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Centralized versus Decentralized Multi-Cell Resource and Power Allocation for Multiuser OFDMA Networks

Mohamad Yassin^{a,b,*}, Samer Lahoud^a, Kinda Khawam^c, Marc Ibrahim^b,
Dany Mezher^b, Bernard Cousin^a

^a*University of Rennes 1, IRISA, Campus de Beaulieu, 35042 Rennes, France*

^b*Saint Joseph University of Beirut, ESIB, CST, Mar Roukoz, Lebanon*

^c*University of Versailles, 45 Av. des Etats-Unis, 78035 Versailles, France*

Abstract

The exponential growth in the usage of mobile networks along with the increasing number of User Equipments (UEs) are exacerbating the scarcity of frequency resources. Dense frequency reuse on the downlink of multiuser Orthogonal Frequency Division Multiple Access networks leads to severe Inter-Cell Interference (ICI) problems. Resource and power allocation techniques are required to alleviate the harmful impact of ICI. Contrarily to the existing techniques that consider single-cell resource and power allocation problem without taking ICI into account, we formulate a centralized downlink multi-cell joint resource and power allocation problem. The objective is to maximize system throughput while guaranteeing throughput fairness between UEs. We demonstrate that the joint problem is separable into two independent problems: a resource allocation problem and a power allocation problem. Lagrange duality theory is used to solve the centralized power allocation problem. We also tackle the resource and power allocation problem differently by addressing it in a decentralized manner. We propose a non-cooperative downlink power allocation approach based on game theory. The players are the base stations, and each base station seeks to maximize its own utility function. We investigate the convergence of our proposed centralized and decentralized approaches, and we

*Corresponding author

Email address: mohamad.yassin@usj.edu.lb (Mohamad Yassin)

compare their performance with that of state-of-the-art approaches.

Keywords: Convex optimization, resource and power allocation, inter-cell interference, ICIC, OFDMA

1. Introduction

Multiuser Orthogonal Frequency Division Multiple Access (OFDMA) networks, such as the Third Generation Partnership Project (3GPP) Long Term Evolution (LTE) [1] and LTE-Advanced (LTE-A) [2] networks, are able to avoid
5 the negative impact of multipath fading and intra-cell interference, by virtue of the orthogonality between subcarrier frequencies. Nevertheless, Inter-Cell Interference (ICI) problems arise on the downlink of dense frequency reuse networks due to simultaneous transmissions on the same frequency resources. System performance is interference-limited, since the achievable throughput is reduced
10 due to ICI.

Fractional Frequency Reuse (FFR) [3] and Soft Frequency Reuse (SFR) [4] were introduced to avoid the harmful impact of ICI on system performance, by applying static rules on Resource Block (RB) usage and power allocation between cell-center and cell-edge users. Heuristic Inter-Cell Interference Co-
15 ordination (ICIC) techniques are proposed to achieve ICI mitigation without severe degradation of the overall system throughput. For instance, authors of [5, 6] propose suboptimal solution for the resource allocation problem. The objective is to minimize ICI by exploiting User Equipment (UE) diversity to maximize system throughput. They propose a two-level algorithm that oper-
20 ates at the evolved-NodeBs (eNodeBs) and at a central controller connected to several eNodeBs. In [7], a heuristic power allocation algorithm is introduced to reduce energy consumption and to improve cell-edge UEs throughput. It has been proven that the proposed algorithm reduces power consumption without reducing the achievable throughput. Moreover, it mitigates ICI and increases
25 the achievable throughput for cell-edge UEs.

Beside heuristic resource and power allocation algorithms [8], convex opti-

mization is used to improve the performance of multiuser OFDMA networks, and to alleviate the negative impact of ICI on UE throughput. Resource and power allocation problem is usually formulated as nonlinear optimization problem, where the objective consists in maximizing system throughput, spectral efficiency, or energy efficiency, with constraints on the minimum throughput per UE or other Quality of Service (QoS) parameters [9]. The exponential growth in the usage of mobile networks along with the increasing number of UEs are exacerbating the scarcity of frequency resources.

The majority of state-of-the-art contributions formulate the resource and power allocation problem for a single cell network [10, 11, 12], or do not consider the impact of ICI on system performance. For instance, the tradeoff between spectral efficiency and energy efficiency is addressed in [12], and a low-complexity suboptimal algorithm is proposed to allocate RBs for practical applications of the tradeoff. However, the system model consists of a single cell OFDMA network, where one subcarrier is assigned to at most one UE. Therefore, ICI problems are not considered. In this article, we formulate the joint resource and power allocation problem for the downlink of multiuser OFDMA networks, as a centralized multi-cell optimization problem. Inter-cell interference is taken into account, and throughput fairness between the different users is guaranteed. We prove that our joint problem is separable into two independent optimization problems: a resource allocation problem and a power allocation problem. Our objective is to maximize system throughput, while satisfying constraints related to resource usage, Signal-to-Interference and Noise Ratio (SINR), and power allocation. We also propose a decentralized power allocation approach that does not rely on centralized controllers. Each base station maximizes its own utility function in a distributed manner. We evaluate the performance of the proposed approaches, and we compare their performance with state-of-the-art resource and power allocation approaches.

The remainder of this article is organized as follows. In section II, we describe the limitations of the existing state-of-the-art approaches. In section III, system model is presented followed by our joint resource and power allocation

problem formulation. The joint problem is decomposed into two independent problems in section IV: a resource allocation problem and a power allocation
60 problem. We also demonstrate the convexity of the formulated problems. In section V, we solve both resource and power allocation problems using the Lagrange duality theory. Our decentralized power allocation approach is introduced in section VI. Then, we investigate the convergence of the centralized and the decentralized approaches in section VII, where we also provide compar-
65 isons with other approaches. Section VIII concludes this article and summarizes our main contributions.

2. Related Work

2.1. State-of-the-Art Contributions

For a given multiuser OFDMA network, resource and power allocation prob-
70 lem is formulated as a centralized optimization problem [10, 11, 12]. Centralized inter-cell coordination is therefore required to find the optimal solution, where the necessary information about SINR, power allocation, and resource usage are sent to a centralized coordination entity.

In [13], the multi-cell optimization problem is decomposed into two dis-
75 tributed optimization problems. The objective of the first problem is to minimize the transmission power allocated for cell-edge UEs, while guaranteeing a minimum throughput for each UE. RB and power are allocated to cell-edge UEs so that they satisfy their minimum required throughput. The remaining RBs and the remaining transmission power are uniformly allocated to cell-center
80 UEs. At this stage, the second problem aims at finding the resource allocation strategy that maximizes the achievable throughput for cell-center UEs. An improved version of this adaptive ICIC technique is proposed in [14], where resource allocation for cell-edge UEs is performed depending on their individual channel conditions. However, the main disadvantage of this adaptive ICIC tech-
85 nique and the proposed improvement is that they do not consider the impact of ICI between adjacent cells when power allocation is performed. Each cell solves

its own optimization problem without requesting additional information from its neighboring cells.

Resource and power allocation for a cluster of coordinated OFDMA cells are studied in [15]. Energy efficiency is maximized under constraints related to the downlink transmission power. However, noise-limited regime is considered, and ICI is neglected. Moreover, energy-efficient resource allocation for OFDMA systems is investigated in [16], where generalized and individual energy efficiencies are defined for the downlink and the uplink of the OFDMA system, respectively. Properties of the energy efficiency objective function are studied, then a low-complexity suboptimal algorithm is introduced to reduce the computational burden of the optimal solution. Subcarrier assignment is made easier using heuristic algorithms. Authors of [17] consider the joint resource allocation, power allocation, and Modulation and Coding Scheme (MCS) selection problem. The joint optimization problem is separated into resource allocation and power allocation problems, and suboptimal algorithms are proposed. Another low complexity suboptimal resource allocation algorithm is proposed in [18]. The objective consists in maximizing the achievable throughput, under constraints related to resource usage in the different cells. Cooperation between adjacent cells is needed. A multi-cell resource allocation approach for OFDMA systems with decode-and-forward relaying is proposed in [19], where an interference constraint is introduced along with time sharing variables. Although this approach guarantees throughput fairness between the different users, the spectral efficiency is reduced since the cells are not allowed to use the available spectrum during 100% of the time due to time sharing between base stations and relays.

Minimizing energy consumption and maximizing spectral efficiency in multiuser OFDMA networks cannot be achieved simultaneously. Energy-bandwidth tradeoff is studied in [20], where authors consider the total energy consumption versus the end-to-end rate in wireless multihop networks. For an arbitrary placement of wireless nodes, resource and power allocation that minimizes the energy level required to achieve a given data rate is found. However, interference-free

resource allocation is considered, and the impact of ICI on system performance is not taken into account.

120 2.2. Our Contributions

The majority of state-of-the-art contributions that formulate spectral efficiency or energy efficiency problems as centralized optimization problems, neglect the impact of ICI on system performance [10, 11, 12], or introduce suboptimal approaches to solve resource and power allocation problems [21, 22, 23].
 125 Moreover, performance comparisons are not made with other distributed heuristic ICIC algorithms, that are usually characterized by a lower computational complexity. In our work, we consider the multi-cell downlink resource and power allocation problem, where the objective is to maximize system throughput while guaranteeing throughput fairness between the different UEs. Moreover, ICI is
 130 taken into account when solving the centralized resource and power allocation problem. We also formulate a decentralized non-cooperative power allocation approach based on game theory. The players are the cells, and each cell seeks maximizing its own utility function independently of the other cells in the network. We investigate the convergence of both centralized and decentralized ap-
 135 proaches, and we compare their performance with that of the frequency reuse-1 model, the frequency reuse-3 model, FFR, and SFR techniques. Our major contributions are summarized as follows:

- Propose an original formulation of the centralized joint resource and power allocation problem: instead of considering a single cell OFDMA network,
 140 we formulate our problem for a multi-cell OFDMA network, taking ICI problems into account. The objective is to maximize the mean rate per UE, and ensure a proportional fair rate for all the active UEs.
- Decompose the joint downlink resource and power allocation problem into two independent problems, and solve the centralized power allocation problem using Lagrange duality theory and subgradient projection
 145 method.

- Formulate a novel decentralized super-modular game for resource and power allocation, and propose a best response algorithm to attain the Nash Equilibrium. Then, solve the decentralized power allocation problem using subgradient projection method.
- Validate the convergence of the proposed centralized and decentralized approaches and evaluate their performance in comparison with broadly adopted state-of-the-art approaches.

3. System Model and Problem Formulation

3.1. System Model

We consider the downlink of a multiuser OFDMA system that consists of I adjacent cells and K active UEs. Let $\mathcal{I} = \{1, 2, \dots, I\}$ denote the the set of cells, and $\mathcal{K} = \{1, 2, \dots, K\}$ the total set of active UEs. We also define $K(i)$ as the number of UEs served by cell i . Thus, we have $\sum_{i=1}^I K(i) = K$. The set of available RBs in each cell is denoted by $\mathcal{N} = \{1, 2, \dots, N\}$.

In OFDMA networks, system spectrum is divided into several channels, where each channel consists of a number of consecutive orthogonal OFDM subcarriers [24]. An RB is the smallest scheduling unit. It consists of 12 consecutive subcarriers in the frequency domain, and seven OFDM symbols with normal cyclic prefix in the time domain [25] (or six OFDM symbols with extended cyclic prefix). Frequency resources are allocated to UEs each Transmit Time Interval (TTI), which is equal to 1 ms. When the frequency reuse-1 model is applied along with homogeneous power allocation, each RB is allocated the same downlink transmission power $\frac{P_{max}}{N}$, where P_{max} denotes the maximum downlink transmission power per cell. The signal to interference and noise ratio for a UE k attached to cell i and allocated RB n is given by:

$$\sigma_{k,i,n} = \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}}, \quad (1)$$

where $\pi_{i,n}$ is the downlink transmission power allocated by cell i to RB n , $G_{k,i,n}$ denotes channel gain for UE k attached to cell i and allocated RB n , and N_0 is the thermal noise power. Indexes i and i' refer to useful and interfering signals respectively. In our work, we rely on perfect channel state information to infer the SINR. Authors of [26] provide models to account for imperfect channel state and study the impact on energy efficiency. Notations, symbols, parameters, and variables used within this document are reported in Table 1.

Table 1: Sets, parameters and variables in the article

i	Index of cell
k	Index of UE
n	Index of RB
\mathcal{I}	Set of cells
\mathcal{K}	Total set of UEs
$\mathcal{K}(i)$	Set of UEs associated to cell i
\mathcal{N}	Set of RBs
$\rho_{k,i,n}$	Peak rate of UE k associated with RB n on cell i
$\pi_{i,n}$	Transmit power of cell i on RB n
$G_{k,i,n}$	Channel gain for UE k over RB n on cell i
N_0	Thermal noise density
$\theta_{k,n}$	Percentage of time UE k is associated with RB n
η	Total system achievable mean rate
$\sigma_{k,i,n}$	SINR for UE k over RB n on cell i
P_{max}	Maximum DL transmission power per cell
π_{min}	Minimum DL transmission power per RB
$\mathcal{I}'(i)$	Set of neighboring cells for cell i

3.2. Problem Formulation

3.2.1. Centralized Multi-Cell Optimization Problem

We define $\theta_{k,n}$ as the percentage of time during which UE k is associated with RB n . $\theta_{k,n}, \forall k \in \mathcal{K}, \forall n \in \mathcal{N}$, and $\pi_{i,n}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$, are the optimization variables of the joint resource and power allocation problem. Our objective is to manage resource and power allocation in a manner that maximizes system throughput and guarantees throughput fairness between the different UEs. The standard approach is to have integer scheduling variables, while in our problem formulation, $\theta_{k,n}$ and $\pi_{i,n}$ are continuous variables. In fact, using continuous variables will decrease the computation time and the complexity of the problem

without losing generality. A simple way of implementing the solution is to extend the Round-Robin scheduler in a way to allocate equal time shares to the
180 users in the cell on each RB.

The peak rate of UE k when associated with RB n on cell i is given by:

$$\rho_{k,i,n} = \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right). \quad (2)$$

Then, the mean rate of UE k is given by:

$$\begin{aligned} \sum_{n \in \mathcal{N}} (\theta_{k,n} \cdot \rho_{k,i,n}) = \\ \sum_{n \in \mathcal{N}} \left(\theta_{k,n} \cdot \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right). \end{aligned} \quad (3)$$

Our centralized multi-cell joint resource and power allocation problem seeks rate maximization in a proportional fair manner. We make use of the logarithmic function that is intimately associated with the concept of proportional fairness [27]. Our problem is formulated in the following:

$$\begin{aligned} \underset{\boldsymbol{\theta}, \boldsymbol{\pi}}{\text{maximize}} \quad & \eta = \\ & \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \log \left(\sum_{n \in \mathcal{N}} \theta_{k,n} \cdot \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right) \end{aligned} \quad (4a)$$

$$\text{subject to} \quad \sum_{k \in \mathcal{K}(i)} \theta_{k,n} \leq 1, \quad \forall n \in \mathcal{N}, \quad (4b)$$

$$\sum_{n \in \mathcal{N}} \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \quad (4c)$$

$$\sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \quad \forall i \in \mathcal{I}, \quad (4d)$$

$$\pi_{i,n} \geq \pi_{min}, \quad \forall i \in \mathcal{I}, \forall n \in \mathcal{N}, \quad (4e)$$

$$0 \leq \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}. \quad (4f)$$

The objective function η ensures a proportional fair rate for all UEs in the network. Constraints (4b) ensure that an RB is used at most 100% of the time,

and constraints (4c) ensure that a UE shares its time on the available RBs. Constraints (4d) guarantee that the total downlink transmission power allocated to the available RBs does not exceed the maximum transmission power P_{max} for each cell i , and constraints (4e) represent the minimum power constraint of the transmit power allocated to each RB. In fact, the power allocated to each RB is larger than a predefined value denoted π_{min} , and the transmit power of cell i is lower than P_{max} . In practice, these bounds are related to hardware limitations. $\theta_{k,n}, \forall k \in \mathcal{K}, \forall n \in \mathcal{N}$, and $\pi_{i,n}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$ are the optimization variables of the joint resource and power allocation problem.

In order to reduce the complexity of the joint resource and power allocation problem (4), we prove that this problem is separable into two independent problems: a resource allocation problem and a power allocation problem. In fact, maximizing the objective function of problem (4) is achieved by maximizing the following term:

$$\sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} (\log(\theta_{k,n}) + \log(\rho_{k,i,n})). \quad (5)$$

The proof of this hypothesis is given in Appendix I.

4. Problem Decomposition

We tackle ICIC as an optimization problem, where we intend to maximize the mean rate of UEs in a multiuser OFDMA system. We consider a system of I cells, having $K(i)$ UEs per cell i . According to (5), and due to the absence of binding constraints, the optimization problem (4) is linearly separable into two independent problems: a power allocation problem and a resource allocation problem.

4.1. Centralized Multi-Cell Power Allocation Problem

In the first problem, the optimization variable $\boldsymbol{\pi}$ is considered, and the problem is formulated as follows:

$$\begin{aligned}
& \underset{\boldsymbol{\pi}}{\text{maximize}} && \eta_1 = \\
& && \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log \left(\log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right) \quad (6a) \\
& \text{subject to} && \sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \quad \forall i \in \mathcal{I}, \quad (6b) \\
& && \pi_{i,n} \geq \pi_{min}, \quad \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (6c)
\end{aligned}$$

Problem (6) consists in finding the optimal power allocation. However, it is not a convex optimization problem as formulated in (6). In the following, we introduce a variable change that allows to formulate problem (6) as a convex optimization problem as follows:

$$\begin{aligned}
& \underset{\boldsymbol{\rho}, \boldsymbol{\pi}}{\text{maximize}} && \eta_1 = \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\rho_{k,i,n}) \quad (7a) \\
& \text{subject to} && \rho_{k,i,n} \leq \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right), \\
& && \forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}, \quad (7b) \\
& && \sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \quad \forall i \in \mathcal{I}, \quad (7c) \\
& && \pi_{i,n} \geq \pi_{min}, \quad \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (7d)
\end{aligned}$$

Let us consider the following variable change:

$$\hat{\rho}_{k,i,n} = \log(\exp(\rho_{k,i,n}) - 1), \quad \forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}, \quad (8a)$$

$$\hat{\pi}_{i,n} = \log(\pi_{i,n}), \quad \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (8b)$$

Hence, the original variables are given by:

$$\rho_{k,i,n} = \log(\exp(\hat{\rho}_{k,i,n}) + 1), \quad \forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}, \quad (9a)$$

$$\pi_{i,n} = \exp(\hat{\pi}_{i,n}), \quad \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (9b)$$

To show that the optimization problem (7) is a convex optimization problem, we need to show that the objective function is concave and the inequality constraint functions define a convex set. After applying the variable change on

peak rate constraints (7b), they can be written as:

$$\begin{aligned}
\rho_{k,i,n} &\leq \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right), \\
\forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N} \\
&\Rightarrow \log(\exp(\hat{\rho}_{k,i,n}) + 1) \leq \log \left(1 + \frac{\exp(\hat{\pi}_{i,n}) G_{k,i,n}}{N_0 + \sum_{i' \neq i} \exp(\hat{\pi}_{i',n}) G_{k,i',n}} \right) \\
&\Rightarrow \exp(\hat{\rho}_{k,i,n}) + 1 \leq 1 + \frac{\exp(\hat{\pi}_{i,n}) G_{k,i,n}}{N_0 + \sum_{i' \neq i} \exp(\hat{\pi}_{i',n}) G_{k,i',n}} \\
&\Rightarrow \frac{\exp(\hat{\rho}_{k,i,n}) \cdot (N_0 + \sum_{i' \neq i} \exp(\hat{\pi}_{i',n}) G_{k,i',n})}{\exp(\hat{\pi}_{i,n}) G_{k,i,n}} \leq 1 \\
&\Rightarrow \log \left(\exp(\hat{\rho}_{k,i,n} - \hat{\pi}_{i,n}) \frac{N_0}{G_{k,i,n}} + \sum_{i' \neq i} \exp(\hat{\rho}_{k,i,n} + \hat{\pi}_{i',n} - \hat{\pi}_{i,n}) \frac{G_{k,i',n}}{G_{k,i,n}} \right) \leq 0, \\
&\forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}.
\end{aligned}$$

These constraints are the logarithmic of the sum of exponential functions. Thus, they are convex functions [28]. When we apply the variable change on power constraints (7c), we get:

$$\begin{aligned}
\sum_{n \in \mathcal{N}} \pi_{i,n} &\leq P_{max}, \forall i \in \mathcal{I} \\
&\Rightarrow \log \left(\sum_{n \in \mathcal{N}} \exp(\hat{\pi}_{i,n}) \right) - \log(P_{max}) \leq 0, \forall i \in \mathcal{I}.
\end{aligned}$$

215 Since $\log(\sum \exp)$ is convex [28], the constraints at hand are therefore convex. Using the variable change, the power allocation problem (7) can be written as follows:

$$\begin{aligned}
&\underset{\hat{\rho}, \hat{\pi}}{\text{maximize}} \quad \eta_1 = \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\log(\exp(\hat{\rho}_{k,i,n}) + 1)) \quad (10a)
\end{aligned}$$

subject to

$$\begin{aligned}
&\log \left(\exp(\hat{\rho}_{k,i,n} - \hat{\pi}_{i,n}) \frac{N_0}{G_{k,i,n}} + \sum_{i' \neq i} \exp(\hat{\rho}_{k,i,n} + \hat{\pi}_{i',n} - \hat{\pi}_{i,n}) \frac{G_{k,i',n}}{G_{k,i,n}} \right) \leq 0, \\
&\forall i \in \mathcal{I}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}, \quad (10b)
\end{aligned}$$

$$\log \left(\sum_{n \in \mathcal{N}} \exp(\hat{\pi}_{i,n}) \right) - \log(P_{max}) \leq 0, \forall i \in \mathcal{I}, \quad (10c)$$

$$\hat{\pi}_{i,n} \geq \log(\pi_{min}), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (10d)$$

The objective function of problem (10) is concave in $\hat{\rho}$ and $\hat{\pi}$, and constraints (10b), (10c), and (10d) are convex functions. Thus, the power allocation
 220 problem is a convex optimization problem.

4.2. Centralized Resource Allocation Problem

The optimization variable θ is considered in the second optimization problem that is given in the following:

$$\underset{\theta}{\text{maximize}} \quad \eta_2 = \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\theta_{k,n}) \quad (11a)$$

$$\text{subject to} \quad \sum_{k \in \mathcal{K}(i)} \theta_{k,n} \leq 1, \quad \forall n \in \mathcal{N}, \quad (11b)$$

$$\sum_{n \in \mathcal{N}} \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \quad (11c)$$

$$0 \leq \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}. \quad (11d)$$

As demonstrated for the power allocation problem (6), we prove that prob-
 225 lem (11) is indeed a convex optimization problem in θ . The objective function (11a) of the resource allocation problem (11) is concave in θ , since the log function is concave for $\theta \in]0; 1]$. Moreover, constraints (11b), (11c), and (11d) are linear and separable constraints. Hence, the resource allocation problem (11) is a convex optimization problem, and it is separable into \mathcal{I} subproblems. For
 230 each cell i , the i th optimization problem is written as follows:

$$\underset{\theta}{\text{maximize}} \quad (\eta_2)_i = \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\theta_{k,n}) \quad (12a)$$

$$\text{subject to} \quad \sum_{k \in \mathcal{K}(i)} \theta_{k,n} \leq 1, \quad \forall n \in \mathcal{N}, \quad (12b)$$

$$\sum_{n \in \mathcal{N}} \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \quad (12c)$$

$$0 \leq \theta_{k,n} \leq 1, \quad \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}. \quad (12d)$$

5. Centralized Multi-Cell Resource and Power Allocation

As stated in the previous section and proven in Appendix I, the joint resource and power allocation problem (4) is separable into two independent convex

optimization problems: a power allocation problem, and a resource allocation
 235 problem. In this section, we solve the resource and power allocation problems
 using Lagrange duality theory and subgradient projection method.

5.1. Solving the Centralized Power Allocation Problem

5.1.1. Lagrange-Based Method

Since the power allocation problem (10) is a convex optimization problem, we
 240 can make use of Lagrange duality properties, which also lead to decomposability
 structures [29]. Lagrange duality theory links the original problem, or *primal*
problem, with a dual maximization problem. The Lagrangian of problem (10)
 is given as follows:

$$\begin{aligned}
 L(\hat{\boldsymbol{\rho}}, \hat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu}) = & \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\log(\exp(\hat{\rho}_{k,i,n}) + 1)) \\
 & - \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \lambda_{k,i,n} (\log(\exp(\hat{\rho}_{k,i,n}) - \hat{\pi}_{i,n}) \frac{N_0}{G_{k,i,n}} \\
 & + \sum_{\substack{i' \in \mathcal{N} \\ i' \neq i}} \exp(\hat{\rho}_{k,i,n} + \hat{\pi}_{i',n} - \hat{\pi}_{i,n}) \frac{G_{k,i',n}}{G_{k,i,n}})) \\
 & - \sum_{i \in \mathcal{I}} \nu_i \left(\log \left(\sum_{n \in \mathcal{N}} \exp(\hat{\pi}_{i,n}) \right) - \log(P_{max}) \right).
 \end{aligned} \tag{13}$$

The optimization variables $\hat{\boldsymbol{\rho}}$ and $\hat{\boldsymbol{\pi}}$ are called the primal variables. $\lambda_{k,i,n}$
 245 and ν_i are the *dual variables* associated with the (k, i, n) th inequality con-
 straint (10b) and with the i th inequality constraint (10c), respectively.

After relaxing the coupling constraints (10b) and (10c), the optimization
 problem separates into two levels of optimization: lower level and higher level.
 At the lower level, $L(\hat{\boldsymbol{\rho}}, \hat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu})$ is the objective function to be maximized. $\hat{\rho}_{k,i,n}$
 and $\hat{\pi}_{i,n}$ are the optimization variables to be found. At the higher level, we have
 the master dual problem in charge of updating the dual variables $\boldsymbol{\lambda}$ and $\boldsymbol{\nu}$ by

solving the dual problem:

$$\underset{\boldsymbol{\lambda}, \boldsymbol{\nu}}{\text{minimize}} \quad \max_{\widehat{\boldsymbol{\rho}}, \widehat{\boldsymbol{\pi}}} (L(\widehat{\boldsymbol{\rho}}, \widehat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu})) \quad (14a)$$

$$\text{subject to} \quad \boldsymbol{\lambda} \geq 0, \quad (14b)$$

$$\boldsymbol{\nu} \geq 0. \quad (14c)$$

In order to solve the primal optimization problem (lower level of optimization), we use the subgradient projection method. It starts with some initial feasible values of $\widehat{\rho}_{k,i,n}$ and $\widehat{\pi}_{i,n}$ that satisfy the constraints (10d). Then, the next iteration is generated by taking a step along the subgradient direction of $\widehat{\rho}_{k,i,n}$ and $\widehat{\pi}_{i,n}$. For the primal optimization variables, iterations of the subgradient projection are given by:

$$\begin{aligned} \widehat{\rho}_{k,i,n}(t+1) &= \widehat{\rho}_{k,i,n}(t) + \delta(t) \times \frac{\partial L}{\partial \widehat{\rho}_{k,i,n}}, \\ \forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}, \end{aligned} \quad (15a)$$

$$\widehat{\pi}_{i,n}(t+1) = \widehat{\pi}_{i,n}(t) + \delta(t) \times \frac{\partial L}{\partial \widehat{\pi}_{i,n}}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (15b)$$

The scalar $\delta(t)$ is a step size that guarantees the convergence of the primal optimization problem [29]. The partial derivatives of the objective function $L(\widehat{\boldsymbol{\rho}}, \widehat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu})$ with respect to $\widehat{\rho}_{k,i,n}$ and $\widehat{\pi}_{i,n}$, are given in the following:

$$\begin{aligned} \frac{\partial L}{\partial \widehat{\rho}_{k,i,n}} &= \frac{\exp(\widehat{\rho}_{k,i,n})}{(\exp(\widehat{\rho}_{k,i,n}) + 1) \log(\exp(\widehat{\rho}_{k,i,n}) + 1)} - \lambda_{k,i,n}, \\ \forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}, \end{aligned} \quad (16a)$$

$$\frac{\partial L}{\partial \widehat{\pi}_{i,n}} = \sum_{k \in \mathcal{K}(i)} \lambda_{k,i,n} - \nu_i \frac{\exp(\widehat{\pi}_{i,n})}{\sum_{n \in \mathcal{N}} \exp(\widehat{\pi}_{i,n})}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}. \quad (16b)$$

The dual function $g(\boldsymbol{\lambda}, \boldsymbol{\nu}) = \max_{\widehat{\boldsymbol{\rho}}, \widehat{\boldsymbol{\pi}}} (L(\widehat{\boldsymbol{\rho}}, \widehat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu}))$ is differentiable. Thus, at the higher optimization level, the master dual problem (14) can be solved using

the following gradient method:

$$\begin{aligned}\lambda_{k,i,n}(t+1) &= \lambda_{k,i,n}(t) + \delta(t) \left(\log \left(\exp(\hat{\rho}_{k,i,n}^* - \hat{\pi}_{i,n}^*) \frac{N_0}{G_{k,i,n}} \right. \right. \\ &\quad \left. \left. + \sum_{\substack{i' \in \mathcal{N} \\ i' \neq i}} \exp(\hat{\rho}_{k,i,n}^* + \hat{\pi}_{i',n}^* - \hat{\pi}_{i,n}^*) \frac{G_{k,i',n}}{G_{k,i,n}} \right) \right), \\ \forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N},\end{aligned}\tag{17a}$$

$$\begin{aligned}\nu_i(t+1) &= \nu_i(t) + \delta(t) \left(\log \left(\sum_{n \in \mathcal{N}} \exp(\hat{\pi}_{i,n}^*) \right) - \log(P_{max}) \right), \\ \forall i \in \mathcal{I}, \forall n \in \mathcal{N},\end{aligned}\tag{17b}$$

where t is the iteration index, and $\delta(t)$ is the step size at iteration t . Appropriate choice of the step size [30] leads to convergence of the dual algorithm. $\hat{\pi}_{i,n}^*$ and $\hat{\rho}_{k,i,n}^*$ denote the solution to the primal optimization problem. When $t \rightarrow \infty$ the dual variables $\lambda(t)$ and $\nu(t)$ converge to the dual optimal λ^* and ν^* , respectively. The difference between the optimal primal objective and the optimal dual objective, called *duality gap*, reduces to zero at optimality, since the problem (10) is convex and the KKT conditions are satisfied. We define $\Delta\hat{\rho}$, $\Delta\hat{\pi}$, $\Delta\lambda$, and $\Delta\nu$ as the differences between the optimization variables obtained at the current iteration and their values at the previous iteration. They are given by:

$$\Delta\hat{\rho}(t+1) = \|\hat{\rho}(t+1) - \hat{\rho}(t)\|,\tag{18a}$$

$$\Delta\hat{\pi}(t+1) = \|\hat{\pi}(t+1) - \hat{\pi}(t)\|,\tag{18b}$$

$$\Delta\lambda(t+1) = \|\lambda(t+1) - \lambda(t)\|,\tag{18c}$$

$$\Delta\nu(t+1) = \|\nu(t+1) - \nu(t)\|.\tag{18d}$$

5.1.2. Iterative Power Allocation Algorithm

The procedure for solving the centralized power allocation problem is described in Algorithm 1. Initially, the primal optimization variables $\hat{\rho}_{k,i,n}$ and $\hat{\pi}_{i,n}$ as well as the dual variables $\lambda_{k,i,n}$ and ν_i start with some initial feasible values. t , t_{primal} , and t_{dual} denote the number of rounds required for the centralized

Algorithm 1 Dual algorithm for centralized power allocation

- 1: Parameters: the utility function $L(\hat{\boldsymbol{\rho}}, \hat{\boldsymbol{\pi}}, \boldsymbol{\lambda}, \boldsymbol{\nu})$, P_{max} , and π_{min} .
 - 2: Initialization: set $t = t_{primal} = t_{dual} = 0$, and $\pi_{i,n} \geq \pi_{min}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$,
such as $\sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \forall i \in \mathcal{I}$. Calculate $\hat{\pi}_{i,n}(0)$ and $\hat{\rho}_{k,i,n}(0)$ accordingly,
 $\forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$.
 - 3: Set $\lambda_{k,i,n}(0)$ and $\nu_i(0)$ equal to some non negative value, $\forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$.
 - 4: $(\hat{\boldsymbol{\pi}}^*(t+1), \hat{\boldsymbol{\rho}}^*(t+1)) \leftarrow \text{PRIMALPROBLEM}(\boldsymbol{\nu}^*(t), \boldsymbol{\lambda}^*(t))$
 - 5: $(\boldsymbol{\nu}^*(t+1), \boldsymbol{\lambda}^*(t+1)) \leftarrow \text{DUALPROBLEM}(\hat{\boldsymbol{\pi}}^*(t+1), \hat{\boldsymbol{\rho}}^*(t+1))$
 - 6: **if** $(\Delta \hat{\boldsymbol{\pi}}^*(t+1) > \epsilon)$ **or** $(\Delta \hat{\boldsymbol{\rho}}^*(t+1) > \epsilon)$ **or** $(\Delta \boldsymbol{\nu}^*(t+1) > \epsilon)$ **or** $(\Delta \boldsymbol{\lambda}^*(t+1) > \epsilon)$ **then**
 - 7: $t \leftarrow t + 1$
 - 8: **go to** 4
 - 9: **end if**
-

power allocation problem to converge, the number of iterations for the primal problem, and the number of iterations for the dual problem, respectively. At each round t , we start by updating the primal optimization variables, using the PRIMALPROBLEM function given in Algorithm 2. The solution to the primal optimization problem at the current round t is denoted by $\hat{\pi}_{i,n}^*(t+1)$ and $\hat{\rho}_{k,i,n}^*(t+1)$. The PRIMALPROBLEM function updates $\hat{\pi}_{i,n}(t_{primal}+1)$ and $\hat{\rho}_{k,i,n}(t_{primal}+1)$, and increments t_{primal} until $\Delta \hat{\boldsymbol{\pi}}(t_{primal}+1)$ and $\Delta \hat{\boldsymbol{\rho}}(t_{primal}+1)$ become less than ϵ .

Then, the solution to the dual optimization problem at the current round t , denoted by $\nu_i^*(t+1)$ and $\lambda_{k,i,n}^*(t+1)$ is calculated using the DUALPROBLEM function given in Algorithm 3. ν_i and $\lambda_{k,i,n}$ are updated using the obtained primal solution $\hat{\pi}_{i,n}^*(t+1)$ and $\hat{\rho}_{k,i,n}^*(t+1)$, until $\Delta \boldsymbol{\nu}(t_{dual}+1)$ and $\Delta \boldsymbol{\lambda}(t_{dual}+1)$ become less than ϵ . An additional round of calculations is performed, and t is incremented as long as $\Delta \hat{\boldsymbol{\pi}}^*(t+1)$ or $\Delta \hat{\boldsymbol{\rho}}^*(t+1)$ or $\Delta \boldsymbol{\nu}^*(t+1)$ or $\Delta \boldsymbol{\lambda}^*(t+1)$ is greater than ϵ . Otherwise, the obtained solution at the current round is the optimal solution to the centralized power allocation problem.

5.2. Solving the Resource Allocation Problem

In this subsection, we search for the optimal solution to the resource allocation problem (12). For each cell i , the problem (12) is a convex optimization problem, as proven previously.

Algorithm 2 Primal problem function

```

1: function PRIMALPROBLEM( $\boldsymbol{\nu}^*(t), \boldsymbol{\lambda}^*(t)$ )
2:   for  $i = 1$  to  $|\mathcal{I}|$  do
3:     for  $n = 1$  to  $|\mathcal{N}|$  do
4:        $\hat{\pi}_{i,n}(t_{\text{primal}} + 1) \leftarrow \max \left( \log(\pi_{\min}); \hat{\pi}_{i,n}(t_{\text{primal}}) + \delta(t) \times \frac{\partial L}{\partial \hat{\pi}_{i,n}} \right)$ 
5:       for  $k = 1$  to  $|\mathcal{K}(i)|$  do
6:          $\hat{\rho}_{k,i,n}(t_{\text{primal}} + 1) \leftarrow \hat{\rho}_{k,i,n}(t_{\text{primal}}) + \delta(t) \times \frac{\partial L}{\partial \hat{\rho}_{k,i,n}}$ 
7:       end for
8:     end for
9:   end for
10:  if  $(\Delta \hat{\pi}(t_{\text{primal}} + 1) > \epsilon)$  or  $(\Delta \hat{\rho}(t_{\text{primal}} + 1) > \epsilon)$  then
11:     $t_{\text{primal}} \leftarrow t_{\text{primal}} + 1$ 
12:    go to 2
13:  end if
14:  return  $\hat{\pi}(t_{\text{primal}} + 1), \hat{\rho}(t_{\text{primal}} + 1)$ 
15: end function

```

Theorem 5.1. *For each cell i , the optimal solution to the resource allocation problem (12) is given by:*

$$\theta_{k,n} = \frac{1}{\max(|\mathcal{K}(i)|, |\mathcal{N}|)}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}. \quad (19)$$

The proof of Theorem 5.1 is given in Appendix II. When the number of active UEs is less than the number of available resources, $\theta_{k,n} = \frac{1}{|\mathcal{N}|}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}$. Thus, the available resources are not fully used over time, and each UE is permanently served. Otherwise, when $|\mathcal{K}(i)| > |\mathcal{N}|$, the optimal solution is: $\theta_{k,n} = \frac{1}{|\mathcal{K}(i)|}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}$. In this case, each RB is fully used over time, while UEs are not permanently served over time.

6. Decentralized Resource and Power Allocation

6.1. Problem Formulation and Decomposition

We have shown that the power allocation problem can be solved optimally in a centralized fashion. The centralized approach is the reference approach for performance comparison, since it finds the optimal resource and power allocation for all the active users. Nevertheless, the computational complexity of the

Algorithm 3 Dual problem function

```

1: function DUALPROBLEM( $\hat{\pi}^*(t+1), \hat{\rho}^*(t+1)$ )
2:   for  $i = 1$  to  $|\mathcal{I}|$  do
3:      $\nu_i(t_{dual} + 1) \leftarrow \max(0; \nu_i(t_{dual}) + \delta(t) \times (\log(\sum_{n \in \mathcal{N}} \exp(\hat{\pi}_{i,n}^*(t+1))) - \log(P_{max})))$ 
4:     for  $n = 1$  to  $|\mathcal{N}|$  do
5:       for  $k = 1$  to  $|\mathcal{K}(i)|$  do
6:          $\lambda_{k,i,n}(t_{dual} + 1) \leftarrow \max(0; \lambda_{k,i,n}(t_{dual}) + \delta(t) \times (\log(\exp(\hat{\rho}_{k,i,n}^*(t+1)) - \hat{\pi}_{i,n}^*(t+1)) \frac{N_0}{G_{k,i,n}} + \sum_{\substack{i' \in \mathcal{N} \\ i' \neq i}} \exp(\hat{\rho}_{k,i',n}^*(t+1) + \hat{\pi}_{i',n}^*(t+1) - \hat{\pi}_{i,n}^*(t+1)) \frac{G_{k,i',n}}{G_{k,i,n}})))$ 
7:       end for
8:     end for
9:   end for
10:  if  $(\Delta \nu(t_{dual} + 1) > \epsilon)$  or  $(\Delta \lambda(t_{dual} + 1) > \epsilon)$  then
11:     $t_{dual} \leftarrow t_{dual} + 1$ 
12:    go to 2
13:  end if
14:  return  $\nu(t_{dual} + 1), \lambda(t_{dual} + 1)$ 
15: end function

```

centralized approach motivates the introduction of low-complexity decentralized approaches. In this section we investigate the decentralized resource and power allocation approach. Base stations of the LTE/LTE-A networks are autonomous entities, and each cell performs resource and power allocation independently of the other cells. Each cell i maximizes its own utility function, which is given by:

$$\begin{aligned}
& \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log \left(\theta_{k,n} \cdot \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right) = \\
& \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\theta_{k,n}) + \\
& \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log \left(\log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right).
\end{aligned} \tag{20}$$

The decentralized joint resource and power allocation problem is separable into two independent problems: a resource allocation problem and a power allocation problem. The resource allocation problem is solved in a distributed manner as proven in the previous section. We propose a decentralized power allocation

285 approach based on game theory, where the cells are the decision makers or
 players of the game. We define a multi-player game \mathcal{G} between the $|\mathcal{I}|$ cells.
 The cells are assumed to make their decisions without knowing the decisions of
 each other.

The formulation of this non-cooperative game $G = \langle \mathcal{I}, S, U \rangle$ can be described
 290 as follows:

- A finite set of cells $\mathcal{I} = (1, \dots, |\mathcal{I}|)$.
- For each cell i , the space of pure strategies is S_i given by what follows:

$$S_i = \{\pi_i \in R^{|\mathcal{N}|} \text{ such as } \pi_{i,n} \geq \pi_{min}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}, \text{ and } \sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \forall i \in \mathcal{I}\}.$$

295 An action of a cell i is the amount of power $\pi_{i,n}$ allocated to the RB n , and
 the strategy chosen by cell i is then $\pi_i = (\pi_{i,1}, \dots, \pi_{i,N})$. A *strategy profile*
 $\pi = (\pi_1, \dots, \pi_{|\mathcal{I}|})$ specifies the strategies of all players and $S = S_1 \times \dots \times S_{|\mathcal{I}|}$
 is the set of all strategies.

- A set of utility functions $U = (U_1(\pi), U_2(\pi), \dots, U_I(\pi))$ that quantify play-
 ers' utility for a given strategy profile π , where a given utility U_i for cell i
 is such as:

$$U_i = \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log \left(\log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}} \right) \right). \quad (21)$$

For every i , U_i is concave w.r.t. π_i and continuous w.r.t. $\pi_l, l \neq i$. Hence, a
 300 Nash Equilibrium (NE) exists [31]. We note that the objective function η_1 of
 the centralized power allocation problem (10) is equivalent to the sum of the
 utility functions U_i of the I cells.

6.2. Super-Modular Games

Super-modular games exhibit strategic complementarity *i.e.*, the marginal
 305 utility for a player in playing a higher strategy increases when the opponents also
 play higher strategy [32]. These games encompass many applied models, and

they are characterized by the existence of pure strategy NE. Before presenting the properties of a super-modular game, we list first the following definition:

Definition 6.1. *If U_i is twice differentiable, it is said to be super-modular if:*

$$\frac{\partial U_i}{\partial \pi_l \partial \pi_i} \geq 0, \forall l \in \mathcal{I} - \{i\}, \forall \pi_i \in S_i.$$

According to [33], a game is super-modular if $\forall i \in \mathcal{I}$:

- 310 1. The strategy space S_i is a compact sublattice of \mathbb{R}^N .
2. The utility function U_i is super-modular.

In [33, 34], proof is given for the following result in a super-modular game:

- If we start with a feasible policy, then the sequence of best responses monotonically converges to an equilibrium: it monotonically increases in all components in the case of maximizing in a super-modular game.

Proposition 6.2. *The game \mathcal{G} is a super-modular game.*

The proof of this proposition is given in Appendix III.

To attain the NE of the game, we implement a best response algorithm where in each round t , cell i strives to find, in parallel for every RB $n \in \mathcal{N}$, the following optimal power level as a response to $\pi_{-i}(t-1)$: $\pi_i^*(t) = \arg \max_{\pi_i} U_i(\pi_i, \pi_{-i}), s.t. \pi_i^* \in S_i$.

The resulting optimization problem for each cell i is as follows:

$$\underset{\pi_i}{\text{maximize}} \quad U_i \tag{22a}$$

$$\text{subject to:} \quad \sum_{n \in \mathcal{N}} \pi_{i,n} \leq P_{max}, \tag{22b}$$

$$\pi_{i,n} \geq \pi_{min}, \forall n \in \mathcal{N}. \tag{22c}$$

6.3. Solving the Decentralized Power Allocation Problem

We use the subgradient projection method to solve the decentralized power allocation problem (22). It is an iterative method that starts with some initial

feasible vector π_i that satisfies constraints (22b) and (22c), and generates the next iteration by taking a step along the subgradient direction of U_i at π_i . For each cell i , iterations of the subgradient projection are given by:

$$\pi_{i,n}(t+1) = \pi_{i,n}(t) + \delta(t) \times \frac{\partial U_i}{\partial \pi_{i,n}}, \forall n \in \mathcal{N}, \quad (23)$$

where the partial derivative of the objective function U_i with respect to $\pi_{i,n}$ is given by:

$$\frac{\partial U_i}{\partial \pi_{i,n}} = \sum_{k \in \mathcal{K}(i)} \frac{G_{k,i,n}}{(N_0 + F_{i,n}) \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + F_{i,n}}\right) \log \left(1 + \frac{\pi_{i,n} G_{k,i,n}}{N_0 + F_{i,n}}\right)}, \quad (24a)$$

$$F_{i,n} = \sum_{\substack{i' \in \mathcal{I} \\ i' \neq i}} \pi_{i',n} G_{k,i',n}, \forall n \in \mathcal{N}. \quad (24b)$$

The scalar $\delta(t) > 0$ is a small step size (*e.g.*, $\delta(t) = 0.001$) chosen appropriately [30] to guarantee the convergence of the decentralized power allocation problem (22). Before updating the variables $\pi_{i,n}(t+1)$, we make sure that $\pi_{i,n}(t+1) \geq \pi_{min}$ in order to satisfy the constraints (22c). Moreover, if constraints (22b) are not satisfied, we perform a projection on the feasible set P_{max} , which is straightforward for a simplex [35]. Then, we calculate the power difference $\Delta\pi_i$, which is the difference between the power allocation vectors of the current and the previous iterations. It is given by:

$$\Delta\pi_i(t+1) = \|\pi_i(t+1) - \pi_i(t)\|. \quad (25)$$

As described in Algorithm 4, each cell i calculates $\pi_{i,n}(t_i+1), \forall n \in \mathcal{N}$, where t_i is the iteration number for cell i . The obtained power values are updated in accordance with the constraints (22b) and (22c). This procedure is repeated and the number of iterations t_i is incremented until $\Delta\pi_i(t_i+1)$ becomes less than ϵ . The number of rounds required for all the cells to converge is denoted by t . An additional round of power calculation is performed for all the cells and t is incremented as long as $\Delta\pi^*(t+1) > \epsilon$, where $\pi^*(t)$ is the

Algorithm 4 Decentralized power allocation

```
1: Parameters: the utility function  $U_i, \forall i \in \mathcal{I}$ , the maximum power per
   cell  $P_{max}$ , and the minimum power per RB  $\pi_{min}$ .
2: Initialization: set  $t = 0, t_i = 0, \forall i \in \mathcal{I}$ , and  $\pi_{i,n}(0)$  to some positive value  $\geq$ 
    $\pi_{min}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$ , such as  $\sum_{n \in \mathcal{N}} \pi_{i,n}(0) \leq P_{max}, \forall i \in \mathcal{I}$ .
3: for  $i = 1$  to  $|\mathcal{I}|$  do
4:   for  $n = 1$  to  $|\mathcal{N}|$  do
5:      $\pi_{i,n}(t_i + 1) \leftarrow \max \left( \pi_{min}; \pi_{i,n}(t_i) + \delta(t_i) \times \frac{\partial U_i}{\partial \pi_{i,n}} \right)$ 
6:   end for
7:   if  $\sum_{n=1}^{|\mathcal{N}|} \pi_{i,n}(t_i + 1) > P_{max}$  then
8:     Perform projection on simplex  $P_{max}$ 
9:   end if
10:  if  $\Delta \pi_i(t_i + 1) > \epsilon$  then
11:     $t_i \leftarrow t_i + 1$ 
12:    go to 4
13:  end if
14:   $\pi_{i,n}^*(t + 1) \leftarrow \pi_{i,n}(t_i + 1), \forall n \in \mathcal{N}$ 
15: end for
16: if  $\Delta \pi^*(t + 1) > \epsilon$  then
17:   $t \leftarrow t + 1$ 
18:  go to 3
19: end if
```

330 power allocation vector obtained at the end of round t .

7. Performance Evaluation

In this section, we evaluate the convergence and the performance of the proposed centralized joint resource and power allocation problem, and the decentralized power allocation approach.

335 7.1. Centralized Resource and Power Allocation

To verify the convergence of the centralized solution, we consider a multi-user OFDMA network, such as LTE/LTE-A networks, that consists of seven adjacent hexagonal cells, with one UE served by each cell. UE positions and radio conditions are randomly generated, and the initial power allocation for
340 each RB equals π_{min} (0.1 W). System bandwidth equals 5 MHz. Thus, 25 RBs are available in each cell. The maximum transmission power per cell P_{max} is

set to 43 dBm or 20 W. At the first iteration, the dual variables $\lambda_{k,i,n}(0), \forall k \in \mathcal{K}(i), \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$, and $\nu_i(0), \forall i \in \mathcal{I}$, are assigned initial positive values. The evolution of $\hat{\pi}_{i,1}$ along with the number of iterations is shown in Fig. 1(a), where $\hat{\pi}_{i,1}$ is the logarithm of the transmission power allocated by the cell i to the RB 1. In addition, the number of primal iterations and the number of dual iterations per round are shown in Fig. 1(b).

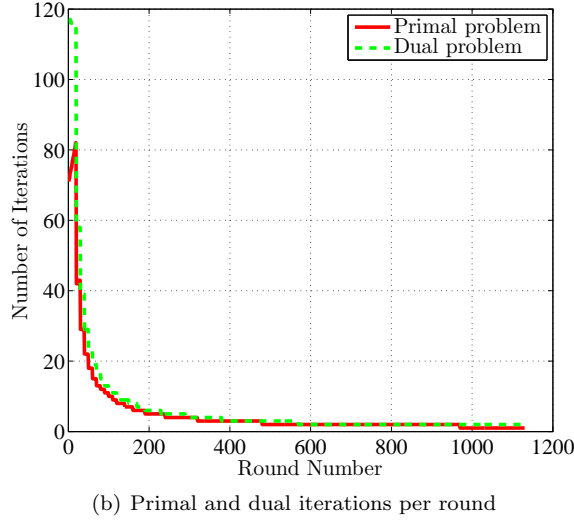
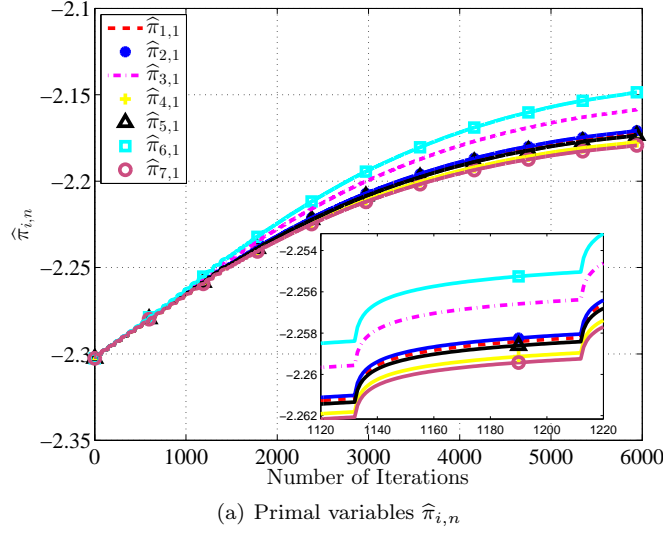


Figure 1: Primal variables and number of iterations

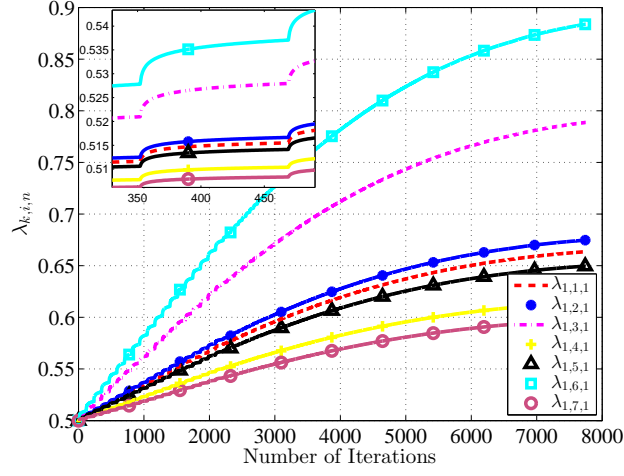
We notice that for the centralized power allocation approach, the primal problem requires approximately 6000 iterations to converge. As shown in Fig. 1(b), 1100 rounds are required to reach the optimal values of the primal and the dual variables. The zoomed box within Fig. 1(a) shows the evolution of $\hat{\pi}_{i,n}$ versus the number of primal iterations for a given round t . The values of $\hat{\pi}_{i,n}$ are calculated using the dual variables obtained at the round $(t-1)$. We also notice that the number of primal iterations and the number of dual iterations decreases with the number of rounds. When t increases, the impact of Lagrange prices $\lambda_{k,i,n}(t)$ and $\nu_i(t)$ on the primal variables calculation is reduced, and the number of primal iterations required for the primal problem to converge becomes lower. The same behavior is noticed for the number of dual iterations when the number of rounds increases.

For the same simulated scenario, we also show the dual variables $\lambda_{k,i,n}$ and ν_i versus the number of dual iterations in Fig. 2(a) and Fig. 2(b), respectively. We notice that approximately 8000 iterations are required for the dual problem to converge. At a given round t , the Lagrange prices $\lambda_{k,i,n}$ and ν_i are updated using the most recent values of the primal variables. The zoomed boxes within Fig. 2(a) and Fig. 2(b) show the evolution of $\lambda_{k,i,n}$ and ν_i versus the number of iterations, respectively. These values are updated until $\Delta\lambda_{k,i,n}$ and $\Delta\nu_i$ become less than ϵ . Convergence of the centralized power allocation problem occurs when two conditions are satisfied: first, the difference between the updated primal variables at round t and their values at round $(t-1)$ is less than ϵ . Second, the difference between the updated primal variables at round t and their values at round $(t-1)$ is less than ϵ .

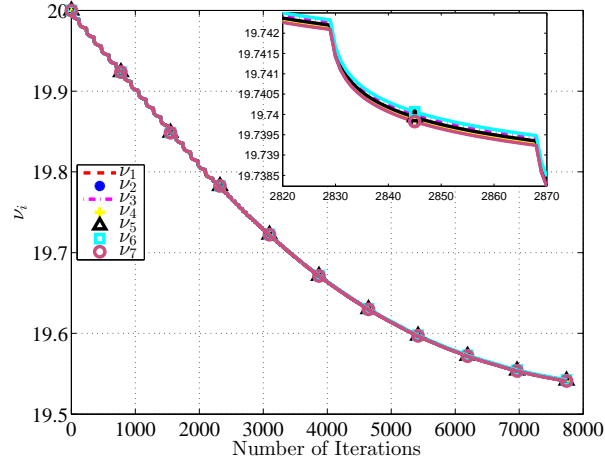
7.2. Decentralized Power Allocation

The same scenario in 7.1 is also simulated in this paragraph to evaluate the performance and convergence of the decentralized power allocation approach. The evolution of the downlink transmission power allocated by all the cells to a given RB is shown in Fig. 3(a).

The initial value of the downlink transmission power allocated to each RB



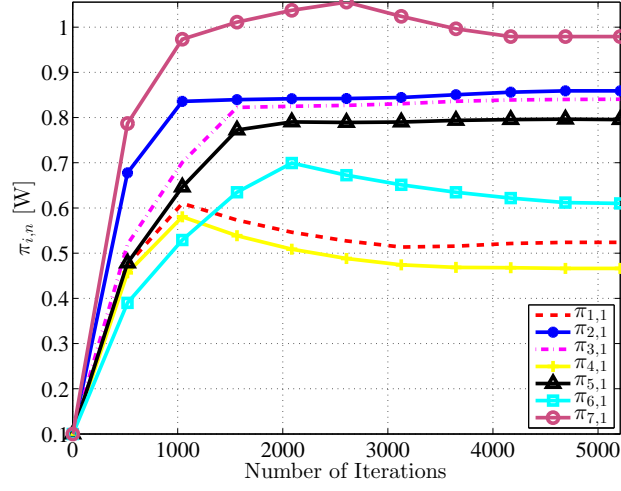
(a) Lagrange prices $\lambda_{k,i,n}$



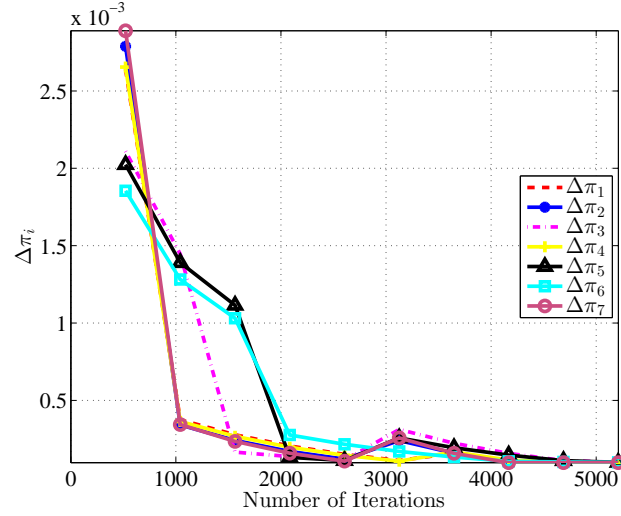
(b) Lagrange prices ν_i

Figure 2: Convergence of the dual variables

equals π_{min} (0.1 W). This allocation satisfies the constraints of the minimal
 downlink transmission power per RB and that of the maximum transmission
 power per cell. Each cell i seeks maximizing its own utility function U_i by ad-
 justing the transmission power allocated to the available RBs. It also estimates
 the interference due to the usage of the same RBs by the neighboring cells. As
 shown in Fig. 3(a), each cell starts increasing the downlink transmission power



(a) $\pi_{i,n}$ versus the number of iterations



(b) $\Delta\pi_i$ versus the number of iterations

Figure 3: Convergence of the transmission power for the decentralized approach

allocated to its RBs, and then the transmission power converges after a given
 385 number of iterations. At convergence, the partial derivative of the objective
 function U_i with respect to $\pi_{i,n}$ becomes negligible. The difference between the
 updated power allocation vector $(\pi_{i,1}, \pi_{i,2}, \dots, \pi_{i,N})$ at iteration $(t + 1)$ and the
 power vector at iteration t becomes less than ϵ .

We also show the evolution of the power vector difference $\Delta\pi_i, \forall i \in \mathcal{I}$, defined
 390 in (25) along with the number of iterations in Fig. 3(b). The obtained curves
 show that $\Delta\pi_i, \forall i \in \mathcal{I}$, decreases when the number of iterations increases. The
 impact of the subgradient projection iterations on the downlink transmission
 power $\pi_{i,n}, \forall i \in \mathcal{I}, \forall n \in \mathcal{N}$, becomes smaller as more iterations are performed.
 Power convergence is achieved when $\Delta\pi_i, \forall i \in \mathcal{I}$, becomes less than ϵ . In fact,
 395 the utility function of each cell i is maximized, and the amount by which the
 downlink transmission power $\pi_{i,n}$ is modified becomes negligible.

7.3. Comparison with State-of-the-Art Resource Allocation Approaches

We also compare the performance of our proposed centralized and decen-
 tralized resource and power allocation approaches with that of state-of-the-art
 400 resource and power allocation approaches [36] such as the frequency reuse-1
 model, the frequency reuse-3 model, FFR, SFR, and a single cell resource and
 power allocation approach [12]. Note that our centralized approach searches for
 the optimal resource and power allocation. It is considered as a reference ap-
 proach when comparing the performance of heuristic algorithms and distributed
 405 approaches.

The frequency reuse-1 model allows the usage of the same frequency spec-
 trum simultaneously in all the network cells. Moreover, homogeneous power
 allocation is performed. In the frequency reuse-3 model, one third of the avail-
 able spectrum is used in each cell in a cluster of three adjacent cells. Interference
 410 problems are eliminated, but the spectral efficiency is reduced. FFR and SFR
 techniques divide each cell into a cell-center and a cell-edge zones, and set restric-
 tions on resource usage and power allocation in each zone. For all the compared
 techniques, resource allocation is performed according to Theorem 5.1.

7.3.1. Spectral Efficiency

415 We investigate the impact of the compared techniques on the spectral effi-
 ciency. Simulation results, including the 95% confidence interval, are shown in
 Fig. 4(a).

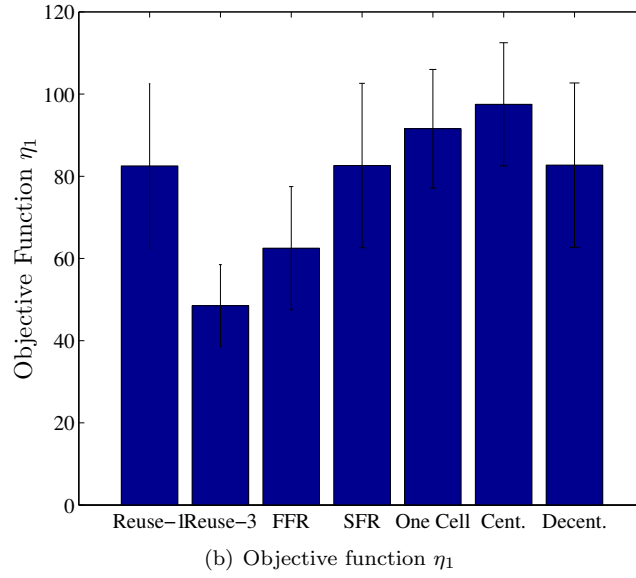
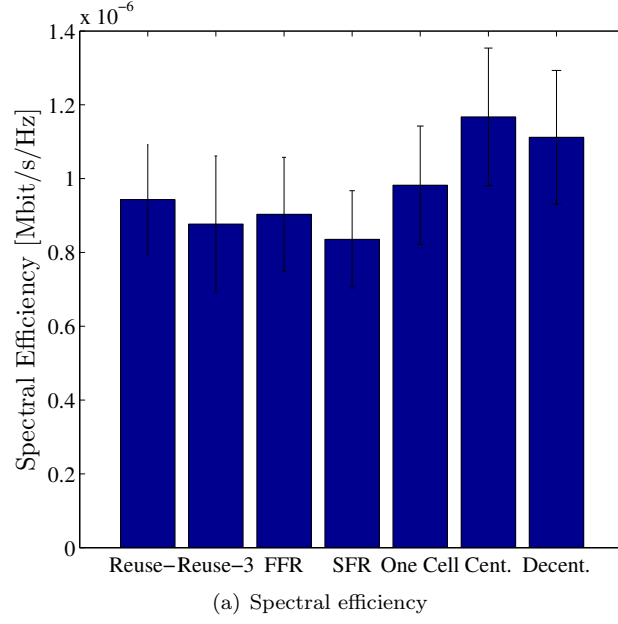


Figure 4: Comparison with state-of-the-art approaches

Our proposed centralized resource allocation approach offers the highest spectral efficiency, since the optimal resource and power allocation is guaran-

420 teed. The spectral efficiency of our decentralized approach is slightly lower than
 that of the centralized approach, due to the lack of information about resource
 usage in the neighboring cells. Nevertheless, the spectral efficiency for both
 the centralized and the decentralized approaches is greater than that of FFR,
 SFR, and the single-cell resource and power allocation approach [12] displayed
 425 as “One Cell” in Fig. 4. In fact, the static resource allocation between cell zones,
 and the quantified transmission power levels do not allow performing flexible
 resource allocation in a manner that satisfies UE needs in each cell. Concerning
 the single-cell approach, it does not take inter-cell interference problems into
 account; however, its spectral efficiency is slightly higher than that of reuse-1
 430 model because it searches for the optimal resource and power allocation locally
 in each cell.

7.3.2. Objective Function

We also compare the objective function η_1 given in (6a) for the different
 resource and power allocation techniques. Simulation results are shown in
 435 Fig. 4(b).

We notice that our centralized approach shows the highest objective function
 η_1 . In fact, it finds the optimal power allocation for the entire system, while
 taking into account constraints related to resource usage and to the maximum
 downlink transmission power per cell. It outperforms the decentralized approach
 440 where each cell strives to maximize its own utility function independently of the
 other cells, and the single-cell approach where inter-cell interference problems
 are not taken into account when solving the resource and power allocation prob-
 lem.

7.4. Centralized Versus Decentralized Complexity Comparison

We evaluate the computational complexity of our centralized and decentral-
 ized resource and power allocation approaches. For the centralized and decen-
 tralized approaches, resource allocation is performed according to Theorem 5.1,
 and it is equivalent to one operation. The complexity of each approach equals

the number of required operations multiplied by the complexity of a single operation, denoted by T_{op} . The complexity of the centralized approach is given by:

$$O[(\text{nb}_{\text{primal}} \times |\mathcal{N}| \times (1 + k) + \text{nb}_{\text{dual}} \times (1 + k \times |\mathcal{N}|)) \times |\mathcal{I}| \times T_{op}]. \quad (26)$$

Similarly, the decentralized approach complexity is given as follows:

$$O(\text{nb}_{\text{iterations}} \times |\mathcal{I}| \times |\mathcal{N}| \times T_{op}), \quad (27)$$

where $\text{nb}_{\text{primal}}$ is the number of primal iterations and nb_{dual} is the number of dual iterations required for convergence of the centralized approach. k is the number of UEs per cell, and $\text{nb}_{\text{iterations}}$ is the number of iterations required for convergence of the decentralized approach.

We notice that the decentralized approach complexity is independent of the number of UEs per cell, contrarily to the centralized approach. The complexity of both techniques depends of the number of cells in the system and the number of RBs available in each cell. Moreover, the computational complexity of these approaches are evaluated under the same simulation scenario as in 7.1. The median number of operations required for the centralized and decentralized approaches are given in Table 2.

Table 2: Median number of operations per approach

Approach	Number of operations
Centralized	$3.02 \cdot 10^8$
Decentralized	$8.84 \cdot 10^5$

According to the results reported in Table 2, the number of operations required for the centralized resource and power allocation approach largely exceeds that of the decentralized approach. In fact, the centralized approach maximizes the objective function for the entire network, contrarily to the decentralized approach where each cell maximizes its objective function independently of the other cells. Therefore, the centralized approach guarantees the optimal solution at the expense of a high computational complexity.

8. Conclusion

Resource and power allocation problem is a challenging problem for nowa-
465 days and future wireless networks. Several state-of-the-art techniques consider
the joint resource and power allocation problem, and formulate it as nonlinear
optimization problems. The objective consists in maximizing system through-
put, spectral efficiency, or energy efficiency under constraints related to the
470 minimum throughput per UE, QoS parameters, and the maximum transmis-
sion power. However, these techniques fall short from considering the impact of
inter-cell interference. Indeed, each cell solves its own resource and power allo-
cation problem without taking into account resource usage and power allocation
in the neighboring cells.

In this article, we formulated the joint resource and power allocation prob-
475 lem for multiuser OFDMA networks as a centralized optimization problem,
where the objective consists in maximizing system throughput while guarantee-
ing throughput fairness between UEs. The joint problem is then decomposed
into two independent problems: a resource allocation problem and a power al-
location problem. Contrarily to the majority of the state-of-the-art approaches,
480 ICI is not neglected, and the impact of the simultaneous transmissions in the
neighboring cells is taken into account when managing the resource and power
allocation. Moreover, we introduced a decentralized power allocation approach
based on game theory. The players are the cells, and each cell aims at max-
imizing its own utility function regardless of the decisions made by the other
485 cells. Simulation results prove the convergence of our proposed approaches, and
show the positive impact of our centralized and decentralized resource alloca-
tion approaches on system performance. In a future work, we will consider
the formulation of a multi-cell and multi-objective resource and power alloca-
tion problem, where we address the compromise between spectral efficiency and
490 energy efficiency.

Appendix I

Upper Bound of the Objective Functions Difference

Given Jensen's inequality and the concavity of the logarithmic function, we have:

$$\log \left(\frac{\sum_{n \in \mathcal{N}} \theta_{k,n} \cdot \rho_{k,i,n}}{|\mathcal{N}|} \right) \geq \frac{\sum_{n \in \mathcal{N}} \log (\theta_{k,n} \cdot \rho_{k,i,n})}{|\mathcal{N}|} \quad (28a)$$

$$\Rightarrow \log \left(\sum_{n \in \mathcal{N}} \theta_{k,n} \cdot \rho_{k,i,n} \right) \geq \frac{\sum_{n \in \mathcal{N}} \log (\theta_{k,n} \cdot \rho_{k,i,n})}{|\mathcal{N}|} + \log (|\mathcal{N}|), \quad (28b)$$

the objective function η can be written as:

$$\begin{aligned} \eta &= \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \log \left(\sum_{n \in \mathcal{N}} \theta_{k,n} \cdot \rho_{k,i,n} \right) \\ &\geq \frac{1}{|\mathcal{N}|} \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log (\theta_{k,n} \cdot \rho_{k,i,n}) + |\mathcal{K}| \cdot \log (|\mathcal{N}|). \end{aligned} \quad (29)$$

Since $\frac{1}{|\mathcal{N}|}$ and $|\mathcal{K}| \cdot \log (|\mathcal{N}|)$ are constant terms, maximizing the objective function of problem (4) is achieved by maximizing the following term:

$$\begin{aligned} \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log (\theta_{k,n} \cdot \rho_{k,i,n}) &= \\ \sum_{i \in \mathcal{I}} \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} (\log (\theta_{k,n}) + \log (\rho_{k,i,n})) &. \end{aligned} \quad (30)$$

In order to decompose the joint problem into two independent problems, we evaluate the gap between the original objective function η and the function given in (30). It is evident that there exists a gap between our objective function η and the function (30) that we will maximize in the following sections. However, maximizing these two functions is equivalent as long as the gap between them is bounded. For this reason, we demonstrate the existence of a finite upper bound on the following difference:

$$0 \leq \log \left(\sum_{n \in \mathcal{N}} \theta_{k,n} \cdot \rho_{k,i,n} \right) - \sum_{n \in \mathcal{N}} \log (\theta_{k,n} \cdot \rho_{k,i,n}) \leq B. \quad (31)$$

Let $\phi_n = \theta_{k,n} \cdot \rho_{k,i,n} > 0$, and suppose that $a \leq \phi_n \leq b$. Thus, there exists $0 \leq \gamma_n \leq 1$ such that $\phi_n = \gamma_n a + (1 - \gamma_n)b$.

$$\begin{aligned}
& \log \left(\sum_{n \in \mathcal{N}} \phi_n \right) - \sum_{n \in \mathcal{N}} \log (\phi_n) \\
&= \log \left(\sum_{n \in \mathcal{N}} (\gamma_n a + (1 - \gamma_n) b) \right) - \sum_{n \in \mathcal{N}} \log (\gamma_n a + (1 - \gamma_n) b) \\
&\leq \log \left(\sum_{n \in \mathcal{N}} (\gamma_n a + (1 - \gamma_n) b) \right) - \sum_{n \in \mathcal{N}} (\gamma_n \log (a) + (1 - \gamma_n) \log (b)) \\
&= \log \left(a \left(\sum_{n \in \mathcal{N}} \gamma_n \right) + b \left(|\mathcal{N}| - \sum_{n \in \mathcal{N}} (\gamma_n) \right) \right) - \log (a) \left(\sum_{n \in \mathcal{N}} \gamma_n \right) \\
&\quad - \log (b) \left(|\mathcal{N}| - \sum_{n \in \mathcal{N}} (\gamma_n) \right).
\end{aligned}$$

Taking $p = \sum_{n \in \mathcal{N}} \gamma_n$ and $q = |\mathcal{N}| - \sum_{n \in \mathcal{N}} \gamma_n$, we have $0 \leq p \leq |\mathcal{N}|$, $0 \leq q \leq |\mathcal{N}|$, and $p + q = |\mathcal{N}|$. Consequently, we have:

$$\begin{aligned}
& \log \left(\sum_{n \in \mathcal{N}} \phi_n \right) - \sum_{n \in \mathcal{N}} \log (\phi_n) \\
&\leq \log (ap + bq) - p \log (a) - q \log (b) \\
&\leq \max_p (\log (ap + b(|\mathcal{N}| - p)) - p \log (a) - (|\mathcal{N}| - p) \log (b)) = B.
\end{aligned}$$

For fixed a and b , let us denote:

$$g(p) = \log (ap + b(|\mathcal{N}| - p)) - p \log (a) - (|\mathcal{N}| - p) \log (b).$$

g is defined, differentiable, and concave on $[0, |\mathcal{N}|]$. Its first derivative is given by:

$$g'(p) = \frac{a - b}{ap + b(|\mathcal{N}| - p)} - \log (a) + \log (b).$$

$g(p)$ attains its maximal value B for a p_0 that satisfies: $g'(p_0) = 0$. Thus, we have $B = g(p_0)$, where p_0 is given by:

$$p_0 = \frac{1}{\log(a) - \log(b)} + \frac{b \cdot |\mathcal{N}|}{b - a}.$$

495 Therefore, the considered difference is bounded by B , which is a finite bound.

Appendix II

Proof of Theorem 5.1

We consider the objective function (12a), which can be written as follows:

$$\begin{aligned} (\eta_2)_i &= \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \log(\theta_{k,n}) \\ &= \log \left(\prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n} \right). \end{aligned} \quad (32)$$

Since the logarithmic function is monotonically increasing, the maximization of $(\eta_2)_i$ becomes equivalent to the maximization of the term $\prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n}$. We
500 consider the following cases:

1. Let us assume that:

$$\sum_{k \in \mathcal{K}(i)} \theta_{k,n} < \sum_{n \in \mathcal{N}} \theta_{k,n}, \quad \forall k \in \mathcal{K}(i), \quad \forall n \in \mathcal{N}. \quad (33)$$

We suppose that $\theta_{k,n}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}$ is an optimal solution to the resource allocation problem (12) *i.e.*, this solution maximizes the objective function (12a). For this solution, we assume that:

$$\exists k \in \mathcal{K}(i) / \sum_{n \in \mathcal{N}} \theta_{k,n} < 1. \quad (34)$$

We define $\epsilon > 0$ as follows:

$$\epsilon = 1 - \sum_{n \in \mathcal{N}} \theta_{k,n},$$

and we demonstrate that this solution is not an optimal solution to problem (12) using the proof by contradiction. In fact, we define a set of $\theta'_{k,n}$ variables as given in the following:

$$\theta'_{k,n} = \begin{cases} \theta_{k,n}, & \forall n \in \mathcal{N}, n \neq n_1, \forall k \in \mathcal{K}(i) \\ \theta_{k,n} + \epsilon, & \text{if } n = n_1, \forall k \in \mathcal{K}(i). \end{cases}$$

Therefore, we have:

$$\prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta'_{k,n} = \prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n} + \epsilon \cdot \prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n} > \prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n},$$

and the assumption made in (34) is false, since it does not maximize the objective function (12a). Consequently, we have:

$$\begin{aligned} \sum_{n \in \mathcal{N}} \theta_{k,n} &= 1, \forall k \in \mathcal{K}(i) \\ \Rightarrow \sum_{k \in \mathcal{K}(i)} \sum_{n \in \mathcal{N}} \theta_{k,n} &= |\mathcal{K}(i)|. \end{aligned}$$

Since the sum of all the $\theta_{k,n}$ variables is constant, the term $\prod_{\substack{k \in \mathcal{K}(i) \\ n \in \mathcal{N}}} \theta_{k,n}$ reaches its maximum when all the variables $\theta_{k,n}$ are equal *i.e.*,

$$\theta_{k,n} = \frac{|\mathcal{K}(i)|}{|\mathcal{K}(i)| \cdot |\mathcal{N}|} = \frac{1}{|\mathcal{N}|}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N},$$

which is an optimal solution to the resource allocation problem (12). Ac-

cording to (33):

$$\begin{aligned}
& \sum_{k \in \mathcal{K}(i)} \theta_{k,n} < \sum_{n \in \mathcal{N}} \theta_{k,n}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N} \\
\Rightarrow & \frac{|\mathcal{K}(i)|}{|\mathcal{N}|} < 1 \\
\Rightarrow & |\mathcal{K}(i)| < |\mathcal{N}|.
\end{aligned}$$

2. Similarly when:

$$\sum_{n \in \mathcal{N}} \theta_{k,n} < \sum_{k \in \mathcal{K}(i)} \theta_{k,n}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}. \quad (35)$$

In this case, the optimal solution is given by:

$$\theta_{k,n} = \frac{|\mathcal{N}|}{|\mathcal{K}(i)| \cdot |\mathcal{N}|} = \frac{1}{|\mathcal{K}(i)|}, \forall k \in \mathcal{K}(i), \forall n \in \mathcal{N}.$$

Appendix III

Proof of Proposition 6.2

To prove the super-modularity of the present game, we need to verify the conditions in 6.2. First, the strategy space S_i is obviously a compact convex set of \mathbb{R}^N . Hence, it suffices to verify the super-modularity of the utility function as there are no constraint policies for \mathcal{G} :

$$\begin{aligned}
& \frac{\partial U_{i,n}}{\partial \pi_{l,n} \partial \pi_{i,n}} = \\
& \frac{1}{\log(1 + \sigma_{k,i,n})} \frac{G_{k,i,n} G_{k,l,n}}{\left(N_0 + \sum_{i' \neq i} \pi_{i',n} G_{k,i',n}\right)^2} \left(\frac{\sigma_{k,i,n}}{\log(1 + \sigma_{k,i,n})} - 1 \right).
\end{aligned}$$

As $\frac{x}{\log(1+x)} > 1$ for $x > 0$, $\frac{\partial U_{i,n}}{\partial \pi_{l,n} \partial \pi_{i,n}} \geq 0$, $\forall l \in \mathcal{I} - \{i\}$ and $\forall n \in \mathcal{N}$.

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