

HYBRID ANALOG-DIGITAL TRANSMITTER WITH PASSIVE ANTENNA ARRAY

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I. INTRODUCTION

Millimeter-wave communication is one of the technologies that enable the high-bandwidth communication. However, due to the small wave length, antenna arrays are often required to compensate for the high loss, which will increase the hardware complexity. In article [1], two methods are investigated, low-resolution ADCs and hybrid analog-digital (HAD) precoding. Receivers equipped low-resolution ADCs have low power consumptions but require more RF chains. On the contrary, systems employing HAD architecture requires less RF chains than the antenna amount, but the performance is limited by the number of RF chains. Low complexity algorithms have also been developed for the HAD systems such as SIC-based hybrid precoding [2]. This precoding method avoid the SVD operations and is more energy efficient than the fully digital systems.

Intelligent reflecting surface (IRS) is a newly proposed reconfigurable communication component [3]. It contains a large number of passive antenna elements whose amplitude/phase is adjustable. Typical applications of IRS includes creating a virtual line-of-sight, protecting the communication from eavesdropper and so on [4]. When the IRS used as a relay, some alternating algorithm can be used with semidefinite relaxation to offer a trade-off between performance and computational complexity [5]. It is also shown in [5] that the use of IRS is more cost efficient than conventional amplify-and-forward relay.

In this article, the IRS is utilized as the analog unit which is placed close to the RF chains in a HAD transmitter. A symbol-level system with this architecture was introduced in [6], where only a single RF chain is included. In that system, the information modulation and precoding

happen at the IRS side. Two designing methods for symbol level point-to-point multiple-input multiple-output (MIMO) precoding in this IRS based transmitter is introduced in [7]. It is shown in this article that although conventional MIMO transmitters such as fully digital ones and fully connected HAD transmitters has better spectral efficiency, their power consumptions are much higher than IRS aided transmitters. A symbol wise precoding scheme is also proposed by [8], which transforms the precoding task into a generalized least-square-error (GLSE) problem to have some control over the power consumptions and solves the problem by iterative optimization.

The above mentioned precoding methods for IRS aided transmitters are all symbol-level precoding. In [9], it is suggested that a block level precoding method can be used for a fully connected HAD transmitter without causing any distortion as long as the number of RF chains is larger than the block length. Moreover, by using block level precoding, the transmitter avoids the symbol by symbol optimization which will reduce the computational loads [10].

This article investigates the precoding scheme in IRS aided HAD systems. The precoding task cannot be formulated as a convex problem. Hence, alternating optimization method is used to optimize the digital unit outputs and IRS configurations. In the meantime, we find out that the optimization of the digital unit output is a convex problem, while the optimization of the IRS configurations is non-convex. We use gradient descent and majorize-minimization (MM) algorithm to reconfigure the IRS. Since the passive antenna elements on the IRS are relative cheap, this article also studies the relation between IRS size and block length.

II. PROBLEM FORMULATION

A. System Model

For down-link transmissions in a multi-user MIMO system consisting of a base station (BS) with M antenna elements and K single antenna user terminals (UTs), the vector of the received signals in this network in the i -th symbol time instance drawn from a block is given by

$$\mathbf{y}_i = \mathbf{H}\mathbf{a}_i + \mathbf{n}_i, \quad (1)$$

where $\mathbf{H} \in \mathbb{C}^{K \times M}$ is the wireless channel matrix from base station to all the UTs if the channel state information (CSI) is known. The received vector is written as

$$\mathbf{y}_i = \begin{bmatrix} y_1 & \dots & y_K \end{bmatrix}^T, \quad (2)$$

where y_k denotes the signal at the k -th UT. Vector $\mathbf{a}_i \in \mathbb{C}^{M \times 1}$ represents the signals transmitted by the BS and \mathbf{n}_i is a K by one noise vector whose entries are i.i.d. drawn from Gaussian distribution with zero mean and variance σ_n^2 .

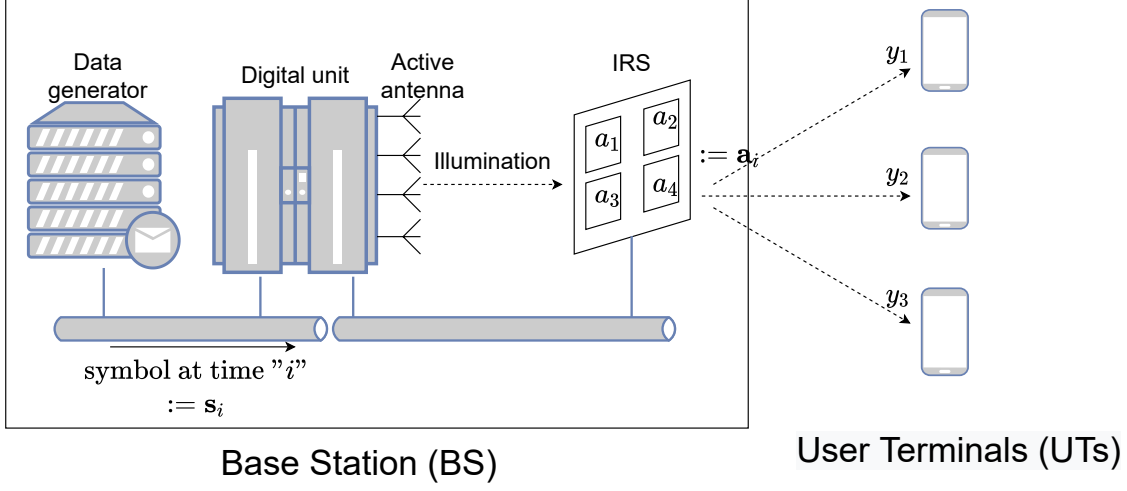


Fig. 1: System Model

The BS employs an IRS-aided HAD transmitter whose architecture is demonstrated in Fig. (1).

Define s_k as the message symbol intended for the k -th UT, and the message vector for all the K UTs at the i -th time interval can be denoted as

$$\mathbf{s}_i = \begin{bmatrix} s_1 & \dots & s_K \end{bmatrix}^T. \quad (3)$$

In this paper, block-wise precoding with block size L is used. This can be achieved by attaching a buffer to the data generator. The message for one entire block is represented as

$$\mathbf{S} = \begin{bmatrix} \mathbf{s}_1 & \dots & \mathbf{s}_L \end{bmatrix}. \quad (4)$$

Similarly, the transmitted signal and the received signal in one block can be modeled respectively as follow,

$$\begin{aligned} \mathbf{A} &= \begin{bmatrix} \mathbf{a}_1 & \dots & \mathbf{a}_L \end{bmatrix} \\ \mathbf{Y} &= \begin{bmatrix} \mathbf{y}_1 & \dots & \mathbf{y}_L \end{bmatrix}. \end{aligned} \quad (5)$$

Therefore, the block-wise system model is described by

$$\mathbf{Y} = \mathbf{H}\mathbf{A} + \mathbf{N}, \quad (6)$$

where \mathbf{N} is a $K \times N$ random matrix standing for the noise. Next, we are going to discuss the transmitter output matrix \mathbf{A} and channel matrix \mathbf{H} as well as the decoder in more detail.

B. Transmitter Architecture

An illustration for the internal structure of the BS with N RF chains and M passive IRS elements is shown as Fig. (2). If the matrix \mathbf{S} contains the intended messages for the UTs

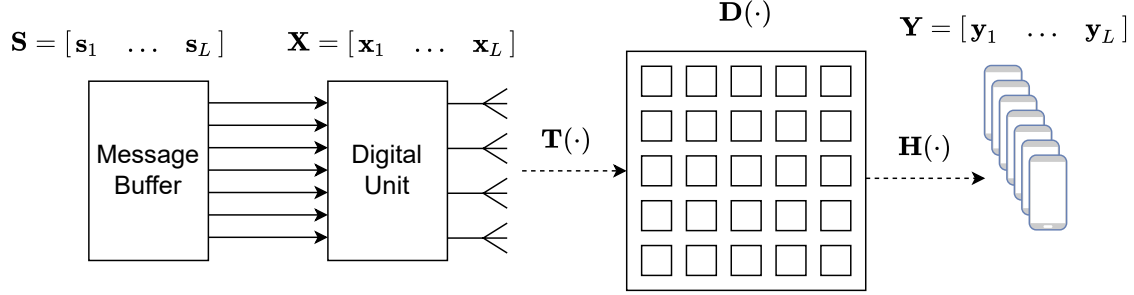


Fig. 2: Signal diagram of the Base Station

during one transmission time block, we can define a matrix $\mathbf{X} \in \mathbb{C}^{N \times L_X}$ to be the RF chain output signals corresponding to that message matrix. Additionally, the phase shifts of the IRS for this message block are contained by $\mathbf{D} = \text{diag}\{e^{j\beta_1}, \dots, e^{j\beta_M}\}$. The IRS is placed in the close proximity of the active antennas, and the illumination channel from the RF chains to the IRS can be considered as fixed during the entire transmission. This fixed channel is denoted as a constant matrix $\mathbf{T} \in \mathbb{C}^{M \times N}$.

Since the UTs are disjoint and don't possess the same computational power as the base station, they can only employ a simple disjoint decoder. The disjoint decoder can be modeled as a diagonal matrix \mathbf{F}_D , i.e.

$$\tilde{\mathbf{Y}} = \mathbf{F}_D \mathbf{Y} = \mathbf{S}. \quad (7)$$

By considering the internal structure of the transmitter, the matrix for the received signals \mathbf{Y} can also be represented as

$$\mathbf{Y} = \mathbf{H} \mathbf{D} \mathbf{T} \mathbf{X} + \mathbf{N}. \quad (8)$$

To make the decoded signal as close as possible to the intended message, \mathbf{Y} and \mathbf{S} must have the same dimension. Therefore, the matrix \mathbf{X} must have as many columns as matrix \mathbf{S} . Thus, we have $L_X = L$ and $\mathbf{X} \in \mathbb{C}^{N \times L}$.

In block-wise precoding, we apply the same phase shifts matrix \mathbf{D} to all the digital-domain vectors contained in the same time block. In this way, the active antennas are updated once every symbol time interval, but the IRS only updates once every time block.

This transmission procedure can be illustrated in Fig. (3). In Fig. (3), the table entries with

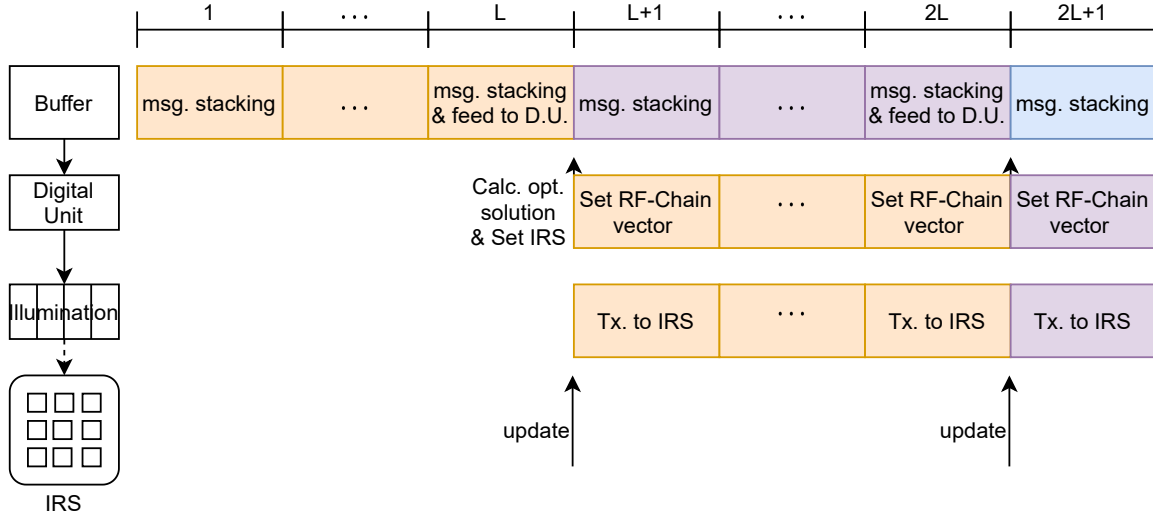


Fig. 3: Transmitter Time Table

the same color indicate that they all correspond to the message symbols of the same block. For example, the RF chain output vectors from time interval $L + 1$ to $2L$ are computed based on the buffered message vectors generated from time interval 1 to L .

If the message vectors are generated at the rate R then the symbol time interval is $T_s = 1/R$. It can be observed that the RF chain output also has the same update rate but with a delay of LT_s relative to the corresponding messages. Compared to symbol-wise transmission, the IRS is reconfigured once per time block, so its update rate is only R/L .

C. Channel Model

According to Eq. (1), the wireless channel from the BS to the k -th UT can be modeled as Eq. (9).

$$y_k = \mathbf{h}_k^T \mathbf{a} + n_k \quad (9)$$

The row vector \mathbf{h}_k^T models the channel coefficients from the IRS to the k -th UT. If we take scattering and antenna spatial correlations into consideration, the channel coefficients can be further modeled in a more detailed form as [11]

$$\mathbf{h}_k^T = \frac{1}{\sqrt{P}} \sum_{p=1}^P h_p \mathbf{h}_t^T(\theta_p^t, \phi_p^t), \quad (10)$$

where $P \in \mathbb{N}$ is the number of effective channel paths, while $h_p \in \mathbb{C}$ is the coefficient that models the path loss and fading of the p -th path. The vector $\mathbf{h}_t(\theta_p^t, \phi_p^t) \in \mathbb{C}^{M \times 1}$ is the steering vector which models the phase shifts of the IRS elements relative to the reference element. The directed angles $\theta_p^t \in [0, \pi]$ and $\phi_p^t \in [0, \pi]$ are the azimuthal and polar angle of the departure beam respectively. Fig. (4) describes this spherical coordinate system closely. The joint points in Fig. (4) represent the IRS reflecting elements.

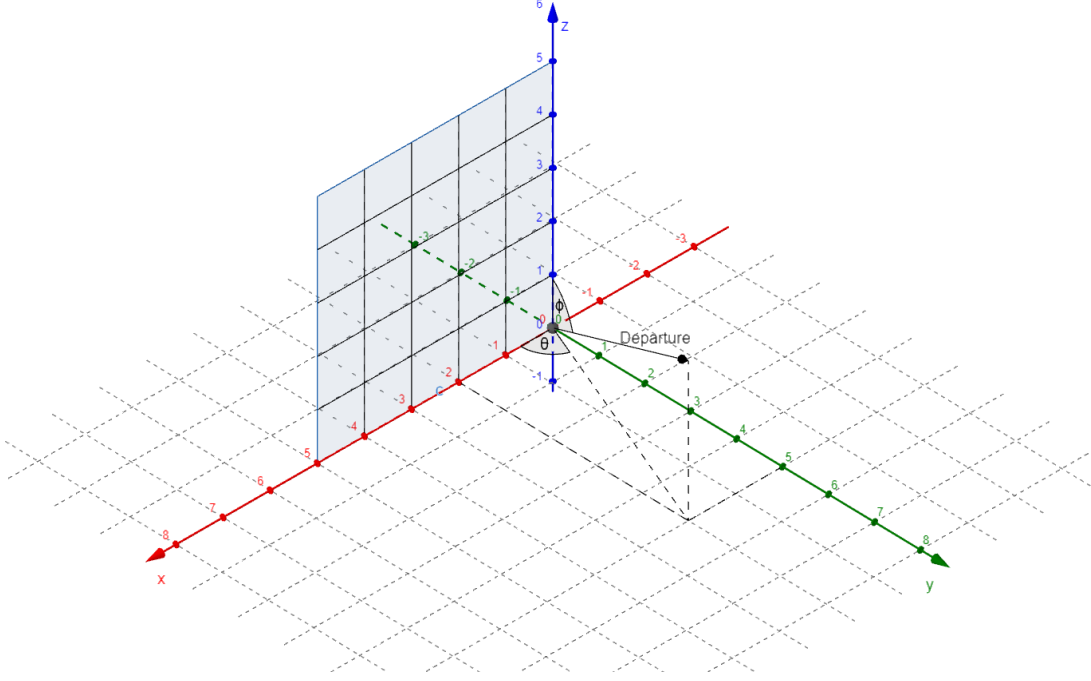


Fig. 4: Angle of departure

W.l.o.g., in the following context, the IRS is placed in the first quadrant of the xz -plane. We also assume that \sqrt{M} is a positive integer. Moreover, we assume that the distances between any two neighboring IRS elements are constant and are denoted as d . The IRS elements are indexed from right to left (along x -axis, 0 to $+\infty$), bottom to top (along z -axis, 0 to $+\infty$). Based on these conventions, if $m \in \mathbb{N} \cap [1, M]$, the m -th IRS element will have the coordinate

$$\begin{bmatrix} x = ((m-1) \bmod \sqrt{M}) \cdot d, \\ y = 0, \\ z = \lfloor (m-1)/\sqrt{M} \rfloor \cdot d \end{bmatrix}. \quad (11)$$

The mod operator in the above equation is the modulo operation which is evaluated to the remainder of the division of $(m-1)$ by \sqrt{M} . The lower half bracket $\lfloor \cdot \rfloor$ is the floor operator

that evaluates to the largest interger that is smaller or equal to the number inside it.

The steering vector $\mathbf{h}_t^T(\theta_p^t, \phi_p^t)$ is described by

$$\mathbf{h}_t(\theta_p^t, \phi_p^t) = \left[e^{j \frac{2\pi d}{\lambda_w} [((m-1) \bmod \sqrt{M}) \sin(\phi_p^t) \cos(\theta_p^t) + \lfloor (m-1)/\sqrt{M} \rfloor \cos(\phi_p^t)]} \right]_{m,1}, \quad (12)$$

where λ_w denotes the wavelength of the transmitted signal.

In Eq. (10), the path loss and the fading is included in the coefficient h_p . According to [11], h_p is modeled as $h_p = \sqrt{\bar{h}_k} \tilde{h}_p$, where \bar{h}_k is the path loss of the k -th UT and \tilde{h}_p is the random fading of the p -th path. Their mathematical descriptions are given by

$$\begin{aligned} \bar{h}_k &= \left(\frac{\lambda_w}{4\pi l_{\text{ref}}} \right)^n \\ \tilde{h}_p &\sim \text{CN}(0, \sigma_f^2). \end{aligned} \quad (13)$$

The variable $n \in \mathbb{R}$ is the path loss exponent and it is greater or equal than two.

D. Precoding via Minimizing RSS

The residual sum of squares (RSS) is adopted here to assess the distortion between the decoded signals and the intended messages. It is defined as

$$d(\mathbf{S}, \tilde{\mathbf{Y}}) = \|\tilde{\mathbf{Y}} - \mathbf{S}\|_{\text{F}}^2 = \text{tr}\{(\tilde{\mathbf{Y}} - \mathbf{S})^H(\tilde{\mathbf{Y}} - \mathbf{S})\}, \quad (14)$$

where $\tilde{\mathbf{Y}} \in \mathbb{C}^{K \times L}$ is the decoded signals of the original received signals \mathbf{Y} in one transmission block. The ideal value of this metric is zero. It is assumed in this article that the distance between BS and UTs are known by UTs. To keep the transmission power relative low, the decoding matrix \mathbf{F}_D is designed to compensate the path loss.

Define $\mathbf{P}_L^{1/2} \in \mathbb{C}^{K \times K}$ to be a diagonal path loss matrix

$$\mathbf{P}_L^{1/2} = \begin{bmatrix} \sqrt{\bar{h}_1} & & \\ & \ddots & \\ & & \sqrt{\bar{h}_K} \end{bmatrix}. \quad (15)$$

The decoder/compensator can then be designed as

$$\mathbf{F}_D = \mathbf{P}_L^{-1/2}. \quad (16)$$

Thus, the compensated received signal is

$$\tilde{\mathbf{Y}} = \mathbf{P}_L^{-1/2} \mathbf{Y}. \quad (17)$$

E. Statistics of the Performance Measure

Most of times, the channel estimation is imperfect. The relation between real channel and estimated channel can be described by the following equation,

$$\mathbf{H} = \hat{\mathbf{H}} + \mathbf{E}, \quad (18)$$

where $\mathbf{H} \in \mathbb{C}^{K \times M}$ stands for the real channel, $\hat{\mathbf{H}} \in \mathbb{C}^{K \times M}$ is the estimated channel and their difference is modeled as a K by M complex random matrix \mathbf{E} . We assume here that each entry of the error matrix are independently and user-wise/row-wise-identically distributed following the two statistics in Eq. (19).

$$\begin{aligned} \mathcal{E}\{e_{k,m}\} &= 0; \\ \mathcal{E}\{e_{k_1,m_1}^* e_{k_2,m_2}\} &= \sigma_{e,k_1}^2 \delta(k_1 - k_2) \delta(m_1 - m_2). \end{aligned} \quad (19)$$

The random variable $e_{k,m} \sim \text{CN}(0, \sigma_{e,k}^2)$ denotes the entry at the k -th row and the m -th column of matrix \mathbf{E} . The $\delta(\cdot)$ is the discrete Dirac function which is evaluated to zero when it has non-zero argument and evaluate to one when its argument is zero.

Taking the estimation error into consideration, the received signals and the decoded/compensated signals can then be written as

$$\begin{aligned} \mathbf{Y} &= (\hat{\mathbf{H}} + \mathbf{E})\mathbf{D}\mathbf{T}\mathbf{X} + \mathbf{N} \\ \tilde{\mathbf{Y}} &= \mathbf{P}_L^{-1/2}[(\hat{\mathbf{H}} + \mathbf{E})\mathbf{D}\mathbf{T}\mathbf{X} + \mathbf{N}]. \end{aligned} \quad (20)$$

To keep the further discussions brief, define the normalized matrices $\tilde{\mathbf{H}}$, $\tilde{\hat{\mathbf{H}}}$, $\tilde{\mathbf{E}}$ and $\tilde{\mathbf{N}}$ as Eq. (21).

$$\begin{aligned} \mathbf{H} &= \mathbf{P}_L^{1/2} \tilde{\mathbf{H}} \\ \hat{\mathbf{H}} &= \mathbf{P}_L^{1/2} \tilde{\hat{\mathbf{H}}} \\ \mathbf{E} &= \mathbf{P}_L^{1/2} \tilde{\mathbf{E}} \\ \mathbf{N} &= \mathbf{P}_L^{1/2} \tilde{\mathbf{N}} \end{aligned} \quad (21)$$

Because of the existence of random terms, statistics of Eq. (14) are used for the designing. Assume that the noise entries are independent with the estimation errors, we then have

$$\begin{aligned} &\mathcal{E}_{\mathbf{E}} \left\{ \mathcal{E}_{\mathbf{N}} \{d(\mathbf{S}, \tilde{\mathbf{Y}})\} \right\} \\ &= \left\| \tilde{\hat{\mathbf{H}}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^H \mathbf{T}^H \mathbf{D}^H \Sigma_{\tilde{\mathbf{E}}} \mathbf{D}\mathbf{T}\mathbf{X}\} + L \text{tr}\{\mathbf{P}_L^{-1}\} \sigma_n^2, \end{aligned} \quad (22)$$

where

$$\Sigma_{\tilde{\mathbf{E}}} = \begin{bmatrix} \sum_{k=1}^K \sigma_{e,k}^2 / \bar{h}_k & & \\ & \ddots & \\ & & \sum_{k=1}^K \sigma_{e,k}^2 / \bar{h}_k \end{bmatrix}. \quad (23)$$

The formula given in Eq. (22) is the expected distance between \mathbf{S} and $\tilde{\mathbf{Y}}$, i.e. $\mathcal{E}_{\mathbf{N}, \mathbf{E}}\{d(\mathbf{S}, \tilde{\mathbf{Y}})\}$. From now on, instead of developing an algorithm that minimize Eq. (14) directly, we try to develop an algorithm that minimize Eq. (22) by optimizing \mathbf{D} and \mathbf{X} .

F. Problem Formulation as a GLSE Precoder

Following [12], if we utilize the peak and total power limitation, the minimization of the expected distance between the intended message and the received symbols leads us to the design of a GLSE precoder. The optimization problem is given by

$$\begin{aligned} \underset{\mathbf{D}, \mathbf{X}}{\text{minimize}} \quad & \left\| \tilde{\mathbf{H}} \mathbf{D} \mathbf{T} \mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^{\text{H}} \mathbf{T}^{\text{H}} \mathbf{D}^{\text{H}} \Sigma_{\tilde{\mathbf{E}}} \mathbf{D} \mathbf{T} \mathbf{X}\} + \lambda \text{tr}\{\mathbf{X} \mathbf{X}^{\text{H}}\} \\ \text{subject to} \quad & \max_{i,j} |x_{i,j}|^2 < P_{\text{peak}}, \end{aligned} \quad (24)$$

where $P_{\text{peak}} \geq 0$ is the peak power constraint. Another variable $\lambda \geq 0$ represents the regularizer that balances the trade-off between the expected distance and total transmission power. With larger λ , the system will have to put more emphasis on constraining the total transmission power. On the contrary, with smaller λ , the system will pay more attention to minimize the expected distortion between received signals and intended signals.

The peak power constraint can be adjusted along with the regularizer to control the peak-to-average-power-ratio (PAPR). To keep the discussion brief, $\text{obj}(\cdot)$ is used to denote the objective function in Eq. (24). It is defined as

$$\text{obj}(\mathbf{D}, \mathbf{X}, \mathbf{S}) = \left\| \tilde{\mathbf{H}} \mathbf{D} \mathbf{T} \mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^{\text{H}} \mathbf{T}^{\text{H}} \mathbf{D}^{\text{H}} \Sigma_{\tilde{\mathbf{E}}} \mathbf{D} \mathbf{T} \mathbf{X}\} + \lambda \text{tr}\{\mathbf{X} \mathbf{X}^{\text{H}}\} \quad (25)$$

The term $L \text{tr}\{\mathbf{P}_{\text{L}}^{-1}\} \sigma_{\text{n}}^2$ is removed from the objective function, because it is a constant term that depends neither on \mathbf{D} nor on \mathbf{X} .

III. GLSE PRECODING VIA ALTERNATING OPTIMIZATION

The joint optimization of the problem stated in (24) is non-trivial since it is generally non-convex. Alternating optimization method is adopted to obtain a suboptimal solution iteratively. In each iteration, the optimal RF chain outputs are first calculated while treating the IRS configurations

as a fixed matrix, and then the system looks for a suitable (may not be optimal) phase shifts configuration while treating \mathbf{X} as a fixed matrix. The pseudo-code in Algorithm (1) depicts the general procedure.

Algorithm 1 General Procedure for finding (\mathbf{D}, \mathbf{X})

- 1: $(\mathbf{D}, \mathbf{X}) \leftarrow$ initial values
 - 2: **while** terminate condition not met **do**
 - 3: $\mathbf{X} \leftarrow \arg \min_{\mathbf{X}} \text{obj}(\mathbf{D}, \mathbf{X}, \mathbf{S})$ s.t. $\max_{i,j} |x_{i,j}|^2 < P_{\text{peak}}$
 - 4: $\mathbf{D} \leftarrow \arg \min_{\mathbf{D}} \text{obj}(\mathbf{D}, \mathbf{X}, \mathbf{S})$
 - 5: **end while**
-

A termination condition for the loop should be chosen accordingly. For example, let the loop be executed until two consequent loops have their cost function difference smaller than a certain threshold or terminates when certain iteration time is reached. The step 3 in the algorithm (optimize with respect to \mathbf{X}) is quadratic programming, which is categorized as convex optimization problem. Therefore, an optimal \mathbf{X} can always be obtained if \mathbf{D} is given.

The domain of \mathbf{D} is not a convex set. This implies that the step 4 in the algorithm is a non-convex problem. Thus, the optimal value in step 4 can generally not be found. This leads to a problem that Algorithm (1) may not converge, because in step 4, the system may find a suboptimal minimal for \mathbf{D} that may lead to a larger cost than the previous iteration.

A. Optimization of the Digital Unit

When \mathbf{D} is treated as a fixed matrix, the problem given by

$$\begin{aligned} & \underset{\mathbf{X}}{\text{minimize}} \quad \left\| \tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^{\text{H}}\mathbf{T}^{\text{H}}\mathbf{D}^{\text{H}}\Sigma_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{X}\} + \lambda \text{tr}\{\mathbf{X}\mathbf{X}^{\text{H}}\} \\ & \text{subject to} \quad \max_{i,j} |x_{i,j}|^2 < P_{\text{peak}} \end{aligned} \quad (26)$$

belongs to the class of convex problem. Moreover, it is worth noticing that the columns (time intervals) of matrix \mathbf{X} can be decoupled according to Eq. (27).

$$\begin{aligned} & \text{tr}\{(\tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S})^{\text{H}}(\tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S})\} + \text{tr}\{\mathbf{X}^{\text{H}}\mathbf{T}^{\text{H}}\mathbf{D}^{\text{H}}\Sigma_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{X}\} + \lambda \text{tr}\{\mathbf{X}^{\text{H}}\mathbf{X}\} \\ & = \sum_{i=1}^L (\tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{x}_i - \mathbf{s}_i)^{\text{H}}(\tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{x}_i - \mathbf{s}_i) + \mathbf{x}_i^{\text{H}}\mathbf{T}^{\text{H}}\mathbf{D}^{\text{H}}\Sigma_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{x}_i + \lambda \mathbf{x}_i^{\text{H}}\mathbf{x}_i. \end{aligned} \quad (27)$$

Therefore, this optimization problem can either be treated as one matrix convex optimization or L independent vector convex optimization problems. This implies that the block length has no influence over the RF chain outputs if \mathbf{D} is fixed.

B. Optimization of the IRS Configuration

When treating \mathbf{X} as a fixed matrix, the optimization problem is then formulated as

$$\underset{\mathbf{D}}{\text{minimize}} \quad \left\| \tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^{\text{H}}\mathbf{T}^{\text{H}}\mathbf{D}^{\text{H}}\Sigma_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{X}\}. \quad (28)$$

It is generally non-convex problem, since the domain of \mathbf{D} is not a convex set. Two methods are used here to obtain a suboptimal point, which are gradient descent and MM algorithm [13]. For further discussions, we denote the objective function in Eq. (28) by

$$\text{obj}_{\mathbf{D}}(\mathbf{D}) = \left\| \tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S} \right\|_{\text{F}}^2 + \text{tr}\{\mathbf{X}^{\text{H}}\mathbf{T}^{\text{H}}\mathbf{D}^{\text{H}}\Sigma_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{X}\}. \quad (29)$$

1) *Gradient descent method:* Gradient descend is a general method to find the local optimal of the functions that maps one or multiple parameters to \mathbb{R} . The local minimal can be obtained by following the opposite gradient of the function iteratively from a random initial point. This can be understood graphically as following the steepest slope on a surface until reaching the trough.

In order to simplify the domain, define the phase vector

$$\boldsymbol{\beta} = \arg \cdot \text{diag}_{\text{vector}}(\mathbf{D}). \quad (30)$$

To keep the notation simple, we declare auxiliary matrices as follow

$$\begin{aligned} \mathbf{U} &= \mathbf{D}\mathbf{T}\mathbf{X}; \\ \mathbf{V} &= \mathbf{S}^{\text{H}}\tilde{\mathbf{H}} \\ \mathbf{Q} &= \tilde{\mathbf{H}}^{\text{H}}\tilde{\mathbf{H}} + \Sigma_{\tilde{\mathbf{E}}}. \end{aligned} \quad (31)$$

Under this convention, the objective function can be expanded as

$$\text{obj}_{\mathbf{D}}(\mathbf{D}) = \text{tr}\{\mathbf{U}^{\text{H}}\mathbf{Q}\mathbf{U} - (\mathbf{U}^{\text{H}}\mathbf{V}^{\text{H}} + \mathbf{V}\mathbf{U}) + \mathbf{S}^{\text{H}}\mathbf{S}\}. \quad (32)$$

Since the differential operation is linear, we first calculate the the derivative of the additive terms in Eq. (32) as

$$\begin{aligned} \frac{\partial \text{tr}\{\mathbf{U}^{\text{H}}\mathbf{Q}\mathbf{U}\}}{\partial \boldsymbol{\beta}} &= -2\text{diag}_{\text{vector}}(\Im(\mathbf{U}\mathbf{U}^{\text{H}}\mathbf{Q})) \\ \frac{\partial \text{tr}\{\mathbf{U}^{\text{H}}\mathbf{V}^{\text{H}} + \mathbf{V}\mathbf{U}\}}{\partial \boldsymbol{\beta}} &= -2\text{diag}_{\text{vector}}(\Im(\mathbf{U}\mathbf{V})). \end{aligned} \quad (33)$$

From Eq. (32) and Eq. (33), the partial derivative of the objective function can be derived as

$$\frac{\partial \text{obj}_{\mathbf{D}}(\mathbf{D})}{\partial \boldsymbol{\beta}} = -2\text{diag}_{\text{vector}}(\Im(\mathbf{U}\mathbf{U}^{\text{H}}\mathbf{Q} - \mathbf{U}\mathbf{V})). \quad (34)$$

The auxiliary vector can be transformed back to the phase shift matrix by Eq. (35).

$$\mathbf{D} = \text{diag}_{\text{matrix}}(\exp(j\boldsymbol{\beta})) \quad (35)$$

Combining with the designing of the digital unit, the pseudo-code of the precoding algorithm with gradient descent is shown as Algorithm (2).

Algorithm 2 Precoding (\mathbf{D} , \mathbf{X}) with gradient descent

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1: ( $\mathbf{D}$ ,  $\mathbf{X}$ )  $\leftarrow$  initial values
2: while terminate condition not met do
3:    $\mathbf{X} \leftarrow \text{convex}_{\text{opt}}\{\text{obj}(\mathbf{D}, \mathbf{X}, \mathbf{S}) \text{ s.t. } \max_{i,j} |x_{i,j}|^2 < P_{\text{peak}}\}$ 
4:   while terminate condition not met do
5:      $\boldsymbol{\beta} \leftarrow \arg \cdot \text{diag}_{\text{vector}}(\mathbf{D})$ 
6:      $\boldsymbol{\beta}_{\text{grad}} \leftarrow -2\text{diag}_{\text{vector}}(\Im(\mathbf{U}\mathbf{U}^H\mathbf{Q} - \mathbf{U}\mathbf{V}))$ 
7:      $\alpha \leftarrow \text{getStepSize}$ 
8:      $\boldsymbol{\beta}' = \boldsymbol{\beta} - \alpha\boldsymbol{\beta}_{\text{grad}}$ 
9:      $\mathbf{D} \leftarrow \text{diag}_{\text{matrix}}(\exp(j\boldsymbol{\beta}'))$ 
10:   end while
11: end while

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In order to avoid the situation where the algorithm diverges, one can design the step size α in the following steps. At the beginning of each gradient descent loop (step 4 - 10 in Algorithm (2)), α is initialized to a relative large value. During each loop, the processor keeps the old value of \mathbf{D} in the memory. If the value of the objective function does not decrease with the new \mathbf{D} , the processor will then load the old value of \mathbf{D} from the memory and redo that gradient descent loop once again with halved step size.

The whole gradient descent loop terminates if the value of α is smaller than a certain threshold. In this way, \mathbf{D} will always be assigned to a value that makes the objective function decrease. Moreover, if the termination threshold for α is small enough, \mathbf{D} will always end up at the close proximity of a critical point after the termination of one entire gradient descent loop.

It is worth noticing that, though we can make the algorithm converge to a suboptimal point, it does not mean that the systems starting with different initial pairs (\mathbf{D}, \mathbf{X}) will converge to the same value.

2) *Majorize-minimization (MM) algorithm*: MM algorithm [13] is another general optimization method that can be used to optimize the IRS configurations. Similar to gradient descent, this algorithm also includes iterative procedures itself. However, when using MM algorithm we don't need to state the step size explicitly.

A special auxiliary function called surrogate function $\overline{\text{obj}}_{\mathbf{D}_0} : \text{domain}(\text{obj}(\cdot)) \rightarrow \mathbb{R}$ is used in MM algorithm. It is a less complicated function that needs to be constructed at any point \mathbf{D}_0 of the objective function. The surrogate function constructed at a point must touch the objective function at that point. Meanwhile, it should always be larger or equal to the objective function at every point within its domain.

The procedure of optimizing \mathbf{D} with MM algorithm can be iteratively described as follows. At the i -th MM iteration for optimizing the IRS, a surrogate function of a point \mathbf{D}_i is constructed. We then find the optimal point \mathbf{D}_{i+1} that minimizes the surrogate function. After that, we proceed to the next MM iteration and optimize a new surrogate function tangent to point \mathbf{D}_{i+1} . We can choose the initial point of \mathbf{D} arbitrarily as long as it lies in the feasible set.

In order to simplify the problem, we wish to construct a linear surrogate function free of quadratic terms. When expanded, the objective function in Eq. (32) is comprised of a quadratic term, two first order terms and one constant term with respect to \mathbf{D} .

By the definition of surrogate function, the objective function is always a surrogate function of itself. The additive property of the surrogate functions can also be proven by the definition. This property states that two surrogate functions are additive if they are both tangent to their objective functions at the same point. Based on these two properties, a linear surrogate function for the objective function Eq. (32) can be constructed as follows.

We first construct a linear surrogate function for the quadratic term of Eq. (32). The first order terms and constant term will serve as surrogate functions of themselves. The sum of those surrogate functions are the surrogate function for the whole objective function described by Eq. (32). Since the constant term does not contribute to the optimization, we can safely discard it.

Next we will give out the equivalent surrogate function.

To find the surrogate function, we use the auxiliary variables defined in Eq. (31)

$$\begin{aligned} \mathbf{U} &= \mathbf{D}^T \mathbf{X}; \\ \mathbf{V} &= \mathbf{S}^H \tilde{\mathbf{H}} \\ \mathbf{Q} &= \tilde{\mathbf{H}}^H \tilde{\mathbf{H}} + \Sigma_{\tilde{\mathbf{E}}}. \end{aligned}$$

Furthermore, define

By substituting the variables according to Eq. (31), the quadratic term can be written as

$$\mathbf{X}^H \mathbf{T}^H \mathbf{D}^H \left(\hat{\mathbf{H}}^H \hat{\mathbf{H}} + \Sigma_{\mathbf{E}} \right) \mathbf{D} \mathbf{T} \mathbf{X} = \mathbf{U}^H \mathbf{Q} \mathbf{U}. \quad (36)$$

To construct the surrogate function for the quadratic term at an arbitrary point \mathbf{D}_0 , we define an auxiliary matrix \mathbf{P} , such that $\mathbf{P} - \mathbf{Q}$ is positive definite. For the tangential point \mathbf{D}_0 , we also define $\mathbf{U}_0 = \mathbf{D}_0 \mathbf{T} \mathbf{X}$. The surrogate for the quadratic term can be obtained by the following inequality [13].

$$\begin{aligned} & \text{tr} \{ \mathbf{U}^H \mathbf{Q} \mathbf{U} \} \\ &= \text{tr} \{ \mathbf{U}_0^H \mathbf{Q} \mathbf{U}_0 + (\mathbf{U} - \mathbf{U}_0)^H \mathbf{Q} \mathbf{U}_0 + \mathbf{U}_0^H \mathbf{Q} (\mathbf{U} - \mathbf{U}_0) + (\mathbf{U} - \mathbf{U}_0)^H \mathbf{Q} (\mathbf{U} - \mathbf{U}_0) \} \\ &\leq \text{tr} \{ \mathbf{U}_0^H \mathbf{Q} \mathbf{U}_0 + (\mathbf{U} - \mathbf{U}_0)^H \mathbf{Q} \mathbf{U}_0 + \mathbf{U}_0^H \mathbf{Q} (\mathbf{U} - \mathbf{U}_0) + (\mathbf{U} - \mathbf{U}_0)^H \mathbf{P} (\mathbf{U} - \mathbf{U}_0) \} \\ &= \text{tr} \{ \mathbf{U}^H \mathbf{P} \mathbf{U} + \mathbf{U}^H (\mathbf{Q} - \mathbf{P}) \mathbf{U}_0 + \mathbf{U}_0^H (\mathbf{Q} - \mathbf{P}) \mathbf{U} + \mathbf{U}_0^H (\mathbf{Q} - \mathbf{P}) \mathbf{U}_0 \} \end{aligned} \quad (37)$$

The newly introduce quadratic term $\mathbf{U}^H \mathbf{P} \mathbf{U}$ will be degraded to a constant term if we specify \mathbf{P} to be $\mathbf{P} = \lambda_{\max}^{\mathbf{Q}} \mathbf{I}$, where $\lambda_{\max}^{\mathbf{Q}}$ is the largest eigenvalue of \mathbf{Q} .

By adding the first order term of the objective function Eq. (29) and removing constant terms, the equivalent surrogate function (not the true surrogate since the constant terms are removed) for the objective function at any point \mathbf{D}_0 reads

$$\overline{\text{obj}_{\mathbf{D}_0}(\mathbf{D})} = \text{tr} \{ \mathbf{U}^H (\mathbf{Q} - \mathbf{P}) \mathbf{U}_0 + \mathbf{U}_0^H (\mathbf{Q} - \mathbf{P}) \mathbf{U} - \mathbf{V} \mathbf{U} - \mathbf{U}^H \mathbf{V}^H \}. \quad (38)$$

For simplicity, define a constant

$$\mathbf{G}_0 = \mathbf{T} \mathbf{X} \mathbf{U}_0^H (\mathbf{Q} - \mathbf{P}) - \mathbf{T} \mathbf{X} \mathbf{V}. \quad (39)$$

By substituting Eq. (39) into Eq. (38), the latter becomes

$$\overline{\text{obj}_{\mathbf{D}_0}(\mathbf{D})} = \text{tr} \{ \mathbf{G}_0 \mathbf{D} + \mathbf{D}^H \mathbf{G}_0^H \} \quad (40)$$

The tangential point \mathbf{D}_1 for starting the next MM iteration is obtained by minimizing the equivalent surrogate function described by Eq. (40). By introducing two auxiliary constant terms $\text{tr} \{ \mathbf{D}^H \mathbf{D} \}$

and $\text{tr}\{\mathbf{G}_0\mathbf{G}_0^H\}$, we have [13]

$$\begin{aligned}
\mathbf{D}_1 &= \arg \min_{\mathbf{D}} (\text{tr}\{\mathbf{G}_0\mathbf{D} + \mathbf{D}^H\mathbf{G}_0^H\}) \\
&= \arg \min_{\mathbf{D}} (\text{tr}\{\mathbf{G}_0\mathbf{D} + \mathbf{D}^H\mathbf{G}_0^H + \mathbf{D}^H\mathbf{D} + \mathbf{G}_0\mathbf{G}_0^H\}) \\
&= \arg \min_{\mathbf{D}} (\text{tr}\{(\mathbf{D} - (-\mathbf{G}_0)^H)^H(\mathbf{D} - (-\mathbf{G}_0)^H)\}) \\
&= \arg \min_{\mathbf{D}} (\|\mathbf{D} - (-\mathbf{G}_0)\|_F^2) \\
&= \exp((\arg(-\mathbf{G}_0))
\end{aligned} \tag{41}$$

The complete pseudo-code for the MM algorithm based precoding is illustrated as Algorithm (3). The convergence of Algorithm (3) is also guaranteed. Due to the definition of surrogate

Algorithm 3 Precoding (\mathbf{D} , \mathbf{X}) with MM algorithm

```

1: ( $\mathbf{D}$ ,  $\mathbf{X}$ )  $\leftarrow$  initial values
2: while terminate condition not met do
3:    $\mathbf{X} \leftarrow \text{convex}_{\text{opt}}\{\|\tilde{\mathbf{H}}\mathbf{D}\mathbf{T}\mathbf{X} - \mathbf{S}\|_F^2 + \text{tr}\{\mathbf{X}^H\mathbf{T}^H\mathbf{D}^H\mathbf{\Sigma}_{\tilde{\mathbf{E}}}\mathbf{D}\mathbf{T}\mathbf{X}\} + \lambda\text{tr}\{\mathbf{X}\mathbf{X}^H\}\}$ 
4:   while terminate condition not met do
5:      $\mathbf{D}_0 \leftarrow \mathbf{D}$ 
6:      $\mathbf{U}_0 \leftarrow \mathbf{D}_0\mathbf{T}\mathbf{X}$ 
7:      $\mathbf{V} \leftarrow \mathbf{S}^H\tilde{\mathbf{H}}$ 
8:      $\mathbf{Q} \leftarrow \tilde{\mathbf{H}}^H\tilde{\mathbf{H}} + \mathbf{\Sigma}_{\tilde{\mathbf{E}}}$ 
9:      $\mathbf{P} \leftarrow \lambda_{\max}^{\mathbf{Q}}\mathbf{I}$ 
10:     $\mathbf{G}_0 \leftarrow \mathbf{T}\mathbf{X}\mathbf{U}_0^H(\mathbf{Q} - \mathbf{P}) - \mathbf{T}\mathbf{X}\mathbf{V}$ 
11:     $\mathbf{D}_1 \leftarrow \exp((\arg(-\mathbf{G}_0))$ 
12:     $\mathbf{D} \leftarrow \mathbf{D}_1$ 
13:   end while
14: end while

```

function, the value of the objective function in the MM iterations (step 4 to step 13 in Algorithm (3)) must form a monotonically decreasing sequence. Therefore, the value of the objective function after the IRS optimization must be smaller than the value before. Thus, Algorithm (3) also converge. However, similar to the gradient descent method, Algorithm (3) may converge to different values if the initial points are chosen differently

l_{ref}	λ_w	σ_f^2	θ_p^t, ϕ_p^t	d	n	P	K
[50, 150]	0.01	1	$[0, \pi]$	$\lambda_w/2$	3	8	8

TABLE I: Parameters for channel H

IV. NUMERICAL INVESTIGATIONS

It is proven in [14] that by using Jensen's inequality, the average rate per user is bounded by

$$\bar{R} \geq \frac{LT_s}{\tau + LT_s} \log \left(1 + \frac{\sigma_s^2}{\sigma_{\tilde{n}_{\max}}^2 + P_{\text{Interference}}} \right), \quad (42)$$

where T_s is symbol time, τ is the duration of guard interval and $P_{\text{Interference}}$ is the average interference power after the decoding/compensation process. The variance $\sigma_{\tilde{n}_{\max}}^2$ is the maximal power of the normalized noise, while σ_s^2 is the power of the intended signal.

Another value of interest is the PAPR. In our simulations, the chain-wise PAPR of every RF chain is calculated by dividing the chain-wise peak power by the average power of transmitted signals in each corresponding chain. The average PAPR is calculated by averaging the chain-wise PAPR over all the RF chains.

A. Simulation Setup

In order to simulate the channel correlations, we place 8 UTs at the range of $50 + 14 \times k$ meters, where $k \in \{0, \dots, 7\}$. Each BS-UT channel consists of 8 effective paths. The intended messages for the UTs have the same power $\sigma_s^2 = 1$. The angles from the BS to the scatterings are constant during the whole transmission.

Since block fading is assumed, the fading parameters are generated once every block. To make the further discussion easier, we assume that the distances between the UTs and the BS are deterministic and known by all the UTs and the BS. The default settings for the channel are concluded as Table (I).

The matrix \mathbf{T} describes the illumination channel from the active antennas to the IRS. In the simulations, N horn-shape active antennas are uniformly placed on a circle parallel to the xz -plane. The center of the the active antenna arrays and the center of the IRS form a line that is also parallel to the y -axis. The radius of the active antenna arrays is denoted as R_r , and the distance between this antenna array and the xz -plane is denoted as R_d . This array setup is exemplified in Fig. (5). Analytically, the illumination channel is given by [11]

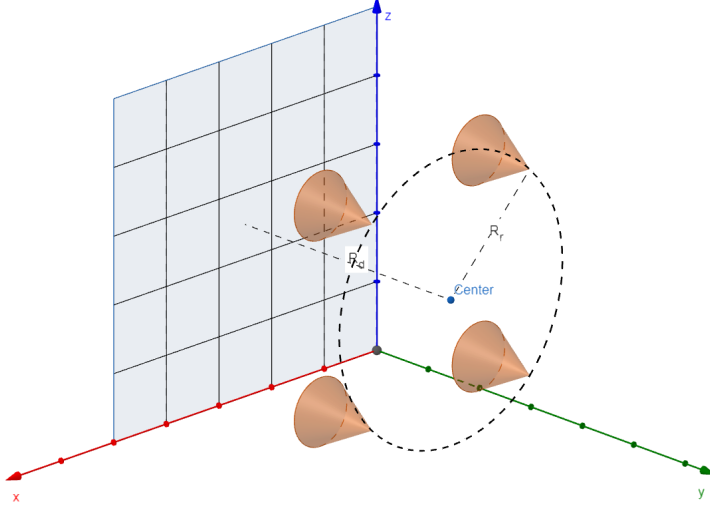


Fig. 5: Placement of the active antennas

$$\mathbf{T} = \left[\frac{\lambda_w \sqrt{\rho G^a(\theta_{m,n}^p, \phi_{m,n}^p) G^p(\theta_{m,n}^a, \phi_{m,n}^a)}}{4\pi r_{m,n}} e^{-j \frac{2\pi r_{m,n}}{\lambda_w}} \right]_{m,n}, \quad (43)$$

where $\rho \in [0, 1]$ denotes the power efficiency of the IRS. The function $G^a(\theta_{m,n}^p, \phi_{m,n}^p)$ and $G^p(\theta_{m,n}^a, \phi_{m,n}^a)$ model the antenna gain of the active antenna and the IRS element respectively. The variable $r_{m,n}$ measures the distance from the m -th IRS element to the n -th active antenna.

We assume here that the IRS elements reflect all the signals in front of them, therefore we have

$$G^p(\theta, \phi) = \begin{cases} 2 & \theta, \phi \in [0, \pi] \\ 0 & \text{otherwise} \end{cases}. \quad (44)$$

The active antenna gain following [11] is assumed to be

$$G^a(\theta, \phi) = \begin{cases} 2(1 + \kappa) \cos^\kappa(\phi) & \phi \in [0, \pi/2] \\ 0 & \text{otherwise} \end{cases}. \quad (45)$$

The κ here is a normalization factor which ensures that the spherical surface integral of $G^a(\theta, \phi)$ is 4π . With larger κ , the beam from the active antenna will be more concentrated to its main direction.

Besides the antenna patterns, the orientations of the active antennas also matter. To investigate the effect of interference on the IRS, two illumination strategies are used.

ρ	R_d	R_r	κ
1	$S1 : \frac{4d\sqrt{M}}{\sqrt{\pi}}, S2 : \frac{4d\sqrt{M}}{\sqrt{N\pi}}$	$S1 : 2d, S2 : \frac{d\sqrt{2M}}{4}$	49

TABLE II: Parameters for fading and path loss

In full illumination (FI), all active antenna illuminates toward the center of IRS. In this way, the IRS receives the most power as well as most interference. This illumination strategy is shown in Fig. (6).

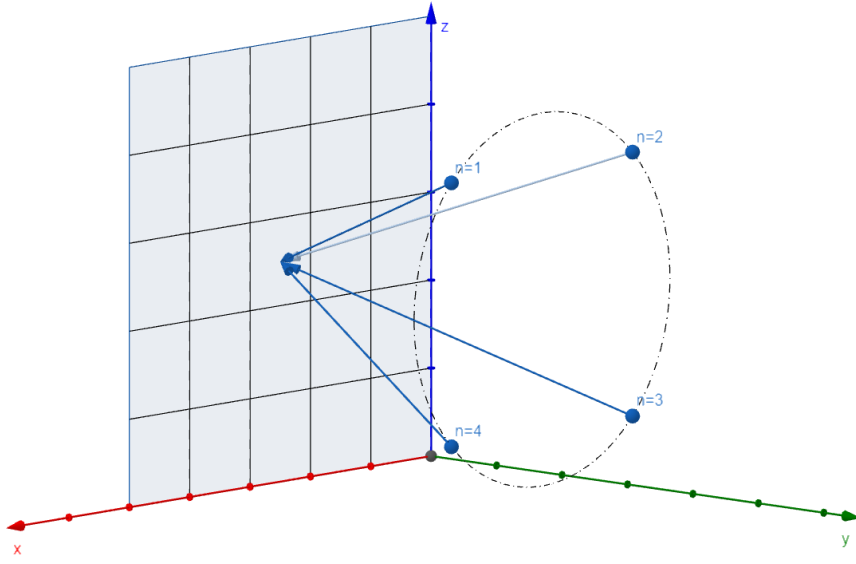


Fig. 6: Full illumination

In partial illumination (PI), we designate N sections on the IRS, such that each section is responsible for one active antenna. Each active antenna illuminates the signals to the center of its designated section. Fig. (7) depicts this scenario.

The illumination parameters are concluded in Table (II) [11].

B. Rate and PAPR

By using the above mentioned setup, and setting IRS antenna element amount to 64, we plot the rate performance for full illumination with gradient descent and MM algorithm. Partial illumination based on MM algorithm is also simulated for comparison. It is assume here that the

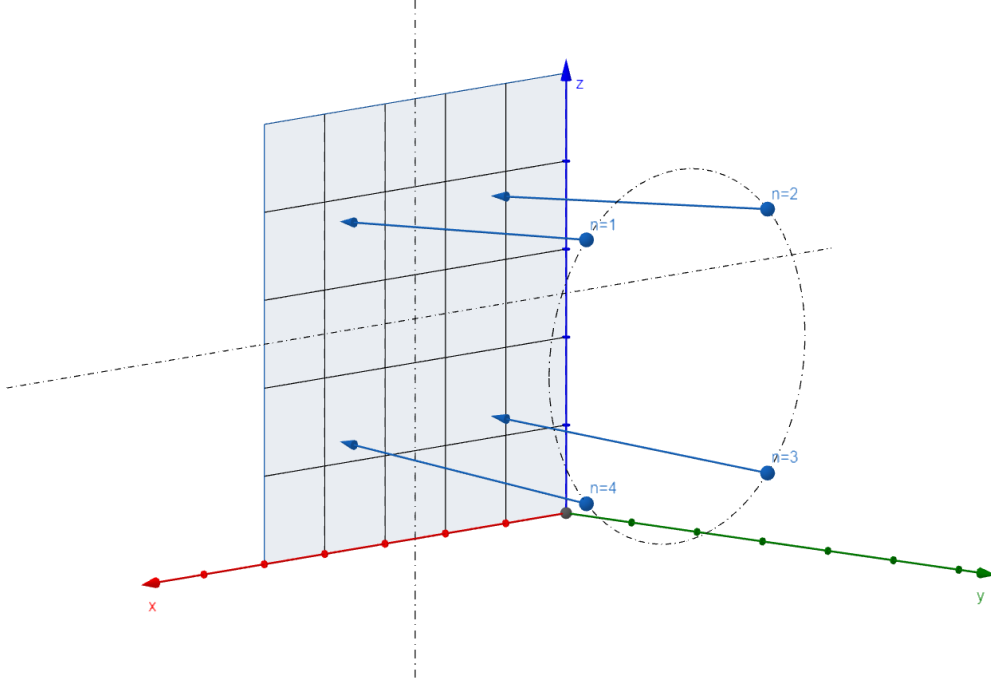


Fig. 7: Partial illumination

estimation error is zero and the maximal normalized noise is set to be $\sigma_{\tilde{n}_{\max}}^2 = 0.001$ such that it has a similar scale as the interference power. The PAPR is adjusted by increasing the peak power constraint while the regularizer λ is fixed to be 0.5. For block length $L = 1$, the result is shown as Fig. (8).

It can be observed by comparing the red line and the yellow line that partial illumination strategy performs slightly worse than the full illumination strategy. One reason is the spillover effect caused by the antenna pattern. In full illumination, the IRS receives more power than the partial illumination strategy. In fact, it can be proven that given enough power and enough IRS size, symbol-wise distortion-free transmission is always achievable.

Theorem 1. *If \mathbf{H} is a complex normal Gaussian random matrix with independent entries and the length of the uni-modular vector \mathbf{d} can be arbitrary long, then it is almost certain that there is a \mathbf{d} such that $\mathbf{H} \cdot \mathbf{d} = \mathbf{0}$.*

Proof. See appendix. □

For the precoding optimization problem with block length $L = 1$ and number of RF chains $N = 1$. Define $\mathbf{H}' = \mathbf{H} \times \text{diag}_{\text{matrix}}\{x\mathbf{t}\}$, and the problem is transformed to finding vector

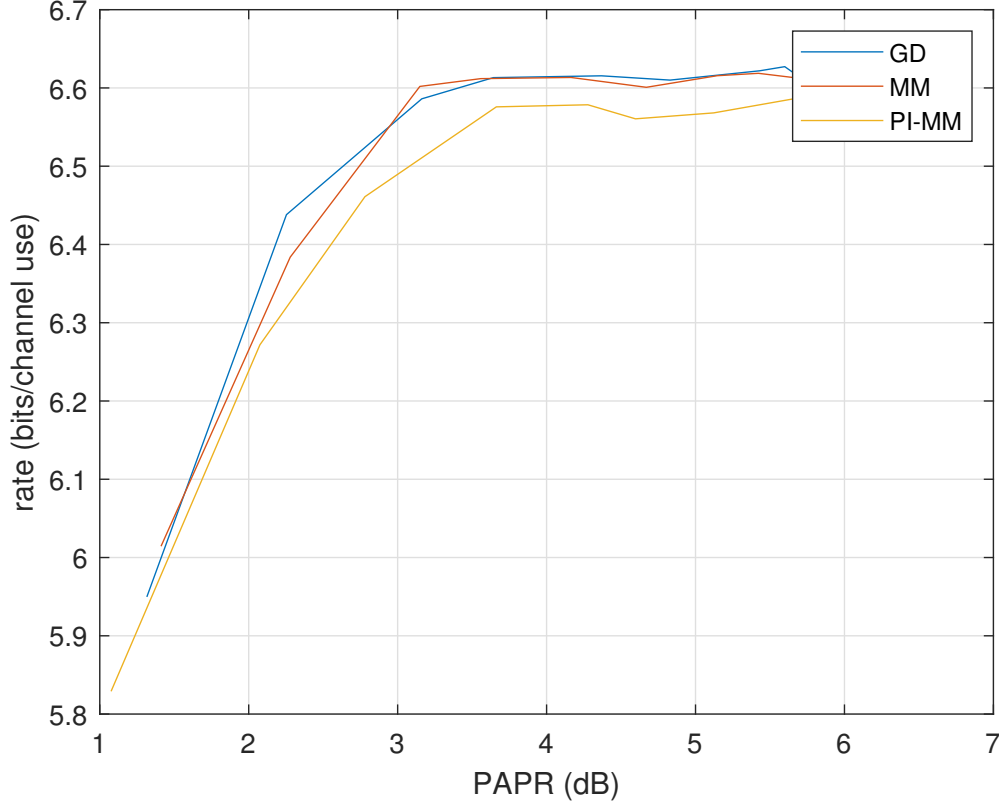


Fig. 8: Rate-PAPR with block length $L = 1$, guard interval $\tau = 0$ and regularizer $\lambda = 0.5$

$\mathbf{d} = \text{diag}_{\text{vector}}\{\mathbf{D}\}$ such that $\mathbf{s} = \mathbf{H}'\mathbf{d}$. To use the theorem, define

$$\begin{aligned} \mathbf{H}'' &= \begin{bmatrix} \mathbf{H}' & \mathbf{s} \end{bmatrix} \\ \mathbf{d}' &= \begin{bmatrix} \mathbf{d}_{\text{aux}} \\ d_{M+1} \end{bmatrix}. \end{aligned} \quad (46)$$

According to the theorem above, a solution for \mathbf{d}' can be found such that $\mathbf{0} = \mathbf{H}''\mathbf{d}'$. Therefore the IRS parameters should be $\mathbf{D} = \text{diag}_{\text{matrix}}\{-d_{M+1}^{-1}\mathbf{d}_{\text{aux}}\}$. For the cases where more RF chains are used $N > 1$, we can always transform it back to the case where $N = 1$ by muting all the RF chains except the first one.

For increased block length, where $L = 2$, the simulation results are shown in Fig. (9).

It can be seen here that with guard interval $\tau = 0$, the systems with larger block lengths tend to have worse rate performance compared to the systems with smaller block lengths. This can be explained by code space density at the receivers side. Let's assume a fixed stream of code and

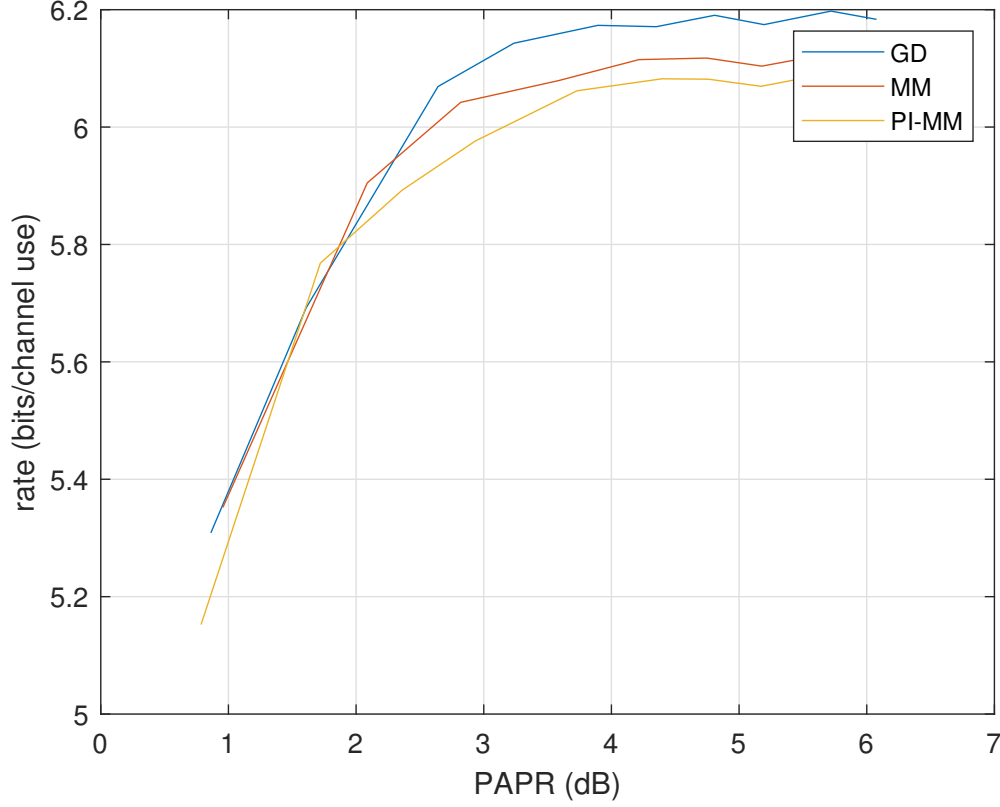


Fig. 9: Rate-PAPR with block length $L = 2$, guard interval $\tau = 0$ and regularizer $\lambda = 0.5$

messages with length L_{tot} . If we have two options for block lengths, L_1 and L_2 . Suppose L_{tot} is dividable by L_1 and L_2 . Following Eq. (27), we can see that the received code at the UTs side is actually generated by any combination of $(\mathbf{x}_i, \dots, \mathbf{x}_{L_{\text{tot}}}, \mathbf{D}_i, \dots, \mathbf{D}_{L_{\text{tot}}})$ such that $\mathbf{D}_i = \mathbf{D}_j$ if i and j belong to the same time block. There are less constraints for the scheme where smaller block length is used. thus, the system which adopts a smaller block length should have a denser code stream space at the receiver side, and it is more likely for the transmitter to find a code stream that is closer to the intended messages.

C. Regularizer

From now on, we set the regularized noise variance to be $\sigma_{\hat{\mathbf{n}}_{\text{max}}}^2 = 0.1$, such that the noise power is relative close to the interference power.

The transmission power is controlled by the regularizer λ . It balances the emphasis between the interference power and the RF chains signal power.

In order to study the behavior of regularizer, we set the peak power constraint to infinity and plot the rate- λ curves in Fig. (10). The default illumination strategy is FI. And the channel estimation error is not included.

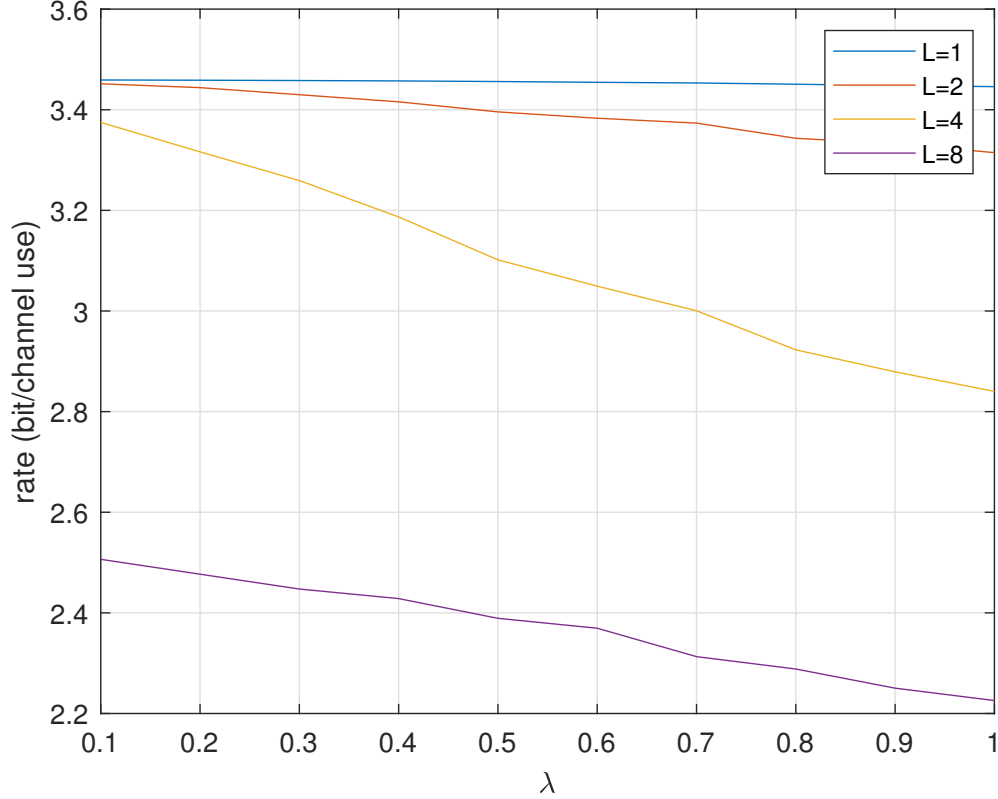


Fig. 10: Rate- λ behavior of MM-algorithm with guard interval $\tau = 0$

Since $\sigma_s^2 = 1$ and $\sigma_{\hat{n}_{\max}}^2 = 0.1$, we would have $rate = \log_2 11 \approx 3.45943$ if the distortion power is zero. Fig. (10) shows that for $L \leq 4$, the curves approach this rate limit as λ tend to zero. Similar phenomenon was researched in [14]. It states that for full rank \mathbf{H} and \mathbf{S} , the interference power $\|\mathbf{H}\mathbf{F}_{\text{GA}}\mathbf{X} - \mathbf{S}\|_{\text{F}}^2$ can never be made exactly to zero if the block length is larger than the number of RF chains. The matrix $\mathbf{F}_{\text{GA}} \in \mathbb{C}^{M \times N}$ in the statement represents the analog unit and it can be any complex matrix of size $M \times N$. Since the IRS structure is a special case of general HAD system, this statement can also be used to explain the purple curve standing for $L = 8$. The behavior of the blue curve $L = 1$ can also be explained by using Theorem (1).

The relation between the average RF chain power and the regularizer lambda is shown as Fig. (11). For relative small block lengths, the system with small L performs better. However, for

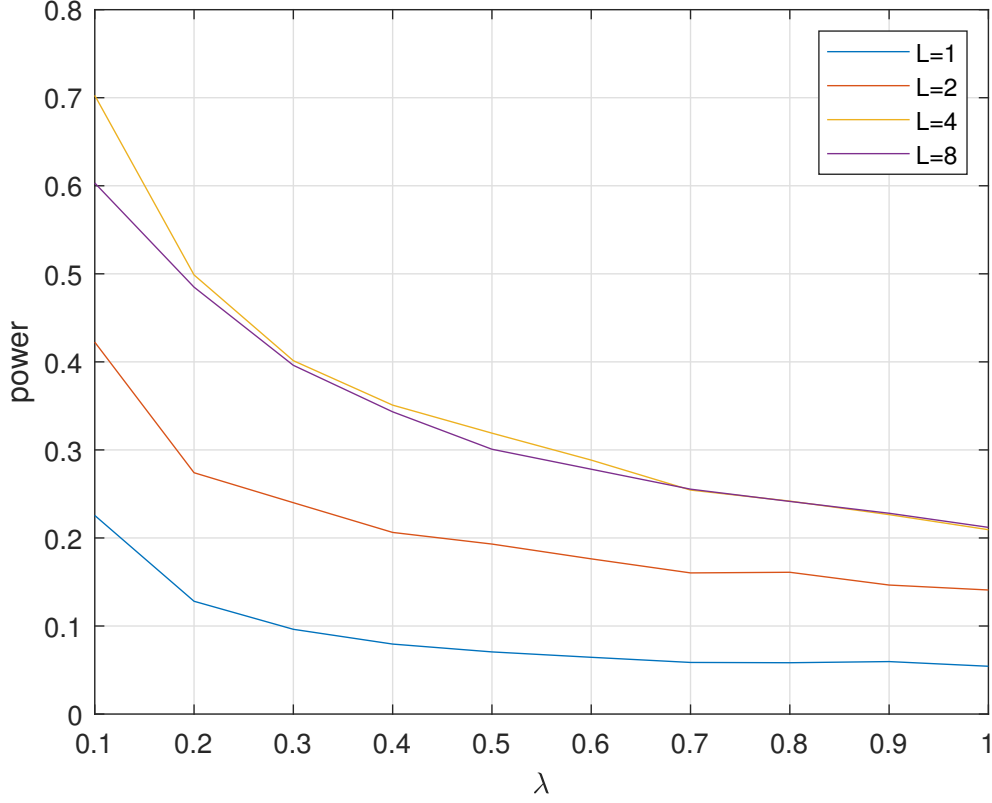


Fig. 11: Average RF chain power- λ with guard interval $\tau = 0$

systems with relative large block lengths, the power curves are close to each other. The reason for this can be explained by analyzing the objective function.

For smaller block lengths, the interference power can be made very small. As a result, the dominant part of the objective function is the term for transmission power. Thus, the system with small L will put more emphasis on reducing the RF chain power. However, systems with larger L will concentrate on minimizing the interference power.

D. Rate, Block Length and Number of IRS Elements

Now, set the guard interval to $\tau = T_s$ and remove the peak power constraint. We assume that the normalized channel estimation error matrix is comprised of i.i.d. entries, i.e.

$$\forall k \in \mathbb{N} \cap [1, K], \sigma_{e,k}^2 / \bar{h}_k = \sigma_e^2, \quad (47)$$

where σ_e^2 is the variance of the entries of the normalized estimation error matrix.

The Fig. (12) shows that the achievable rate increases as the number of IRS elements increases and the rate decreases as the variance of the estimation error increases, which are expected.

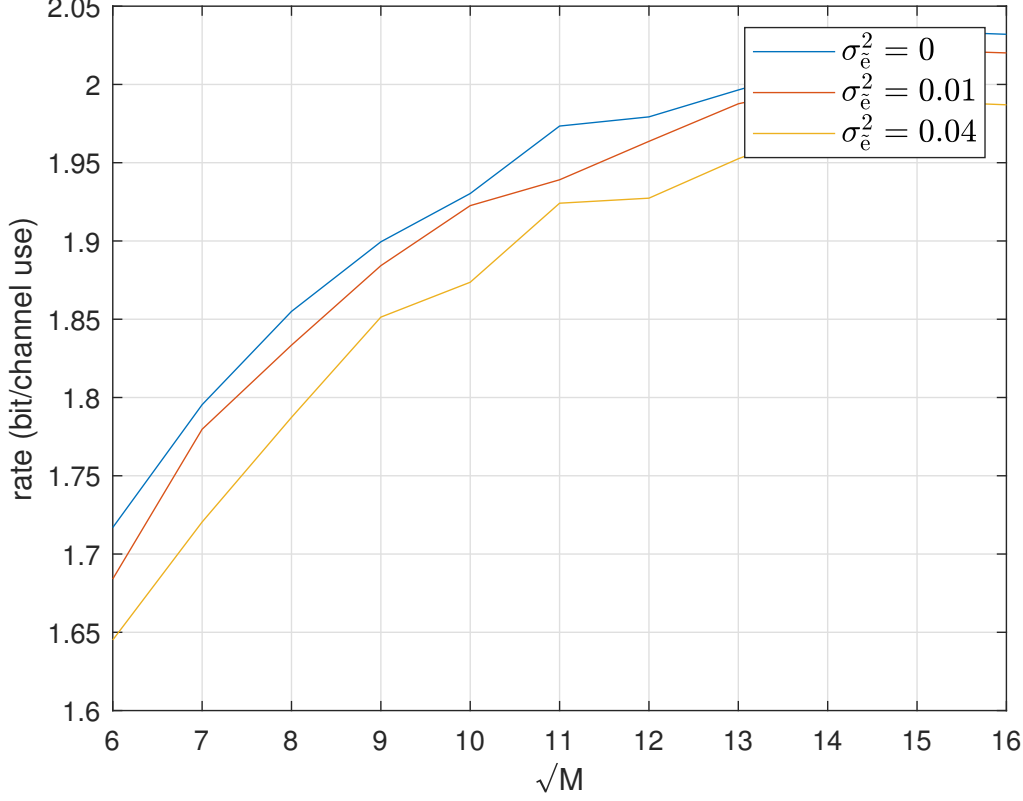


Fig. 12: Rate- \sqrt{M} with guard interval $\tau = T_s$ and $L = 10$

Next, we want to study the rate-block length behavior. The number of IRS elements is set to $M = 144$, and the simulation results are plotted in Fig. (13). In this figure, the rate first increases and then decreases. This phenomenon is caused by the non-zero guard interval.

For smaller block length, the interference is small relative to the noise power. As a result, the dominant part is the guard interval. It can be seen that the at $L = 1$ the rate of the estimation error free curve is 1.73(bits/channeluse) which is one half of $\log_2 11$. This is consistent with Theorem (1). As the block length increases, the term $\frac{LT_s}{\tau + LT_s}$ tends to 1 and the dominant part becomes the term $\log(1 + SINR)$.

Based on these two points, we can predict that for larger τ , the peak of the rate-L plot will shift to right since the term $\frac{LT_s}{\tau + LT_s}$ monotonically increases as L increases while the term

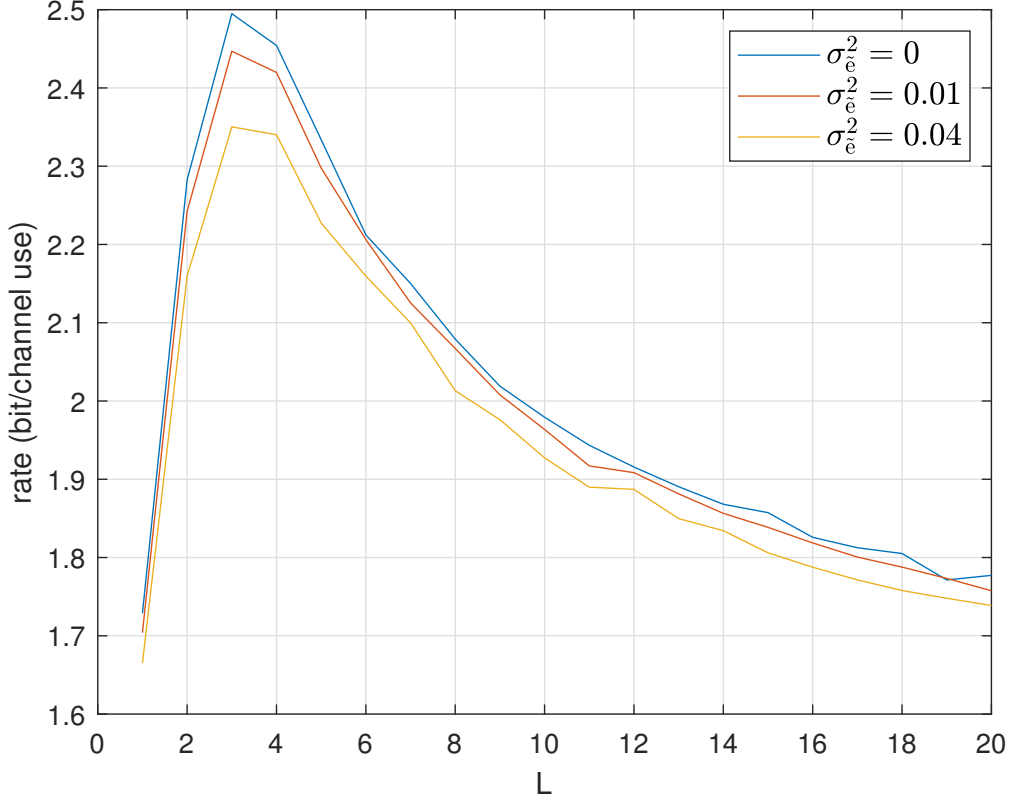


Fig. 13: Rate- L with guard interval $\tau = T_s$ and $M = 144$

$\log(1 + SINR)$ decreases as L increases. The complete relation of rate, \sqrt{M} and L are shown as a mesh surface in Fig. (14).

At last it would also be interesting to know the minimal required IRS elements to achieve a certain rate with increasing block length. Instead of directly plotting by reading Fig. (14), we calculate the rate of each transmission block, and use this value to obtain the averaged minimal IRS size for reaching a certain rate.

For rate threshold 1.8(bits/channeluse), the averaged minimal M versus L is plotted as Fig. (15).

V. APPENDIX

A. Proof for Theorem 1

For the simple case where \mathbf{H} is a row vector, namely when there is only one user $K = 1$, define $\mathbf{h}^T = \mathbf{H}$. We then try to find out the lower bound of probability $\Pr(\exists \mathbf{d}, (\mathbf{h}^T \mathbf{d} = 0))$.

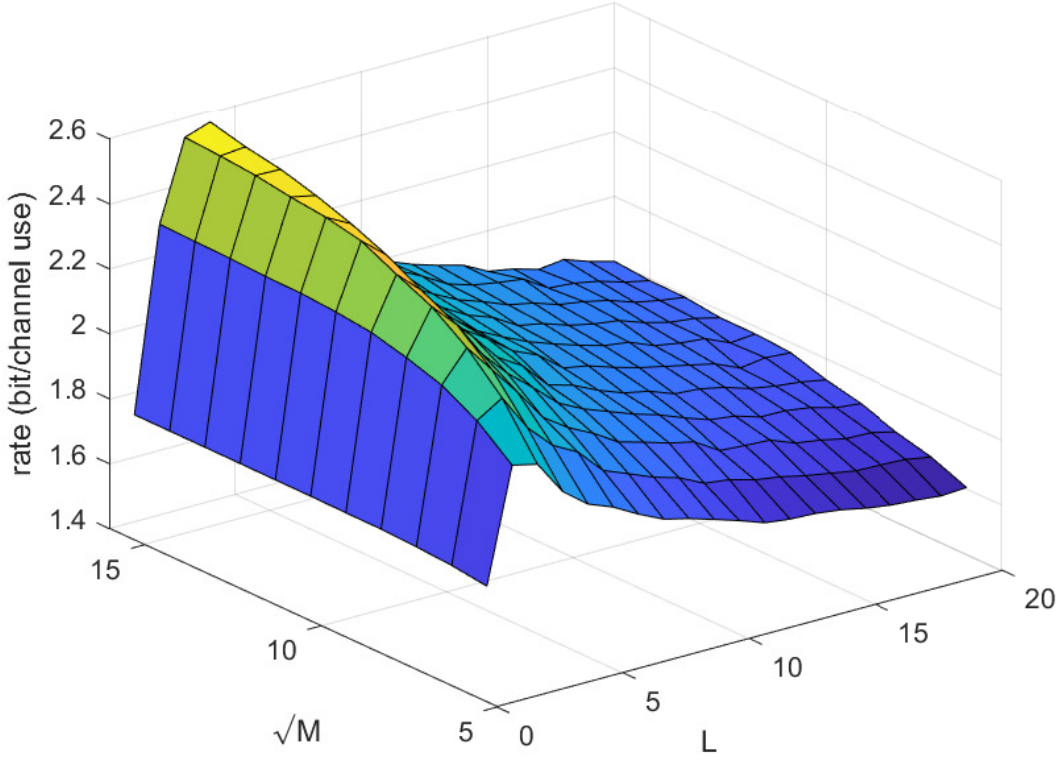


Fig. 14: Impact of IRS size and Block size ($\tau = 1$)

This can be translated into geometry representations. Each complex entry in vector \mathbf{h}^T is an 2-D vector whose length is a random variable following standard Gaussian distribution and phase is uniformly distributed between 0 to 2π . From this point of view, the uni-modular vector \mathbf{d} represents the rotation operations which are applied to the columns of \mathbf{h}^T with controllable angles. Therefore, the proposition $\exists \mathbf{d}, (\mathbf{h}^T \mathbf{d} = 0)$ can be interpreted as finding an rotation angle for each of the entries in vector \mathbf{h}^T such that these rotated entries add up to zero. Following triangular inequality, we can say that there exists a vector \mathbf{d} , such that $\mathbf{h}^T \mathbf{d} = 0$. This means

$$\max(|h_1|, \dots, |h_M|) \leq |h_1| + \dots + |h_M| - \max(|h_1|, \dots, |h_M|). \quad (48)$$

It can also be observed that $|h_1| > |h_2| + \dots + |h_M|$ and $|h_2| > |h_1| + \dots + |h_M|$ will never happen simultaneously. Therefore, we can write the above probability in negation form while

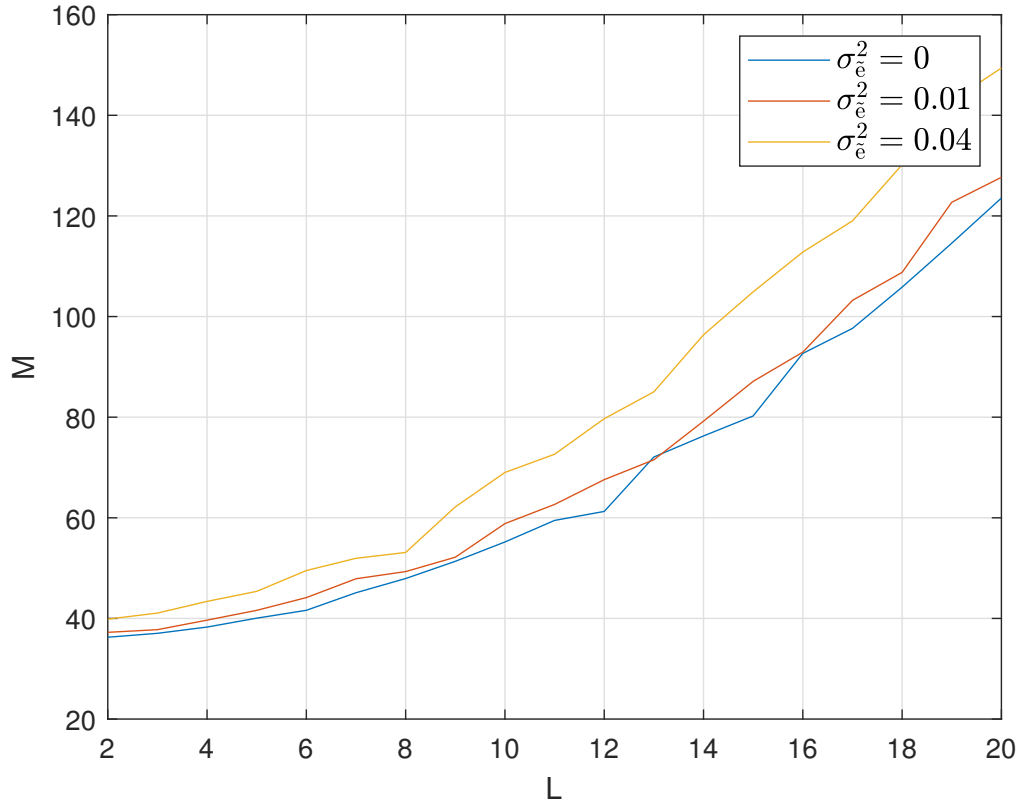


Fig. 15: Averaged minimal M - L with rate_threshold= 1.8

taking the independence assumption into consideration

$$\begin{aligned}
 P_1 &= \Pr(\exists \mathbf{d}, (\mathbf{h}^T \mathbf{d} = 0)) \\
 &= 1 - (M \times \Pr(|h_1| > |h_2| + \dots + |h_M|))
 \end{aligned} \tag{49}$$

The properties of F-distribution can be used to determine the upper bound of $\Pr(|h_1| > |h_2| +$

$\dots + |h_M|).$

$$\begin{aligned}
& M \times \Pr(|X| > |X_1| + \dots + |X_{M-1}|) \\
& \leq M \times \Pr(|X|^2 > |X_1|^2 + \dots + |X_{M-1}|^2) \\
& \leq M \times \Pr(|X|^2 + |Y|^2 > |X_1|^2 + \dots + |X_{M-1}|^2) \\
& = M \times \Pr\left(\frac{2}{M-1} > \frac{(|X_1|^2 + \dots + |X_{M-1}|^2)/(M-1)}{(|X|^2 + |Y|^2)/2}\right) \\
& = M \times F\left(\frac{2}{M-1}; M-1, 2\right) = M \times I_{\frac{1}{2}}\left(\frac{M-1}{2}, 1\right) \\
& = M \times \left(\frac{1}{2}\right)^{\frac{M-1}{2}},
\end{aligned} \tag{50}$$

where $F(x; z_1, z_2)$ is the cumulative distribution function with z_1 and z_2 degrees of freedom and $I_x(a, b)$ is the regularized incomplete beta function. The first less equal sign holds because $|X|^2 > (|X_1|^2 + \dots + |X_{M-1}|^2)^2 > |X_1|^2 + \dots + |X_{M-1}|^2$.

For the multi-user case where $K \neq 1$, the matrix \mathbf{H} has more than one rows. A solution for \mathbf{d} can be constructed recursively.

We first try to construct a sufficient condition for the existence of such vector \mathbf{d} . Then we try to prove that the probability for this sufficient condition not to hold tends to zero.

For sake of simplicity, we denote the channel matrix as $\mathbf{H}^{(k)}$ if $K = k$. Suppose we know how to construct \mathbf{d} if $K = k-1$ such that $\mathbf{H}^{(k-1)} \cdot \mathbf{d} = \mathbf{0}$. This also implies that $(\mathbf{H}^{(k-1)} \cdot \mathbf{d}) \cdot e^{j\varphi} = \mathbf{0}$, where φ is an arbitrary number.

Based on these assumptions and discussions, for the case where $K = k$, we can sectorize $\mathbf{H}^{(k)}$ as Fig. (16) and sectorize \mathbf{d} into several vectors correspondingly.

We only need to find one solution for this problem, and the condition for that solution to exist can serve as sufficient condition for $\exists \mathbf{d}, (\mathbf{H}^{(k)} \mathbf{d} = \mathbf{0})$.

One instance of \mathbf{d} may be constructed via the following steps. We first find $\mathbf{d}_1, \dots, \mathbf{d}_B$ such that $\forall i \in \{1, \dots, B\}, \mathbf{H}_i^{(k-1)} \mathbf{d}_i = \mathbf{0}$. Then for the last row of matrix $\mathbf{H}^{(k)}$, we know that $\mathbf{h}_i^T \cdot \mathbf{d}_i$ is a scalar value. At the same time, for each subvector \mathbf{d}_i , we can still apply an additional rotation $e^{j\varphi_i}$ to each of the vectors \mathbf{d}_i .

Now the problem is transformed back to the simple case, where we want to find rotation

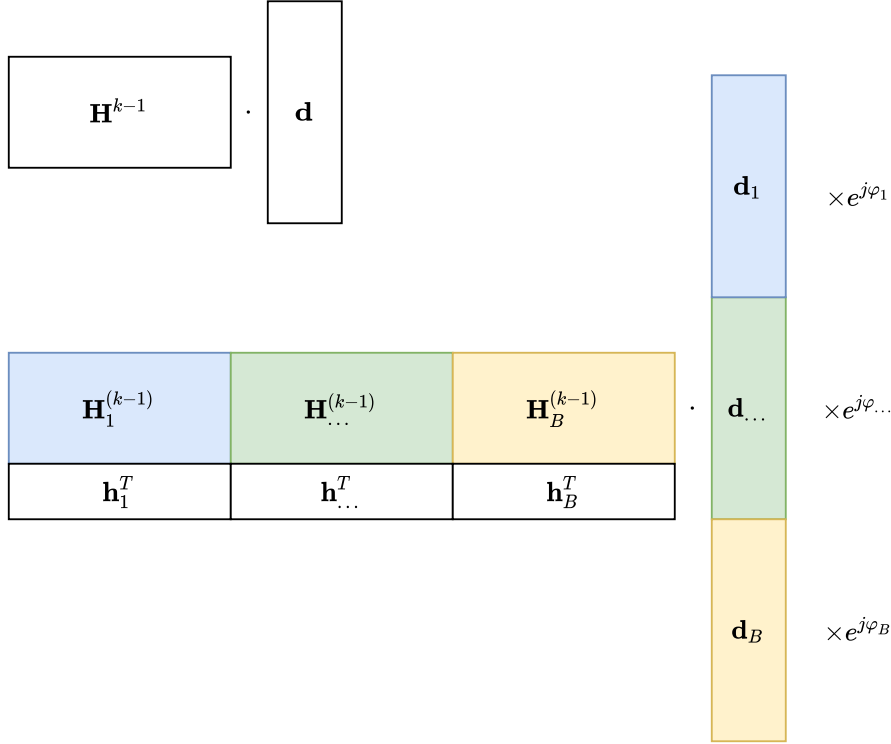


Fig. 16: Idea of construction

angles $\varphi_1, \dots, \varphi_B$ such that

$$\begin{bmatrix} \mathbf{h}_1^T \mathbf{d}_1 & \dots & \mathbf{h}_B^T \mathbf{d}_B \end{bmatrix} \cdot \begin{bmatrix} e^{j\varphi_1} \\ \vdots \\ e^{j\varphi_B} \end{bmatrix} = 0 \quad (51)$$

By finding appropriate \mathbf{d}_i and φ_i , the rotation vector \mathbf{d} is obtained. This is the general idea for the proof. Next, we will provide a detailed proof for this theorem.

Suppose $\mathbf{H}^{(k)}$ has M columns. We can find at least one $E \in \mathbb{N}$, such that $M \in [E^k, (E+1)^k]$. We then sectorize the matrix $\mathbf{H}^{(k)}$ with the scheme shown as Fig. (17). It is worth noticing that the sectorization method is not unique, the only requirement is that the number of columns in each sub-matrix lies within the range of $[E^{k-1}, (E+1)^{k-1}]$. After the sectorization, there are either E or $E+1$ blocks in total.

If the rotation vector \mathbf{d} cannot be found with the idea illustrated in Fig (16), then one of the two sufficient conditions does not hold, i.e., either one of the colored submatrix $\mathbf{H}_i^{(k-1)}$ cannot be rotated to zero or the last row of $\mathbf{H}^{(k)}$ cannot be made zero. Recursive method can be used

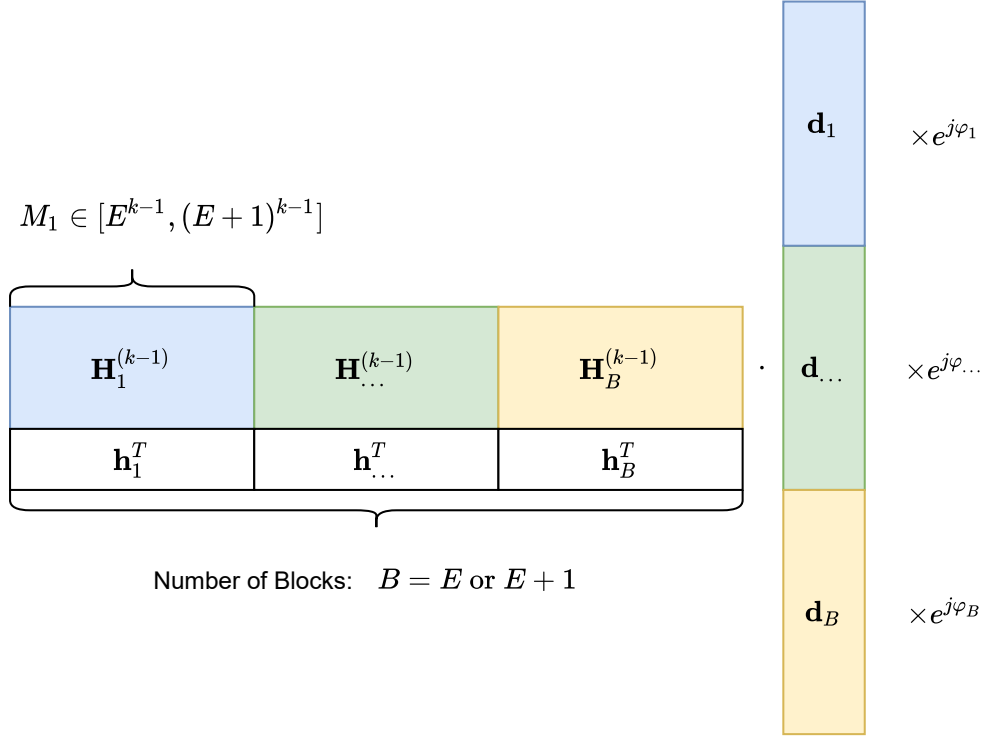


Fig. 17: Submatrix Construction

to calculate the probability that one of the colored sub-matrices cannot be made zero. We focus on analyzing the last row of matrix $\mathbf{H}^{(k)}$ first. Define

$$B \in \{E, E + 1\} : \text{number of blocks} \quad (52)$$

$$M_{1\dots B} \in [E^{k-1}, (E + 1)^{k-1}] : \text{number of columns in one block.}$$

as is also illustrated in Fig. (17). Each block \mathbf{h}_i^T in the last row will be multiplied with correspondent vector block \mathbf{d}_i , which creates a complex Gaussian random variable $\mathbf{h}_i^T \mathbf{d}_i \sim \mathcal{CN}(0, M_i)$. The F-distribution is defined on the normal Gaussian variables. Thus, we should first normalize the variance to 1. This is done in Eq. (53) and Eq. (54) Define

$$\begin{aligned} \forall b \in \{1, \dots, B\}, Y_b &= \mathbf{h}_b^T \mathbf{d}_b \sim \mathcal{CN}(0, M_b) \\ \forall b \in \{1, \dots, B\}, X_b &\sim \mathcal{CN}(0, 1). \end{aligned} \quad (53)$$

Then we have

$$\begin{aligned}
& \sum_{b=1}^B \Pr(|Y_b| > |Y_1| + \dots + |Y_{b-1}| + |Y_{b+1}| + \dots + |Y_B|) \\
& \leq \sum_{b=1}^B \Pr\left(|Y_b| > \sqrt{M_{\min}} \left(\frac{|Y_1|}{\sqrt{M_1}} + \dots + \frac{|Y_{b-1}|}{\sqrt{M_{b-1}}} + \frac{|Y_{b+1}|}{\sqrt{M_{b+1}}} + \dots + \frac{|Y_B|}{\sqrt{M_B}} \right)\right) \\
& = \sum_{b=1}^B \Pr\left(\frac{\sqrt{M_b}}{\sqrt{M_{\min}}} |Y_b| > |X_1| + \dots + |X_{B-1}|\right) \\
& \leq B \times \Pr\left(\frac{\sqrt{M_{\max}}}{\sqrt{M_{\min}}} |X| > |X_1| + \dots + |X_{B-1}|\right) \\
& \leq B \times \left(\frac{M_{\max}}{M_{\max} + M_{\min}}\right)^{\frac{B-1}{2}}.
\end{aligned} \tag{54}$$

For larger M , we have

$$\lim_{M \rightarrow +\infty} \frac{M_{\max}}{M_{\max} + M_{\min}} \leq \lim_{E \rightarrow +\infty} \frac{(E+1)^{k-1}}{2 \times E^{k-1}} = \frac{1}{2} \leq \frac{2}{3}. \tag{55}$$

Therefore,

$$\Pr\left(\#(\varphi_1, \dots, \varphi_B) \cdot \begin{bmatrix} \mathbf{h}_1^T \mathbf{d}_1 & \dots & \mathbf{h}_B^T \mathbf{d}_B \end{bmatrix} \cdot \begin{bmatrix} e^{j\varphi_1} \\ \vdots \\ e^{j\varphi_B} \end{bmatrix} = 0\right) \leq (N+1) \times \left(\frac{2}{3}\right)^{\frac{E-1}{2}}. \tag{56}$$

The bound for the complete probability can then be obtained

$$\begin{aligned}
P_k &= \Pr(\# \mathbf{d}, (\mathbf{H}^{(k)})^T \mathbf{d} = 0) \\
&\leq (E+1) \times P_{k-1} + (E+1) \times \left(\frac{2}{3}\right)^{\frac{E-1}{2}}.
\end{aligned} \tag{57}$$

Based on this inequality, one can define an auxiliary sequence for the upper bound of the probabilities iteratively

$$\begin{aligned}
P'_1 &= (E+1) \times \left(\frac{2}{3}\right)^{\frac{E-1}{2}} \\
P'_k &= (E+1) \times P_{k-1} + (E+1) \times \left(\frac{2}{3}\right)^{\frac{E-1}{2}}.
\end{aligned} \tag{58}$$

By using mathematical induction, it can be proven that $\forall k \in \mathbb{N}, P_k < P'_k$. This sequence can also be written in analytic form by solving difference equation as

$$P'_k = \frac{(E+1)^{k+1} - (E+1)}{E} \left(\frac{2}{3}\right)^{\frac{E-1}{2}}. \tag{59}$$

It can be observed that this upper bound will converge to zero if E tend to infinity.

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