

A Geometrical Approach for Power Optimization in Relay-Based Cellular Networks

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Abstract—Relays are considered as a feasible way to increase coverage and capacity of cellular networks. However, optimization of relay-enhanced networks is still a matter of research. This paper proposes a simple geometrical approach to derive the optimal long-term relay downlink transmit power with the aim of maximizing the average spectral efficiency in a relay-based cellular network. In our model, a User Equipment (UE) is served by the network node from which it receives the highest power. The spectral efficiency at the UE location is maximized when this serving node is also the one that provides the best Signal to Interference Ratio (SIR). Our approach thus consists in setting the relay node (RN) power such that its *serving area* and its *SIR area* coincide as much as possible. Results show that our algorithm provides a fine approximation of the optimal RN transmit power under a wide range of RN location scenarios and propagation parameters.

I. INTRODUCTION

The deployment of Relay Nodes (RN) is a central feature of next generation cellular networks (LTE-A and IEEE 802.16j). In this paper, we consider decode-and-forward non-cooperative RNs [1] and we focus on the downlink. Our goal is to optimize the RNs transmit power on the long-term in order to maximize the cell average spectral efficiency.

Several papers in the literature tackle the problem of optimal transmit power in relay-assisted cellular networks. In [2], authors come up with a convex weighted power minimization problem under rate, bandwidth and power constraints and give some insights on the optimal solution. In [3], a system capacity maximization problem is formulated under adaptive sub-channel, path and

power allocation. A heuristic is proposed to approximately solve the mixed integer programming problem. Authors of [4] maximize the system goodput with appropriate sub-band, power and rate allocation, while taking into account fairness and imperfect channel information. The main drawbacks of these approaches (and other in the literature, e.g. [5]) are the computational complexity and the induced protocol overhead. Proposed solutions indeed require to be regularly updated on a frame-by-frame basis in order to adapt RN and Base-Station (BS) powers to variable traffic, UEs channel conditions, etc. On the contrary, setting a constant transmit power has the advantage of simplicity and is a common assumption in multi-cell performance evaluations [6], [7]. In this case, the need arises for optimizing capacity by appropriately setting, at the network deployment, the RN power for a long term.

In this paper, we thus derive an approximation of the long-term optimal RN transmit power. Our approach is based on matching the *serving area* of a RN (the set of locations served by the relay) with its *SIR area* (the area where the relay provides the best SIR) as closely as possible. We use the fluid model [8] as a means to approximate the average SIR. The implementation of our method improves network performance without requiring any additional protocols overhead. Moreover, no extra computational burden for BSs is involved, and RNs transmit power does not need to be regularly updated. We show that our algorithm yields accurate results, as compared to the actual optimal long-term RNs transmit power, in numerous simulation scenarios. Algorithm complexity decreases when fluid model is used to compute SIR, as well as results accuracy. This

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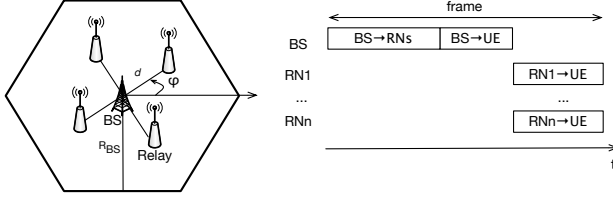


Fig. 1. Relay location (left) and radio resource partitioning (right).

tradeoff is exhaustively discussed.

The paper is organized as follows: Section II describes the system model. Section III formulates the optimization problem and Section IV presents our maximization approach. Section V gives numerical results, while Section VI concludes the paper.

II. SYSTEM MODEL

We consider the downlink of a single-frequency hexagonal cellular network in an urban environment. Let the cell radius be less than 1km. One BS is located in the center of each cell. The BS is assisted by N_{RN} RN, placed on a circle of radius d around the BS. RN are regularly spaced on the circle (see e.g. [2] [9]). Let ϕ be the angle shift of the first RN with respect to the horizontal axis (see Fig. 1, left). BSs and RNs transmit respectively with power P_{BS} and P_{RN} , with $P_{RN} < P_{BS}$ [1]. The set of network nodes is denoted with $\mathcal{S} = \mathcal{S}_R \cup \mathcal{S}_B$, where \mathcal{S}_R represents the set of RNs and \mathcal{S}_B is the set of BSs of cardinality $|\mathcal{S}_B| = B$. Both BS and RN are equipped with a single omni-directional antenna and they transmit in division mode of operations, i.e., using orthogonal time-frequency resources (see e.g. [10]). An example of time division is shown in Fig. 1 (right).

A. Serving and SIR areas

Received power at location z from node i is given by:

$$P_i(z) = P_{tx,i} K d_{i,z}^{-\eta}, \quad (1)$$

where $P_{tx,i} \in \{P_{BS}, P_{RN}\}$ is the transmit power of i , K is a constant, $d_{i,z}$ is the distance between i and z and η is the propagation exponent. UEs are supposed to be distributed over the cell area according to a uniform probability distribution (homogeneous spatial Poisson process). UEs are served by the node (the *best server*) providing the highest received power as e.g. in [7]. We define the *serving area* of i as the set of locations where i is the serving station.

We consider thermal noise to be negligible compared to the sum of in-cell and out-of-cell interference, in agreement with [8], which starts from the same assumptions made in this paper. Hence, Signal to Interference Ratio (SIR) is chosen as radio channel performance measure. At location z , the SIR provided by node i is given by:

$$\Gamma_i(z) = \frac{P_{tx,i} K d_{i,z}^{-\eta}}{\sum_{k \in \mathcal{S}_i \setminus \{i\}} P_{tx,k} K d_{k,z}^{-\eta}} = \frac{d_{i,z}^{-\eta}}{\sum_{k \in \mathcal{S}_i \setminus \{i\}} d_{k,z}^{-\eta}}, \quad (2)$$

where \mathcal{S}_i is the set of nodes of the same kind of i (BS or RN). We define the *SIR area* of i as the set of locations where i provides the best SIR.

Fig. 2 illustrates the notions of serving and SIR areas. It focuses on a BS i and a RN j in a cell \mathcal{A}_c . $\mathcal{A}_{ij}^P(P_{RN})$ denotes the area (delimited by the curve $\pi_{ij}(P_{RN})$), where the power received by i is higher than the one received by j . The serving area of i is the intersection of all \mathcal{A}_{ij}^P , where j is the index of a RN of the cell. \mathcal{A}_{ij}^S is the area, where the SIR obtained from i is higher than the SIR obtained from j and the SIR area is the intersection of such areas in the cell.

According to (1) and (2), the SIR area is independent of the nodes transmit power, whereas the serving area depends on P_{RN} , and can be smaller or greater than the SIR area according to it. We denote $\mathcal{D}_{ij}(P_{RN})$ the symmetric difference of the two sets in the cell, i.e.,

$$\mathcal{D}_{ij}(P_{RN}) \triangleq \mathcal{A}_c \cap (\mathcal{A}_{ij}^S \triangle \mathcal{A}_{ij}^P), \quad (3)$$

where the symmetric difference \triangle is defined for any two sets A and B as follows: $A \triangle B \triangleq (A \setminus B) \cup (B \setminus A)$. Let $\mathcal{D}_i \triangleq \cup_{j \in \mathcal{S}_R} \mathcal{D}_{ij}$. \mathcal{D}_i represents the area in the cell, where UEs are not served by the best station.

B. Fluid Model

Under our assumptions $\Gamma_i(z)$ can be approximated by means of the fluid model [8] for the regular hexagonal network formed by the BSs. The fluid model is a powerful tool for simplifying SIR formulas in a wireless network. The main idea is to replace a discrete set of transmitters by a continuum and thus transform discrete sums into integrals.

If $i \in \mathcal{S}_B$ and the network is large we have:

$$\Gamma_i(z) \approx \frac{\eta - 2}{2\pi\rho_{BS}d_{i,z}^\eta(2R_{BS} - d_{i,z})^{2-\eta}}, \quad (4)$$

where ρ_{BS} is the BS density and R_{BS} is the half inter-BS distance. Strictly speaking, the fluid model is not

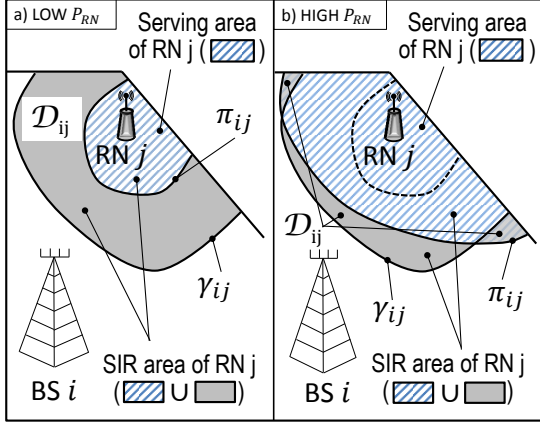


Fig. 2. SIR and serving areas of RN j .

applicable to the generally non-regular RN network. However, we show that in a number of scenarios this method still provides a fair approximation of $\Gamma_i(z)$, $i \in \mathcal{S}_R$ (see Section V) provided that R_{BS} is replaced by R_{RN} , the minimum half inter-RN distance, and ρ_{BS} by ρ_{RN} , the RN density.

C. Spectral Efficiency

The spectral efficiency at location z is defined as follows:

$$C(z, P_{RN}, P_{BS}) = \log_2(1 + \Gamma_{b(z, P_{RN}, P_{BS})}(z)), \quad (5)$$

where $b(z, P_{RN}, P_{BS})$ is the serving station of z (as we focus on P_{RN} optimization, the dependence on P_{BS} is skipped in the following developments). We also define $C_H(z)$ as the spectral efficiency in z associated to the station providing the best SIR in z .

III. PROBLEM FORMULATION

Our aim is to maximize the cell average spectral efficiency on the downlink, via a proper tuning of the RN transmit power. The optimization is focused on a generic cell \mathcal{A}_c of area A_c , centered at BS i . We formulate the problem as follows:

$$P_{RN}^* = \operatorname{argmax}_{P_{RN}} \{\bar{C}(P_{RN})\}, \quad (6)$$

under the constraint:

$$0 < P_{RN} < P_{BS}, \quad (7)$$

where $\bar{C}(P_{RN}) = \frac{1}{A_c} \int_{\mathcal{A}_c} C(z, P_{RN}) dz$ is the cell average spectral efficiency.

Ideally, the highest $\bar{C}(P_{RN})$ is reached for $\mathcal{D}_i \equiv \emptyset$, i.e., when all UEs are served by the station which can

guarantee them the highest SIR. However, this condition cannot be achieved in general, as there is normally no P_{RN} for which $\mathcal{A}_{ij}^P \equiv \mathcal{A}_{ij}^S$.

Let us work out (6) as follows:

$$\begin{aligned} P_{RN}^* &= \operatorname{argmin}_{P_{RN}} \{-\bar{C}(P_{RN})\} \\ &= \operatorname{argmin}_{P_{RN}} \{K - \bar{C}(P_{RN})\} \\ &= \operatorname{argmin}_{P_{RN}} \left\{ \frac{1}{A_c} \int_{\mathcal{A}_c} (C_H(z) - C(z, P_{RN})) dz \right\}, \end{aligned} \quad (8)$$

where $K \in \mathbb{R}$ is an arbitrary constant. The second equality is valid because $\frac{1}{A_c} \int_{\mathcal{A}_c} C_H(z) dz$ is a constant w.r.t. P_{RN} . Now, the difference in the integrand is zero outside \mathcal{D}_i . Hence (8) is equivalent to

$$P_{RN}^* = \operatorname{argmin}_{P_{RN}} \left\{ \frac{1}{A_c} \int_{\mathcal{D}_i} (C_H(z) - C(z, P_{RN})) dz \right\}. \quad (9)$$

Solving (9) requires finding the borders of \mathcal{D}_i and computing the integral for every candidate value of P_{RN} . However, it is possible to restrict the range of P_{RN} to be explored. This range being very tight, it is actually not even needed to evaluate the integral, as it will be shown in the next section. Solution is proposed for the simple system model hypothesized in this paper. However, our approach can still be applied with few modifications to more complex scenarios, i.e., when dealing with more complex path-loss or non-hexagonal BS network.

IV. PROPOSED SOLUTION

In this section, we first derive the borders of the area \mathcal{D}_{ij} and then show that the spectral efficiency is first increasing and then decreasing with P_{RN} . This allows us to propose a fast algorithm to numerically solve the optimization problem.

A. Derivation of $\pi_{ij}(P_{RN})$ and γ_{ij}

Let us again focus on a pair of BS i and RN j in a cell \mathcal{A}_c . The expression of $\pi_{ij}(P_{RN})$ is found by equating the received powers from i and j . With the axis system shown on Fig. 2 and using (1), we can write for $(x, y) \in \mathcal{A}_c$:

$$\pi_{ij}(P_{RN}) : (x - x_\delta)^2 + (y - y_\delta)^2 = r^2, \quad (10)$$

where $r = \sqrt{\Omega(x_{RN}(j)^2 + y_{RN}(j)^2)/(\Omega - 1)^2}$, $\Omega = (P_{RN}/P_{BS})^{2/\eta}$ and $(x_{RN}(j), y_{RN}(j))$ is the location

of RN j . Hence, $\pi_{ij}(P_{RN})$ is a part of a circle of radius r , increasing with P_{RN} . The circle is centered at $(x_\delta, y_\delta) = (x_{RN}/(1-\Omega), y_{RN}/(1-\Omega))$, along the line connecting i and j . It is easy to verify that increasing P_{RN} has the effect shown on Fig. 2, i.e., the set \mathcal{A}_{ij}^P is decreasing. Similarly to (10), γ_{ij} is found by equating the SIR expressions related to i and j . If we apply the fluid model, we obtain for $(x, y) \in \mathcal{A}_c$:

$$\gamma_{ij} : \frac{\rho_{BS}}{\rho_{RN}} \left(\frac{d_i(x, y)}{d_j(x, y)} \right)^\eta = \left(\frac{2R_{RN} - d_j}{2R_{BS} - d_i} \right)^{2-\eta}, \quad (11)$$

where $d_i(x, y) = \sqrt{x^2 + y^2}$ and $d_j(x, y) = \sqrt{(x - x_{RN}(j))^2 + (y - y_{RN}(j))^2}$. If the fluid model does not provide a sufficient approximation, the SIR can be directly computed with (2). Note again that varying P_{RN} does not affect the equation of γ_{ij} .

B. Derivation of Bounds on P_{RN} Search Range

We see from Fig. 2 that starting from a low RN power and increasing its value has the effect to first decrease and then increase the area \mathcal{D}_{ij} , while average spectral efficiency first increases and then decreases. This intuition is formalized as follows.

Lemma 1: (i) Suppose that for every RN j in the cell, $\mathcal{A}_{ij}^S \subseteq \mathcal{A}_{ij}^P(P_{RN})$, then $\bar{C}(P_{RN})$ is increasing in P_{RN} . (ii) Suppose that for every RN j in the cell, $\mathcal{A}_{ij}^P(P_{RN}) \subseteq \mathcal{A}_{ij}^S$, then $\bar{C}(P_{RN})$ is decreasing in P_{RN} .

Proof: For (i) consider two RN powers $P_2 > P_1$ for which the condition is valid. Using (10), it is easy to verify that $\mathcal{A}_{ij}^P(P_2) \subseteq \mathcal{A}_{ij}^P(P_1)$. Using now the definition of \mathcal{D}_{ij} in (3) and the lemma first condition, we obtain $\mathcal{D}_{ij}(P_2) \subseteq \mathcal{D}_{ij}(P_1)$ for every j . For $P_{RN} = P_2$ integral (9) is hence computed on a subset of the area associated to P_1 and the integrand is positive, which implies $\bar{C}(P_2) \geq \bar{C}(P_1)$. For (ii) the proof is similar, except that now $\mathcal{D}_{ij}(P_1) \subseteq \mathcal{D}_{ij}(P_2)$. ■

According to lemmas 1, P_{RN}^* must thus lie between a lower and upper bound: $P_{RN}^L \leq P_{RN}^* \leq P_{RN}^U$, where P_{RN}^L and P_{RN}^U are respectively the highest and the lowest P_{RN} for which the conditions $\mathcal{A}_{ij}^S \subseteq \mathcal{A}_{ij}^P(P_{RN})$ and $\mathcal{A}_{ij}^P(P_{RN}) \subseteq \mathcal{A}_{ij}^S$ are valid for all j . Note that in general we cannot draw any conclusion about spectral efficiency when $P_{RN}^L \leq P_{RN} \leq P_{RN}^H$. Bounds can be interpreted as follows:

- P_{RN}^L is the RN power for which $\mathcal{A}_{ij}^S \subseteq \mathcal{A}_{ij}^P(P_{RN})$ and $\pi_{ij}(P_{RN})$ is tangent to γ_{ij} for some j .

Algorithm 1 Optimization Algorithm

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1: Initialization:  $\mathcal{P}_i = \emptyset$ 
2: Set up a grid  $\mathcal{G}$  of  $N$  points paving the cell area
3: for each relay  $j$  in the cell do
4:   Find a set of points  $p_n^{ij} \in \gamma_{ij}$  using (2) or (4) on  $\mathcal{G}$ 
5:   for each  $p_n^{ij} = (x_n^{ij}, y_n^{ij}) \in \gamma_{ij}$  do
6:     Compute  $\tilde{P}_{RN}(p_n^{ij})$  using (12)
7:      $\mathcal{P}_i \leftarrow \mathcal{P}_i \cup \{\tilde{P}_{RN}(p_n^{ij})\}$ 
8:   end for
9: end for
10:  $(P_{RN}^L, P_{RN}^U) \leftarrow (\min \mathcal{P}_i, \max \mathcal{P}_i)$ 
11: Return  $P_{RN}^* = (P_{RN}^L + P_{RN}^U)/2$ 

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- P_{RN}^U is the RN power for which $\mathcal{A}_{ij}^P(P_{RN}) \subseteq \mathcal{A}_{ij}^S$ and $\pi_{ij}(P_{RN})$ is tangent to γ_{ij} for some j .

The bounds cannot be derived analytically since (11) has no explicit solution. We can however obtain approximations in three steps:

- 1) We find numerically a set of points $\{p_n^{ij}\}_{n \in [1, N_{ij}]} \in \gamma_{ij}$. For that, we build a regular grid of N points on the cell area and, using (2) or (4), we find the p_n^{ij} by interpolation.
- 2) For each $p_n^{ij} = (x_n^{ij}, y_n^{ij})$, we find the associated RN power $\tilde{P}_{RN}(p_n^{ij})$, defined as the power for which x_n^{ij} and y_n^{ij} satisfy the equation of $\pi_{ij}(P_{RN})$. It is found by substituting x_n^{ij} and y_n^{ij} into (10), and solving for P_{RN} , obtaining:

$$\tilde{P}_{RN}(p_n^{ij}) = P_{BS} (l_{nj}^2 / l_{n0}^2)^{\frac{\eta}{2}}, \quad (12)$$

where $l_{nj}^2 = (x_n^{ij} - x_{RN}(j))^2 + (y_n^{ij} - y_{RN}(j))^2$ is the squared distance to RN j and $l_{n0}^2 = x_n^{ij2} + y_n^{ij2}$ is the squared distance to the BS. We repeat the operation for all j and we collect the obtained RN power values in a set $\mathcal{P}_i = \cup_j \{\tilde{P}_{RN}(p_n^{ij})\}$.

- 3) We compute: $P_{RN}^L = \min \mathcal{P}_i$, $P_{RN}^U = \max \mathcal{P}_i$.

The accuracy of the approach improves with N . Given that the bounds are usually tight, we propose to approximate P_{RN}^* by: $P_{RN}^* \approx (P_{RN}^L + P_{RN}^U)/2$. Hence, using this approximation, we do not need to compute integral (9) for several values of P_{RN} . Our approach is summarized in Algorithm 1.

V. NUMERICAL RESULTS

A. Simulation Settings

Two alternatives are compared for our proposal, which differ at step 4 of the algorithm: use either the fluid model ((4) and (11)) or the exact SIR formulation (2).

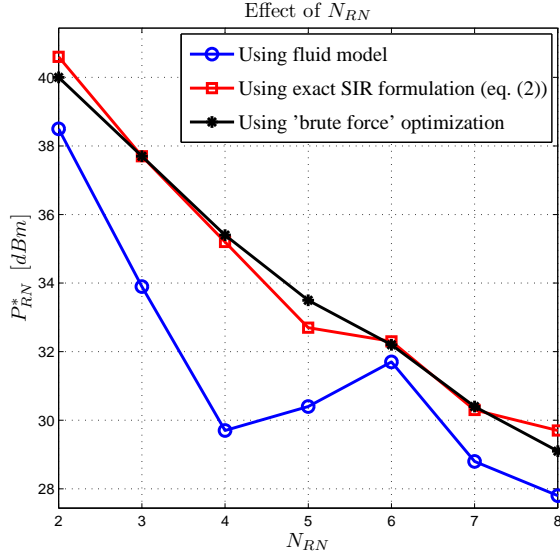


Fig. 3. Effect of N_{RN} ($d=3/8D$, $\phi=0$, $\eta=3.5$)

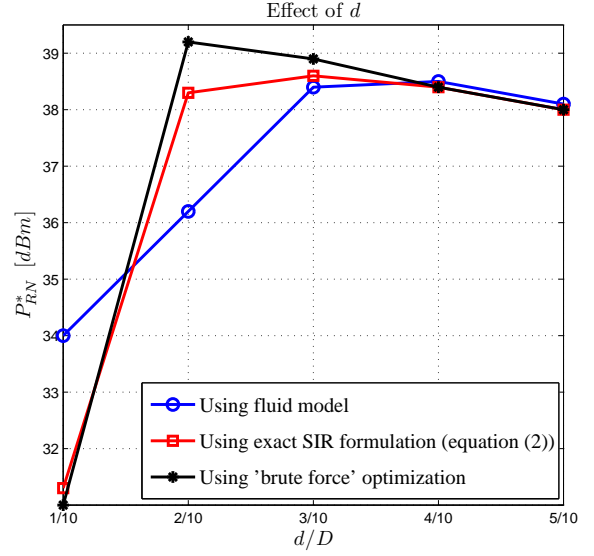


Fig. 5. Effect of d ($N_{RN}=3$, $\eta=3$, $\phi=30$)

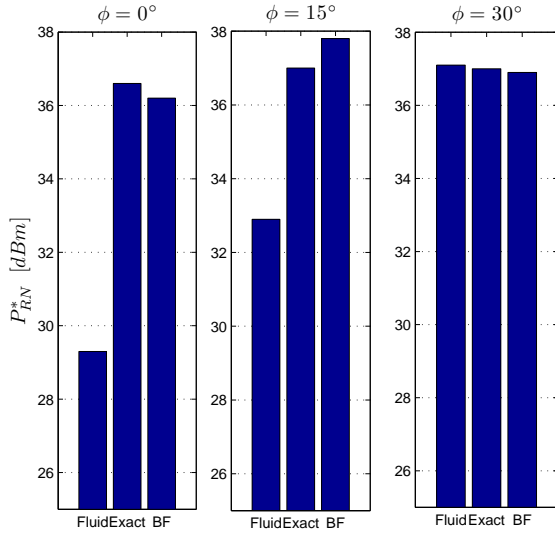


Fig. 4. Effect of ϕ ($N_{RN}=3$, $\eta=3.5$ and $d=9/20D$)

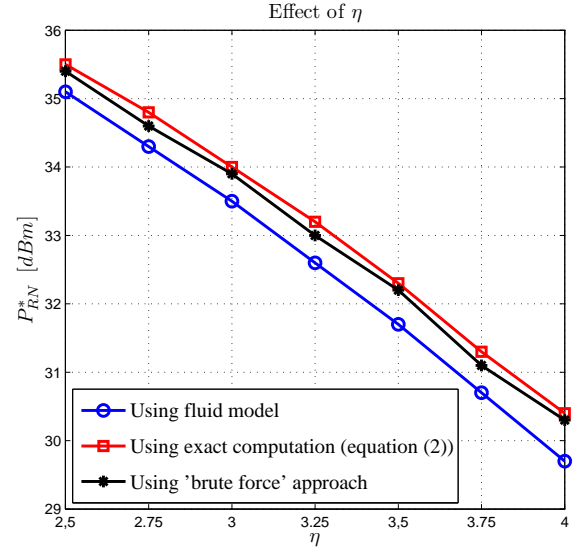


Fig. 6. Effect of η ($N_{RN}=6$, $d=3/8D$, $\phi=0$)

B. Results Analysis

We consider a network of 4 rings of cells around a central cell. We set the inter-BS distance to $D = 2R_{BS} = 500$ m. Computations are performed on a grid of points of \mathcal{A}_c of resolution 10 m. We set $P_{BS} = 46$ dBm [1]. Results are compared with the value of P_{RN}^* obtained by means of brute force optimization: SIR and spectral efficiency are computed for each point of a grid of resolution equal to 10 m, and for each considered RN power between 25 and 40 dBm with a 0.1 dB step.

Figure 3 shows the effect of N_{RN} . As expected, P_{RN}^* is a decreasing function of N_{RN} because when the RN density increases, their SIR areas shrink. In all cases, the optimal power found using Eq. (2) is within 1 dBm of the optimal (in most cases, even within 0.5 dBm). On Figures 4 and 5, we see that ϕ and d have also little impact on the accuracy of P_{RN}^* . The fluid model provides faster results at the price of reduced accuracy. It is known that the fluid model is more accurate when the

network is regular or close to be regular. This explains why results are better when N_{RN} increases and when RNs are not too close from the BS. When $\phi = 0$ and $d = 9/20D$, RNs from different cells are clustered close to the cell corners so that the obtained RN network is very irregular. In this case, the fluid model applied to our algorithm yields a very inaccurate result. The precision improves for growing ϕ because RNs tend to be more regularly spaced, and for $\phi = 30$ deg the difference with the optimal value is approximately 0.2 dBm.

Figure 6 shows that η has little influence on the accuracy of our approach. In the chosen scenario, the fluid model provides also a good approximation.

We have also performed simulations considering log-normal shadowing, and compared our approach with the result of Monte Carlo simulations, finding out that our algorithm is still within 0.3 dBm of the optimal value even for shadowing standard deviation $\sigma = 10dB$.

Finally our algorithm is computationally faster compared to brute force optimization. The latter requires $O(N_P N B N_{RN})$ operations. We notice that $|\mathcal{P}_i|$ is negligible compared to N and to $B N_{RN}$, so that Algorithm 1 using (2) requires $O(N B N_{RN})$ operations, while using (4) it requires $O(B N_{RN} + N)$ operations. We have also observed that N can be kept sensibly lower in our algorithm, compared to brute force. As an illustrative example, using Matlab for the presented scenarios, the savings in terms of computational time are of several orders of magnitude (from few seconds for our algorithm to several hours for the brute force).

VI. CONCLUSION

In this paper we have proposed a fast method to compute, for a given RN placement pattern, the RN transmit power optimizing the average cell spectral efficiency. This method is based on matching the SIR area of stations with their serving area, as closely as possible. Fluid model is proposed as a means for computing the SIR, because of the lower computational burden associated with it, as compared to straightforward SIR computation. However, the increased speed comes at the price of an accuracy loss. Results show that our method yields accurate results, particularly when an exact SIR formulation is adopted. However, accuracy is still high on a number of scenarios, when fluid model is used. The tradeoff between computational speed and precision is

investigated. Future work may study the impact of bursty traffic and backhaul link on performance.

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