

NONNEGATIVE MATRIX INEQUALITIES AND THEIR APPLICATION TO NONCONVEX POWER CONTROL OPTIMIZATION*

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Abstract. Maximizing the sum rates in a multiuser Gaussian channel by power control is a nonconvex NP-hard problem that finds engineering application in code division multiple access (CDMA) wireless communication network. In this paper, we extend and apply several fundamental nonnegative matrix inequalities initiated by Friedland and Karlin in a 1975 paper to solve this nonconvex power control optimization problem. Leveraging tools such as the Perron–Frobenius theorem in nonnegative matrix theory, we (1) show that this problem in the power domain can be reformulated as an equivalent convex maximization problem over a closed unbounded convex set in the logarithmic signal-to-interference-noise ratio domain, (2) propose two relaxation techniques that utilize the reformulation problem structure and convexification by Lagrange dual relaxation to compute progressively tight bounds, and (3) propose a global optimization algorithm with ϵ -suboptimality to compute the optimal power control allocation. A byproduct of our analysis is the application of Friedland–Karlin inequalities to inverse problems in nonnegative matrix theory.

Key words. nonconvex optimization, convex relaxation, maximization of convex functions, nonnegative matrix theory, spectral radii of irreducible nonnegative matrices, wireless networks

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1. Introduction. We study the problem of total data throughput maximization using power control in a code division multiple access (CDMA) wireless communication network, where interference is a major source of performance impairment. Due to the broadcast nature of the wireless medium, the data rates in a wireless network are affected by interference when all the users transmit simultaneously over the same frequency band using CDMA. Power control is used to mitigate the effect of multiuser interference on performance and maximize the total data rates of all the users [4]. The CDMA wireless network can be modeled by an information-theoretic interference channel that treats multiuser interference as additive Gaussian noise [4]. A widely studied problem is to find the optimal power allocation that maximizes the sum rates over this multiuser Gaussian channel, and this requires solving a nonconvex problem [22], [5], [4], [7], [15]. This nonconvex problem also finds applications in the throughput maximization for digital subscriber line (DSL) wireline systems [24], [16], [17].¹

The complexity of an exhaustive search is prohibitively expensive, since this optimization problem is NP-hard, and may even be hard to approximate [16]. The authors in

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¹The sum rate maximization problem that we study in this paper is a special case of the problem formulation in the DSL setting, whereby each user allocates power only in a single frequency. On the other hand, each user in a DSL system allocates their power over more than one frequency.

[5], [22] formulated the problem as a signomial program, and used a successive convex approximation method based on geometric programming. In [7], the solution to a two-user special case was analyzed. The authors in [16] showed the NP-hardness of the problem, and used the Lyapunov theorem in functional analysis to deduce a zero duality gap result between a related primal of a continuous problem formulation in the DSL setting and its dual. The authors in [17] estimated the size of this duality gap for the finite problem in the DSL setting using Lagrangian dual relaxation combined with a linear program. The authors in [20] proposed approximation algorithms to solve the problem with individual power constraints (applicable to a CDMA uplink system). The authors in [21] solved the problem with a single total power constraint (applicable to a CDMA downlink system) under low to medium interference conditions.

We now state briefly the sum rate maximization problem with individual power constraints. We refer the readers to section 2 for all definitions, notations, and motivations. Let $F = [f_{ij}]_{i,j=1}^L$ and $\mathbf{v} = (v_1, \dots, v_L)^\top$ be an $L \times L$ matrix with zero diagonal and positive off-diagonal elements and a positive vector, respectively. Let $\bar{\mathbf{p}} = (\bar{p}_1, \dots, \bar{p}_L)^\top$ and $\mathbf{w} = (w_1, \dots, w_L)^\top$ be a given positive vector and a given probability vector, respectively. The sum rate maximization problem is given by

$$(1.1) \quad \max_{0 \leq p_l \leq \bar{p}_l \forall l} \sum_{l=1}^L w_l \log \left(1 + \frac{p_l}{\sum_{j \neq l} f_{lj} p_j + v_l} \right).$$

This is a nonconvex optimization problem that has a nonlinear-fractional objective function of positive variables over a simple box constraint set. The exact solution to this problem is also known to be (strongly) NP-hard [16]. An often used technique to tackle nonconvexity is the standard Lagrange dual relaxation of (1.1) in the power domain. However, the shortcoming of this approach is that there can exist a positive duality gap between the global optimal primal and optimal dual value of (1.1) [16]. Also, finding an optimal primal solution given an optimal dual solution, or vice versa, is in general difficult.

We adopt a reformulation-relaxation approach to tackle (1.1). Our reformulation possesses certain desirable properties, which enable the application of nonnegative matrix theory, especially the Friedland–Karlin inequalities stated in [10], to find the global optimal solution and motivate efficient relaxation techniques. In particular, we utilize the problem structure to develop suitable fast computational procedures for solving and computing useful bounds to the sum rate maximization problem. Furthermore, analytical solution to both the sum rate maximization problem and its relaxed problem can also be characterized by the spectra of specially crafted nonnegative matrices. A byproduct of our analysis is a refinement of the Friedland–Karlin inequalities in [10] and its application to an inverse problem in nonnegative matrix theory. From an engineering perspective, our algorithms operate in the logarithmic signal-to-interference-noise ratio domain or, equivalently, the dB domain that is lingua franca in existing wireless technology.

Overall, the contributions of the paper are as follows:

1. We study a reformulation of the sum rate maximization problem showing that it is equivalent to a convex maximization problem on a closed unbounded convex set.
2. Exploiting the structure of the reformulated problem, we propose two relaxation techniques that find progressively tighter bounds on the global optimal value. The first one is a convex relaxation technique that uses Lagrange duality (and its connection to convex envelope relaxation) and the Friedland–Karlin

inequalities (basic inequalities that characterize the spectral radius of a nonnegative matrix) to solve a sequence of linear programs. The second method exploits the spectra of specially crafted nonnegative matrices in a successive convex approximation method.

3. Utilizing the relaxation techniques, we propose a global optimization algorithm (with ϵ -suboptimality) to solve the sum rate maximization problem.
4. We give new applications of the Friedland–Karlin inequalities to inverse problems in nonnegative matrix theory.
5. Numerical examples illustrate the performance of our techniques, and include a comparison between our relaxation techniques and the standard Lagrange dual relaxation.

This paper is organized as follows. In section 2, we state definitions, notations, and a short motivation. We give a characterization of the image of the multidimensional box $[\mathbf{0}, \bar{\mathbf{p}}] \subset \mathbb{R}_+^L$ by a mapping in terms of the spectral radii of a set of nonnegative matrices. In section 3, we study the problem of sum rate maximization by power control in a wireless network. We give necessary and sufficient conditions for an extremal point $\mathbf{p} \in [\mathbf{0}, \bar{\mathbf{p}}]$ to be a local optimal value. In section 4, we exploit the structure of the reformulated problem to study two relaxation techniques to find useful upper bounds to the global optimal value. In section 5, we propose a global optimization algorithm to solve the sum rate maximization problem. In section 6, we evaluate the performance of our algorithms. In section 7, we conclude our paper. In Appendix A, we restate some useful results from [10], and give their extensions and applications to inverse problems in nonnegative matrix theory, which are needed in this paper.

2. Notations and preliminary results. Throughout the paper, we use the following notations. Let $\mathbb{R}^{m \times n} \supset \mathbb{R}_+^{m \times n}$ denote the set of $m \times n$ matrices and its subset of nonnegative matrices. For $A, B \in \mathbb{R}^{m \times n}$, we denote $A \preceq B$ if $B - A \in \mathbb{R}_+^{m \times n}$. We denote $A \leq B$, $A < B$ if $B - A$ is a nonzero nonnegative and positive matrix, respectively. We denote the entries of a matrix $A \in \mathbb{R}^{m \times n}$ by the small letters; i.e., $A = [a_{ij}]_{i,j=1}^{m,n}$. Identify $\mathbb{R}^m = \mathbb{R}^{m \times 1}$, $\mathbb{R}_+^m = \mathbb{R}_+^{m \times 1}$.

A column vector is denoted by the **bold** letter $\mathbf{x} = (x_1, \dots, x_L)^\top \in \mathbb{R}^L$. We denote $e^{\mathbf{x}} := (e^{x_1}, \dots, e^{x_m})^\top$. For $\mathbf{x} > \mathbf{0}$, we let $\mathbf{x}^{-1} := (\frac{1}{x_1}, \dots, \frac{1}{x_m})^\top$ and $\log \mathbf{x} = (\log x_1, \dots, \log x_L)^\top$. Let $\mathbf{x} \circ \mathbf{y}$ denote the Schur product of the vectors \mathbf{x} and \mathbf{y} ; i.e., $\mathbf{x} \circ \mathbf{y} = [x_1 y_1, \dots, x_L y_L]^\top$. Let $\mathbf{1} = (1, \dots, 1)^\top \in \mathbb{R}^L$. For $\mathbf{p} \leq \bar{\mathbf{p}} \in \mathbb{R}^L$, denote by $[\mathbf{p}, \bar{\mathbf{p}}]$ the set of all $\mathbf{x} \in \mathbb{R}^L$ satisfying $\mathbf{p} \leq \mathbf{x} \leq \bar{\mathbf{p}}$. For a vector $\mathbf{y} = (y_1, \dots, y_L)^\top$, denote by $\text{diag}(\mathbf{y})$ the diagonal matrix $\text{diag}(y_1, \dots, y_L)$. We let $(g(\mathbf{y}))_l$ denote the l th element of a function vector $g(\mathbf{y}) : \mathbb{R}^L \rightarrow \mathbb{R}^L$. For example, when $B \in \mathbb{R}_+^{L \times L}$ and $\mathbf{y} \in \mathbb{R}_+^L$, $(B\mathbf{y})_l$ denotes the l th element of the vector $B\mathbf{y}$. The Perron–Frobenius eigenvalue of a nonnegative matrix F is denoted as $\rho(F)$, and the Perron (right) and left eigenvector of F associated with $\rho(F)$ are denoted by $\mathbf{x}(F)$ and $\mathbf{y}(F)$ (or simply \mathbf{x} and \mathbf{y} when the context is clear), respectively. Assume that F is an irreducible nonnegative matrix. Then, $\rho(F)$ is simple and positive, and both $\mathbf{x}(F)$ and $\mathbf{y}(F)$ are positive [2]. We will assume the normalization $\mathbf{x}(F) \circ \mathbf{y}(F)$ is a probability vector. The superscript $(\cdot)^\top$ denotes transpose. For a positive integer n , denote by $\langle n \rangle$ the set $\{1, \dots, n\}$. Let $P : X \rightarrow Y$ be a mapping from the space X to the space Y . For a subset $Z \subset X$, we denote by $P(Z)$ the image of the set Z .

Consider an interference channel with L transmitter–receiver pairs (each transmitter–receiver pair is also called a user). The data transmission in this system with L users can be modeled as a Gaussian interference channel given by the following baseband signal model:

$$(2.1) \quad \mathbf{y}_l = \mathbf{h}_l \mathbf{x}_l + \sum_{j \neq l} \mathbf{h}_{lj} \mathbf{x}_j + \mathbf{z}_l,$$

where $\mathbf{y}_l \in \mathbb{C}^{1 \times 1}$ is the received signal of the l th user, $\mathbf{h}_{lj} \in \mathbb{C}^{1 \times 1}$ is the channel coefficient between the transmitter of the j th user and the receiver of the l th user, $(\mathbf{x}_1, \dots, \mathbf{x}_L)^\top \in \mathbb{C}^{L \times 1}$ is the transmitted (information carrying) signal vector, and \mathbf{z}_l 's are the independent and identically distributed (i.i.d.) additive white Gaussian noise (AWGN) coefficient with variance $n_l/2$ on each of its real and imaginary components. The first term on the right-hand side of (2.1) represents the desired signal, whereas the second term represents the interference signals from the other users. At each transmitter, the signal is constrained by an average power constraint, i.e., $\mathbb{E}[|\mathbf{x}_l|^2] = p_l$, which we assume to be upper bounded by \bar{p}_l for all l .

The vector $\mathbf{p} = (p_1, \dots, p_L)^\top$ is the transmit power vector and is the optimization variable of interest in this paper. Let $G = [g_{lj}]_{l,j=1}^L > 0_{L \times L}$ be an $L \times L$ channel gain matrix, where $g_{lj} = |\mathbf{h}_{lj}|^2$ is the channel gain from the j th transmitter to the l th receiver, and $\mathbf{n} = (n_1, \dots, n_L)^\top > \mathbf{0}$, where n_l is the noise power at the l th receiver. Assuming a linear matched-filter receiver at each user (treating multiuser interference as additive Gaussian noise), the signal-to-interference-noise ratio (SINR) for the l th receiver is defined as the ratio of the received signal power $g_{ll}p_l$ to the sum of interference signal power and additive Gaussian noise power $\sum_{j \neq l} g_{lj}p_j + n_l$. We denote the SINR of the l th receiver by γ_l , and consider it as a scalar nonnegative function of \mathbf{p} as follows. Let us first define

$$(2.2) \quad F = [f_{ij}]_{i,j=1}^L, \quad \text{where } f_{ij} = \begin{cases} 0 & \text{if } i = j, \\ \frac{g_{ji}}{g_{ii}} & \text{if } i \neq j, \end{cases}$$

and

$$(2.3) \quad \mathbf{g} = (g_{11}, \dots, g_{LL})^\top, \quad \mathbf{n} = (n_1, \dots, n_L)^\top, \quad \mathbf{v} = \left(\frac{n_1}{g_{11}}, \frac{n_2}{g_{22}}, \dots, \frac{n_L}{g_{LL}} \right)^\top.$$

For $\mathbf{p} = (p_1, \dots, p_L)^\top \geq \mathbf{0}$, we define the following transformation: $\mathbf{p} \mapsto \boldsymbol{\gamma}(\mathbf{p})$, where

$$(2.4) \quad \gamma_l(\mathbf{p}) := \frac{g_{ll}p_l}{\sum_{j \neq l} g_{lj}p_j + n_l}, \quad l = 1, \dots, L,$$

and we denote the vector $\boldsymbol{\gamma}(\mathbf{p}) = (\gamma_1(\mathbf{p}), \dots, \gamma_L(\mathbf{p}))^\top = \mathbf{p} \circ (F\mathbf{p} + \mathbf{v})^{-1}$.

We state the following result that characterizes the mapping in (2.4), which was first established in [9]. We include a new proof on this result for completeness.

LEMMA 2.1. *Let \mathbf{p} be a nonnegative vector. Assume that $\boldsymbol{\gamma}(\mathbf{p})$ is defined by (2.4). Then, $\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))F) < 1$, where F is defined by (2.2). Hence, for $\boldsymbol{\gamma} = \boldsymbol{\gamma}(\mathbf{p})$,*

$$(2.5) \quad \mathbf{p} = P(\boldsymbol{\gamma}) := (I - \text{diag}(\boldsymbol{\gamma})F)^{-1} \text{diag}(\boldsymbol{\gamma})\mathbf{v}.$$

Vice versa, if $\boldsymbol{\gamma}$ is in the set

$$(2.6) \quad \boldsymbol{\Gamma} := \{\boldsymbol{\gamma} \geq \mathbf{0}, \rho(\text{diag}(\boldsymbol{\gamma})F) < 1\},$$

then the vector \mathbf{p} defined by (2.5) is nonnegative. Furthermore, $\boldsymbol{\gamma}(P(\mathbf{p})) = \boldsymbol{\gamma}$. That is, $\boldsymbol{\gamma}: \mathbb{R}_+^L \rightarrow \boldsymbol{\Gamma}$ and $P: \boldsymbol{\Gamma} \rightarrow \mathbb{R}_+^L$ are inverse mappings.

Proof. Observe that (2.4) is equivalent to the equality

$$(2.7) \quad \mathbf{p} = \text{diag}(\boldsymbol{\gamma})F\mathbf{p} + \text{diag}(\boldsymbol{\gamma})\mathbf{v}.$$

First, let us assume that \mathbf{p} is a positive vector; i.e., $\mathbf{p} > \mathbf{0}$. Hence, $\boldsymbol{\gamma}(\mathbf{p}) > \mathbf{0}$. Since all off-diagonal entries of F are positive, it follows that the matrix $\text{diag}(\boldsymbol{\gamma})F$ is irreducible. As $\mathbf{v} > \mathbf{0}$, we deduce that $\max_{l \in [1, n]} \frac{(\text{diag}(\boldsymbol{\gamma})F\mathbf{p})_l}{p_l} < 1$. The min-max characterization of Wielandt of $\rho(\text{diag}(\boldsymbol{\gamma})F)$ (for example, see [2] and [11, equation (38), p. 64]) implies that $\rho(\text{diag}(\boldsymbol{\gamma})F) < 1$. Hence, $\boldsymbol{\gamma}(\mathbf{p}) \in \boldsymbol{\Gamma}$. Assume that $\mathbf{p} \geq \mathbf{0}$. Note that $p_l > 0 \Leftrightarrow \gamma_l(\mathbf{p}) > 0$. So $\mathbf{p} = \mathbf{0} \Leftrightarrow \boldsymbol{\gamma}(\mathbf{p}) = \mathbf{0}$. Clearly, $\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{0}))F) = \rho(0_{L \times L}) = 0 < 1$. Assume that $\mathbf{p} \succeq \mathbf{0}$. Let $\mathcal{A} = \{l: p_l > 0\}$. Denote by $\boldsymbol{\gamma}(\mathbf{p})(\mathcal{A})$ the vector composed of positive entries of $\boldsymbol{\gamma}(\mathbf{p})$. Let $F(\mathcal{A})$ be the principal submatrix of F with rows and columns in \mathcal{A} . It is straightforward to see that $\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))F) = \rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p})(\mathcal{A}))F(\mathcal{A}))$. The arguments above imply that

$$\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))F) = \rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p})(\mathcal{A}))F(\mathcal{A})) < 1.$$

Assume that $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$. Then,

$$(2.8) \quad (I - \text{diag}(\boldsymbol{\gamma})F)^{-1} = \sum_{k=0}^{\infty} (\text{diag}(\boldsymbol{\gamma})F)^k \geq 0_{L \times L}.$$

Hence, $P(\boldsymbol{\gamma}) \geq \mathbf{0}$. The definition of $P(\boldsymbol{\gamma})$ implies that $\boldsymbol{\gamma}(P(\boldsymbol{\gamma})) = \boldsymbol{\gamma}$. \square

LEMMA 2.2. *The set $\boldsymbol{\Gamma} \subset \mathbb{R}_+^L$ is monotonic with respect to the order \geq . That is, if $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$ and $\boldsymbol{\gamma} \geq \boldsymbol{\beta} \geq \mathbf{0}$, then $\boldsymbol{\beta} \in \boldsymbol{\Gamma}$. Furthermore, the function $P(\boldsymbol{\gamma})$ is monotone on $\boldsymbol{\Gamma}$.*

$$(2.9) \quad P(\boldsymbol{\gamma}) \geq P(\boldsymbol{\beta}) \quad \text{if } \boldsymbol{\gamma} \in \boldsymbol{\Gamma} \quad \text{and} \quad \boldsymbol{\gamma} \geq \boldsymbol{\beta} \geq \mathbf{0}.$$

Equality holds if and only if $\boldsymbol{\gamma} = \boldsymbol{\beta}$.

Proof. Clearly, if $\boldsymbol{\gamma} \geq \boldsymbol{\beta} \geq \mathbf{0}$, then $\text{diag}(\boldsymbol{\gamma})F \geq \text{diag}(\boldsymbol{\beta})F$, which implies $\rho(\text{diag}(\boldsymbol{\gamma})F) \geq \rho(\text{diag}(\boldsymbol{\beta})F)$. Hence, $\boldsymbol{\Gamma}$ is monotonic. Next, we use the Neumann expansion (2.8) to deduce the monotonicity of P . The equality case is straightforward. \square

Note that $\boldsymbol{\gamma}(\mathbf{p})$ is not monotonic in \mathbf{p} . Indeed, if one increases only the l th coordinate of \mathbf{p} , then one increases the l th coordinate of $\boldsymbol{\gamma}(\mathbf{p})$ and decreases all other coordinates of $\boldsymbol{\gamma}(\mathbf{p})$. As usual, let $\mathbf{e}_l = (\delta_{1l}, \dots, \delta_{ll})^\top$, $l = 1, \dots, L$, be the standard basis in \mathbb{R}^L . We have the following result.

THEOREM 2.1. *Let $l \in [1, L]$ be an integer and $a > 0$. Denote by $[0, a]_l \times \mathbb{R}_+^{L-1}$ the set of all $\mathbf{p} = (p_1, \dots, p_L)^\top \in \mathbb{R}_+^L$ satisfying $p_l \leq a$. Then, the image of the set $[0, a]_l \times \mathbb{R}_+^{L-1}$ by the map $\boldsymbol{\gamma}$ in (2.4) is given by*

$$(2.10) \quad \rho(\text{diag}(\boldsymbol{\gamma})(F + (1/a)\mathbf{v}\mathbf{e}_l^\top)) \leq 1, \quad \boldsymbol{\gamma} \geq \mathbf{0}.$$

Furthermore, $\mathbf{p} = (p_1, \dots, p_L) \in \mathbb{R}_+^L$ satisfies the condition $p_l = a$ if and only if $\boldsymbol{\gamma} = \boldsymbol{\gamma}(\mathbf{p})$ satisfies

$$(2.11) \quad \rho(\text{diag}(\boldsymbol{\gamma})(F + (1/a)\mathbf{v}\mathbf{e}_l^\top)) = 1.$$

Proof. Suppose that $\boldsymbol{\gamma}$ satisfies (2.10). We claim that $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$. Suppose first that $\boldsymbol{\gamma} > \mathbf{0}$. Then, $\text{diag}(\boldsymbol{\gamma})(F + t_1\mathbf{v}\mathbf{e}_l^\top) \preceq \text{diag}(\boldsymbol{\gamma})(F + t_2\mathbf{v}\mathbf{e}_l^\top)$ for any $t_1 < t_2$. [11, Lemma 2, section 2, Chapter XIII] yields

$$(2.12) \quad \begin{aligned} \rho(\text{diag}(\boldsymbol{\gamma})F) &< \rho(\text{diag}(\boldsymbol{\gamma})(F + t_1\mathbf{v}\mathbf{e}_l^\top)) < \rho(\text{diag}(\boldsymbol{\gamma})(F + t_2\mathbf{v}\mathbf{e}_l^\top)) \\ &< \rho(\text{diag}(\boldsymbol{\gamma})(F + (1/a)\mathbf{v}\mathbf{e}_l^\top)) \leq 1 \quad \text{for } 0 < t_1 < t_2 < 1/a. \end{aligned}$$

Thus, $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$. Combine the above argument with the arguments of the proof of Lemma 2.1 to deduce that $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$ for $\boldsymbol{\gamma} \geq \mathbf{0}$.

We now show that $(P(\boldsymbol{\gamma}))_l \leq a$. The continuity of P implies that it suffices to consider the case $\boldsymbol{\gamma} > \mathbf{0}$. Combine the Perron–Frobenius theorem (see, e.g., [2]) with (2.12) to deduce

$$(2.13) \quad 0 < \det(I - \text{diag}(\boldsymbol{\gamma})(F + t\mathbf{v}\mathbf{e}_l^\top)) \quad \text{for } t \in [0, a^{-1}).$$

We now expand the right-hand side of the above inequality. Let $B = \mathbf{xy}^\top \in \mathbb{R}^{L \times L}$ be a rank one matrix. Then, B has $L - 1$ zero eigenvalues and one eigenvalue equal to $\mathbf{y}^\top \mathbf{x}$. Hence, $I - \mathbf{xy}^\top$ has $L - 1$ eigenvalues equal to 1 and one eigenvalue is $(1 - \mathbf{y}^\top \mathbf{x})$. Therefore, $\det(I - \mathbf{xy}^\top) = 1 - \mathbf{y}^\top \mathbf{x}$. Since $\boldsymbol{\gamma} \in \boldsymbol{\Gamma}$, $(I - \text{diag}(\boldsymbol{\gamma})F)$ is invertible. Thus, for any $t \in \mathbb{R}$,

$$(2.14) \quad \begin{aligned} & \det(I - \text{diag}(\boldsymbol{\gamma})(F + t\mathbf{v}\mathbf{e}_l^\top)) \\ &= \det(I - \text{diag}(\boldsymbol{\gamma})F) \det(I - t((I - \text{diag}(\boldsymbol{\gamma})F)^{-1} \text{diag}(\boldsymbol{\gamma})\mathbf{v})\mathbf{e}_l^\top) \\ & \det(I - \text{diag}(\boldsymbol{\gamma})F)(1 - t\mathbf{e}_l^\top(I - \text{diag}(\boldsymbol{\gamma})F)^{-1} \text{diag}(\boldsymbol{\gamma})\mathbf{v}). \end{aligned}$$

Combine (2.13) with the above identity to deduce that

$$(2.15) \quad 1 > t\mathbf{e}_l^\top(I - \text{diag}(\boldsymbol{\gamma})F)^{-1} \text{diag}(\boldsymbol{\gamma})\mathbf{v} = tP(\boldsymbol{\gamma})_l \quad \text{for } t \in [0, a^{-1}).$$

Letting $t \nearrow a^{-1}$, we deduce that $(P(\boldsymbol{\gamma}))_l \leq a$. Hence, the set of $\boldsymbol{\gamma}$ defined by (2.10) is a subset of $\boldsymbol{\gamma}([0, a]_l \times \mathbb{R}_+^{L-1})$.

Let $\mathbf{p} \in [0, a]_l \times \mathbb{R}_+^{L-1}$ and denote $\boldsymbol{\gamma} = \boldsymbol{\gamma}(\mathbf{p})$. We show that $\boldsymbol{\gamma}$ satisfies (2.10). Lemma 2.1 implies that $\rho(\text{diag}(\boldsymbol{\gamma})F) < 1$. Since $\mathbf{p} = P(\boldsymbol{\gamma})$ and $p_l \leq a$, we deduce (2.15). Use (2.14) to deduce (2.13). As $\rho(\text{diag}(\boldsymbol{\gamma})F) < 1$, the inequality (2.13) implies that $\rho(\text{diag}(\boldsymbol{\gamma})(F + t\mathbf{v}\mathbf{e}_l^\top)) < 1$ for $t \in (0, a^{-1})$. Hence, (2.10) holds.

It remains to show that the condition (2.11) holds if and only if $(P(\boldsymbol{\gamma}))_l = a$. Assume that $\mathbf{p} = (p_1, \dots, p_L)^\top \in \mathbb{R}_+^L$, $p_l = a$ and let $\boldsymbol{\gamma} = \boldsymbol{\gamma}(\mathbf{p})$. We claim that equality holds in (2.10). Assume to the contrary that $\rho(\text{diag}(\boldsymbol{\gamma})(F + (1/a)\mathbf{v}\mathbf{e}_l^\top)) < 1$. Then, there exists $\boldsymbol{\beta} > \boldsymbol{\gamma}$ such that $\rho(\text{diag}(\boldsymbol{\beta})(F + (1/a)\mathbf{v}\mathbf{e}_l^\top)) < 1$. Since P is monotonic, $(P(\boldsymbol{\beta}))_l > p_l = a$. On the other hand, since $\boldsymbol{\beta}$ satisfies (2.10), we deduce that $(P(\boldsymbol{\beta}))_l \leq a$. This contradiction yields (2.11). Similarly, if $\boldsymbol{\gamma} \geq \mathbf{0}$ and (2.11) holds, then $(P(\boldsymbol{\gamma}))_l = a$. \square

COROLLARY 2.2. *Let $\bar{\mathbf{p}} = (\bar{p}_1, \dots, \bar{p}_L)^\top$ be a given positive vector. Then, $\boldsymbol{\gamma}([\mathbf{0}, \bar{\mathbf{p}}])$, the image of the set $[\mathbf{0}, \bar{\mathbf{p}}]$ by the map $\boldsymbol{\gamma}$ (2.4), is given by*

$$(2.16) \quad \rho(\text{diag}(\boldsymbol{\gamma})(F + (1/\bar{\mathbf{p}}_l)\mathbf{v}\mathbf{e}_l^\top)) \leq 1 \quad \text{for } l = 1, \dots, L, \quad \text{and } \boldsymbol{\gamma} \in \mathbb{R}_+^L.$$

In particular, any $\boldsymbol{\gamma} \in \mathbb{R}_+^L$ satisfying the conditions (2.16) satisfies the inequalities

$$(2.17) \quad \boldsymbol{\gamma} \leq \bar{\boldsymbol{\gamma}} = (\bar{\gamma}_1, \dots, \bar{\gamma}_L)^\top, \quad \text{where } \bar{\gamma}_l = \frac{\bar{p}_l}{v_l}, \quad i = 1, \dots, L.$$

Proof. Theorem 2.1 yields that $\boldsymbol{\gamma}([\mathbf{0}, \bar{\mathbf{p}}])$ is given by (2.16). Using (2.4), we have

$$\gamma_l(\mathbf{p}) = \frac{p_l}{(F\mathbf{p})_l + v_l} \leq \frac{p_l}{v_l} \leq \frac{\bar{p}_l}{v_l} \quad \text{for } \mathbf{p} \in [\mathbf{0}, \bar{\mathbf{p}}].$$

It is easy to see that equality holds for $\mathbf{p} = \bar{p}_l \mathbf{e}_l$. \square

Remark 1. Corollary 2.2 shows that the (nonconvex) set (2.16) is contained in a rectangular set (2.17). This fact is later used in our relaxation techniques.

3. The sum rate maximization problem. We assume the use of single-user decoder at each receiver, i.e., treating interference as additive Gaussian noise, and all users have perfect channel state information at the receiver. We also assume that the coherence time of the channel is less than the duration of the whole transmission by any user. This assumption is valid, for example, when fading occurs sufficiently slowly in the channel, i.e., flat-fading, so that the channel can be considered essentially fixed during transmission. We further assume that all users employ random Gaussian codes for transmission. In practice, Gaussian codes can be replaced by finite-order signal constellations such as the use of quadrature-amplitude modulation (QAM) or other practical (suboptimal) coding schemes. Assuming a fixed bit error rate (BER) at the receiver, the Shannon capacity formula can be used to deduce the achievable data rate (maximum information rate) of the l th user as [6]

$$(3.1) \quad r_l = \log \left(1 + \frac{\gamma_l(\mathbf{p})}{\Gamma} \right) \quad \text{nats/symbol,}$$

where Γ is the SINR gap to capacity, which is always greater than 1. In this paper, we absorb $(1/\Gamma)$ into g_{ll} for all l , and instead write the achievable data rate as $r_l = \log(1 + \gamma_l(\mathbf{p}))$.

Let $\mathbf{w} = (w_1, \dots, w_L)^\top$ be a given probability vector, where w_l is a positive weight assigned to the l th link to reflect priority (a larger weight reflects a higher priority). The problem of maximizing the sum rate can be stated as the following optimization problem:

$$(3.2) \quad \begin{aligned} & \text{maximize } \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p})) = \sum_{l=1}^L w_l \log(1 + \gamma_l(\mathbf{p})) \\ & \text{subject to } \mathbf{0} \leq \mathbf{p} \leq \bar{\mathbf{p}}, \\ & \text{variables: } \mathbf{p} = (p_1, \dots, p_L)^\top \in \mathbb{R}_+^L. \end{aligned}$$

Let $\mathbf{p}^* = (p_1^*, \dots, p_L^*)^\top$ be a global optimal solution to (3.2). We first derive necessary conditions obtained by straightforward differentiation for an optimal solution \mathbf{p}^* of (3.2).

LEMMA 3.1. *Denote the gradient of $\Phi_{\mathbf{w}}(\boldsymbol{\gamma})$ by*

$$\nabla \Phi_{\mathbf{w}}(\boldsymbol{\gamma}) = \left(\frac{w_1}{1 + \gamma_1}, \dots, \frac{w_L}{1 + \gamma_L} \right)^\top = \mathbf{w} \circ (\mathbf{1} + \boldsymbol{\gamma})^{-1}.$$

Let $\boldsymbol{\gamma}(\mathbf{p})$ be defined as in (2.4). Then, $\mathbf{H}(\mathbf{p}) = [\frac{\partial \gamma_l}{\partial p_j}]_{l,j=1}^L$, the Hessian matrix of $\boldsymbol{\gamma}(\mathbf{p})$, is given by

$$\mathbf{H}(\mathbf{p}) = \text{diag}((F\mathbf{p} + \mathbf{v})^{-1})(-\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))F + I).$$

In particular,

$$\nabla_{\mathbf{p}} \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p})) = \mathbf{H}(\mathbf{p})^\top \nabla \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p})).$$

COROLLARY 3.1. *Divide the set $\langle L \rangle = \{1, \dots, L\}$ into the following three disjoint sets S_{\max} , S_{in} , and S_0 :*

$$S_{\max} = \{l \in \langle L \rangle, p_l^* = \bar{p}_l\}, \quad S_{\text{in}} = \{l \in \langle L \rangle, p_l^* \in (0, \bar{p}_l)\}, \quad S_0 = \{l \in \langle L \rangle, p_l^* = 0\}.$$

Then, the following conditions hold:

$$(3.3) \quad \begin{aligned} (\mathbf{H}(\mathbf{p}^*)^\top \nabla \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}^*)))_l &\geq 0 \quad \text{for } l \in S_{\max}, \\ (\mathbf{H}(\mathbf{p}^*)^\top \nabla \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}^*)))_l &= 0 \quad \text{for } l \in S_{\text{in}}, \\ (\mathbf{H}(\mathbf{p}^*)^\top \nabla \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}^*)))_l &\leq 0 \quad \text{for } l \in S_0. \end{aligned}$$

Proof. Assume that $p_l^* = \bar{p}_l$. Then, $\frac{\partial}{\partial p_l} \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}))(\mathbf{p}^*) \geq 0$. Assume that $0 < p_l^* < \bar{p}_l$. Then, $\frac{\partial}{\partial p_l} \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}))(\mathbf{p}^*) = 0$. Assume that $p_l^* = 0$. Then, $\frac{\partial}{\partial p_l} \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}))(\mathbf{p}^*) \leq 0$. \square

Instead of solving (3.2) by dealing with the powers directly, we now turn to a reformulation-relaxation approach that solves and provides useful bounds to (3.2) indirectly. We first need the following lemma.

LEMMA 3.2. *Let \mathbf{w} be a probability vector, and assume that $\mathbf{p}^* = (p_1^*, \dots, p_L^*)^\top$ is an optimal solution to (3.2). Then, $p_l^* = \bar{p}_l$ for some l . Furthermore if $w_j = 0$, then $p_j^* = 0$.*

Proof. Assume to the contrary that $\mathbf{p}^* < \bar{\mathbf{p}}$. Let $\boldsymbol{\gamma}^* = \boldsymbol{\gamma}(\mathbf{p}^*)$. Since P is continuous on Γ , there exists $\boldsymbol{\gamma} \in \Gamma$ such that $\boldsymbol{\gamma} > \boldsymbol{\gamma}^*$ such that $P(\boldsymbol{\gamma}) < \bar{\mathbf{p}}$. Clearly, $\Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}^*)) < \Phi_{\mathbf{w}}(\boldsymbol{\gamma})$. As $\boldsymbol{\gamma} = \boldsymbol{\gamma}(P(\boldsymbol{\gamma}))$, we deduce that \mathbf{p}^* is not an optimal solution to (3.2), contrary to our assumptions.

Suppose that $w_j = 0$. For $\mathbf{p} = (p_1, \dots, p_L)^\top$, let \mathbf{p}_j be obtained from \mathbf{p} by replacing the j th coordinate in \mathbf{p} by 0. Assume that $p_j > 0$. Then, $\gamma_l(\mathbf{p}) < \gamma_l(\mathbf{p}_j)$ for $l \neq j$. Since $w_j = 0$, it follows that $\Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p})) < \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}_j))$. \square

We combine the above lemma with Theorem 2.1 and Corollary 2.2 to deduce an alternative formulation of (3.2).

THEOREM 3.2. *Problem (3.2) is equivalent to the following optimization problem:*

$$(3.4) \quad \begin{aligned} &\text{maximize } \Phi_{\mathbf{w}}(\boldsymbol{\gamma}) \\ &\text{subject to } \rho(\text{diag}(\boldsymbol{\gamma})(F + (1/\bar{p}_l)\mathbf{ve}_l^\top)) \leq 1 \quad \forall l, \\ &\text{variables: } \boldsymbol{\gamma} = (\gamma_1, \dots, \gamma_L)^\top \in \mathbb{R}_+^L, \end{aligned}$$

where $\boldsymbol{\gamma}^*$ is an optimal solution of the above problem if and only if $P(\boldsymbol{\gamma}^*)$ is an optimal solution \mathbf{p}^* of the problem (3.2). In particular, any optimal solution $\boldsymbol{\gamma}^*$ satisfies the equality (2.16) for some integer $l \in [1, L]$.

Remark 2. Note that (3.4) is a nonconvex problem having a strictly concave objective function and a set of nonconvex spectral radius constraints.

We now show that the optimization problem (3.4) can be restated as an optimization problem with a convex objective function on a closed unbounded convex domain. For $\boldsymbol{\gamma} = (\gamma_1, \dots, \gamma_L)^\top > 0$, we define the logarithmic mapping

$$(3.5) \quad \tilde{\boldsymbol{\gamma}} = \log \boldsymbol{\gamma};$$

i.e., $\boldsymbol{\gamma} = e^{\tilde{\boldsymbol{\gamma}}}$. Recall that for an irreducible nonnegative matrix $B \in \mathbb{R}_+^{L \times L}$, $\log \rho(\text{diag}(e^{\mathbf{x}})B)$ is a convex function [14]. This is the log-convexity property of the

Perron–Frobenius eigenvalue [3]. Furthermore, $\log(1 + e^t)$ is a strictly convex function in $t \in \mathbb{R}$. Hence, the optimization problem in (3.4) is equivalent to the problem

$$(3.6) \quad \begin{aligned} & \text{maximize } \Phi_{\mathbf{w}}(e^{\tilde{\gamma}}) \\ & \text{subject to } \log(\text{diag}(e^{\tilde{\gamma}})(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)) \leq 0 \quad \forall l, \\ & \text{variables: } \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L. \end{aligned}$$

The unboundedness of the convex set in (3.6) is due to the identity $0 = e^{-\infty}$. In view of Lemma 3.2, it suffices to consider the optimization problem (3.2) in the case where $\mathbf{w} > \mathbf{0}$. Using the reformulation in (3.6), we deduce the following result that any solution satisfying (3.3) in (3.1) is also locally optimal to (3.2).

THEOREM 3.3. *Consider the optimization problem (3.2). Then, any point \mathbf{p}^* satisfying $\mathbf{0} \leq \mathbf{p}^* \leq \bar{\mathbf{p}}$ and the conditions (3.3) is a local optimal solution.*

Proof. Since $\mathbf{w} > \mathbf{0}$, $\Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}})$ is a strictly convex function in $\tilde{\boldsymbol{\gamma}} \in \mathbb{R}^L$. Hence, the optimal value of (3.6) is achieved exactly on the extreme points of the closed unbounded set specified in (3.6). It may happen that some but not all the coordinates of the extreme point are $-\infty$. Translating this observation to the optimization problem (3.2), we deduce the theorem. \square

Although the reformulation in (3.6) is a convex maximization problem over a closed unbounded convex set, we choose not to rehash standard global optimization methods for solving a standard convex maximization problem (cf. [23], [13]). Rather, we choose to exploit the problem structure of (3.6) to first compute good bounds to (3.6) (cf. section 4) and then to propose a global optimization algorithm (with ϵ -suboptimality) to solve (3.6) (cf. section 5). The global optimization algorithm is motivated by the relaxation techniques and the problem structure (namely the log-convexity of the Perron–Frobenius eigenvalue; see, e.g., [14], [3], and the separability in the objective function), and it differs from the standard global optimization technique found in the literature, e.g., [23], [13]. In comparison to our recent work in [19], a cutting plane method (which is an outer approximation method [23], [13]) was proposed in [19] to solve (3.6) asymptotically. This cutting plane method, however, requires vertex enumeration (a time-consuming procedure for large problem size) at each iteration of the method and does not yield an ϵ -suboptimal solution.

We now give simple lower and upper bounds on the value of (3.2).

LEMMA 3.3. *Consider the optimization problem (3.2). Denote $R = \max_{l \in \langle L \rangle} \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)$. Let $\bar{\boldsymbol{\gamma}}$ be defined by (2.17). Then,*

$$\Phi_{\mathbf{w}}((1/R)\mathbf{1}) \leq \max_{\mathbf{p} \in [\mathbf{0}, \bar{\mathbf{p}}]} \Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p})) \leq \Phi_{\mathbf{w}}(\bar{\boldsymbol{\gamma}}).$$

Proof. By Corollary 2.2, $\boldsymbol{\gamma}(\mathbf{p}) \leq \bar{\boldsymbol{\gamma}}$ for $\mathbf{p} \in [\mathbf{0}, \bar{\mathbf{p}}]$. Hence, the upper bound holds. Clearly, for $\boldsymbol{\gamma} = (1/R)\mathbf{1}$, we have that $\rho(\text{diag}(\boldsymbol{\gamma})(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)) \leq 1$ for $l \in \langle L \rangle$. Then, from Theorem 3.2, $\Phi_{\mathbf{w}}((1/R)\mathbf{1})$ yields the lower bound. Equality is achieved in the lower bound when $\mathbf{p}^* = t\mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = \arg \max_{l \in \langle L \rangle} \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)$ for some $t > 0$. \square

The upper bound in Lemma 3.3 is trivial and can be too loose to be useful (as it disregards the interference power and the number of interferers at each receiver). The lower bound is obtained when all the users have a common SINR value. Necessary conditions under which this lower bound is tight are given later (cf. Corollary 5.2). We

will examine how to exploit the problem structure of (3.6) to obtain progressively tighter bounds in section 4.

4. Relaxations and convex approximations. In this section, we use two different approaches that exploit the problem structure of (3.4) or equivalently (3.6) to construct several relaxed versions of (3.4), which can compute useful upper bounds to (3.4). The first relaxation approach leverages the Friedland–Karlin inequalities, the separability of the objective function, and its convex envelope over box constraints (a sum of readily computed functions) to construct a linear program whose optimal value upper bounds that of (3.4). Progressively tighter bounds are then obtained by successive partitioning of the box constraints. The second relaxation approach replaces the L spectral constraints in (3.4) by a single one, which has three different versions depending on the choice of a nonnegative matrix. Each relaxed version is still nonconvex, but necessary conditions under which the relaxations are tight are given. Further, simpler algorithms derived from this approach are shown numerically to solve optimally these relaxations, thus providing useful upper bounds to (3.2). Figure 4.1 gives an overview of the development of these two relaxation approaches as well as a global optimization approach based on these relaxation techniques (see section 5 later).

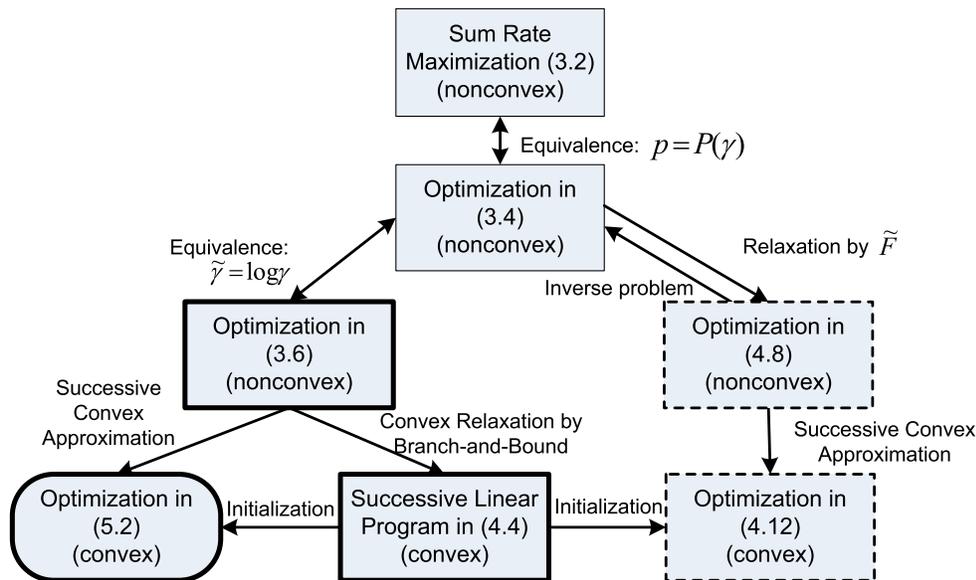


FIG. 4.1. Overview of the two relaxation techniques and a global optimization technique used on the sum rate maximization problem: The relaxation techniques are (1) a convex relaxation with branch-and-bound method, and (2) a relaxation by three different versions of a matrix \tilde{F} with successive convex approximation method and its connection to the inverse problem given in the appendix (see Theorems 4.2 and 4.3). The global optimization technique relies on the first relaxation technique to find a good initial point (within ϵ -suboptimality) that is then combined with the successive convex approximation method to solve the sum rate maximization problem. The key optimization problems (whether convex or nonconvex) and their relationships are also highlighted. Key steps that lead to the algorithms developed will be highlighted in the main text by the blocks with different line boundary patterns in the figure.

4.1. Convex relaxation. We replace the (convex) spectral radius constraint set in (3.6) with a larger set by exploiting the Friedland–Karlin inequalities. We thus consider the following optimization problem:

$$\begin{aligned}
 & \text{maximize } \Phi_{\mathbf{w}}(e^{\tilde{\gamma}}) \\
 & \text{subject to } \sum_{j=1}^L (\mathbf{x}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \circ \mathbf{y}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top))_j \tilde{\gamma}_j \\
 & \qquad \leq -\log \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \quad \forall l, \\
 & \qquad -K \leq \tilde{\gamma}_l \leq \log \bar{\gamma}_l \quad \forall l, \\
 (4.1) \quad & \text{variables: } \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L.
 \end{aligned}$$

Note that (4.1) is a convex maximization problem with a polyhedron constraint set (still a nonconvex and NP-hard problem). Observe that the constraint set of (4.1) consists of a polyhedron and a box constraint set. In particular, this constraint structure allows us to compute useful upper bounds to (4.1) by exploiting several results in [8] that connect the relationship between relaxation via convexification and the Lagrange dual relaxation. More precisely, the optimal Lagrange dual of (4.1) (which upper bounds (4.1)) can be computed by considering the convex envelope of (4.1).

To compute the convex envelope of a separable function over a box constraint set, it is sufficient to compute the convex envelope of the individual summand of the function over their respective domains (cf. Theorem 2.3 in [8]). For any l , the convex envelope of $w_l \log(1 + e^{\tilde{\gamma}_l})$ over a constraint set $\check{\gamma}_l \leq \tilde{\gamma}_l \leq \hat{\gamma}_l$ is a linear function in $\tilde{\gamma}_l$ given by

$$(4.2) \quad w_l \left(\frac{\log(1 + e^{\hat{\gamma}_l}) - \log(1 + e^{\check{\gamma}_l})}{\hat{\gamma}_l - \check{\gamma}_l} (\tilde{\gamma}_l - \check{\gamma}_l) + \log(1 + e^{\check{\gamma}_l}) \right).$$

Now, using (4.2) and letting $\check{\gamma}_l = -K$, $\hat{\gamma}_l = \log \bar{\gamma}_l$ for all l , we replace the objective function of (4.1) by its convex envelope over the box constraint set $\{-K \leq \tilde{\gamma}_l \leq \log \bar{\gamma}_l, \forall l\}$ to obtain the following linear program:²

$$\begin{aligned}
 & \text{maximize } \Phi_{\mathbf{w}}^c(e^{\tilde{\gamma}}) := \sum_{l=1}^L w_l \left(\frac{\log(1 + \bar{\gamma}_l) - \log(1 + e^{-K})}{\log \bar{\gamma}_l + K} (\tilde{\gamma}_l + K) + \log(1 + e^{-K}) \right) \\
 & \text{subject to } \sum_{j=1}^L (\mathbf{x}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \circ \mathbf{y}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top))_j \tilde{\gamma}_j \\
 & \qquad \leq -\log \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \quad \forall l, \\
 & \qquad -K \leq \tilde{\gamma}_l \leq \log \bar{\gamma}_l \quad \forall l, \\
 (4.3) \quad & \text{variables: } \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L.
 \end{aligned}$$

Interestingly, the optimal Lagrange dual of (4.3) (equivalent to the optimal value of (4.3) and also obtained by a dual linear program) is equal to the optimal Lagrange dual of (4.1) [8]. In other words, (4.3) is the “dual of the dual” of (4.1). For a large enough K (to approximate $K \rightarrow \infty$), the optimal Lagrange dual of (4.3) upper bounds the optimal

²Due to separability, the convex envelope of $\Phi_{\mathbf{w}}(e^{\tilde{\gamma}})$, denoted by $\Phi_{\mathbf{w}}^c(e^{\tilde{\gamma}})$, is the sum of the convex envelope of its constituents.

Lagrange dual of (3.6). Thus, the optimal value of (4.3) gives an upper bound to (3.6).

Although the upper bound obtained by solving (4.3) may be loose, tighter bounds to (3.6) can be obtained iteratively by combining this convex relaxation approach with a branch-and-bound method in [1], [8], [13] that subdivides the set $\{-K \leq \tilde{\gamma}_l \leq \log \bar{\gamma}_l \forall l\}$ into successively smaller subsets (the rectangular method in [1]; also see, e.g., Chapter 7 in [13]). The search for the global optimal solution is performed over the subdivided sets organized in a binary tree data structure. More precisely, using the branch-and-bound method, (4.3) is solved in the first iteration (at the root of the binary tree). In subsequent iterations (lower levels of the binary tree), the set of lower and upper bounds on $\tilde{\boldsymbol{\gamma}}$ in (4.3) is replaced by a subdivided set, and the objective function of (4.3) is then replaced with a reweighted function using (4.2), i.e., the convex envelope of $\Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}})$ over the subdivided set. In particular, at the k th iteration of the branch-and-bound algorithm, we consider a subdivided box $\{\tilde{\gamma}_l | \check{\gamma}_l(k) \leq \tilde{\gamma}_l \leq \hat{\gamma}_l(k) \forall l\}$ and solve

$$\begin{aligned} & \text{maximize} \quad \sum_{l=1}^L w_l \left(\frac{\log(1 + e^{\hat{\gamma}_l(k)}) - \log(1 + e^{\check{\gamma}_l(k)})}{\hat{\gamma}_l(k) - \check{\gamma}_l(k)} (\tilde{\gamma}_l - \check{\gamma}_l(k)) + \log(1 + e^{\check{\gamma}_l(k)}) \right) \\ & \text{subject to} \quad \sum_{j=1}^L (\mathbf{x}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \circ \mathbf{y}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top))_j \tilde{\gamma}_j \\ & \quad \leq -\log \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \quad \forall l, \\ & \quad \check{\gamma}_l(k) \leq \tilde{\gamma}_l \leq \hat{\gamma}_l(k) \quad \forall l, \end{aligned}$$

(4.4) variables: $\tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L$

at one of the two nodes at the k th iteration of the branch-and-bound algorithm. We denote $\tilde{\boldsymbol{\gamma}}^{\text{LP}_k}$ as the optimal solution to (4.4). A feasible power vector can be obtained from $\mathbf{p}^{\text{LP}_k} = \min\{P(e^{\tilde{\boldsymbol{\gamma}}^{\text{LP}_k}}), \bar{\mathbf{p}}\}$, and a lower bound to (3.2) is then given by $\Phi_{\mathbf{w}}(\boldsymbol{\gamma}(\mathbf{p}^{\text{LP}_k}))$. Rules to subdivide a rectangular box constraint and branch into a selected subdivided set as well as the convergence of this branch-and-bound method can be found in [1].

In this method, at the k th iteration of the branch-and-bound algorithm, taking the maximum over all the lower bound at each child node across all the levels in the binary tree (denote this maximum value as L_{bb}^k and its corresponding solution as $\tilde{\boldsymbol{\gamma}}^{\text{BB}_k}$) gives a global lower bound on the optimal value of (3.2). Likewise, taking the maximum over all the upper bound at each child node across all the levels in the binary tree (denote this maximum value as U_{bb}^k) gives a global upper bound on the optimal value of (3.2). The difference between these two bounds is nonincreasing with k . Suppose $U_{bb}^k - L_{bb}^k \leq \epsilon$ for some positive ϵ ; then we have $\Phi_{\mathbf{w}}(\boldsymbol{\gamma}^*) \leq \Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}^{\text{BB}_k}}) + \epsilon$. Intuitively speaking, this relaxation method systematically narrows down the SINR region that contains the global optimal solution of (3.6); i.e., locate an ϵ -suboptimal neighborhood of $\tilde{\boldsymbol{\gamma}}^*$. This first relaxation method is denoted by the two blocks with thick boundary lines in Figure 4.1.

4.2. Relaxation by nonnegative matrices. In this section, we study the second relaxation method that uses specially constructed nonnegative matrices to find useful upper bounds to (3.2). Conditions under which the relaxations are tight are stated, and simpler (lower complexity and faster) algorithms are proposed to compute the upper bounds.

Now, we consider a general matrix \tilde{F} that is used to denote one of the following three matrices:

$$\begin{aligned}
& 1) F + \text{diag}(\tilde{\boldsymbol{\gamma}})^{-1}, \\
& 2) F + (1/\mathbf{1}^\top \tilde{\mathbf{p}})\mathbf{v}\mathbf{1}^\top, \\
(4.5) \quad & 3) F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top, \quad i = \arg \max_l \rho(F + (1/\tilde{p}_l)\mathbf{v}\mathbf{e}_l^\top).
\end{aligned}$$

Observe that the entries of \tilde{F} are functions of all the problem parameters of (3.2). In the following, we consider a relaxation to (3.4) that has only a single spectral radius constraint in \tilde{F} . We then utilize the spectra of \tilde{F} to find useful upper bounds to (3.4), which in turn upper bounds (3.2).

LEMMA 4.1. *Let $\mathbf{0} \leq \mathbf{p} \leq \tilde{\mathbf{p}}$. Assume that $\boldsymbol{\gamma}(\mathbf{p})$ is given by (2.4) and \tilde{F} is given by any of the three matrices in (4.5). Then,*

$$(4.6) \quad \mathbf{p} \geq \text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}\mathbf{p},$$

and

$$(4.7) \quad \rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}) \leq 1.$$

Proof. The assumption that $0 \leq p_l \leq \tilde{p}_l$ implies that $\frac{2u}{g_{ii}} \geq \tilde{f}_{ii}p_l$. From the definition of $\boldsymbol{\gamma}(\mathbf{p})$, we deduce that $p_l = \gamma_l(\mathbf{p})(\sum_{j \neq l} f_{lj}p_j + v_l)$, which together with the definition of \tilde{F} and the above observation implies (4.6). The inequality (4.7) is a consequence of Wielandt's characterization of the spectral radius of an irreducible nonnegative matrix [11]. Indeed, if $\mathbf{p} > \mathbf{0}$ —i.e., all the coordinates of \mathbf{p} are positive—then $\boldsymbol{\gamma}(\mathbf{p}) > \mathbf{0}$. Hence, $\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}$ is a positive matrix. Then, using Wielandt's characterization, we have

$$\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}) \leq \max_{l=1, \dots, L} \frac{(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}\mathbf{p})_l}{p_l} \leq 1.$$

Observe next that if $p_l = 0$, then $(\boldsymbol{\gamma}(\mathbf{p}))_l = 0$. So if some of $p_l = 0$, then $\rho(\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F})$ is the spectral radius of the maximal positive submatrix of $\text{diag}(\boldsymbol{\gamma}(\mathbf{p}))\tilde{F}$. By applying Wielandt's characterization to this positive submatrix, we deduce (4.7). \square

Lemma 4.1 shows that any feasible \mathbf{p} satisfies (4.7). This leads to the following relaxation of (3.4) that has only a single constraint involving \tilde{F} in (4.5).

LEMMA 4.2. *The optimal value of*

$$\begin{aligned}
& \text{maximize } \Phi_{\mathbf{w}}(\boldsymbol{\gamma}) \\
& \text{subject to } \rho(\text{diag}(\boldsymbol{\gamma})\tilde{F}) \leq 1, \\
& \quad \boldsymbol{\gamma} \leq \tilde{\boldsymbol{\gamma}}, \\
(4.8) \quad & \text{variables: } \boldsymbol{\gamma} = (\gamma_1, \dots, \gamma_L)^\top \in \mathbb{R}_+^L
\end{aligned}$$

is not less than the optimal value of (3.2). Further, using $P(\boldsymbol{\gamma})$, the optimal solution of (4.8) expressed in the power domain is given by $\mathbf{x}(\text{diag}(\boldsymbol{\gamma}')\tilde{F})$, where $\boldsymbol{\gamma}'$ solves (4.8). In particular, $P(\boldsymbol{\gamma}) = \mathbf{p}^$ if $P(\boldsymbol{\gamma}')$ satisfies (3.3) in Corollary 3.1.*

Proof. In view of (4.7), we see that the optimal value in (4.8) is achieved on a bigger set than the optimal value in (3.2). In view of (4.6), the optimal solution to (4.8) satisfies $\mathbf{p} = \text{diag}(\boldsymbol{\gamma}')\tilde{F}\mathbf{p}$. Together with (4.7), this implies that $\mathbf{p} = \mathbf{x}(\text{diag}(\boldsymbol{\gamma}')\tilde{F})$. \square

Remark 3. Note that the constraint $\boldsymbol{\gamma} \leq \tilde{\boldsymbol{\gamma}}$ has been included explicitly in (4.8), because the spectral radius constraint $\rho(\text{diag}(\boldsymbol{\gamma})\tilde{F}) \leq 1$ does not imply $\boldsymbol{\gamma} \leq \tilde{\boldsymbol{\gamma}}$ (cf. Corollary 2.2 and the reformulation in (3.4)).

The second relaxation method is to solve (4.8) by considering all the three choices of \tilde{F} in (4.5), and to find the tightest relaxation to (3.2) among the three choices of \tilde{F} . Note that, in (4.5), the first two nonnegative matrices have positive diagonals, whereas the third nonnegative matrix has only a single positive diagonal element. This fact will be important in characterizing the optimal solution of the relaxed problems based on the inverse problem given in the appendix (see Theorems 4.2 and 4.3 later). From a computational viewpoint, solving (4.8) is also useful when L is large (as the computational time to solve (3.4) increases with L).

COROLLARY 4.1. *We have $\rho(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) = 1$ in (4.8), where $\boldsymbol{\gamma}'$ solves (4.8) optimally.*

Proof. Corollary 4.1 is easily proved by noting that both the objective function and the spectral radius function in (4.8) increase with $\boldsymbol{\gamma}$. \square

Using the logarithmic mapping in (3.5), solving (4.8) is thus equivalent to solving

$$\begin{aligned}
 & \text{maximize } \Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{y}}}) \\
 & \text{subject to } \log \rho(\text{diag}(e^{\tilde{\boldsymbol{y}}})\tilde{F}) \leq 0, \\
 & \quad \tilde{\boldsymbol{y}} \leq \log \tilde{\boldsymbol{\gamma}}, \\
 (4.9) \quad & \text{variables: } \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L.
 \end{aligned}$$

Still, (4.9), or equivalently (4.8), is nonconvex and hard to solve. In the following, we give conditions that relate the optimal power \mathbf{p}^* and the solution of (4.9) for different \tilde{F} . These conditions are also necessary when the solution of (4.9) solves (3.6); i.e., (4.9) is a tight relaxation of (3.6).

Corollary 4.1 implies that if the optimizer of (4.8) $\boldsymbol{\gamma}'$ satisfies $P(\boldsymbol{\gamma}') \leq \bar{\mathbf{p}}$, then $P(\boldsymbol{\gamma}')$ is also the global optimizer of (3.2). Hence, $P(\boldsymbol{\gamma}') \leq \bar{\mathbf{p}}$ is a necessary and sufficient condition for the relaxation to be tight. Weaker necessary conditions can, however, be obtained by checking that the following holds:

$$(4.10) \quad \text{diag}(e^{\tilde{\boldsymbol{y}}})\tilde{F}\mathbf{p}^* = \mathbf{p}^*$$

for some \mathbf{p}^* . A summary of the necessary conditions on \mathbf{p}^* satisfying (4.10) for the three different versions of \tilde{F} is given in Table 4.1, whereby the relaxed problem (4.9) solves (3.6).

This second relaxation method is denoted by the two blocks with dotted boundary lines in Figure 4.1. Now, we consider using a successive convex approximation method to solve (4.9) directly. This method is motivated by the inverse problem given in the appendix. This is given in the following algorithm to solve (4.9) and also yield a feasible solution to (3.2).

TABLE 4.1

A comparison of the different versions of \tilde{F} in the second relaxation method, wherein the optimal solution in (4.8) is given by $\boldsymbol{\gamma}'$. Necessary conditions under which the relaxed problem (4.9) solves (3.6), equivalently (3.2), are given.

\tilde{F}	Necessary condition for $P(\boldsymbol{\gamma}') = \mathbf{p}^*$
$F + \text{diag}(\tilde{\boldsymbol{\gamma}})^{-1}$	$p_i^* = \{0, \bar{p}_i\}, \forall i$
$F + (1/\mathbf{1}^\top \bar{\mathbf{p}})\mathbf{v}\mathbf{1}^\top$	$\mathbf{p}^* = \bar{\mathbf{p}}$
$F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top,$ $i = \arg \max_i \rho(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$	$\mathbf{p}^* = \mathbf{x}$ with $p_i^* = \bar{p}_i,$ $\mathbf{w} = \mathbf{x} \circ \mathbf{y}$ (cf. Corollary 5.2)

ALGORITHM 1 (ITERATIVELY REWEIGHTED RELAXATION ALGORITHM).

1. Compute the weight $\mathbf{m}(k+1)$:

$$(4.11) \quad \mathbf{m}(k+1) = \frac{\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(k)}) \circ e^{\tilde{\mathbf{r}}(k)}}{\mathbf{1}^\top (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(k)}) \circ e^{\tilde{\mathbf{r}}(k)})}.$$

2. Obtain $\tilde{\boldsymbol{\gamma}}(k+1)$ as the optimal solution to

$$(4.12) \quad \begin{aligned} & \text{maximize} \quad \sum_{l=1}^L m_l(k+1) \tilde{\gamma}_l \\ & \text{subject to} \quad \log \rho(\text{diag}(e^{\tilde{\mathbf{r}}}) \tilde{F}) \leq 0, \tilde{\gamma}_l \leq \log \bar{\gamma}_l \quad \forall l, \\ & \text{variables:} \quad \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L. \end{aligned}$$

3. Compute the power

$$(4.13) \quad \mathbf{p}(k+1) = \min \{P(e^{\tilde{\mathbf{r}}(k+1)}), \bar{\mathbf{p}}\}.$$

LEMMA 4.3. *For any $\tilde{\boldsymbol{\gamma}}(0)$ in a sufficiently close neighborhood of $\tilde{\boldsymbol{\gamma}}'$, $\tilde{\boldsymbol{\gamma}}(k)$ in Algorithm 1 converges to the optimal solution of (4.8).*

Proof. We use the fact that in a sufficiently close neighborhood of $\tilde{\boldsymbol{\gamma}}'$, the domain set is convex, and the objective function $\Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}})$ is twice continuously differentiable. We then use a successive convex approximation technique to compute $\tilde{\boldsymbol{\gamma}}'$ assuming that the initial point is sufficiently close to $\tilde{\boldsymbol{\gamma}}'$. The convergence conditions for such a technique are given in [18], [5]. Instead of solving (4.8) directly, we replace the objective function of (4.8) in a neighborhood of a feasible point $\tilde{\boldsymbol{\gamma}}(0)$ by its Taylor series (up to the first order terms): $\Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(0)}) + (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(0)}) \circ e^{\tilde{\mathbf{r}}(0)})^\top (\tilde{\boldsymbol{\gamma}} - \tilde{\boldsymbol{\gamma}}(0))$. Assume a feasible $\tilde{\boldsymbol{\gamma}}(0)$ that is close to $\tilde{\boldsymbol{\gamma}}'$. We then compute a feasible $\tilde{\boldsymbol{\gamma}}(k+1)$ by solving the $(k+1)$ th approximation problem:

$$(4.14) \quad \begin{aligned} & \text{maximize} \quad (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(k)}) \circ e^{\tilde{\mathbf{r}}(k)}) / \mathbf{1}^\top (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}(k)}) \circ e^{\tilde{\mathbf{r}}(k)})^\top (\tilde{\boldsymbol{\gamma}} - \tilde{\boldsymbol{\gamma}}(k)) \\ & \text{subject to} \quad \log \rho(\text{diag}(e^{\tilde{\mathbf{r}}}) \tilde{F}) \leq 0, \quad \tilde{\gamma}_l \leq \log \bar{\gamma}_l \quad \forall l, \\ & \text{variables:} \quad \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_n)^\top \in \mathbb{R}^L, \end{aligned}$$

where $\tilde{\boldsymbol{\gamma}}(k)$ is the optimal solution of the k th approximation problem. This inner approximation technique converges to a local optimal solution [18], [5]. In addition, if $\tilde{\boldsymbol{\gamma}}(0)$ is sufficiently close to $\tilde{\boldsymbol{\gamma}}'$, then $\lim_{k \rightarrow \infty} \tilde{\boldsymbol{\gamma}}(k) = \tilde{\boldsymbol{\gamma}}'$.

Next, we leverage Corollary A.6 in the appendix to solve (4.14). At the global optimality of (4.8), we have the necessary condition of (4.8): $\mathbf{x}(\text{diag}(e^{\tilde{\mathbf{r}}}) \tilde{F}) \circ \mathbf{y}(\text{diag}(e^{\tilde{\mathbf{r}}}) \tilde{F}) = \nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}}) \circ e^{\tilde{\mathbf{r}}} / \mathbf{1}^\top (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\mathbf{r}}}) \circ e^{\tilde{\mathbf{r}}})$. \square

Remark 4. Now, (4.12) can be reformulated as an equivalent geometric program by utilizing a basic result in nonnegative matrix theory, which states that the Perron–Frobenius eigenvalue is given by $\rho(A) = \inf\{\lambda | A\mathbf{v} \leq \lambda \mathbf{v} \text{ for some positive } \mathbf{v}\}$ and that the infimum is achieved; e.g., see Chapter 4 of [3]. This fact can be used to transform the spectral radius constraint in (4.12) so that (4.12) can be readily solved using standard convex optimization solvers (e.g., those in [12]).

Interestingly, Algorithm 1 can be viewed as an iteratively reweighted method that produces better estimates with each iteration. The weights in (4.12) are determined by the previous solution. Now, suppose the positive weight vector $\nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}'}) \circ e^{\tilde{\gamma}'}/\mathbf{1}^\top(\nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}'}) \circ e^{\tilde{\gamma}'})$ is used as the weight input to (4.12). Then, (4.12) yields $\tilde{\gamma}'$. Intuitively speaking, if the weight vector is approximately proportional to $\nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}'}) \circ e^{\tilde{\gamma}'}/\mathbf{1}^\top(\nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}'}) \circ e^{\tilde{\gamma}'})$, then Algorithm 1 should converge to a unique $\tilde{\gamma}'$. Based on the inverse problem given in the appendix (see Theorems 4.2 and 4.3), we quantify this in the following for the different choices of \tilde{F} (and the diagonals of \tilde{F} matter in a unique $\tilde{\gamma}'$).

THEOREM 4.2. *Suppose that \tilde{F} is given by (1) $F + \text{diag}(\tilde{\gamma})^{-1}$ or (2) $F + (1/\mathbf{1}^\top\tilde{\mathbf{p}})\mathbf{v}\mathbf{1}^\top$ in (4.5). Let $\mathbf{m} = (m_1, \dots, m_L)^\top$ be a positive probability vector. Then,*

$$(4.15) \quad \max_{\boldsymbol{\gamma}=(\gamma_1, \dots, \gamma_L)^\top > \mathbf{0}, \rho(\text{diag}(\boldsymbol{\gamma})\tilde{F}) \leq 1} \sum_{l=1}^L m_l \log \gamma_l = \sum_{l=1}^L m_l \log \gamma'_l,$$

where $\boldsymbol{\gamma}' = (\gamma'_1, \dots, \gamma'_L)^\top > \mathbf{0}$ is the unique vector satisfying the following conditions: $\rho(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) = 1$ and $\mathbf{x}(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) \circ \mathbf{y}(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) = \mathbf{m}$.

Proof. We use Theorem A.3 and Corollary A.6 in the appendix to prove Theorem 4.2. \square

Combining Theorem A.3 and Corollary A.10 in the appendix, we deduce the following result.

THEOREM 4.3. *Suppose that \tilde{F} is given by (3) $F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top$, $i = \arg \max_l \rho(F + (1/\tilde{p}_l)\mathbf{v}\mathbf{e}_l^\top)$ in (4.5). Let $\mathbf{m} = (m_1, \dots, m_L)^\top$ be a positive probability vector satisfying the condition*

$$(4.16) \quad \sum_{j \neq l} m_j > m_l \quad \forall l \in \langle L \rangle.$$

Then,

$$(4.17) \quad \max_{\boldsymbol{\gamma}=(\gamma_1, \dots, \gamma_L)^\top > \mathbf{0}, \rho(\text{diag}(\boldsymbol{\gamma})\tilde{F}) \leq 1} \sum_{l=1}^L m_l \log \gamma_l = \sum_{l=1}^L m_l \log \gamma'_l,$$

where $\boldsymbol{\gamma}' = (\gamma'_1, \dots, \gamma'_L)^\top > \mathbf{0}$ is a vector satisfying the following conditions: $\rho(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) = 1$ and $\mathbf{x}(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) \circ \mathbf{y}(\text{diag}(\boldsymbol{\gamma}')\tilde{F}) = \mathbf{m}$.

The above last two theorems enable us to choose \mathbf{m} for which we know the solution to the optimization problems (4.15) and (4.17). Namely, choose $\boldsymbol{\beta}_1, \boldsymbol{\beta}_2 > \mathbf{0}$ such that $A_1 = \text{diag}(\boldsymbol{\beta}_1)\tilde{F}$, $A_2 = \text{diag}(\boldsymbol{\beta}_2)\tilde{F}$ have spectral radius one. Let $\mathbf{m}_i = \mathbf{x}(A_i) \circ \mathbf{y}(A_i)$ for $i = 1, 2$. Then, for \mathbf{m}_1 , (4.15) has the unique optimal solution $\boldsymbol{\gamma}^* = \boldsymbol{\beta}_1$. For \mathbf{m}_2 , (4.17) has an optimal solution $\boldsymbol{\gamma}^* = \boldsymbol{\beta}_2$. In view of Theorem A.3, \mathbf{m}_2 does not have to satisfy the condition $\sum_{j \neq l} m_j > m_l$ for all $l \in \langle L \rangle$.

4.3. Relaxation with improved initialization. Observe that it is viable to apply the first relaxation technique, i.e., the convex relaxation and branch-and-bound method, to (4.9), and obtain upper bounds to (4.9). The bounds obtained will be looser than that employed on (3.6). On the other hand, Algorithm 1 requires an initial point that is sufficiently close to the optimal solution. We now propose a natural procedure of finding such a good initial point. The basic idea is to employ the first relaxation technique, i.e., solve (4.4) iteratively by a branch-and-bound method to locate a point $\log \boldsymbol{\gamma}(\min\{P(e^{\tilde{\gamma}^{\text{LP}_k}}), \tilde{\mathbf{p}}\})$ close enough to the optimal solution $\tilde{\boldsymbol{\gamma}}^*$ —i.e., a point in an ϵ -suboptimal region—and then input it as the initial point in Algorithm 1.

5. Global optimization algorithm. We now state a global optimization algorithm that combines the relaxation techniques and the improved initialization in section 4 to solve (3.6) and equivalently to yield an optimal solution of (3.2) to within a prescribed accuracy on the suboptimality. This approach is denoted by the extreme left-hand-side block with round corners in Figure 4.1.

ALGORITHM 2 (ITERATIVELY REWEIGHTED OPTIMAL ALGORITHM).

Initial phase: Set a prescribed accuracy ϵ . Find an ϵ -suboptimal region of (3.6) using the branch-and-bound technique in section 4 and return as output the solution that yields the tightest lowest bound L_{bb} : $\tilde{\boldsymbol{\gamma}}^{BB}$. Set $\tilde{\boldsymbol{\gamma}}(0) = \tilde{\boldsymbol{\gamma}}^{BB}$.

1. Compute the weight $\mathbf{m}(k+1)$:

$$(5.1) \quad \mathbf{m}(k+1) = \frac{\nabla \Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}(k)}) \circ e^{\tilde{\boldsymbol{\gamma}}(k)}}{\mathbf{1}^\top (\nabla \Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}(k)}) \circ e^{\tilde{\boldsymbol{\gamma}}(k)})}.$$

2. Obtain $\tilde{\boldsymbol{\gamma}}(k+1)$ as the optimal solution to

$$(5.2) \quad \begin{aligned} & \text{maximize} \quad \sum_{l=1}^L m_l(k+1) \tilde{\gamma}_l \\ & \text{subject to} \quad \log \rho(\text{diag}(e^{\tilde{\boldsymbol{\gamma}}}(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top))) \leq 0 \quad \forall l, \\ & \text{variables:} \quad \tilde{\boldsymbol{\gamma}} = (\tilde{\gamma}_1, \dots, \tilde{\gamma}_L)^\top \in \mathbb{R}^L. \end{aligned}$$

3. Compute the power

$$(5.3) \quad \mathbf{p}(k+1) = \min \{P(e^{\tilde{\boldsymbol{\gamma}}(k+1)}), \bar{\mathbf{p}}\}.$$

THEOREM 5.1. *Suppose we have $U_{bb} - L_{bb} \leq \epsilon$ at the completion of the initial phase in Algorithm 2, and the initial phase yields an output $\tilde{\boldsymbol{\gamma}}^{BB}$ that is a feasible solution to (3.6). If $\tilde{\boldsymbol{\gamma}}(0) = \tilde{\boldsymbol{\gamma}}^{BB}$, then $\tilde{\boldsymbol{\gamma}}(k)$ in Algorithm 2 converges to a point $\tilde{\boldsymbol{\gamma}}_\epsilon$ in an ϵ -suboptimal neighborhood of the optimal solution of (3.6); i.e., $\Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}_\epsilon}) - \Phi_{\mathbf{w}}(e^{\tilde{\boldsymbol{\gamma}}_\epsilon}) \leq \epsilon$.*

Furthermore, $\tilde{\boldsymbol{\gamma}}(k)$ in Algorithm 2 converges to $\tilde{\boldsymbol{\gamma}}^$ for a sufficiently small ϵ .*

Proof. Theorem 5.1 can be proved by combining the previous proofs in section 4. \square

The following result demonstrates a special case in which the optimal solution of (3.2) is given analytically and can be computed by Algorithm 2 using a simpler initial point and in only one iteration.

COROLLARY 5.2. *If $\tilde{\gamma}_l^*$ is equal for all l , then $\mathbf{w} = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = \arg \max_l \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)$. In this special case, $\tilde{\boldsymbol{\gamma}}(k)$ in Algorithm 2 converges to $-\log \rho(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ in only one iteration from any initial point $\tilde{\boldsymbol{\gamma}}(0)$ such that $\tilde{\gamma}_l(0)$ are equal for all l . Moreover, $\mathbf{p}(k)$ in Algorithm 2 converges to the optimal solution of (3.2) given by $\mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ (up to a scaling factor).*

Proof. Suppose that $\tilde{\gamma}_l^*$ is equal (to a value $\tilde{\gamma}^*$) for all l . At optimality, the constraint set of (3.6) reduces to $\tilde{\gamma}^* + \log \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top) \leq 0$ for all l , and since at least one of the spectral radius constraints in (3.6) is tight, $\tilde{\gamma}^* = -\log \rho(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = \arg \max_l \rho(F + (1/\bar{p}_l)\mathbf{v}\mathbf{e}_l^\top)$ (cf. the third matrix in (4.5)). Now, from Corollary A.6, we also have the optimality condition

$$(5.4) \quad \mathbf{x}(\text{diag}(e^{\tilde{\gamma}^*})(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)) \circ \mathbf{y}(\text{diag}(e^{\tilde{\gamma}^*})(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top))) = \nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}^*}) \circ e^{\tilde{\gamma}^*} / \mathbf{1}^\top (\nabla\Phi_{\mathbf{w}}(e^{\tilde{\gamma}^*}) \circ e^{\tilde{\gamma}^*}).$$

Using the fact that $\tilde{\gamma}_l^*$ is equal for all l , (5.4) reduces to

$$\mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) = \mathbf{w}.$$

Hence, $\mathbf{w} = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ only if $\tilde{\gamma}_l^*$ is equal for all l .

To prove the convergence of $\tilde{\boldsymbol{\gamma}}(k)$, any initial point $\tilde{\boldsymbol{\gamma}}(0)$ such that $\tilde{\gamma}_l(0)$ are equal for all l yields $\mathbf{m}(k) = \mathbf{w}$ for all k . Hence, the solution of (5.2) is always $-\log \rho(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$. Since the optimality of (3.2) and (3.6) implies

$$(5.5) \quad \text{diag}(e^{\tilde{\boldsymbol{\gamma}}^*})(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)\mathbf{p}^* = \mathbf{p}^* = \rho(\text{diag}(e^{\tilde{\boldsymbol{\gamma}}^*})(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top))\mathbf{p}^*,$$

\mathbf{p}^* can be interpreted as the right eigenvector of $\text{diag}(e^{\tilde{\boldsymbol{\gamma}}^*})(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$. Together with the assumption that $\tilde{\boldsymbol{\gamma}}_l^*$ is equal for all l , this implies that $\mathbf{p}^* = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ (up to a scaling factor). This proves Corollary 5.2. \square

An example with a geometrical illustration of Corollary 5.2 is given in the following.

Example 5.1. We give a simple illustrative example for the two-user case. The channel gains are given by $G_{11} = 0.73$, $G_{12} = 0.04$, $G_{21} = 0.03$, and $G_{22} = 0.89$. The AWGN for the first and second user are 0.1 and 0.3, respectively. The individual maximum power vector $\bar{\mathbf{p}}$ is $(1, 50)^\top$. We then set $\mathbf{w} = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = 1$ in (3.2). The rate of the two users evaluated at the solution of (3.2) given by $\mathbf{p}^* = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ (up to a scaling factor) is then plotted on the achievable rate region (showing that maximizing the minimum rate coincides with the weighted sum

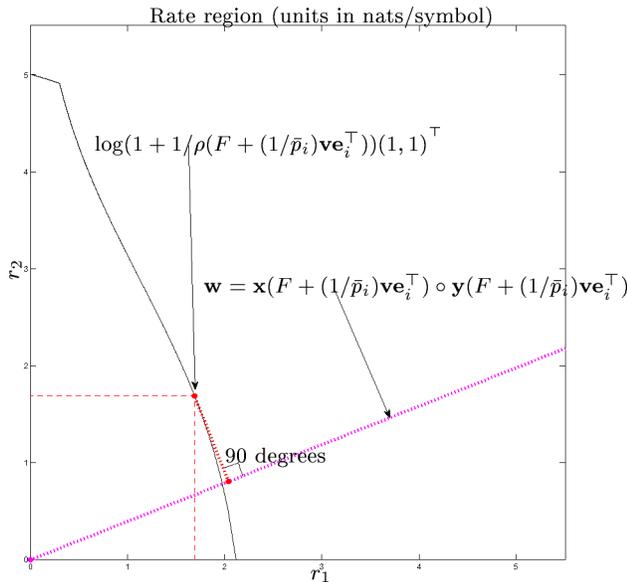


FIG. 5.1. Achievable rate region for a 2-user interference channel. From a geometrical perspective, the weighted sum rate point (with the weight vector $\mathbf{w} = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = 1$, superimposed on the rate region) evaluated at the optimal solution $\boldsymbol{\gamma}^* = -\log \rho(F + (1/\bar{p}_i)\mathbf{v}\mathbf{e}_i^\top)(1, 1)^\top$ finds the largest hypercube that is contained inside the achievable rate region.

rate). Algorithm 2 converges to this point in one iteration starting from any positive initial point such that $\tilde{\boldsymbol{\gamma}}(k) = \mathbf{1}$ (up to a scaling factor). Figure 5.1 illustrates this example that gives the geometrical perspective: the weighted sum rate point (with the weight vector $\mathbf{w} = \mathbf{x}(F + (1/\bar{p}_i)\mathbf{ve}_i^\top) \circ \mathbf{y}(F + (1/\bar{p}_i)\mathbf{ve}_i^\top)$, where $i = 1$, superimposed on the rate region) evaluated at the optimal solution $\boldsymbol{\gamma}^* = -\log \rho(F + (1/\bar{p}_i)\mathbf{ve}_i^\top)(1, 1)^\top$ finds the largest hypercube that is contained inside the achievable rate region.

6. Numerical examples. In this section, we evaluate the performance of our global optimization algorithm, the relaxation techniques, Algorithm 1 and Algorithm 2. In the branch-and-bound technique, we choose the rectangular set with the largest upper bound and use the rule that splits the rectangular set with the longest edge [1]. We use nats per symbol for the sum rate unit.

6.1. Experiment 1 (comparison with Lagrange dual relaxation). We compare the upper bounds obtained by our relaxation techniques and the Lagrange dual relaxation. We consider the two-user example in [16]:

$$(6.1) \quad \max_{\mathbf{0} \leq p_i \leq 2, l=1,2} \frac{1}{2} \log \left(1 + \frac{p_1}{p_2 + 1} \right) + \frac{1}{2} \log \left(1 + \frac{p_2}{p_1 + 1} \right).$$

In [16], the optimal value of (6.1) and its optimal Lagrange dual value are computed explicitly as $(\log 3)/2$ and $(\log 5)/2$, respectively, i.e., a positive duality gap value of $\log(5/3)/2$. The optimal solution \mathbf{p}^* is either $(2, 0)^\top$ or $(0, 2)^\top$ (respectively, $\boldsymbol{\gamma}^*$ is either $(2, e^{-K})^\top$ or $(e^{-K}, 2)^\top$ in (3.4)) because, for the two-user case with $w_1 = w_2$, it suffices to check the extreme points of the feasible set of (6.1) to solve (6.1) [7].

Figure 6.1 illustrates how the first method by convex relaxation computes an upper bound to (6.1). In this example, an upper bound very close to the optimal value of $(\log 3)/2$ within an acceptable accuracy can be obtained after solving two linear programs. An upper bound better than the optimal Lagrange dual value $(\log 5)/2$ is obtained after solving ten linear programs, and it takes another twelve linear programs to certify that $(\log 3)/2$ is the global optimal value. The binary tree in Figure 6.1 has a

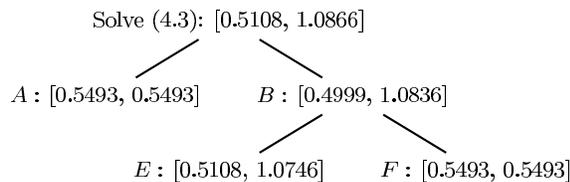


FIG. 6.1. Solving (6.1) using the first relaxation technique: Successive linear program with a branch-and-bound algorithm (the rectangular method in [1]; also see, e.g., Chapter 7 in [13]). We use a rectangular set $[-K, \log(2)]^2$ with $K = 100$. The lower and upper bounds are depicted in brackets next to the subdivided set. At the root of the tree, (4.3) is solved (original rectangle). We then have $L_{bb} = 0.5108$ and $U_{bb} = 1.0866$. In the second iteration, the rectangular set is partitioned into two: A and B (A is the set $\{\tilde{\boldsymbol{\gamma}}_1 \in [-K, \log(2)/2], \tilde{\boldsymbol{\gamma}}_2 \in [-K, \log 2]\}$ and B is the set $\{\tilde{\boldsymbol{\gamma}}_1 \in [\log(2)/2, \log 2], \tilde{\boldsymbol{\gamma}}_2 \in [-K, \log 2]\}$). We then have $L_{bb} = 0.5493$ and $U_{bb} = 1.0836$. At the third iteration, we partition the set B to obtain the bottom leaf children C and D (C is the set $\{\tilde{\boldsymbol{\gamma}}_1 \in [\log(2)/2, \log 2], \tilde{\boldsymbol{\gamma}}_2 \in [\log(2)/2, \log 2]\}$ and D is the set $\{\tilde{\boldsymbol{\gamma}}_1 \in [\log(2)/2, \log 2], \tilde{\boldsymbol{\gamma}}_2 \in [-K, \log(2)/2]\}$). We then have $L_{bb} = 0.5493$ and $U_{bb} = 1.0746$. Observe that a lower and upper bound of 0.5493 within acceptable accuracy to the optimal primal value $(\log 3)/2$ is obtained after solving two linear programs (at the node with set A containing the global optimal solution $(-K, \log 2)^\top$). Note that set F containing the other global optimal solution $(\log 2, -K)^\top$ also yields the same lower and upper bound of 0.5493 within acceptable accuracy after solving three linear programs.

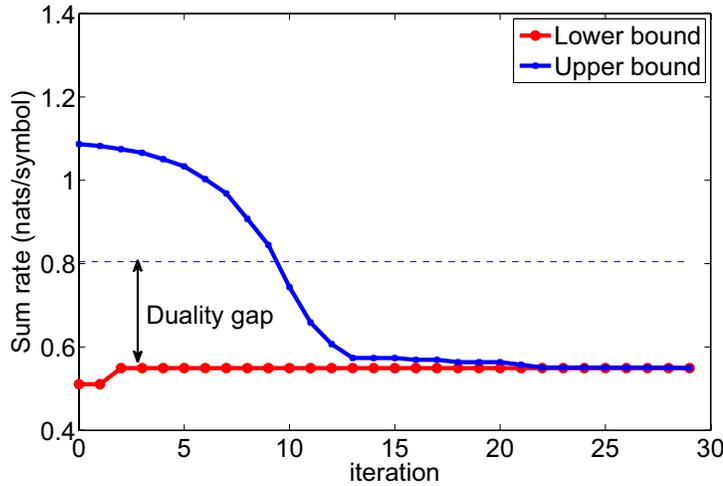


FIG. 6.2. Upper and lower bounds on the sum rate computed by the first relaxation technique that uses successive linear programs and branch-and-bound in Experiment 1.

total of fourteen levels (only the root and the first two levels are shown in Figure 6.1). Figure 6.2 shows the upper and lower bounds computed by the branch-and-bound method, and the branch-and-bound method converges after twenty-two iterations ($\epsilon = 1.5 \times 10^{-3}$).

Next, we apply the second relaxation method that uses Algorithm 1 on (6.1), and we use an initial vector $(-0.172, -1.061)^\top$ (not too close to $\tilde{\gamma}^*$). The optimal value computed by Algorithm 1 when we use: (1) $\tilde{F} = F + \text{diag}(\tilde{\gamma})$, (2) $\tilde{F} = F + (1/\mathbf{1}^\top \tilde{\mathbf{p}})\mathbf{v}\mathbf{1}^\top$, and (3) $\tilde{F} = F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top$ is given by $(\log 3)/2$, $(\log 5)/2$, and $(\log 3)/2$, respectively. The first and third versions of \tilde{F} yield the global optimal primal value (as well as a feasible power vector $(2, 0)^\top$), whereas the second version yields the optimal dual value (with an infeasible power vector $(4, 0)^\top$). This shows that Algorithm 1 can find the global optimal power solution of (3.2) (cf. the necessary condition in the second row of Table 4.1).

Lastly, using the feasible $\tilde{\gamma}$ obtained by the first relaxation method $(\log(2/3), \log(2/3))^\top$ and $(\log 2, -K)^\top$ that yields L_{bb} at the root and first level of the binary tree in Figure 6.1, respectively, as initial points, Algorithm 2 converges to the global optimal solution $(\log 2, -K)^\top$.

6.2. Experiment 2 (convergence of Algorithm 2). The channel gains are given by $G_{11} = 0.73$, $G_{12} = 0.04$, $G_{21} = 0.03$, $G_{22} = 0.89$, and the AWGN for the first and second user are 0.1 and 0.1, respectively. The individual maximum power vector $\tilde{\mathbf{p}}$ is $(1.8, 20.5)^\top$. We then set $\mathbf{w} = \mathbf{x}(F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top) \circ \mathbf{y}(F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top)$, where $i = 1$ in (3.2), and the optimal power is $\mathbf{p}^* = \mathbf{x}(F + (1/\tilde{p}_i)\mathbf{v}\mathbf{e}_i^\top)$ in (3.2) (hence the optimal solution to (3.4) is $8.334(1, 1)^\top$). Using the two different outputs obtained from the branch-and-bound algorithm (at the root level and at the 10th iteration) shown in Figure 6.3(a) as initial points, the convergence of Algorithm 2 is shown in Figure 6.3(b). As observed in Figure 6.3(b), Algorithm 2 converges to the optimal solution with the two different initial points as input after four iterations.

6.3. Experiment 3 (larger examples of branch-and-bound relaxation). Figure 6.4 illustrates the convergence of the branch-and-bound relaxation method

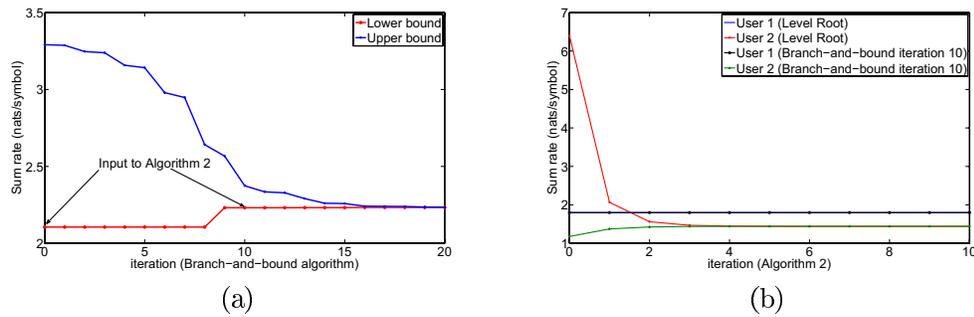


FIG. 6.3. (a) Illustration of the convergence of the branch-and-bound relaxation with the feasible point of the tightest lower bound used as input to Algorithm 2. (b) Illustration of the convergence of Algorithm 2 with an initial point obtained at the root level and the 10th iteration of the branch-and-bound method.

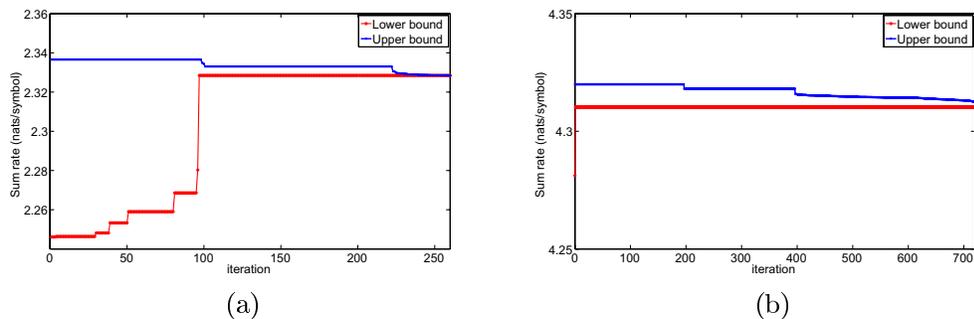


FIG. 6.4. Illustration of the convergence of the branch-and-bound relaxation for (a) 100 users and (b) 200 users in Experiment 3.

for randomly generated problems with a larger number of users. As observed from Figure 6.4, when $L = 100$, the optimal solution can be computed after solving 190 linear programs, but it takes another 310 linear programs to certify that the optimal solution is within a (1×10^{-3}) -suboptimal region. When $L = 200$, the optimal solution can be computed after solving 4 linear programs, but it takes approximately another 1400 linear programs to certify that the optimal solution is within a (1×10^{-3}) -suboptimal region.

7. Conclusion. We studied the nonconvex sum rate maximization problem that finds applications in power control of CDMA wireless networks. Using tools from non-negative matrix theory, in particular the Perron–Frobenius theorem and the Friedland–Karlin inequalities, we showed that this problem can be reformulated as an equivalent convex maximization problem on a closed unbounded convex set. Utilizing the reformulation problem structure, we studied two relaxation techniques to compute progressively tight bounds to the nonconvex problem. One was based on convex relaxation by convex envelope and the other was based on successive convex approximation. We showed that the optimal solution to the sum rate maximization and its relaxed problem can be analytically characterized by the spectra of specially crafted nonnegative matrices. By exploiting the relaxation and approximation techniques, we proposed a global optimization algorithm with ϵ -suboptimality to solve the sum rate maximization problem. We also gave new applications of the Friedland–Karlin inequalities to inverse problems

in nonnegative matrix theory. As future work, we plan to extend our results in this paper to a multiple-frequency channel that has applications in a DSL multiuser system.

Appendix A. Results related to Friedland–Karlin inequalities. In this section, we recall some results from [10] and state the extensions of these results, and then illustrate their applications in this paper. We first state the following extension of [10, Theorem 3.1].

THEOREM A.1. *Let $A \in \mathbb{R}_+^{L \times L}$ be an irreducible nonnegative matrix. Assume that $\mathbf{x}(A) = (x_1(A), \dots, x_L(A))^\top$, $\mathbf{y}(A) = (y_1(A), \dots, y_L(A))^\top > \mathbf{0}$ are left and right Perron–Frobenius eigenvectors of A , normalized such that $\mathbf{x}(A) \circ \mathbf{y}(A)$ is a probability vector. Suppose $\boldsymbol{\gamma}$ is a nonnegative vector. Then,*

$$(A.1) \quad \rho(A) \prod_{l=1}^L \gamma_l^{(\mathbf{x}(A) \circ \mathbf{y}(A))_l} \leq \rho(\text{diag}(\boldsymbol{\gamma})A).$$

If $\boldsymbol{\gamma}$ is a positive vector, then equality holds if and only if all γ_l are equal. Furthermore, for any positive vector $\mathbf{z} = (z_1, \dots, z_L)^\top$, the following inequality holds:

$$(A.2) \quad \rho(A) \leq \prod_{l=1}^L \left(\frac{(A\mathbf{z})_l}{z_l} \right)^{(\mathbf{x}(A) \circ \mathbf{y}(A))_l}.$$

If A is an irreducible nonnegative matrix with positive diagonal elements, then equality holds in (A.2) if and only if $\mathbf{z} = t\mathbf{x}(A)$ for some positive t .

Proof. Theorem 3.1 in [10] makes the following assumptions. First, in (A.1), it assumes that $\boldsymbol{\gamma} > \mathbf{0}$. Second, in (A.2), it assumes that $\rho(A) = 1$. Third, the equality case in (A.2) for $\mathbf{z} > \mathbf{0}$ is stated for a positive matrix A . We now show how to deduce the stronger version of Theorem 3.1 claimed here.

First, by using the continuity argument, we deduce the validity of (A.1) for any $\boldsymbol{\gamma} \geq \mathbf{0}$. Second, by replacing A by tA , where $t > 0$, we deduce that it suffices to show (A.2) in the case $\rho(A) = 1$.

Third, to deduce the equality case in (A.2) for $\mathbf{z} > \mathbf{0}$, we need to examine the proof of Lemma 3.2 in [10]. The proof of the Lemma 3.2 applies if the following condition holds. For any sequence of probability vectors $\mathbf{z}_i = (z_{1,i}, \dots, z_{L,i})^\top, i = 1, \dots$, which converges to a probability vector $\boldsymbol{\zeta} = (\zeta_1, \dots, \zeta_L)^\top$, where $\boldsymbol{\zeta}$ has at least one zero coordinate, the function $\prod_{l=1}^L \left(\frac{(A\mathbf{z}_i)_l}{z_{l,i}} \right)^{(\mathbf{x}(A) \circ \mathbf{y}(A))_l}$ tends to ∞ on the sequence $\mathbf{z}_i, i = 1, \dots$. Assume that $\mathcal{A} = \{l \in \langle L \rangle, \zeta_l = 0\}$. Note that the complement of \mathcal{A} in $\langle L \rangle$, denoted by \mathcal{A}^c , is nonempty.

Since $A = [a_{ij}]$ has positive diagonal entries, it follows that $\frac{(A\mathbf{z}_i)_l}{z_{l,i}} \geq a_{ll} > 0$ for each $l \in \langle L \rangle$. Since A is irreducible, there exist $l \in \mathcal{A}$ and $m \in \mathcal{A}^c$ such that $a_{lm} > 0$. Hence, $\lim_{i \rightarrow \infty} \frac{(A\mathbf{z}_i)_l}{z_{l,i}} = \infty$. This shows that the unboundedness condition holds. \square

The following result gives an interpretation of (A.1) in terms of the supporting hyperplane of the convex function $\log \rho(\text{diag}(e^\boldsymbol{\xi})B)$, where $B \in \mathbb{R}_+^{L \times L}$ is irreducible and $\boldsymbol{\xi} \in \mathbb{R}^L$.

THEOREM A.2. *Let $B \in \mathbb{R}_+^{L \times L}$ be an irreducible nonnegative matrix. Let $\boldsymbol{\eta} = (\eta_1, \dots, \eta_L)^\top \in \mathbb{R}^L$ satisfy the condition $\rho(\text{diag}(e^\boldsymbol{\eta})B) = 1$. Denote $A = \text{diag}(e^\boldsymbol{\eta})B$ and assume that $\mathbf{x}(A) = (x_1(A), \dots, x_L(A))^\top$, $\mathbf{y}(A) = (y_1(A), \dots, y_L(A))^\top > \mathbf{0}$ are left and right Perron–Frobenius eigenvectors of A , normalized such that $\mathbf{x}(A) \circ \mathbf{y}(A)$ is a probability vector. Let*

$$(A.3) \quad H(\xi) = \sum_{l=1}^L x_l(A)y_l(A)(\xi_l - \eta_l).$$

Then, $H(\xi) \leq 0$ is the unique supporting hyperplane to the convex set $\log \rho(\text{diag}(e^\xi)B) \leq 0$ at $\xi = \eta$.

Proof. Let $\xi \in \mathbb{R}^L$. Then, $\text{diag}(e^\xi)B = \text{diag}(e^{\xi-\eta})A$. Theorem A.1 implies that $H(\xi) \leq \log \rho(\text{diag}(e^\xi)B)$. Thus, $H(\xi) \leq 0$ if $\log \rho(\text{diag}(e^\xi)B) \leq 0$. Clearly, $H(\eta) = 0$. Hence, $H(\xi) \leq 0$ is a supporting hyperplane of the convex set $\log \rho(\text{diag}(e^\xi)B) \leq 0$. Since the function $\log \rho(\text{diag}(e^\xi)B)$ is a smooth function of ξ , it follows that $H(\xi) \leq 0$ is unique. \square

We now give an application of (A.2) in Theorem A.1.

THEOREM A.3. *Let $B \in \mathbb{R}_+^{L \times L}$ be an irreducible nonnegative matrix. Let $\eta = (\eta_1, \dots, \eta_L)^\top \in \mathbb{R}^L$. Let $\mathbf{m} = \mathbf{x}(\text{diag}(e^\eta)B) \circ \mathbf{y}(\text{diag}(e^\eta)B) = (m_1, \dots, m_L)^\top$ be a probability vector. Then, for any positive vector $\mathbf{z} = (z_1, \dots, z_L)^\top$,*

$$(A.4) \quad \sum_{l=1}^L m_l \log \frac{z_l}{(B\mathbf{z})_l} \leq -\log \rho(\text{diag}(e^\eta)B) + \sum_{l=1}^L m_l \eta_l.$$

If B has a positive diagonal, then equality holds if and only if $\mathbf{z} = t\mathbf{x}(\text{diag}(e^\eta)B)$ for some $t > 0$.

Proof. Let $A = \text{diag}(e^\eta)B$. Then,

$$\sum_{l=1}^L m_l \log \frac{(B\mathbf{z})_l}{z_l} = \sum_{l=1}^L m_l \log \frac{(A\mathbf{z})_l}{z_l} - \sum_{l=1}^L m_l \eta_l.$$

Use (A.2) to deduce (A.4). The equality case follows from the equality case in (A.2). \square

We now turn to applying Theorem A.1 to solve the following inverse problem.

Problem A.4. Let $B \in \mathbb{R}_+^{L \times L}$, $\mathbf{m} \in \mathbb{R}_+^L$ be a given irreducible nonnegative matrix and a positive probability vector, respectively. When does there exist $\eta \in \mathbb{R}^L$ such that $\mathbf{x}(\text{diag}(e^\eta)B) \circ \mathbf{y}(\text{diag}(e^\eta)B) = \mathbf{m}$? If such η exists, when is it unique up to an addition $t\mathbf{1}$?

To solve the inverse problem, we recall Theorem 3.2 in [10] (a consequence of Theorem A.1, i.e., Theorem 3.1 in [10]) that is reproduced in the following.

THEOREM A.5. *Let $A \in \mathbb{R}_+^{L \times L}$, $\mathbf{u}, \mathbf{v} \in \mathbb{R}_+^L$ be given, where A is irreducible with positive diagonal elements and \mathbf{u}, \mathbf{v} are positive. Then, there exists $D_1, D_2 \in \mathbb{R}_+^{L \times L}$ such that*

$$(A.5) \quad D_1 A D_2 \mathbf{u} = \mathbf{u}, \quad \mathbf{v}^\top D_1 A D_2 = \mathbf{v}^\top, \quad D_1 = \text{diag}(\mathbf{f}), \quad D_2 = \text{diag}(\mathbf{g}), \quad \text{and} \quad \mathbf{f}, \mathbf{g} > \mathbf{0}.$$

The pair (D_1, D_2) is unique to the change $(tD_1, t^{-1}D_2)$ for any $t > 0$. There exist $\eta \in \mathbb{R}^L$ such that $\mathbf{x}(\text{diag}(e^\eta)B) \circ \mathbf{y}(\text{diag}(e^\eta)B) = \mathbf{m}$. Furthermore, η is unique up to an addition $t\mathbf{1}$.

COROLLARY A.6. *Let $B \in \mathbb{R}_+^{L \times L}$, $\mathbf{m} \in \mathbb{R}_+^L$ be a given irreducible nonnegative matrix with positive diagonal elements and a positive probability vector, respectively. Then, there exists $\eta \in \mathbb{R}^L$ such that $\mathbf{x}(\text{diag}(e^\eta)B) \circ \mathbf{y}(\text{diag}(e^\eta)B) = \mathbf{m}$. Furthermore, η is unique up to an addition of $t\mathbf{1}$. In particular, this η can be computed by solving the following convex optimization problem:*

$$\begin{aligned}
 & \text{maximize } \mathbf{m}^\top \boldsymbol{\eta} \\
 & \text{subject to } \log \rho(\text{diag}(e^\boldsymbol{\eta})B) \leq 0, \\
 \text{(A.6)} \quad & \text{variables: } \boldsymbol{\eta} = (\eta_1, \dots, \eta_L)^\top \in \mathbb{R}^L.
 \end{aligned}$$

Proof. Let $\mathbf{u} = \mathbf{1}$, $\mathbf{v} = \mathbf{m}$. Then, from Theorem A.5, there exists D_1, D_2 two diagonal matrices with positive diagonal entries such that $D_1BD_2\mathbf{1} = \mathbf{1}$, $\mathbf{m}^\top D_1BD_2 = \mathbf{m}^\top$. Consider the matrix $D_2D_1B = D_2(D_1BD_2)D_2^{-1}$. It is straightforward to see that $\mathbf{x}(D_2D_1B) \circ \mathbf{y}(D_2D_1B) = \mathbf{m}$. Hence, $\boldsymbol{\eta}$ is the unique solution of $\text{diag}(e^\boldsymbol{\eta}) = D_2D_1$.

Assume that $\boldsymbol{\zeta} \in \mathbb{R}^L$ satisfies $\mathbf{x}(\text{diag}(e^{\boldsymbol{\zeta}})B) \circ \mathbf{y}(\text{diag}(e^{\boldsymbol{\zeta}})B) = \mathbf{m}$. By considering $\tilde{\boldsymbol{\zeta}} = \boldsymbol{\zeta} + t\mathbf{1}$, we may assume that $\rho(\text{diag}(e^{\boldsymbol{\zeta}})B) = 1$. Let $D_4 = \text{diag}(\mathbf{x}(\text{diag}(e^{\boldsymbol{\zeta}})B))$. Then, $(D_4^{-1}\text{diag}(e^{\boldsymbol{\zeta}})BD_4)\mathbf{1} = \mathbf{1}$. Let $D_3 = D_4^{-1}\text{diag}(e^{\boldsymbol{\zeta}})$. Hence, $\mathbf{y}(D_3BD_4) = \mathbf{m}$. In view of Theorem A.5, $\text{diag}(e^{\boldsymbol{\zeta}}) = D_4D_3 = D_2D_1 = \text{diag}(e^\boldsymbol{\eta})$.

Next, we show that (A.6) computes the required $\boldsymbol{\eta}$. Since (A.6) is convex, we apply the Karush–Kuhn–Tucker conditions to (A.6). The stationarity of the Lagrangian yields $\mathbf{x}(\text{diag}(e^\boldsymbol{\eta})B) \circ \mathbf{y}(\text{diag}(e^\boldsymbol{\eta})B) = \mathbf{m}$, thus proving the corollary. \square

Remark 5. The convex optimization problem in (A.6) can be reformulated as an equivalent geometric program by utilizing a basic result in nonnegative matrix theory, which states that the Perron–Frobenius eigenvalue is given by $\rho(A) = \inf\{\lambda | A\mathbf{v} \leq \lambda\mathbf{v} \text{ for some positive } \mathbf{v}\}$ and that the infimum is achieved; e.g., see Chapter 4 of [3]. The geometric program can then be solved numerically using an interior-point solver (see, for example, the cvx software package [12]), whose solution can be transformed back to that of (A.6).

We illustrate the necessity of the irreducible nonnegative matrix having positive diagonal elements in Corollary A.6 by the following example.

Example A.7. Let us look at the matrix F defined in (2.2) having zero diagonal entries and positive off-diagonal entries. For $L = 2$, it is easy to show that $\mathbf{x}(F) \circ \mathbf{y}(F) = (\frac{1}{2}, \frac{1}{2})^\top$. In particular, for $L = 2$, Problem A.4 is not solvable for $\mathbf{m} \neq (\frac{1}{2}, \frac{1}{2})^\top$. Similarly, given positive $\mathbf{u}, \mathbf{v} \in \mathbb{R}^2$ such that $\mathbf{u} \circ \mathbf{v} \neq t(1, 1)$ for any positive t , (A.5) does not hold for $A = F$. For $L \geq 3$, the situation is different, and is illustrated in the following result.

THEOREM A.8. *Let $L \geq 3$, $A \in \mathbb{R}_+^{L \times L}$, $\mathbf{u} = (u_1, \dots, u_L)^\top$, $\mathbf{v} = (v_1, \dots, v_L)^\top \in \mathbb{R}_+^L$ be given, where A is a matrix with zero diagonal entries and positive off-diagonal elements, and \mathbf{u}, \mathbf{v} are positive. Assume that $\mathbf{m} = \mathbf{u} \circ \mathbf{v}$ is a probability vector satisfying the condition*

$$\text{(A.7)} \quad \sum_{j \neq l} m_j > m_l \quad \forall l \in \langle L \rangle.$$

Then, there exists $D_1, D_2 \in \mathbb{R}_+^{L \times L}$ such that (A.5) holds.

Proof. Let $A_i = A + (1/i)I$, $i = 1, \dots$, where I is the $L \times L$ identity matrix. Theorem A.5 implies

$$\begin{aligned}
 B_i &= D_{1,i}A_iD_{2,i}, \quad D_{1,i} = \text{diag}(\mathbf{f}_i), \quad D_{2,i} = \text{diag}(\mathbf{g}_i), \quad B_i\mathbf{u} = \mathbf{u}, \quad \mathbf{v}^\top B_i = \mathbf{v}^\top, \\
 \mathbf{f}_i &= (f_{1,i}, \dots, f_{L,i})^\top, \quad \mathbf{g}_i = (g_{1,i}, \dots, g_{L,i})^\top, \quad s_i = \max_{j \in \langle L \rangle} f_{j,i} = \max_{j \in \langle L \rangle} g_{j,i}, \quad i = 1, \dots
 \end{aligned}$$

Note that each entry of B_i is bounded by $\frac{\max_j u_j}{\min_j u_j}$. By passing to the subsequence $B_{i_k}, \mathbf{f}_{i_k}, \mathbf{g}_{i_k}$, $1 \leq i_1 < i_2 < \dots$, we can assume that the first subsequence converges to B , and the last two subsequences converge in a generalized sense:

$$\begin{aligned} \lim_{k \rightarrow \infty} B_{i_k} &= B = [b_{jl}] \in \mathbb{R}_+^{L \times L}, & \lim_{k \rightarrow \infty} \mathbf{f}_{i_k} &= \mathbf{f} = (f_1, \dots, f_L)^\top, \\ \lim_{k \rightarrow \infty} \mathbf{g}_{i_j} &= \mathbf{g} = (g_1, \dots, g_L)^\top, \\ f_j, g_j &\in [0, \infty], & j &= 1, \dots, L, \\ \lim_{k \rightarrow \infty} s_{i_k} &= s = \max_{j \in \langle L \rangle} f_j = \max_{j \in \langle L \rangle} g_j \in [0, \infty]. \end{aligned}$$

Note that

$$(A.8) \quad B\mathbf{u} = \mathbf{u}, \quad \mathbf{v}^\top B = \mathbf{v}^\top.$$

Assume first that $s < \infty$. Then, $B = \text{diag}(\mathbf{f})A \text{diag}(\mathbf{g})$. In view of (A.8), $\mathbf{f} \circ \mathbf{g} > \mathbf{0}$. This proves the theorem in this case.

Assume now that $s = \infty$. Let

$$\begin{aligned} \mathcal{F}_\infty &= \{j \in \langle L \rangle, f_j = \infty\}, & \mathcal{F}_+ &= \{j \in \langle L \rangle, f_j \in (0, \infty)\}, & \mathcal{F}_0 &= \{j \in \langle L \rangle, f_j = 0\}, \\ \mathcal{G}_\infty &= \{j \in \langle L \rangle, g_j = \infty\}, & \mathcal{G}_+ &= \{j \in \langle L \rangle, g_j \in (0, \infty)\}, & \mathcal{G}_0 &= \{j \in \langle L \rangle, g_j = 0\}. \end{aligned}$$

Since the off-diagonal entries of A are positive, and $B \in \mathbb{R}_+^{L \times L}$, it follows that $\mathcal{F}_\infty = \mathcal{G}_\infty = \{l\}$ for some $l \in \langle L \rangle$. Furthermore, $\mathcal{F}_+ = \mathcal{G}_+ = \emptyset$. So $\mathcal{F}_0 = \mathcal{G}_0 = \langle L \rangle \setminus \{l\}$. Assume first that $l = 1$. Then, the principal submatrix $[b_{jl}]_{j=l=2}^L$ is zero, and (A.8) yields that

$$b_{j1} = \frac{u_j}{u_1}, \quad b_{1j} = \frac{v_j}{v_1} \quad \text{for } j = 2, \dots, L, \quad b_{11}u_1v_1 + \sum_{j=2}^L u_jv_j = u_1v_1.$$

Since $b_{11} \geq 0$, the above last equality contradicts the condition (A.7) for $l = 1$. Similar argument implies the impossibility of $\mathcal{F}_\infty = \mathcal{G}_\infty = \{l\}$ for any $l \geq 2$. Hence, $s < \infty$ and we conclude the theorem. \square

We do not know whether, under the conditions of Theorem A.8, the diagonal matrices (D_1, D_2) are unique up to the transformation $(tD_1, t^{-1}D_2)$. We now generalize the above theorem.

THEOREM A.9. *Let*

$$L \geq 2, \quad A = [a_{jl}]_{j=l=1}^L \in \mathbb{R}_+^{L \times L}, \quad \mathbf{0} < \mathbf{u} = (u_1, \dots, u_L)^\top, \quad \mathbf{v} = (v_1, \dots, v_L)^\top \in \mathbb{R}_+^L$$

be given. Assume that A has positive off-diagonal elements, and $\mathbf{m} = \mathbf{u} \circ \mathbf{v}$ is a probability vector satisfying the condition

$$(A.9) \quad \sum_{\forall j \neq l} m_j > m_l$$

for each l such that $a_{ll} = 0$. Then, there exists $D_1, D_2 \in \mathbb{R}_+^{L \times L}$ such that (A.5) holds.

Proof. Assume first that $L \geq 3$. In view of Theorems A.5 and A.8, it suffices to assume that A has positive and zero diagonal entries. We then apply the proof of Theorem A.8 and the following observation: If $\mathcal{F}_\infty = \mathcal{G}_\infty = \{l\}$, then $a_{ll} = 0$.

Assume now that $L = 2$. Note that if A has a zero diagonal, then the condition (A.7) cannot hold. Assume now that A has at least one positive diagonal element. Then, the above arguments for $L \geq 3$ apply. \square

Using Theorem A.9, we obtain the following corollary that is used to prove Theorem 4.3 in section 4.

COROLLARY A.10. Let $B = [b_{jl}]_{j,l=1}^L \in \mathbb{R}_+^{L \times L}$, $\mathbf{m} \in \mathbb{R}_+^L$ be a given matrix with positive off-diagonal elements and a positive probability vector, respectively. Assume that $L \geq 2$ and \mathbf{m} satisfies the conditions (A.9) for each l such that $b_{ll} = 0$. Then, there exists $\boldsymbol{\eta} \in \mathbb{R}^L$ such that $\mathbf{x}(\text{diag}(e^\boldsymbol{\eta})B) \circ \mathbf{y}(\text{diag}(e^\boldsymbol{\eta})B) = \mathbf{m}$.

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