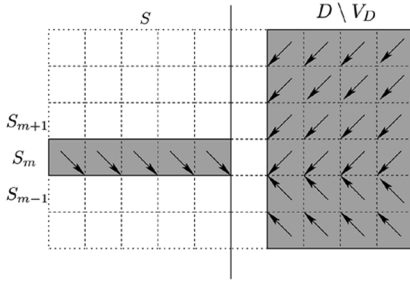


Fig. 7. The displacement of the nodes inside the squarelets to squarelet vertices, indicated by arrows.

For the second term in (10), we simply have

$$\begin{aligned} & \max_{\tilde{Q}(\tilde{H}) \geq 0} \mathbb{E} \left(\text{Tr}(\tilde{H}\tilde{Q}(\tilde{H})\tilde{H}^*)1_{B_{n,\varepsilon}^c} \right) \\ & \mathbb{E}(\text{Tr}\tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n) \\ & \leq \max_{\tilde{Q}(\tilde{H}) \geq 0} \mathbb{E} \left(\|\tilde{H}\|^2 \text{Tr}\tilde{Q}(\tilde{H})1_{B_{n,\varepsilon}^c} \right) \\ & \mathbb{E}(\text{Tr}\tilde{Q}(\tilde{H})) \leq P_{\text{tot}}(n) \\ & \leq n^\varepsilon P_{\text{tot}}(n). \end{aligned}$$

The last thing that needs therefore to be checked is the scaling of $P_{\text{tot}}(n)$ stated in Lemma 5.2.

Let us divide the network area into n squarelets of area 1. By Part a) of Lemma 5.1, there are no more than $\log n$ nodes in each squarelet with high probability. Let us consider grouping the squarelets of S into \sqrt{n} rectangular areas S_m of height 1 and width \sqrt{n} , as shown in Fig. 7. Thus, $S = \bigcup_{m=1}^{\sqrt{n}} S_m$. We are interested in bounding above

$$P_{\text{tot}}(n) = PG \sum_{k \in S} d_k = PG \sum_{m=1}^{\sqrt{n}} \sum_{k \in S_m} d_k.$$

Let us consider

$$\sum_{k \in S_m} d_k = \sum_{k \in S_m, i \in D \setminus V_D} r_{ik}^{-\alpha} \quad (12)$$

for a given m . Note that if we move the nodes that lie in each squarelet of S_m together with the nodes in the squarelets of $D \setminus V_D$ onto the squarelet vertex as indicated by the arrows in Fig. 7, all the (positive) terms in the summation in (12) can only increase since the displacement can only decrease the Euclidean distance between the nodes involved. Note that the modification results in a regular network with at most $\log n$ nodes at each squarelet vertex on the left and at most $2 \log n$ nodes at each squarelet vertex on the right. Considering the same reasoning for all rectangular slabs $S_m, m = 1, \dots, \sqrt{n}$ allows to conclude that $P_{\text{tot}}(n)$ for the random network is with high probability less than the same quantity computed for a regular network with $\log n$ nodes at each left-hand side vertex and $2 \log n$ nodes at each right-hand side vertex.

The most convenient way to index the node positions in the resulting regular network is to use double indices. The left-hand side nodes are located at positions $(-k_x + 1, k_y)$ and those on

the right at positions (i_x, i_y) where $k_x, k_y, i_x, i_y = 1, \dots, \sqrt{n}$, so that

$$\tilde{H}_{ik} = \frac{e^{j\theta_{ik}}}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/4}} \frac{1}{\sqrt{d_{k_x, k_y}}}$$

and

$$d_{k_x, k_y} = \sum_{i_x, i_y=1}^{\sqrt{n}} \frac{1}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/2}} \quad (13)$$

which yields the following upper bound for $P_{\text{tot}}(n)$ of the random network

$$P_{\text{tot}}(n) \leq 2(\log n)^2 PG \sum_{k_x, k_y=1}^{\sqrt{n}} d_{k_x, k_y}. \quad (14)$$

The following lemma establishes the scaling of d_{k_x, k_y} defined in (13).

Lemma 5.4: There exist constants $K'_2, K'_3 > 0$ independent of k_x, k_y , and n such that

$$d_{k_x, k_y} \leq \begin{cases} K'_2 \log n, & \text{if } \alpha = 2 \\ K'_2 k_x^{2-\alpha}, & \text{if } \alpha > 2 \end{cases}$$

and

$$d_{k_x, k_y} \geq K'_3 k_x^{2-\alpha}, \quad \text{for } \alpha \geq 2.$$

The rigorous proof of the lemma is given at the end of Appendix III. A heuristic way of thinking about the approximation

$$d_{k_x, k_y} \approx k_x^{2-\alpha}$$

can be obtained through Laplace's principle. The summation in d_{k_x, k_y} scales the same as the maximum term in the sum times the number of terms which have roughly this maximum value. The maximum term is of the order of $1/k_x^\alpha$. The terms that take on roughly this value are those for which i_x runs from 1 to the order of k_x and i_y runs from k_y to k_y plus or minus the order of k_x . There are roughly k_x^2 such terms. Hence $d_{k_x, k_y} \approx 1/k_x^\alpha \cdot k_x^2 = k_x^{2-\alpha}$.

We can now use the upper bound given in the above lemma to yield

$$\sum_{k_x, k_y=1}^{\sqrt{n}} d_{k_x, k_y} \leq \begin{cases} K'_4 n \log n & \text{if } \alpha = 2 \\ K'_4 n^{2-\alpha/2}, & \text{if } 2 < \alpha \leq 3 \\ K'_4 \sqrt{n} \log n, & \text{if } \alpha = 3 \\ K'_4 \sqrt{n}, & \text{if } \alpha > 3 \end{cases}$$

for another constant $K'_4 > 0$ independent of n . This upper bound combined with (14) completes the Proof of Lemma 5.2. \square

VI. DISCUSSIONS ON THE MODEL AND THE RESULTS

In this section, we point out the scope and limitations of the model and the results.

A. Scaling Laws Versus Performance Analysis

We should emphasize that the focus in this paper, as well as in [3] and the follow-up works, is on scaling laws, i.e., scaling of the aggregate throughput in the limit when the number of users gets large. The main advantage of studying scaling laws is to highlight qualitative and architectural properties of the system

without getting bogged down by too many details. For example, the linear scaling law in dense networks we derived highlights the fact that interference limitation in the Gupta–Kumar scaling is not a fundamental one and can be alleviated by more complicated physical-layer processing.

It is important to distinguish between such scaling law study and the design and performance analysis of a scheme for a network with a given number of users. While scaling law results provide some architectural guidelines on how to design schemes that scale well, detailed design and performance analysis would involve tuning of many parameters and improvements of the scheme to optimize the pre-constant in the system throughput. For example, our scheme quantizes the received analog signal at each node and forward the bits to the final destination, but the quantized bits are correlated across the receive nodes and hence a reduction in the overhead can be achieved by doing some Slepian–Wolf coding (as proposed in [15]). Such work is beyond the scope of the present paper.

We studied two different scaling laws in this paper, one for dense and one for extended networks. Given a network with a specific number of nodes occupying a specific area, a natural question is: is this network best described by the dense scaling regime or the extended scaling regime? What our results say is that a better delineation is in terms of whether we are in the degree-of-freedom-limited or power (coverage)-limited regime, because this is what will have architectural implications for the communication scheme (for example, whether bursty transmissions are required). To get a sense of the operating regime a given network is in, our results suggest a rule-of-thumb quantity that can be calculated: the total received SNR per node, total over all the transmit powers of the nodes in the network. If this quantity is much larger than 0 dB, then the network is in the degree-of-freedom-limited regime; otherwise it is in the power-limited regime.

B. Far-Field Assumption in Dense Scaling

One potential concern with the dense scaling is that the far-field assumption will eventually break down as the number of nodes gets too large. In practice, the typical separation between nodes is so much larger than the carrier wavelength that the number of nodes for which the far-field assumption fails has to be extremely large, i.e., there is a clear separation between the large and the small spatial scales. Consider the following numerical example: suppose the area of interest is 1 km², well within the communication range of many radio devices. With a carrier frequency of 3 GHz, the carrier wavelength is 0.1 m. Even with a very large system size of $n = 10000$ nodes, the typical separation between nearest neighbors is 10 m, very much in the far field. Under free-space propagation and assuming unit transmit and receive antenna gains, the attenuation given by Friis' formula (2) is about 10^{-6} , much smaller than unity. At the same time, the total received SNR per node (assuming transmit power P of 1 mW per node, thermal noise N_0 at -174 dBm, a bandwidth W of 10 MHz, and noise figure $NF = 10$ dB) is 84 dB, very much in the degree-of-freedom-limited regime.⁵ (Looking at even only one point-to-point link at distance 1 km,

the received SNR is 34 dB). Hence, this example gives evidence that there are networks for which simultaneously the number of nodes is large, the far-field assumption holds, and the received SNR across the network is high. However, a careful performance analysis of the pre-constants is required to confirm that linear scaling of our scheme has already kicked in and our scheme indeed outperforms multihop in this parameter range. Nevertheless, we do believe that the linear scaling obtained here also applies for a relatively small number of nodes. The intuition for this is that our strategy relies on the use of MIMO communication, whose linear capacity scaling has never been disputed in the range of a small number of antennas.

C. Transport Capacity

Let us finally mention that a more general measure of network performance has been introduced in [3]: the *transport capacity* of a network, defined as the maximum number of bits exchanged in the network per second, weighted by their traveled distances. From an upper bound on transport capacity, one can easily deduce an upper bound on the aggregate throughput for the special traffic model considered in this paper, where nodes are paired up one-to-one into source–destination pairs that communicate at a common rate, which is the traffic requirement considered in the current paper. But the interest in an *upper bound* on transport capacity lies in the fact that it applies to more general communication scenarios. Reciprocally, it has been shown recently in [16] that for a network with a random placement of nodes, there is a natural way to deduce an upper bound on transport capacity from an upper bound on throughput, by studying cut-set bounds over multiple cuts (as first suggested in [5]). Applying this technique to the present result leads to the following conclusion: the transport capacity $T_c(n)$ of the extended network is upper-bounded by

- $T_c(n) \leq K'n^{2.5-\alpha/2+\varepsilon}$, for $2 \leq \alpha \leq 3$;
- $T_c(n) \leq K'n^{1+\varepsilon}$, for $\alpha > 3$;

for any $\varepsilon > 0$, where $K' > 0$ is a constant independent of n . Note that these scaling laws for the transport capacity are also achievable within a factor of n^ε .

VII. CONCLUSION

In point-to-point communication, performance is limited by either the power or the degrees of freedom (bandwidth and number of antennas) available, depending on whether the link is operating at low or high SNR. In a network with multiple source–destination pairs, performance can further be limited by the interference between simultaneous transmission of information. In this paper, we have shown that by achieving near global MIMO cooperation between nodes without introducing significant cooperation overhead, interference can be successfully removed as a limitation, at least as far as scaling laws are concerned. Moreover, such near-global MIMO cooperation also allows the maximum transfer of energy between all source–destination pairs, provided that the path loss across the network is not too much. This implies that in degrees-of-freedom-limited scenarios, such as in dense networks or extended networks with path loss exponent $\alpha = 2$, the full degrees of freedom in the network can be shared among all nodes and a linear capacity scaling can be achieved. In power-limited scenarios but with

⁵ $\text{SNR}_{\text{dB}} = P_{\text{dBm}} + 10 \log_{10} n + \text{path loss}_{\text{dB}} - (N_0)_{\text{dBm}} - 10 \log_{10} W - NF_{\text{dB}}$.

low attenuation, such as extended networks with α between 2 and 3, our scheme achieves the optimal (power-limited) capacity scaling law.

The key ideas behind our scheme are as follows:

- using MIMO for long-range communication to achieve spatial multiplexing;
- using local transmit and receive cooperation to maximize spatial reuse;
- setting up the intra-cluster cooperation such that it is yet another digital communication problem, but in a smaller network, thus enabling a hierarchical cooperation architecture.

We have focused on the two-dimensional (2D) setting, where the nodes are on the plane, but our results generalize naturally to networks where nodes live in d -dimensional space. For dense networks, linear scaling is achievable whenever $\alpha > d$, i.e., whenever spatial reuse is possible. For extended networks, the scaling exponent $e_d^*(\alpha)$ is given by

$$e_d^*(\alpha) = \begin{cases} 2 - \frac{\alpha}{d}, & d \leq \alpha \leq d + 1 \\ 1 - \frac{1}{d}, & \alpha > d + 1. \end{cases}$$

For α between d and $d + 1$, hierarchical MIMO achieves the optimal scaling, and for $\alpha > d + 1$, nearest neighbor multihop is optimal.

APPENDIX I
LINEAR SCALING LAW FOR THE MIMO GAIN
UNDER FAST FADING ASSUMPTION

Proof of Lemma 4.3: The $M \times M$ MIMO channel between two clusters S and D is given by $Y = HX + Z$, where H_{ik} are given in (1). Recall that $Z = (Z_k)$ is uncorrelated background noise plus interference at the receiver nodes. Assume that the transmitted signals $X = (X_i)$ are from an i.i.d. $\sim \mathcal{N}_{\mathbb{C}}(0, \sigma^2)$ randomly chosen codebook with

$$\sigma^2 = \frac{P(r_{SD})^\alpha}{M}.$$

It is well known that the achievable mutual information is lower-bounded by assuming that the interference-plus-noise Z is i.i.d. Gaussian. (See, for example, [17, Theorem 5] for a precise statement and proof of this in the MIMO case.) With our transmission strategy in the MIMO phase, there exists $b > a > 0$ with a and b independent of n , such that $r_{ik}^{-\alpha/2} = r_{SD}^{-\alpha/2} \rho_{ik}$, where all ρ_{ik} lie in the interval $[a, b]$ both in the cases when S and D are neighboring clusters or not.

By assuming perfect channel state information at the receiver side, the mutual information of the above MIMO channel is given by

$$\begin{aligned} I(X; Y, H) &\geq \mathbb{E} \left(\log \det \left(I + \frac{\sigma^2}{N} H H^* \right) \right) \\ &= \mathbb{E} \left(\log \det \left(I + \frac{\text{SNR}}{M} F F^* \right) \right) \end{aligned} \quad (15)$$

where

$$\text{SNR} = \frac{GP}{N} \quad (N = \text{total interference-plus-noise power})$$

and $F_{ik} = \rho_{ik} \exp(j\theta_{ik})$. Let λ be chosen uniformly among the M eigenvalues of $\frac{1}{M} F F^*$. The above mutual information may be rewritten as

$$\begin{aligned} I(X; Y, H) &\geq M \mathbb{E}(\log(1 + \text{SNR} \lambda)) \\ &\geq M \log(1 + \text{SNR} t) \mathbb{P}(\lambda > t) \end{aligned}$$

for any $t \geq 0$. By the Paley–Zygmund inequality, if $0 \leq t < \mathbb{E}(\lambda)$, we have

$$\mathbb{P}(\lambda > t) \geq \frac{(\mathbb{E}(\lambda) - t)^2}{\mathbb{E}(\lambda^2)}.$$

We therefore need to compute both $\mathbb{E}(\lambda)$ and $\mathbb{E}(\lambda^2)$. We have

$$\begin{aligned} \mathbb{E}(\lambda) &= \frac{1}{M} \mathbb{E} \left(\text{Tr} \left(\frac{1}{M} F F^* \right) \right) \\ &= \frac{1}{M^2} \sum_{i,k=1}^M \mathbb{E}(|F_{ik}|^2) \\ &= \frac{1}{M^2} \sum_{i,k=1}^M \rho_{ik}^2 \geq a^2 \end{aligned}$$

and

$$\begin{aligned} \mathbb{E}(\lambda^2) &= \frac{1}{M} \mathbb{E} \left(\text{Tr} \left(\frac{1}{M^2} F F^* F F^* \right) \right) \\ &= \frac{1}{M^3} \sum_{iklm=1}^M \mathbb{E}(F_{ik} \overline{F_{ik}} F_{lm} \overline{F_{lm}}) \\ &\leq \frac{2}{M^3} \sum_{ikm=1}^M \mathbb{E}(|F_{ik}|^2) \mathbb{E}(|F_{im}|^2) \\ &= \frac{2}{M^3} \sum_{ikm=1}^M \rho_{ik}^2 \rho_{im}^2 \leq 2b^4 \end{aligned}$$

so $\mathbb{E}(\lambda) \geq a^2$ and $\mathbb{E}(\lambda^2) \leq 2b^4$. This leads us to the conclusion that for any $t < a$, we have

$$I(X; Y, H) \geq M \log(1 + \text{SNR} t) \frac{(a^2 - t)^2}{2b^4}. \quad (16)$$

Choosing, e.g., $t = a/2$ shows that $I(X; Y, H)$ grows at least linearly with M . \square

Lemma 1.1 (Paley–Zygmund Inequality): Let X be a non-negative random variable such that $\mathbb{E}(X^2) < \infty$. Then for any $t \geq 0$ such that $t < \mathbb{E}(X)$, we have

$$\mathbb{P}(X > t) \geq \frac{(\mathbb{E}(X) - t)^2}{\mathbb{E}(X^2)}$$

Proof: By the Cauchy–Schwarz inequality, we have for any $t \geq 0$

$$\mathbb{E}(X 1_{X>t}) \leq \sqrt{\mathbb{E}(X^2) \mathbb{P}(X > t)}$$

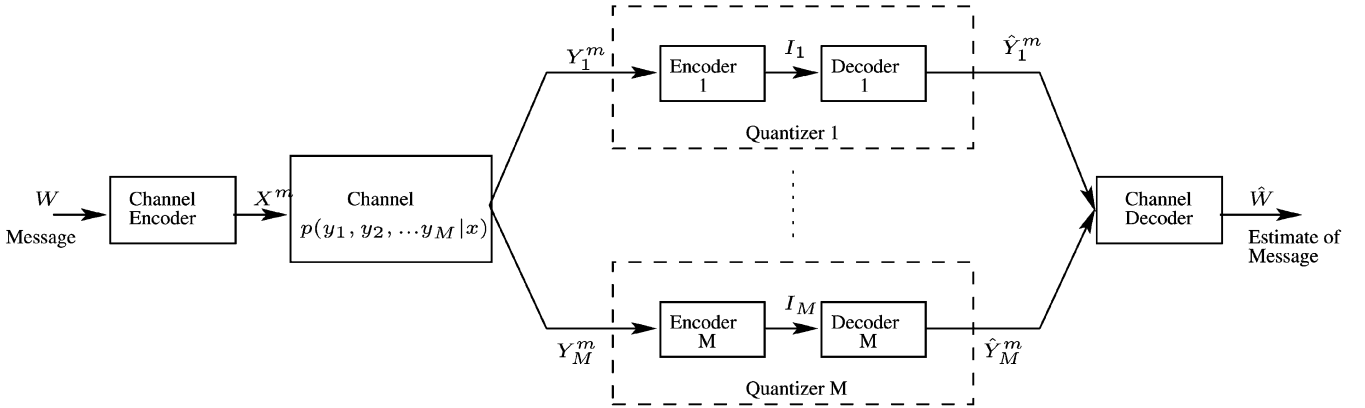


Fig. 8. The quantized channel problem.

and also, if $t < \mathbb{E}(X)$

$$\mathbb{E}(X1_{X>t}) = \mathbb{E}(X) - \mathbb{E}(X1_{X\leq t}) \geq \mathbb{E}(X) - t > 0$$

Therefore

$$\mathbb{P}(X > t) \geq \frac{(\mathbb{E}(X) - t)^2}{\mathbb{E}(X^2)}. \quad \square$$

Note that the achievability results in this paper can be extended to the slow fading case, provided that Lemma 4.3 can be proved in the slow fading setting. In that case, one would need to show that the expression inside the expectation in (15) concentrates around its mean exponentially fast in M . However, another difficulty might arise from the lack of averaging of the phases in the interference term, which leads to a non spatially decorrelated noise term Z . Although proving the result might require some technical effort, we believe it holds true, due to the self-averaging effect of a large number of independent random variables.

APPENDIX II

ACHIEVABLE RATES ON QUANTIZED CHANNELS

In order to conclude the discussion on the throughput achieved by our scheme, we need to show that the quantized MIMO channel achieves the same spatial multiplexing gain as the MIMO channel. In Theorem II.1, we give a simple achievability region for general quantized channels. Note that a stronger result is established in [18, Theorem 3] that implies Theorem II.1 as a special case. The required result for the quantized MIMO channel is then found as an easy application of Theorem II.1. We start by formally defining the general quantized channel problem in a form that is of interest to us and proceed with several definitions that will be needed in the sequel.

Let us consider a discrete-time memoryless channel with single input of alphabet \mathcal{X} and M outputs of respective alphabets $\mathcal{Y}_1, \dots, \mathcal{Y}_M$. The channel is statistically described by a conditional probability distribution $p(y_1, \dots, y_M | x)$ for each $y_1 \in \mathcal{Y}_1, \dots, y_M \in \mathcal{Y}_M$ and $x \in \mathcal{X}$. The outputs of

the channel are to be followed by quantizers which independently map the output alphabets $\mathcal{Y}_1, \dots, \mathcal{Y}_M$ to the respective reproduction alphabets $\hat{\mathcal{Y}}_1, \dots, \hat{\mathcal{Y}}_M$. The aim is to recover the transmitted information through the channel by observing the outputs of the quantizers. Communication over the channel takes place in the following manner: a message W , drawn from the index set $\{1, 2, \dots, L\}$, is encoded into a codeword $X^m(W) \in \mathcal{X}^m$, which is received as M random sequences $(Y_1^m, \dots, Y_M^m) \sim p(y_1^m, \dots, y_M^m | x^m)$ at the outputs of the channel. The quantizers themselves consist of encoders and decoders, where the j th encoder describes its corresponding received sequence Y_j^m by an index $I_j(Y_j^m) \in \{1, 2, \dots, L_j\}$, and decoder j represents Y_j^m by an estimate $\hat{Y}_j^m(I_j) \in \hat{\mathcal{Y}}_j^m$. The channel decoder then observes the reconstructed sequences $\hat{Y}_1^m, \dots, \hat{Y}_M^m$ and guesses the index W by an appropriate decoding rule $\hat{W} = g(\hat{Y}_1^m, \dots, \hat{Y}_M^m)$. An error occurs if \hat{W} is not the same as the index W that was transmitted. The complete model under investigation is shown in Fig. 8. An $(L; L_1, \dots, L_M; m)$ code for this channel is a joint (L, m) channel code and M quantization codes $(L_1, m), \dots, (L_M, m)$; more specifically, it is two sets of encoding and decoding functions, the first set being the channel encoding function $X^m : \{1, 2, \dots, L\} \rightarrow \mathcal{X}^m$ and the channel decoding function $g : \hat{\mathcal{Y}}_1^m \times \dots \times \hat{\mathcal{Y}}_M^m \rightarrow \{1, 2, \dots, L\}$, and the second set consists of the encoding functions $I_j : \mathcal{Y}_j^m \rightarrow \{1, 2, \dots, L_j\}$ and decoding functions $\hat{Y}_j^m : \{1, 2, \dots, L_j\} \rightarrow \hat{\mathcal{Y}}_j^m$ for $j = 1, \dots, M$ used for the quantizations. We define the (average) probability of error for the $(L; L_1, \dots, L_M; m)$ code by

$$P_e^m = \frac{1}{L} \sum_{k=1}^L \mathbb{P}(\hat{W} \neq k | W = k).$$

A set of rates $(R; R_1, \dots, R_M)$ is said to be achievable if there exists a sequence of $(2^{mR}, 2^{mR_1}, \dots, 2^{mR_M}; m)$ codes with $P_e^m \rightarrow 0$ as $m \rightarrow \infty$. Note that determining achievable rates $(R; R_1, \dots, R_M)$ is not a trivial problem, since there is tradeoff between maximizing R and minimizing R_1, \dots, R_M .

Theorem II.1 (Achievability for the Quantized Channel Problem): Given a probability distribution $q(x)$ on \mathcal{X} and M conditional probability distributions $q_j(\hat{y}_j | y_j)$ where $y_j \in \mathcal{Y}_j$,

$\hat{y}_j \in \hat{\mathcal{Y}}_j$, and $j = 1, \dots, M$, all rates $(R; R_1, \dots, R_M)$ such that $R < I(X; \hat{Y}_1, \dots, \hat{Y}_M)$ and $R_j > I(Y_j; \hat{Y}_j)$ are achievable. Specifically, given any $\delta > 0$, $q(x)$, and $q_j(\hat{y}_j | y_j)$, together with rates $R < I(X; \hat{Y}_1, \dots, \hat{Y}_M)$ and $R_j > I(Y_j; \hat{Y}_j)$ for $j = 1, \dots, M$; there exists a $(2^{mR}; 2^{mR_1}, \dots, 2^{mR_M}; m)$ code such that $P_e^m < \delta$.

Proof: The proof of the theorem for discrete finite-size alphabets relies on a random coding argument based on the idea of joint (*strong*) typicality. For the idea of strong typicality and properties of typical sequences, see [19, Ch. 13.6]. The proof can be outlined as follows. Given $q(x)$ generate a random channel codebook \mathcal{C}_c with 2^{mR} codewords, each of length m , independently from the distribution

$$q(x^m) = \prod_{k=1}^m q(x^m(k))$$

and call them $X^m(1), X^m(2), \dots, X^m(2^{mR})$. Also generate M quantization codebooks $\mathcal{C}_j, j = 1, \dots, M$, each codebook \mathcal{C}_j consisting of 2^{mR_j} codewords drawn independently from

$$p_j(\hat{y}_j^m) = \prod_{k=1}^m \sum_{\substack{x \in \mathcal{X} \\ y_1 \in \mathcal{Y}_1, \dots, y_M \in \mathcal{Y}_M}} q(x) p(y_1, \dots, y_M | x) q_j(\hat{y}_j^m(k) | y_j).$$

and index them as $\hat{Y}_j^m(1), \hat{Y}_j^m(2), \dots, \hat{Y}_j^m(2^{mR_j})$. Given the message w , send the codeword $X^m(w)$ through the channel. The channel will yield Y_1^m, \dots, Y_M^m . Given the channel output Y_j^m at the j th quantizer, choose i_j such that $(Y_j^m, \hat{Y}_j^m(i_j))$ are jointly typical. If there exists no such i_j , declare an error. If the number of codewords in the quantization codebook 2^{mR_j} is greater than $2^{mI(Y_j; \hat{Y}_j)}$, the probability of finding no such i_j decreases to zero exponentially as m increases. The probability of failing to find such an index in at least one of the M quantizers is bounded above by the union bound with the sum of M exponentially decreasing probabilities in m . Given $\hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M)$ at the channel decoder, choose the unique \hat{w} such that $(X^m(\hat{w}), \hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M))$ are jointly typical. The fact that $(X^m(w), \hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M))$ will be jointly typical with high probability can be established by identifying the Markov chains in the problem and applying Markov lemma [19, Lemma 14.8.1] repeatedly. Observing that $(Y_1^m, \dots, Y_M^m, \hat{Y}_1^m, \dots, \hat{Y}_j^m) - Y_{j+1}^m - \hat{Y}_{j+1}^m$ forms a Markov chain and recursively applying Markov lemma, we conclude that $(Y_1^m, \dots, Y_M^m, \hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M))$ are jointly typical with probability approaching 1 as m increases. Observing that $X^m - (Y_1^m, \dots, Y_M^m) - (\hat{Y}_1^m, \dots, \hat{Y}_M^m)$ forms another Markov chain, again by Markov lemma we have $(X^m(w), \hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M))$ jointly typical with high probability. If there are more than one codewords X^m that are jointly typical with $(\hat{Y}_1^m(i_1), \dots, \hat{Y}_M^m(i_M))$, we declare an error. The probability of having more than one such sequence will decrease exponentially to zero as m increases, if the number of channel codewords 2^{mR} is less than $2^{mI(X; \hat{Y}_1, \dots, \hat{Y}_M)}$. Hence, if $R < I(X; \hat{Y}_1, \dots, \hat{Y}_M)$ and $R_j > I(Y_j; \hat{Y}_j)$, the probability of error averaged over all codes decreases to zero as $m \rightarrow \infty$. This shows the existence of a code that achieves rates

$(R; R_1, \dots, R_M)$ with arbitrarily small probability of error. The result can be readily extended to memoryless channels with discrete-time and continuous alphabets by standard arguments (see [20, Ch.7]). \square

Proof of Lemma 4.4: Now we turn to our original problem. We need to show that it is possible to encode the observations at the outputs of the MIMO channel at a fixed rate, while preserving the spatial multiplexing gain of the MIMO channel. This is a direct consequence of Theorem II.1: Consider the conditional probability densities

$$q_j(\hat{y}_j | y_j) \sim \mathcal{N}_{\mathbb{C}}(y_j, \Delta^2)$$

for the quantization process. From Theorem II.1 we know that for any distribution $p(x)$ on the input space, all rate pairs $(R; R_1, \dots, R_M)$ are simultaneously achievable if

$$R_j > I(Y_j; \hat{Y}_j), \quad j = 1, \dots, M \quad \text{and} \\ R < I(X; \hat{Y}_1, \dots, \hat{Y}_M)$$

where now R_j is the encoding rate of the j th stream and R is the total transmission rate over the MIMO channel. Using Lemma 4.5, we have that $I(Y_j; \hat{Y}_j) \leq \log(1 + \frac{P_2}{\Delta^2})$ for any probability distribution $p(x)$ on the input space. So if we choose

$$R_j = \log\left(1 + \frac{P_2}{\Delta^2}\right) + \varepsilon, \quad \forall j = 1, \dots, M$$

for some $\varepsilon > 0$, all rates

$$R \leq I(X; \hat{Y}_1, \dots, \hat{Y}_M)$$

are achievable on the quantized MIMO channel for any input distribution $p(x)$. Note that now the channel from X to $\hat{Y}_1, \dots, \hat{Y}_M$ is given by

$$\hat{Y} = HX + Z + D$$

where $D \sim \mathcal{N}_{\mathbb{C}}(0, \Delta^2 I)$. Obviously, this channel has the same spatial multiplexing gain as the original MIMO channel. \square

Proof of Lemma 4.6: Consider the case where the MIMO signals are corrupted by interference of increasing power $K_I \log M$. In this case, the power received by the destination nodes is not bounded any more and increases as $P_2 + K_I \log M$ with increasing M . In order to apply the technique employed in the proof of Lemma 4.4, one can first normalize the received signal by multiplying it by $q = \sqrt{\frac{P_2}{P_2 + K_I \log M}}$ and then do the quantization as before. Note that the resultant *scaled* quantized MIMO channel is given by

$$\hat{Y} = q(HX + Z) + D$$

where again $D \sim \mathcal{N}_{\mathbb{C}}(0, \Delta^2 I)$ and $Z = (z_k)$ is the background noise plus interference vector independent of the signal with uncorrelated entries of power $\mathbb{E}[z_k^2] \leq N_0 + K_I \log M$. Thus, we can apply the result of Lemma 4.3. Note that the resultant signal-to-noise-ratio $\text{SNR} \geq \frac{K}{\log M}$ for a constant $K > 0$. Plugging this SNR expression into (16) yields $M/\log M$ capacity scaling for the resultant channel. \square

APPENDIX III
LARGEST EIGENVALUE BEHAVIOR OF THE EQUALIZED
CHANNEL MATRIX \tilde{H}

In this appendix, we give the proofs of Lemmas 5.3 and 5.4. We start with Lemma 5.3. The proof of the second lemma is given at the end of the section.

Proof of Lemma 5.3: Let us start by considering the $2m$ th moment of the spectral norm of \tilde{H} given by (see [21, Ch. 5])

$$\|\tilde{H}\|^{2m} = \rho(\tilde{H}^* \tilde{H})^m = \lim_{l \rightarrow \infty} \{\text{Tr}((\tilde{H}^* \tilde{H})^l)\}^{m/l}.$$

By dominated convergence theorem and Jensen's inequality, we have

$$\mathbb{E}(\|\tilde{H}\|^{2m}) \leq \lim_{l \rightarrow \infty} \{\mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^l))\}^{m/l}.$$

In the subsequent paragraphs, we will prove that the following upper bound holds with high probability:

$$\mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^l)) \leq t_l n (K'_1 \log n)^{3l} \quad (17)$$

where $t_l = \frac{(2l)!}{l!(l+1)!}$ are the Catalan numbers and $K'_1 > 0$ is a constant independent of n . By Chebyshev's inequality, this allows to conclude that for any m

$$\begin{aligned} \mathbb{P}(B_{n,\varepsilon}) &\leq \frac{\mathbb{E}(\|\tilde{H}\|^{2m})}{n^{m\varepsilon}} \leq \frac{1}{n^{m\varepsilon}} \lim_{l \rightarrow \infty} (t_l m (K'_1 \log n)^{3l})^{m/l} \\ &\leq \frac{(4(K'_1 \log n)^3)^m}{n^{m\varepsilon}} \end{aligned}$$

since $\lim_{l \rightarrow \infty} t_l^{1/l} = 4$. For any $\varepsilon > 0$, choosing m sufficiently large shows therefore that $\mathbb{P}(B_{n,\varepsilon})$ decays polynomially with arbitrary exponent as $n \rightarrow \infty$, which is the result stated in Lemma 5.3.

There remains to prove the upper bound in (17). Expanding the expression gives

$$\begin{aligned} \mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^l)) &= \sum_{\substack{i_1, \dots, i_l \in D \setminus V_D \\ k_1, \dots, k_l \in S}} \mathbb{E} \left(\overline{\tilde{H}_{i_1 k_1}} \tilde{H}_{i_1 k_2} \overline{\tilde{H}_{i_2 k_2}} \tilde{H}_{i_2 k_3} \cdots \overline{\tilde{H}_{i_l k_l}} \tilde{H}_{i_l k_1} \right). \end{aligned} \quad (18)$$

Recall that the random variables \tilde{H}_{ik} are independent and zero-mean, so the expectation is only nonzero when the terms in the product form conjugate pairs. Let us consider the case $l = 2$ as an example. We have

$$\begin{aligned} \mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^2)) &= \sum_{\substack{i_1, i_2 \in D \setminus V_D \\ k_1, k_2 \in S}} \mathbb{E} \left(\overline{\tilde{H}_{i_1 k_1}} \tilde{H}_{i_1 k_2} \overline{\tilde{H}_{i_2 k_2}} \tilde{H}_{i_2 k_1} \right) \\ &= \sum_{\substack{i_1, i_2 \in D \setminus V_D \\ k \in S}} |\tilde{H}_{i_1 k}|^2 |\tilde{H}_{i_2 k}|^2 + \sum_{\substack{i \in D \setminus V_D \\ k_1 \neq k_2 \in S}} |\tilde{H}_{i k_1}|^2 |\tilde{H}_{i k_2}|^2 \end{aligned} \quad (19)$$

since the expectation is nonzero only when either $k_1 = k_2 = k$ or $i_1 = i_2 = i$. Note that we have removed the expectations

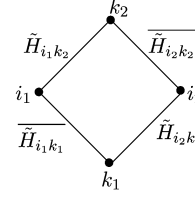


Fig. 9. The product in (19) illustrated as a ring.

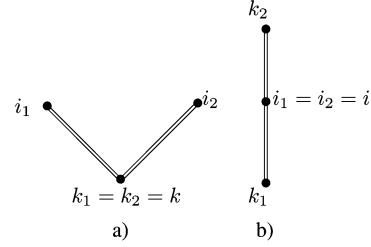


Fig. 10. Two possible graphs corresponding to the nonzero terms in (19).

in (20) since $|\tilde{H}_{ik}|^2$ is a deterministic quantity in our case. The expression can be bounded above by

$$\begin{aligned} \mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^2)) &\leq \sum_{\substack{i_1, i_2 \in D \setminus V_D \\ k \in S}} |\tilde{H}_{i_1 k}|^2 |\tilde{H}_{i_2 k}|^2 + \sum_{\substack{i \in D \setminus V_D \\ k_1, k_2 \in S}} |\tilde{H}_{i k_1}|^2 |\tilde{H}_{i k_2}|^2 \end{aligned} \quad (21)$$

where we now double-count the terms with $i_1 = i_2 = i$ and $k_1 = k_2 = k$, that is, the terms of the form $|\tilde{H}_{ik}|^4$.

The nonvanishing terms in the sum in (19) can also be determined by the following approach, which generalizes to larger l : let each index be associated to a vertex and each term in the product in (19) to an edge between its corresponding vertices. Note that the resulting graph is in general a ring with four edges as depicted in Fig. 9. A term in the summation in (19) is only nonzero if each edge of its corresponding graph has even multiplicity. Such a graph can be obtained from the ring in Fig. 9 by merging some of the vertices, thus equating their corresponding indices. For example, merging the vertices k_1 and k_2 into a single vertex k gives the graph in Fig. 10(a); on the other hand, merging i_1 and i_2 into a single vertex i gives Fig. 10(b). Note that in the first figure i_1, i_2 can take values in $D \setminus V_D$ and k can take values in S , thus the sum of all such terms yields

$$\sum_{\substack{i_1, i_2 \in D \setminus V_D \\ k \in S}} |\tilde{H}_{i_1 k}|^2 |\tilde{H}_{i_2 k}|^2. \quad (22)$$

Similarly, the terms of the form in Fig. 10(b) sum up to

$$\sum_{\substack{i \in D \setminus V_D \\ k_1, k_2 \in S}} |\tilde{H}_{i k_1}|^2 |\tilde{H}_{i k_2}|^2. \quad (23)$$

Note that another possible graph composed of edges with even multiplicity can be obtained by further merging the vertices i_1 and i_2 into a single vertex i in Fig. 10(a), or equivalently merging k_1 and k_2 into k in Fig. 10(b). This will result in a graph with only two vertices k and i and a quadruple edge in between which corresponds to terms of the form $|\tilde{H}_{ik}|^4$ with

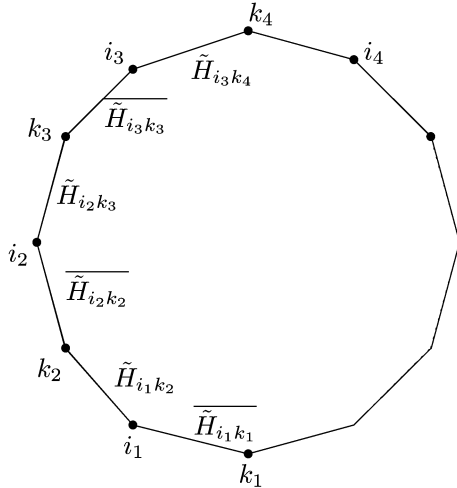


Fig. 11. The product in (18) illustrated as a ring.

$i \in D \setminus V_D$ and $k \in S$. Note, however, that such terms have already been considered in both (22) and (23) since we did not exclude the case $i_1 = i_2$ in (22) and $k_1 = k_2$ in (23). In fact, terms corresponding to any graph with number of vertices less than three are already accounted for in either one of the sums in (22) and (23), or simultaneously in both. Hence, the sum of (22) and (23) is an upper bound for (19) yielding again (21).

In the general case with $l \geq 2$, considering (18) leads to a larger ring with $2l$ edges, as depicted in Fig. 11. Similarly to the case $l = 2$, the nonvanishing terms in (18) are those that correspond to a graph having only edges of even multiplicity. Since each edge can have at least double multiplicity, such graphs can have at most l edges. In turn, a graph with l edges can have at most $l + 1$ vertices which is the case of a tree. Hence, let us first start by considering such trees; namely, planar trees with l branches that are rooted (at k_1) and planted, implying that rotating asymmetric trees around the root results in a new tree. See Fig. 12 which depicts the five possible trees with $l = 3$ branches where we relabel the resultant $l + 1 = 4$ vertices as p_1, \dots, p_4 . In general, the number of different planar, rooted, planted trees with l branches is given by the l th Catalan number t_l [22]. In each of these trees, the $l + 1$ vertices p_1, \dots, p_{l+1} take values in either $D \setminus V_D$ or S . Hence, each tree T_i^l corresponds to a group of nonzero terms

$$T_i^l = \sum_{p_1, \dots, p_{l+1}} f_{T_i^l}(p_1, \dots, p_{l+1}), \quad i = 1, \dots, t_l. \quad (24)$$

Note that if a nonvanishing term in (18) corresponds to a graph with less than $l + 1$ vertices, then the corresponding graph possesses either edges with multiplicity larger than 2 or cycles, and this term is already accounted for in either one or more of the terms in (24). This fact can be observed by noticing that both edges with large multiplicity as well as cycles can be untied to get trees with l branches, with some of the $l + 1$ indices constrained however to share the same values (see Fig. 13). Note that such cases are not excluded in the summations in (24), thus we have

$$\mathbb{E}(\text{Tr}((\tilde{H}^* \tilde{H})^l)) \leq \sum_{i=1}^{t_l} T_i^l.$$

Below we show that

$$T_i^l \leq n(K'_1 \log n)^l, \quad \forall i \quad (25)$$

in a regular network and

$$T_i^l \leq n(K'_1 \log n)^{3l}, \quad \forall i \quad (26)$$

with high probability in a random network. We first concentrate on regular networks in order to reveal the proof idea in the simplest setting. A binning argument then allows to extend the result to random networks.

a) *Regular Network*: Recall that in the regular case, the nodes on the left half are located at positions $(-k_x + 1, k_y)$ and those on the right half at (i_x, i_y) for $k_x, k_y, i_x, i_y = 1, \dots, \sqrt{n}$. In this case, the matrix elements of \tilde{H} are given by

$$\tilde{H}_{ik} = \frac{e^{j\theta_{ik}}}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/4}} \frac{1}{\sqrt{d_{k_x, k_y}}}$$

and

$$d_{k_x, k_y} = \sum_{i_x, i_y=1}^{\sqrt{n}} \frac{1}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/2}}.$$

In the following discussion, we will need an upper bound on the scaling of $\sum_{i=1}^n |\tilde{H}_{ik}|^2$ and $\sum_{k=1}^n |\tilde{H}_{ik}|^2$. By Lemma 5.4, we have

$$d_{k_x, k_y} \geq K'_3 k_x^{2-\alpha}$$

for a constant $K'_3 > 0$ independent of n which, in turn, yields the upper bound

$$\begin{aligned} |\tilde{H}_{ik}|^2 &\leq \frac{1}{K'_3} \frac{k_x^{\alpha-2}}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/2}} \\ &\leq \frac{1}{K'_3} \frac{1}{(i_x + k_x - 1)^2 + (i_y - k_y)^2}. \end{aligned}$$

Summing over either i or k , and using the upper bound in Lemma 5.4 for $\alpha = 2$ yields

$$\sum_{i=1}^n |\tilde{H}_{ik}|^2, \quad \sum_{k=1}^n |\tilde{H}_{ik}|^2 \leq K'_1 \log n \quad (27)$$

where $K'_1 = \frac{K'_2}{K'_3}$ with K'_2 and K'_3 being the constants appearing in the lemma.

Let us first consider the simplest case where the tree is composed of l height 1 branches and denote it by T_1^l (see Fig. 14). We have

$$\begin{aligned} T_1^l &= \sum_{p_1, \dots, p_{l+1}=1}^n f_{T_1^l}(p_1, \dots, p_{l+1}) \\ &= \sum_{p_1, \dots, p_{l+1}=1}^n \left| \tilde{H}_{p_2 p_1} \right|^2 \left| \tilde{H}_{p_3 p_1} \right|^2 \cdots \left| \tilde{H}_{p_{l+1} p_1} \right|^2 \\ &= \sum_{p_1=1}^n \left(\sum_{p_2=1}^n \left| \tilde{H}_{p_2 p_1} \right|^2 \right)^l \\ &\leq n(K'_1 \log n)^l \end{aligned} \quad (28)$$

which follows from the upper bound (27).

Now let us consider the general case of an arbitrary tree T_i^l having s leaves, where $1 \leq s \leq l$ (see Fig. 15). Let the indices

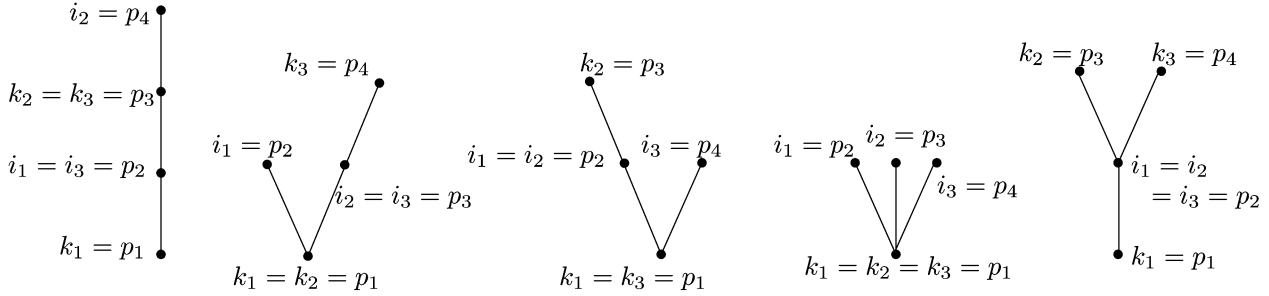


Fig. 12. Planar rooted planted trees with three branches. Note that each edge is actually a double edge in our case, although depicted with a single line in the figure.

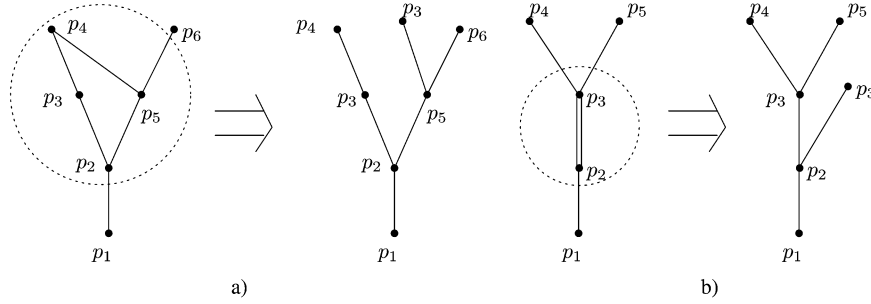


Fig. 13. The graphs with a) cycles and b) edges with large multiplicity can be united to get trees where some vertices are constrained to share the same indices.

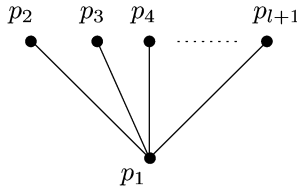


Fig. 14. A simple tree with l branches.

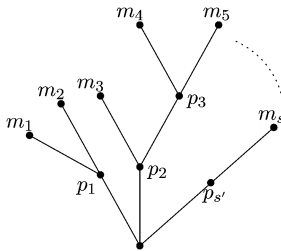


Fig. 15. A tree with leaves m_1, m_2, \dots, m_s .

corresponding to these leaves be m_1, \dots, m_s . Let us denote the “parent” vertices of these leaves by $p_1, \dots, p_{s'}$ and assume that p_1 is the common parent vertex of leaves m_1, \dots, m_{d_1} ; p_2 is the common parent vertex of leaves $m_{(d_1+1)}, \dots, m_{d_2}$ etc. and finally $p_{s'}$ is the parent of $m_{(d_{s'+1})}, \dots, m_s$. The term T_i^l corresponding to this tree is given by

$$\begin{aligned}
 T_i^l &= \sum_{\substack{m_1, \dots, m_s \\ p_1, \dots, p_{(l+1-s)}}} f_{T_i^l}(p_1, \dots, p_{l+1}) \\
 &= \sum_{p_1, \dots, p_{(l+1-s)}=1}^n f_{T_i^{l-s}}(p_1, \dots, p_{(l+1-s)}) \\
 &\quad \times \sum_{m_1, \dots, m_s=1}^n |\tilde{H}_{m_1 p_1}|^2 \dots |\tilde{H}_{m_{d_1} p_1}|^2 |\tilde{H}_{m_{(d_1+1)} p_2}|^2 \\
 &\quad \times \dots |\tilde{H}_{m_{d_2} p_2}|^2 \dots |\tilde{H}_{m_{(d_{s'+1})} p_{s'}}|^2 \dots |\tilde{H}_{m_s p_{s'}}|^2 \quad (29) \\
 &\leq T_i^{l-s} (K'_1 \log n)^s \quad (30)
 \end{aligned}$$

where T_i^{l-s} corresponds to a smaller (and shorter) tree $T_i^{(l-s)}$ with $l-s$ branches.⁶ The argument above decreases the height of the tree by 1, hence, it can be applied recursively to get a simple tree composed only of height 1 branches in which case the upper bound in (28) applies. Thus, given T_i^l let h be the number of recursions to get a simple tree and s_1, \dots, s_h denote the number of leaves in the trees observed at each step of the recursion. We have

$$\begin{aligned}
 T_i^l &\leq (K'_1 \log n)^{s_1} (K'_1 \log n)^{s_2} \dots (K'_1 \log n)^{s_h} T_1^{l-s_1 \dots -s_h} \\
 &\leq n (K'_1 \log n)^l
 \end{aligned}$$

since $T_1^{l-s_1 \dots -s_h} \leq n (K'_1 \log n)^{l-s_1 \dots -s_h}$ by (28). Thus, (25) follows.

b) Random Network: We denote the locations of the nodes to the left of the cut by $a_k = (-a_k^x, a_k^y)$ where a_k^x is the x -coordinate and a_k^y is the y -coordinate of node $k \in S$ and those to the right of the cut are similarly denoted by $b_i = (b_i^x, b_i^y)$ for $i \in D \setminus V_D$. In this case, the matrix elements of \tilde{H} are given by

$$\tilde{H}_{ik} = \frac{e^{j\theta_{ik}}}{\left((b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2 \right)^{\alpha/4} \sqrt{d_k}}$$

and

$$d_k = \sum_{i \in D \setminus V_D} \frac{1}{\left((b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2 \right)^{\alpha/2}}.$$

In parallel to the regular case, we will need an upper bound on $\sum_{i \in D \setminus V_D} |\tilde{H}_{ik}|^2$ and $\sum_{k \in S} |\tilde{H}_{ik}|^2$. The upper bound can be obtained in two steps by first showing that

$$d_k \geq K'_3 \frac{(a_k^x)^{2-\alpha}}{\log n} \quad (31)$$

⁶Note that the term corresponding to a leaf m can be either $|\tilde{H}_{mp}|^2$ or $|\tilde{H}_{pm}|^2$ depending on whether the height of the leaf is even or odd. However, in (29), we ignore this issue in order to simplify the notation since the upper bound (30) applies in both cases.

with high probability for a constant $K'_3 > 0$ independent of n , which leads to

$$\begin{aligned} |H_{ik}|^2 &\leq \frac{1}{K'_3} \log n \frac{(a_k^x)^{\alpha-2}}{\left((b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2\right)^{\alpha/2}} \\ &\leq \frac{1}{K'_3} \log n \frac{1}{(b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2} \end{aligned} \quad (32)$$

for all i, k . This, in turn yields

$$\sum_{k \in S} |\tilde{H}_{ik}|^2, \sum_{i \in D \setminus V_D} |\tilde{H}_{ik}|^2 \leq K'_1 (\log n)^3 \quad (33)$$

with high probability for another constant $K'_1 > 0$ independent of n . Recalling the leaf removal argument discussed for regular networks immediately leads to (26).

Both the lower bound in (31) and the upper bound in (33) regarding random networks can be proved using binning arguments that provide the connection to regular networks. In order to prove the lower bound, we consider Part b) of Lemma 5.1, while the upper bound (33) is proved using Part a) of the same lemma.

Let us first consider dividing the right half network into squarelets of area $2 \log n$. Given a left-hand side node k located at $(-a_k^x, a_k^y)$, let us move the nodes inside each right-hand side squarelet onto the squarelet vertex that is farthest to k . Since this displacement can only increase the Euclidean distance between the nodes involved, and since by Part b) of Lemma 5.1, we know that there is at least one node inside each squarelet, we have the equation at the bottom of the page by using the lower bound in Lemma 5.4.

Now having (32) in hand, in order to show (33), we divide the network into n squarelets of area 1. By Part a) of Lemma 5.1, there are at most $\log n$ nodes inside each squarelet. Considering the argument in Section V and the displacement of the nodes as illustrated in Fig. 7 yields a regular network with at most $2 \log n$ nodes at each vertex in the right half network

$$\begin{aligned} \sum_{i \in D \setminus V_D} |\tilde{H}_{ik}|^2 &\leq \frac{2}{K'_3} \log n \sum_{i \in D \setminus V_D} \frac{1}{(b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2} \\ &\leq \frac{4}{K'_3} (\log n)^2 \sum_{i_x, i_y=1}^{\sqrt{n}} \frac{1}{(i_x + k_x)^2 + (i_y - k_y)^2} \\ &\leq 4K'_1 (\log n)^3 \end{aligned}$$

by employing the upper bound in Lemma 5.4 for $\alpha = 2$. The same bound follows similarly for $\sum_{k \in S} |\tilde{H}_{ik}|^2$, thus the desired result in (33). \square

Proof of Lemma 5.4: Both the lower and upper bound for d_{k_x, k_y} can be obtained by straightforward manipulations. Recall that

$$d_{k_x, k_y} = \sum_{i_x, i_y=1}^{\sqrt{n}} \frac{1}{((i_x + k_x - 1)^2 + (i_y - k_y)^2)^{\alpha/2}}.$$

The upper bound can be obtained as follows:

$$\begin{aligned} d_{k_x, k_y} &= \sum_{y=1-k_y}^{\sqrt{n}-k_y} \sum_{x=k_x}^{k_x+\sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} \\ &\leq \sum_{y=1-k_y}^{\sqrt{n}-k_y} \left(\frac{1}{(k_x^2 + y^2)^{\alpha/2}} \right. \\ &\quad \left. + \int_{k_x}^{k_x+\sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} dx \right) \\ &\leq k_x^{-\alpha} + \int_{k_x}^{k_x+\sqrt{n}-1} \frac{1}{x^\alpha} dx \\ &\quad + \int_{1-k_y}^{\sqrt{n}-k_y} \frac{1}{(k_x^2 + y^2)^{\alpha/2}} dy \\ &\quad + \int_{1-k_y}^{\sqrt{n}-k_y} \int_{k_x}^{k_x+\sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} dx dy \\ &\leq k_x^{-\alpha} + (1 + \pi) k_x^{1-\alpha} + \int_{-\pi/2}^{\pi/2} \int_{k_x}^{3\sqrt{n}} \frac{1}{r^\alpha} r dr d\theta \end{aligned}$$

So

$$\begin{aligned} d_{k_x, k_y} &= \begin{cases} k_x^{-\alpha} + (1 + \pi) k_x^{1-\alpha} + \pi \log r \Big|_{k_x}^{3\sqrt{n}}, & \text{if } \alpha = 2 \\ k_x^{-\alpha} + (1 + \pi) k_x^{1-\alpha} + \frac{\pi}{(2-\alpha)} r^{2-\alpha} \Big|_{k_x}^{3\sqrt{n}}, & \text{if } \alpha > 2 \end{cases} \\ &\leq \begin{cases} K'_2 \log n, & \text{if } \alpha = 2 \\ K'_2 k_x^{2-\alpha}, & \text{if } \alpha > 2 \end{cases} \end{aligned} \quad (34)$$

for a constant $K'_2 > 0$ independent of n , since the dominating terms in (34) are the third ones. The lower bound follows similarly

$$\begin{aligned} d_{k_x, k_y} &= \sum_{y=1-k_y}^{\sqrt{n}-k_y} \sum_{x=k_x}^{k_x+\sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} \\ &\geq \sum_{y=1-k_y}^{\sqrt{n}-k_y} \int_{k_x}^{k_x+\sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} dx \end{aligned}$$

$$\begin{aligned} d_k &= \sum_{i \in D \setminus V_D} \frac{1}{\left((b_i^x + a_k^x)^2 + (b_i^y - a_k^y)^2\right)^{\alpha/2}} \\ &\geq \sum_{i_x, i_y=1}^{\sqrt{n/2 \log n}} \frac{1}{\left((i_x \sqrt{2 \log n} + a_k^x)^2 + (i_y \sqrt{2 \log n} - a_k^y)^2\right)^{\alpha/2}} \geq K'_3 \frac{(a_k^x)^{2-\alpha}}{2 \log n} \end{aligned}$$

$$d_{k_x, k_y} = \begin{cases} \arctan\left(\frac{1}{2}\right) \log r \Big|_{\sqrt{2}k_x}^{k_x + \sqrt{n}-1} + \frac{1}{\alpha-1} x^{1-\alpha} \Big|_{k_x}^{k_x + \sqrt{n}-1}, & \text{if } \alpha = 2 \\ \arctan\left(\frac{1}{2}\right) \frac{1}{2-\alpha} r^{2-\alpha} \Big|_{\sqrt{2}k_x}^{k_x + \sqrt{n}-1} + \frac{1}{\alpha-1} x^{1-\alpha} \Big|_{k_x}^{k_x + \sqrt{n}-1}, & \text{if } \alpha > 2 \end{cases}$$

$$\geq K'_3 k_x^{2-\alpha} \quad (35)$$

$$\begin{aligned} &\geq \int_{1-k_y}^{\sqrt{n}-k_y} \int_{k_x}^{k_x + \sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} dx dy \\ &\quad - \int_{k_x}^{k_x + \sqrt{n}-1} \frac{1}{x^\alpha} dx \\ &\geq \int_0^{\sqrt{n}} \int_{k_x}^{k_x + \sqrt{n}-1} \frac{1}{(x^2 + y^2)^{\alpha/2}} dx dy \\ &\quad + \frac{x^{1-\alpha}}{\alpha-1} \Big|_{k_x}^{k_x + \sqrt{n}-1} \\ &\geq \int_0^{\arctan(1/2)} \int_{\sqrt{2}k_x}^{k_x + \sqrt{n}-1} \frac{1}{r^\alpha} r dr d\theta + \frac{x^{1-\alpha}}{\alpha-1} \Big|_{k_x}^{k_x + \sqrt{n}-1} \end{aligned}$$

So for all $\alpha \geq 2$, we have (35) at the top of the page, where $K'_3 > 0$ is a constant independent of n , since the dominating terms in (35) are the first ones. This concludes the proof of the lemma. \square

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