Efficient evaluation of the error probability for pilot-assisted URLLC with Massive MIMO

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Abstract—We propose a numerically efficient method for evaluating the random-coding union bound with parameter s on the error probability achievable in the finite-blocklength regime by a pilot-assisted transmission scheme employing Gaussian codebooks and operating over a memoryless block-fading channel. Our method relies on the saddlepoint approximation, which, differently from previous results reported for similar scenarios, is performed with respect to the number of fading blocks (a.k.a. diversity branches) spanned by each codeword, instead of the number of channel uses per block. This different approach avoids a costly numerical averaging of the error probability over the realizations of the fading process and of its pilot-based estimate at the receiver and results in a significant reduction of the number of channel realizations required to estimate the error probability accurately. Our numerical experiments for both single-antenna communication links and massive multiple-input multiple-output (MIMO) networks show that, when two or more diversity branches are available, the error probability can be estimated accurately with the saddlepoint approximation with respect to the number of fading blocks using a numerical method that requires about two orders of magnitude fewer Monte-Carlo samples than with the saddlepoint approximation with respect to the number of channel uses per block.

Index Terms—Pilot-assisted transmission, finite-blocklength information theory, saddlepoint approximation, ultra-reliable low-latency communication, massive MIMO system

I. INTRODUCTION

Next-generation wireless communication systems will support mission-critical links operating under stringent reliability and latency constraints. Denoted as ultra-reliable low-latency communications (URLLC), this type of links will enable applications such as vehicle-to-everything communication, factory automation [2], autonomous driving [3], and haptic communications [4].

One crucial characteristic of the URLLC traffic is that it often involves small information payloads combined with short packets, i.e., packets consisting of a small number of coded symbols. To understand why short packets are needed, it is worth recalling that the length of a data packet depends on the product of the available bandwidth and the signal duration. In URLLC, the signal duration is limited because of the latency constraint of the targeted applications (e.g., control of automated factories, critical internet-of-things services). The bandwidth is often also limited, because of the need to orthogonalize the transmission of different users to avoid multiuser interference, which has a negative impact on the packet error probability. As pointed out in, e.g., [5], the classic asymptotic performance metrics used to design communication systems, i.e., the ergodic and the outage rates, are unsuitable in the short-packet regime. Thus, a much more precise characterization of the tradeoff between transmission rate and error probability is required.

Finite-blocklength information theory, a field whose relevance to URLLC has become apparent after the seminal works in [6], [7], provides a precise characterization of such tradeoffs, in terms of nonasymptotic upper (achievability) and lower (converse) bounds on the smallest error probability compatible with a given SNR, transmission rate and packet size.

To satisfy the reliability requirements over fading channels in URLLC, under the above-mentioned diversity limitations in both time and frequency, it becomes crucial to leverage on the spatial diversity offered by multiple antennas. A promising approach is to use massive multiple-input multiple-output (MIMO)—a wireless cellular network architecture in which a base station (BS) with a large number of active antennas serves multiple users on the same time-frequency resources [8]. The benefits of massive MIMO are well understood [9], and this technology has been incorporated into the 5G standard.

Focusing on communication over memoryless block-fading channels, we present in this paper a numerically efficient method to evaluate information-theoretic upper bounds on the finite-blocklength error probability achievable in practically relevant scenarios, including massive MIMO deployments. Methods such as the one presented in this paper are necessary since evaluating most of the available information-theoretic error-probability bounds and approximations that are accurate for scenarios of interest for URLLC [10]–[12] is—as we shall see—extremely time consuming. This prevents the use of such expressions within URLLC optimization routines such as resource-allocation and scheduling algorithms.

State of the art: Throughout the paper, we will focus on the upper bound on the error probability obtained by using the random-coding union bound with parameter s (RCUs) proposed in [13]. As discussed in, e.g., [10], this bound...
is particularly suited for transmission over fading channels because it provides achievable results that hold both for the optimal noncoherent maximum-likelihood (ML) decoder, as well as for more practically relevant transmission schemes that rely on pilot-assisted transmission (PAT). For example, PAT schemes include the case in which the acquired channel estimate at the receiver is treated as perfect via the use of a mismatched scaled nearest-neighbor (SNN) decoder [14]. In fact, for this setup, it appears that the RCUs bound is the only known tractable bound on the error probability. Moreover, the normal and saddlepoint approximations seem to be the only reasonable ways to approximate the RCUs bound. The RCUs bound involves the computation of a certain tail probability, which is not known in closed form and needs to be evaluated numerically. If performed naively, this step is time consuming because of the low error probabilities of interest in URLLC. A common approach to circumvent this issue encompasses the following two steps. One starts by noting that, given the realization of the fading channel and of its estimate at the receiver, the random variable whose tail probability is of interest can be written as a sum of independent random variables. Then, one uses the central-limit theorem to approximate this tail probability by a Gaussian tail probability. The resulting approximation, which is typically referred to as normal approximation (see, e.g., [7, Sec.IV]), is, however, not accurate for the error probabilities of interest in the URLLC regime [12], [13]. Furthermore, this approach still requires one to perform a Monte-Carlo averaging over the channel realizations and their estimate at the receiver, which is time consuming.

As shown in, e.g., [13], [15], a much more accurate approximation can be obtained by using the so-called saddlepoint method [16]. Consider a memoryless block-fading channel where each packet is assumed to span $n_b$ fading blocks. Assume that, within each fading block, $n_s$ coded symbols are transmitted. For this scenario, the saddlepoint method can be applied in two different ways: we can fix $n_b$ and perform a saddlepoint expansion with respect to (w.r.t.) $n_s$, i.e., perform an expansion that is accurate when the number of symbols per block is large. Alternatively, we can fix $n_s$ and perform a saddlepoint expansion w.r.t $n_b$, i.e., perform an expansion that is accurate when the number of fading blocks (a.k.a. diversity branches) is large.

For PAT transmission and SNN decoding, the first approximation has been recently studied in [10], [12] for the special case $n_b = 1$. As discussed in [12], this approach yields an approximation on the conditional error probability given the channel and its estimate, which needs then to be averaged w.r.t. the channel realizations and their estimates at the receiver. A different approach to evaluate this conditional error probability is described in [17].

The second approximation was studied in [11], but only for the case of optimal ML decoder. This approximation pertains the unconditional tail probability, and, hence, does not require an additional averaging over the realizations of the channel and its estimate. As we shall see, this makes the numerical evaluation of this approximation computationally efficient.

Contributions: Focusing on independent and identically distributed (i.i.d.) Gaussian codebooks, we present in this paper two saddlepoint approximations on the RCUs for the practically relevant case of PAT and SNN decoding: the one w.r.t. $n_s$ generalizes the one reported in [12] to arbitrary $n_b$ values; the one w.r.t. to $n_b$ generalizes the one reported in [11] to PAT and SNN decoding. Considering the URLLC regime, we then provide a detailed analysis of the accuracy and the computational complexity of both approximations in: i) a single-input single-output (SISO) setup; ii) the uplink of a two-user single-cell massive MIMO network; iii) the uplink of a multi-user multi-cell massive MIMO network. This progression allows us to understand the impact of the number of BS and users in the network on both the accuracy and the numerical complexity of the considered approximations. Our analysis shows that, despite being developed under the assumption of large $n_b$, the saddlepoint w.r.t. $n_b$ is accurate for $n_b$ values as small as 2 in both SISO and multi-user MIMO scenarios. Furthermore, it entails a much smaller computational complexity than the saddlepoint w.r.t. $n_s$. Specifically, whenever $n_b \geq 2$, the number of samples required to evaluate the saddlepoint w.r.t. $n_b$ via Monte-Carlo simulation is typically around 2 orders of magnitude smaller than the number of samples required to evaluate the saddlepoint approximation w.r.t. $n_s$, once the averaging over the channel and its estimate is accounted for. We also show that, for the scenarios considered in the paper, the normal approximation is typically not accurate.

Notation: We denote random vectors and random scalars by upper-case boldface letters such as $X$ and upper-case standard letters, such as $X$, respectively. Their realizations are indicated by lower-case letters of the same font. We use upper-case letters of two special fonts to denote deterministic matrices (e.g., $Y$) and random matrices (e.g., $\mathcal{Y}$). To avoid ambiguities, we use another font, such as $\mathbb{R}$ for rate, to denote constants that are typically capitalized in the literature. The identity matrix of size $a \times a$ is written as $I_a$.

The circularly-symmetric Gaussian distribution is denoted by $CN(0, \sigma^2)$, where $\sigma^2$ denotes the variance. The superscripts $(\cdot)^T$, $(\cdot)^H$, and $(\cdot)^*$ denote transposition, Hermitian transposition, and complex conjugation, respectively. We write $\log(\cdot)$ to denote the natural logarithm, $\|\cdot\|$ stands for the $\ell^2$-norm, $\mathbb{P}[\cdot]$ for the probability of an event, $E[\cdot]$ for the expectation operator, $\mathbb{V}[\cdot]$ for the variance of a random variable, and $Q(\cdot)$ for the Gaussian Q-function. Finally, for two functions $f(n)$ and $g(n)$, the notation $f(n) = o(g(n))$ means that $\lim_{n \to \infty} f(n)/g(n) = 0$ and the notation $f(n) = O(g(n))$ means that $\limsup_{n \to \infty} |f(n)/g(n)| < \infty$.

Organization of the paper: In Section II, we present a finite-blocklength upper bound on the error probability for the SISO Rayleigh block-fading channel. We then introduce different methods to evaluate this bound in the URLLC regime. The extension of our framework to multicell, multiuser massive MIMO networks is presented in Section III. In Section IV, we discuss the accuracy and computational complexity of the methods introduced in Section II with the help of numerical examples. Concluding remarks are provided in Section V.
II. A NON-ASYMPTOTIC UPPER BOUND ON THE ERROR PROBABILITY

A. The SISO System Model

We start by considering a SISO memoryless block-fading channel. Specifically, the channel is assumed to stay constant over the transmission of a block of $n_c$ channel uses and to change independently across blocks. Each transmitted packet spans $n_b$ such fading blocks. Hence, each packet consists of $n_b n_c$ complex-valued symbols. We assume that the first $n_p$ symbols within each block are pilots known to the receiver and the remaining $n_s = n_c - n_p$ symbols are data symbols, as illustrated in Fig. 1. Pilot symbols are used to estimate the fading channel within the corresponding block. The input-output relation corresponding to the pilot transmission phase within block $\ell$ is scaled by the available channel estimate. Mathematically, given the received vector $[y_{1\ell}^T, \ldots, y_{n_s\ell}^T]^T$ and the channel estimates $\{h_1, \ldots, h_{n_b}\}$, the decoded codeword $\hat{x} = [\hat{x}_{1\ell}^T, \ldots, \hat{x}_{n_s\ell}^T]^T$ is determined as follows:

$$\hat{x} = \arg \min_{x} \sum_{\ell=1}^{n_b} \|y_{\ell} - \hat{h}_{\ell} x_{\ell}\|^2.$$  (3)

This decoder, which is known as mismatched SNN decoder, coincides with the ML decoder only when the receiver has perfect channel-state information, i.e., $\hat{h}_{\ell} = h_{\ell}$ for $\ell = 1, \ldots, n_b$. The attractive feature of this decoder is that information-theoretic bounds on its error probability can be approached in practice using good channel codes for the nonfading AWGN channel [10]. In contrast, approaching information-theoretic error-probability bounds for the optimal ML decoder considered in [11] with low-complexity coding schemes is still an open problem (note, however, the recent progress reported in [18]).

B. The RCUs Finite-Blocklength Bound

Like most of the achievability results in information theory, the RCUs bound [13] we shall focus on in this paper is obtained by means of a random-coding argument. Specifically, one evaluates the average error probability averaged over a randomly constructed ensemble of codebooks. In this paper, we consider the i.i.d. Gaussian ensemble, in which each symbol of each codeword is drawn independently from a $CN(0, \rho^2)$ distribution, where $\rho$ models the average transmit power per symbol in the data phase (same power as in the pilot phase). Although suboptimal at finite blocklength [19] in the nonfading SISO case, the i.i.d. Gaussian ensemble is often used in the literature because it leads tractable expressions when applied to PAT, SNN decoding, and multiuser MIMO scenarios.

Specialized to our setup, the RCUs bound results in the following upper bound $\epsilon_{\text{ub}}$ on the packet error probability $\epsilon$:

$$\epsilon \leq \epsilon_{\text{ub}} = P \left[ \log \frac{U}{n_s n_b} + \frac{1}{n_c n_b} \sum_{\ell=1}^{n_b} t_s(X_{\ell}; Y_{\ell}, \hat{H}_{\ell}) \leq R \right].$$  (4)

Here, $U$ is a random variable that is uniformly distributed on $[0, 1]$ and independent of all other quantities,

$$t_s(X_{\ell}; Y_{\ell}, \hat{H}_{\ell}) = \sum_{k=1}^{n_s} t_s(X_{k,\ell}; Y_{k,\ell}, \hat{H}_{\ell})$$  (5)

where $X_{k,\ell}$ and $Y_{k,\ell}$ are the $k$th element of $X_{\ell}$ and $Y_{\ell}$ respectively, and $t_s(x; y, \hat{h})$ is the so-called generalized information density, which, for the case of i.i.d. $CN(0, \rho)$ codebooks and SNN decoding, is given by [12, App. A]

$$t_s(x; y, \hat{h}) = -s \left| y - \hat{h} x \right|^2 + \frac{s |y|^2}{1 + s \hat{h}^2} + \log \left( 1 + s \rho \left| \hat{h} \right|^2 \right).$$  (6)
Finally, \( s > 0 \) is an optimization parameter that can be used to tighten the bound.

In general, no closed-form expression is available for (4). Hence, this probability needs to be evaluated with numerical methods. A naïve implementation of this step results in time-consuming simulations, given the low target error probabilities of interest in URLLC. We next discuss two approaches to compute (4) efficiently: one is based on asymptotic expansions of (4) applied w.r.t. the number of data symbols per block \( n_s \), and the other is based on asymptotic expansions of (4) applied w.r.t. the number of blocks \( n_b \). For each approach, we will present an expansion based on the central-limit theorem, which will result in the so-called normal approximation, and an expansion based on the saddlepoint method.

### C. Asymptotic Expansion w.r.t. \( n_s \)

The idea behind this approach, which for the case \( n_b = 1 \) has been explored in [12], is to analyze first a conditional version of the probability in (4), in which the channel and its estimate within each block are given. Specifically, one focuses on

\[
\epsilon_{ab}\left(\mathbf{h}, \hat{\mathbf{h}}\right) = P\left[\frac{\log U}{n_c n_b} + \frac{1}{n_c n_b} \sum_{\ell=1}^{n_b} \sum_{k=1}^{n_s} t_s(X_{k,\ell}; Y_{k,\ell}, \hat{h}_\ell) \leq R | \mathbf{H} = \mathbf{h}, \hat{\mathbf{H}} = \hat{\mathbf{h}}\right]
\]

(7)

where \( \mathbf{H} = [H_1, \ldots, H_{n_b}]^T \) and \( \hat{\mathbf{H}} = [\hat{H}_1, \ldots, \hat{H}_{n_b}]^T \). Then one seeks an asymptotic approximation to this conditional probability. The approximations that will be presented in this section are easy to evaluate numerically because they depend on quantities that can be evaluated in closed form. To obtain the desired estimate of \( \epsilon_{ab} \), one still needs to perform the averaging

\[
\epsilon_{ab} = \mathbb{E}_{\mathbf{H}, \hat{\mathbf{H}}} [\epsilon_{ab}(\mathbf{H}, \hat{\mathbf{H}})]
\]

over the channel and its estimate numerically.

#### 1) Normal Approximation w.r.t. \( n_s \): One way to numerically approximate (7) is to perform a normal approximation w.r.t. \( n_s \) based on the Berry–Esseen central-limit theorem [20, Ch. XVI.5]. Specifically, note that given \( \mathbf{H} \) and \( \hat{\mathbf{H}} \), the \( n_b n_s \) random variables \( \{t_s(X_{k,\ell}; Y_{k,\ell}, \hat{h}_\ell)\} \) in (7) are conditionally independent and identically distributed within each fading block. Let

\[
I_s(h_\ell, \hat{h}_\ell) = \mathbb{E}[t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell)] \quad \text{and} \quad V_s(h_\ell, \hat{h}_\ell) = \mathbb{V}[t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell)]
\]

denote the mean and the variance of the information density, respectively. Then, conditioned on \( \mathbf{H} = \mathbf{h} \) and \( \hat{\mathbf{H}} = \hat{\mathbf{h}} \) it follows that

\[
\mathbb{E}\left[\sum_{\ell=1}^{n_b} \sum_{k=1}^{n_s} t_s(X_{k,\ell}; Y_{k,\ell}, \hat{h}_\ell)\right] = n_s \sum_{\ell=1}^{n_b} \mathbb{E}\left[t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell)\right]
\]

(9)

and

\[
\mathbb{V}[\sum_{\ell=1}^{n_b} \sum_{k=1}^{n_s} t_s(X_{k,\ell}; Y_{k,\ell}, \hat{h}_\ell)] = n_s \sum_{\ell=1}^{n_b} \mathbb{V}[t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell)]
\]

(11)

Note that in (9) and (11) we set \( k = 1 \) without loss of generality. We now apply the Berry–Esseen central-limit theorem [20, Ch. XVI.5] to the tail probability in (7) to obtain

\[
\epsilon_{ab}(\mathbf{h}, \hat{\mathbf{h}}) = Q\left(\frac{n_s \sum_{\ell=1}^{n_b} I_s(h_\ell, \hat{h}_\ell) - n_c n_b R}{\sqrt{n_s \sum_{\ell=1}^{n_b} V_s(h_\ell, \hat{h}_\ell)}}\right) + \mathcal{O}\left(\frac{1}{n_s}\right).
\]

(13)

Here, both \( I_s(h_\ell, \hat{h}_\ell) \) and \( V_s(h_\ell, \hat{h}_\ell) \) are available in closed-form as

\[
I_s(h_\ell, \hat{h}_\ell) = \left(1 + s \rho |\hat{h}_\ell|^2\right) + (\beta_\ell - \alpha_\ell)
\]

(14)

\[
V_s(h_\ell, \hat{h}_\ell) = (\beta_\ell - \alpha_\ell)^2 + 2\alpha_\ell \beta_\ell (1 - \nu_\ell)
\]

(15)

where

\[
\alpha_\ell = s \left(\rho |\hat{h}_\ell|^2 + \sigma^2\right)
\]

(16)

\[
\beta_\ell = \frac{s}{1 + s \rho |\hat{h}_\ell|^2} \left(\rho |\hat{h}_\ell|^2 + \sigma^2\right)
\]

(17)

\[
\nu_\ell = \frac{s^2 |\rho |\hat{h}_\ell|^2 + \sigma^2 - h_\ell^T \hat{h}_\ell f|^2}{\alpha_\ell \beta_\ell (1 + s \rho |\hat{h}_\ell|^2)}.
\]

(18)

Strictly speaking, according to [20, Ch. XVI.5], for (13) to hold, we need to verify that the third central moment of \( t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell) \) exists for every \( \ell \in \{1, \ldots, n_b\} \). Otherwise, the error term in (13) does not vanish as \( n_s \) grows, and the normal approximation is not applicable. We will show in Section II-C2 that the third central moment of \( t_s(X_{1,\ell}; Y_{1,\ell}, \hat{h}_\ell) \) indeed exists. The normal approximation w.r.t. \( n_s \) is finally obtained by ignoring the \( \mathcal{O}(\cdot) \) term in (13) and by averaging the resulting approximation over \( \mathbf{H} \) and \( \hat{\mathbf{H}} \). This is typically done via Monte-Carlo simulations.

#### 2) Saddlepoint Approximation w.r.t. \( n_s \): Since it is based on a central-limit theorem, the normal approximation is typically accurate only in the regime in which the target rate \( R \) is close to the mean of the information density [11]. However, this regime may be of limited interest in URLLC, since it may correspond to packet error probability values above the URLLC target (see, e.g., [12, Fig. 1]).

A more refined approximation can be obtained by using the so-called saddlepoint method. It results in an error probability expansion given in terms of a leading factor that
decays exponentially with $n_s$ and captures the behavior of the error probability in the large-deviation regime, and a sub-exponential factor, which is obtained by applying a refined normal approximation, and which makes the resulting approximation accurate in the short-packet regime. This method allows one to obtain an approximation that is accurate for a large range of target error probabilities and rates, including the ones relevant in URLLC scenarios.

We now state this approximation. Consider again the conditional probability given in (7). We fix again $k = 1$, without loss of generality, and let $\kappa(\cdot)$ be the cumulant generating function (CGF) of the random variable $-\sum_{\ell=1}^{n_s} t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)$:

$$\kappa(\cdot) = \log \mathbb{E} \left[ e^{-\zeta \sum_{\ell=1}^{n_s} t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)} \right]$$

(19)

$$= \sum_{\ell=1}^{n_s} \log \mathbb{E} \left[ e^{-\zeta t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)} \right].$$

(20)

Note that $\kappa(\cdot)$ depends on $\hat{h}$ and $h$, but we choose not to make this dependence explicit, to keep the notation compact. Each term in (20) admits a closed-form expression. Specifically, let

$$g(\zeta, h_\ell, \hat{h}_\ell) = \mathbb{E} \left[ e^{-\zeta t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)} \right]$$

be the moment generating function (MGF) of the random variable $-t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)$. Then [12, Eq. (56)]

$$g(\zeta, h_\ell, \hat{h}_\ell) = \frac{(1 + (\beta_\ell - \alpha_\ell) \zeta - \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta^2)^{-1}}{1 + s \rho \hat{h}_\ell^2}$$

(21)

where $\alpha_\ell$, $\beta_\ell$, and $\nu_\ell$ were defined in (16), (17), and (18).

Note that, by substituting (22) into (20), one can obtain a closed-form expression not only for $\kappa(\cdot)$, but also for its first and second derivatives, which we shall denote as $\kappa'(\cdot)$ and $\kappa''(\cdot)$, and which we shall need shortly. Specifically, let $\kappa(\cdot) = \log g(\zeta, h_\ell, \hat{h}_\ell)$, so that $\kappa(\cdot) = \sum_{\ell=1}^{n_s} \kappa_\ell(\cdot)$. We have that (see [12, Eqs. (16)–(18)]

$$\kappa_\ell(\zeta) = -\zeta \log(1 + s \rho \hat{h}_\ell^2) - \log(1 + (\beta_\ell - \alpha_\ell) \zeta - \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta^2)$$

(23)

$$\kappa'_\ell(\zeta) = -\log(1 + s \rho \hat{h}_\ell^2) - \frac{(\beta_\ell - \alpha_\ell) - 2 \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta}{1 + (\beta_\ell - \alpha_\ell) \zeta - \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta^2}$$

(24)

$$\kappa''_\ell(\zeta) = \left[ \frac{(\beta_\ell - \alpha_\ell) - 2 \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta}{1 + (\beta_\ell - \alpha_\ell) \zeta - \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta^2} \right]^2 + \frac{2 \alpha_\ell \beta_\ell (1 - \nu_\ell)}{1 + (\beta_\ell - \alpha_\ell) \zeta - \alpha_\ell \beta_\ell (1 - \nu_\ell) \zeta^2}.$$  

(25)

A saddlepoint expansion can be established that the third derivative of the MGF of $-t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)$ exists in a neighborhood of zero for every $\ell \in \{1, \ldots, n_s\}$. Indeed, the saddlepoint expansions presented next depend on the third derivative of the MGF, albeit this term is included in the $O(\cdot)$ and $O(\cdot)$ terms and, hence, does not appear explicitly in the expansions. Specifically, for every $\ell \in \{1, \ldots, n_s\}$, we require that there exist two values $\zeta_\ell < \zeta < \zeta_\ell$, such that

$$\sup_{\zeta < \zeta < \zeta} \left| \frac{d^3}{dz^3} g(\zeta, h_\ell, \hat{h}_\ell) \right| < \infty.$$  

(26)

As shown in [12, Appendix B], this condition holds in our setup with

$$\zeta_\ell = -\sqrt{\frac{(\beta_\ell - \alpha_\ell)^2 + 4 \alpha_\ell \beta_\ell (1 - \nu_\ell)}{2 \alpha_\ell \beta_\ell (1 - \nu_\ell)}}$$

(27)

$$\zeta_\ell = -\sqrt{\frac{(\beta_\ell - \alpha_\ell)^2 + 4 \alpha_\ell \beta_\ell (1 - \nu_\ell)}{2 \alpha_\ell \beta_\ell (1 - \nu_\ell)}} - \alpha_\ell + \beta_\ell.$$  

(28)

This implies in particular that the third moment of $-t_s(X_1, \ell; Y_1, \ell, \hat{h}_\ell)$, which can be obtained by evaluating the third derivative in (26) at $\zeta = 0$, exists—a condition we required to establish the normal approximation in Section II-C1.

By taking $\zeta = \max\{\zeta_1, \ldots, \zeta_{n_s}\}$ and $\zeta = \min\{\zeta_1, \ldots, \zeta_{n_s}\}$, we ensure that (26) holds simultaneously for every $\ell \in \{1, \ldots, n_s\}$. The saddlepoint expansion w.r.t. $n_s$ is stated in the following theorem.

**Theorem 1**: Assume that there exists a $\zeta \in [\zeta, \zeta]$ satisfying $R = -\kappa'(\zeta)n_s/(n_1n_s)$. If $\zeta \in [0, 1]$, then

$$\epsilon_{ab}(\hat{h}, h) = e^{n_s(\kappa(\zeta) - \kappa'(\zeta))} \left[ \Psi_{n,\zeta}(\zeta) + \Psi_{n,\zeta}(1 - \zeta) + O\left(\frac{1}{\sqrt{n_s}}\right) \right]$$

(29)

where

$$\Psi_{b,\zeta}(u) = e^{b \frac{2}{\zeta} \kappa''(\zeta)} Q\left(u \sqrt{b n''(\zeta)}\right).$$

(30)

If $\zeta > 1$, then

$$\epsilon_{ab}(\hat{h}, h) = e^{n_s(\kappa(1) - \kappa'(1))} \left[ \tilde{\Psi}_{n,1}(1, 1) + \tilde{\Psi}_{n,0}(0, 1) + o\left(\frac{1}{\sqrt{n_s}}\right) \right]$$

(31)

where

$$\tilde{\Psi}_{b,a_1, a_2} = e^{b a_1 \left[ -\kappa'(1) - R + \frac{b(\kappa'(1))}{2} \right]} \times Q\left(a_1 \sqrt{b n''(1)} - a_2 \frac{b(\kappa'(1) + R)}{\sqrt{b n''(1)}}\right).$$

(32)

If $\zeta < 0$, then

$$\epsilon_{ab}(\hat{h}, h) = 1 - e^{n_s(\kappa(\zeta) - \kappa'(\zeta))} \left[ \Psi_{n,\zeta}(\zeta) - \Psi_{n,\zeta}(1 - \zeta) + o\left(\frac{1}{\sqrt{n_s}}\right) \right].$$

(33)

**Proof**: Although a direct proof of this theorem is not available in the literature, the desired expansions can be established following steps similar to the ones reported in [20, App. E] for the case of abstract channels and generic mismatch decoding rules and in [11, App. I] for the case of memoryless block-fading channels and ML decoding rule.

We obtain the desired saddlepoint approximation of $\epsilon_{ab}(\hat{h}, h)$ in (7) by omitting the $O(\cdot)$ and the $O(\cdot)$ terms in (29), (31), and (33).
D. Asymptotic Expansion w.r.t. $n_b$

We next present a different approach, which avoids the conditioning w.r.t. $H$ and $H$ and the associated, often time-consuming, Monte-Carlo step. The idea is to exploit directly that the random variables $\{t_s(X_\ell; Y_\ell, \hat{H}_\ell)\}$ in (4) are i.i.d. across the block index $\ell$, and perform an asymptotic expansion of the tail probability in (4) w.r.t. the number of blocks $n_b$.

1) Normal Approximation w.r.t. $n_b$: Proceeding as in Section II-C1, we can obtain an asymptotic expansion—this time directly of $\epsilon_{ub}$ in (4)—by applying the Berry-Esseen central-limit theorem. Since the random variables $\{t_s(X_\ell; Y_\ell, \hat{H}_\ell)\}$ are i.i.d. in $\ell$, it follows that

$$
\mathbb{E}\left[\sum_{\ell=1}^{n_b} t_s(X_\ell; Y_\ell, \hat{H}_\ell)\right] = n_b \mathbb{E}\left[t_s(X_1; Y_1, \hat{H}_1)\right]
$$

and

$$
\text{Var}\left[\sum_{\ell=1}^{n_b} t_s(X_\ell; Y_\ell, \hat{H}_\ell)\right] = n_b \text{Var}\left[t_s(X_1; Y_1, \hat{H}_1)\right].
$$

Here, we have fixed $\ell = 1$ without loss of generality. Furthermore, since $t_s(X_1; Y_1, \hat{H}_1) = \sum_{k=1}^{n_s} t_s(X_{k,1}; Y_{k,1}, \hat{H}_1)$ and since conditioned on $(H_1, \hat{H}_1)$, the random variables $\{t_s(X_{k,1}; Y_{k,1}, \hat{H}_1)\}$ are i.i.d. in $k$, we conclude that

$$
\mathbb{E}\left[\sum_{\ell=1}^{n_b} t_s(X_\ell; Y_\ell, \hat{H}_\ell)\right] = n_b n_s \mathbb{E}\left[\mathbb{E}\left[t_s(X_{1,1}; Y_{1,1}, \hat{H}_1)\bigg| H_1, \hat{H}_1\right]\right]
$$

and

$$
\text{Var}\left[\sum_{\ell=1}^{n_b} t_s(X_\ell; Y_\ell, \hat{H}_\ell)\right] = n_b \left( \mathbb{E}\left[\text{Var}\left[t_s(X_1; Y_1, \hat{H}_1)\bigg| H_1, \hat{H}_1\right]\right]
\right. + \left. \text{Var}\left[\mathbb{E}\left[t_s(X_1; Y_1, \hat{H}_1)\bigg| H_1, \hat{H}_1\right]\right]\right).
$$

It then follows that

$$
\epsilon_{ub} = Q \left( \frac{\sqrt{n_b} \mathbb{E}\left[I_s(H_1, \hat{H}_1)\right] - n_b R}{\sqrt{n_s \text{Var}\left[V_s(H_1, \hat{H}_1)\right] + n_s^2 \text{Var}\left[I_s(H_1, \hat{H}_1)\right]}}, + O\left(\frac{1}{\sqrt{n_b}}\right) \right).
$$

The normal approximation is obtained by neglecting the $O(\cdot)$ term.

Note that, differently from the normal approximation provided in (13), the one provided in (40) applies directly to $\epsilon_{ub}$ and not to the conditional probability $\epsilon_{ub}(h, h)$. Hence, no Monte-Carlo averaging step is required at the end. On the negative side, although $I_s(\cdot, \cdot)$ and $V_s(\cdot, \cdot)$ are available in closed form (see (14) and (15)), the terms $\mathbb{E}\left[I_s(H_1, \hat{H}_1)\right]$, $\mathbb{E}\left[V_s(H_1, \hat{H}_1)\right]$, and $\text{Var}\left[I_s(H_1, \hat{H}_1)\right]$ need to be evaluated numerically, e.g., Monte-Carlo methods. Similarly, the existence of the third central moment of $t_s(X_1; Y_1, \hat{H}_1)$, which is required for (40) to hold, needs to be ensured with numerical methods.

2) Saddlepoint Approximation w.r.t. $n_b$: We now proceed as in Section II-C2 and obtain a saddlepoint approximation of $\epsilon_{ub}$ w.r.t. $n_b$. As pointed out in Section II-C2, to establish a saddlepoint asymptotic expansion, we need that the third derivative of the MGF of the random variables at hand exists in a neighborhood of zero. Specifically, to establish a saddlepoint approximation of $\epsilon_{ub}$ w.r.t. $n_b$, we shall require that, for some $\zeta < 0 < \zeta^*$,

$$
\sup_{\zeta < \zeta^*} \left| \frac{d^3}{d\zeta^3} \mathbb{E}\left[e^{-\zeta t_s(X_1; Y_1, \hat{H}_1)}\right] \right| < \infty.
$$

Unfortunately, differently from (26), the moment-generating function $\mathbb{E}\left[e^{-\zeta t_s(X_1; Y_1, \hat{H}_1)}\right]$ is not known in closed form. Hence, no closed-form expressions for $\zeta$ and $\zeta^*$ are available and these quantities need to be estimated with numerical methods.

To state the saddlepoint approximation, we shall need the CGF $\gamma(\zeta)$ of the random variable $-t_s(X_1; Y_1, \hat{H}_1)$

$$
\gamma(\zeta) = \log \mathbb{E}\left[e^{-\zeta t_s(X_1; Y_1, \hat{H}_1)}\right]
$$

and its first and second derivatives $\gamma'(\zeta)$ and $\gamma''(\zeta)$. Note that, given $H_1$ and $\hat{H}_1$, the random variable $t_s(X_1; Y_1, \hat{H}_1)$ can be decomposed into the sum of $n_s$ conditionally i.i.d. terms (see (5)). Hence,

$$
\gamma(\zeta) = \log \mathbb{E}_{H_1, \hat{H}_1} \left[ \prod_{k=1}^{n_s} \mathbb{E}\left[e^{-\zeta t_s(X_{k,1}; Y_{k,1}, \hat{H}_1)}\bigg| H_1, \hat{H}_1\right]\right]
$$

where the function $g$ was defined in (21). Let $p(\zeta)$ be the MGF of $-t_s(X_1; Y_1, \hat{H}_1)$. Then $p(\zeta)$ and its first and second derivatives with respect to $\zeta$ are given as

$$
p(\zeta) = \mathbb{E}_{H_1, \hat{H}_1} \left[ g(\zeta, H_1, \hat{H}_1)\right]^{n_s}
$$

$$
p'(\zeta) = n_s \mathbb{E}_{H_1, \hat{H}_1} \left[ g(\zeta, H_1, \hat{H}_1)\right]^{n_s-1} g'(\zeta, H_1, \hat{H}_1)
$$

$$
p''(\zeta) = n_s \mathbb{E}_{H_1, \hat{H}_1} \left[ (n_s-1)g(\zeta, H_1, \hat{H}_1)\right]^{n_s-2} g''(\zeta, H_1, \hat{H}_1)^2
$$

$$
\quad + g(\zeta, H_1, \hat{H}_1)^{n_s-1} g''(\zeta, H_1, \hat{H}_1)
$$

where $g'$, $g''$, $p'$ and $p''$ denote the first and second derivatives of $p(\zeta)$ with respect to $\zeta$ of the functions $g$ and $p$, respectively. We can then write $\gamma(\zeta)$ and its first and second derivatives as

$$
\gamma(\zeta) = \log p(\zeta)
$$

$$
\gamma'(\zeta) = \frac{p'(\zeta)}{p(\zeta)}
$$

$$
\gamma''(\zeta) = \frac{p''(\zeta)p(\zeta) - p'(\zeta)^2}{p(\zeta)^2}.\quad(49)
$$
We are now ready to state the saddlepoint expansion w.r.t. $n_b$.

**Theorem 2:** Assume that there exists a $\zeta \in [\zeta, \zeta]$ satisfying $R = -\gamma'(\zeta)/n_c$. If $\zeta \in [0, 1]$ then
\[
\epsilon_{ub} = e^{n_b[\gamma(\zeta) - \zeta \gamma'(\zeta)]} \times \left[ \Phi_{n_b, \zeta}(\zeta) + \Phi_{n_b, \zeta}(1 - \zeta) + o\left( \frac{1}{\sqrt{n_b}} \right) \right]
\] (50)
where
\[
\Phi_{b, \zeta}(u) = e^{b u^2/2} [1 - \gamma''(1 - R + \gamma''(1) + O\left( \frac{1}{\sqrt{n_b}} \right))]
\] (51).

If $\zeta > 1$, then
\[
\epsilon_{ub} = e^{n_b[\gamma(1) - \gamma'(\zeta)]} \times \left[ \Phi_{n_b, 1}(1) + \Phi_{n_b, 0}(1 - 1) + O\left( \frac{1}{\sqrt{n_b}} \right) \right]
\] (52)
where
\[
\Phi_b(a_1, a_2) = e^{b a_1 [1 - \gamma'(1) + R + \gamma''(1)/2]}
\times \left[ 1 - b(\gamma'(1) + R) - b(\gamma'(1) + R) + O\left( \frac{1}{\sqrt{n_b}} \right) \right].
\] (53)

Finally, if $\zeta < 0$,
\[
\epsilon_{ub} = 1 - e^{n_b[\gamma(\zeta) - \zeta \gamma'(\zeta)]} \times \left[ \Phi_{n_b, \zeta}(-\zeta) - \Phi_{n_b, \zeta}(1 - \zeta) + o\left( \frac{1}{\sqrt{n_b}} \right) \right].
\] (54)

**Proof:** The proof follows by combining the steps in the proofs of [11, App. I] and [20, App. E]. Note that the saddlepoint approximation in [11] was developed for Rayleigh SISO block-fading channels but under the assumption of ML decoding. Furthermore, this expansion was provided only for the case $\zeta \in [0, 1]$. The saddlepoint derived in [20] holds only for channels whose input and output belong to finite-cardinality alphabets, but applies to arbitrary mismatched decoding rules and arbitrary values of $\zeta$. Our result is obtained by carefully combining the proof techniques used in these two papers.

We obtain the desired saddlepoint approximation of $\epsilon_{ub}$ in (4) w.r.t. $n_b$ by neglecting the $O(\cdot)$ and the $o(\cdot)$ terms in (50), (52), and (54). Note that, differently from the asymptotic expansion proved in Theorem 1, the one provided in Theorem 2 applies directly to $\epsilon_{ub}$, and not to the conditional probability $\epsilon_{ub}(h, \tilde{h})$. Hence, no Monte-Carlo averaging step is required at the end. On the negative side, the function $\gamma(\cdot)$ and its first and second derivatives are not available in closed form and one needs to resort to numerical methods, such as Monte-Carlo averaging, to evaluate them. Specifically, one needs to evaluate numerically, the expectation over the channel $H_1$ and its estimate $\hat{H}_1$ appearing in the definition of $p(\zeta)$ and of its first and second derivatives in (44), (45), and (46). This is all one needs to evaluate numerically, in order to compute both the normal approximation and the saddlepoint approximation w.r.t. $n_b$. Indeed, according to (47), (48), (49), from $p(\zeta)$ and its first and second derivatives, one obtains $\gamma(\zeta)$ and its first and second derivatives, which are the required quantities to evaluate the saddlepoint approximation in Theorem 1. Furthermore, since $\gamma'(0) = n_c \mathbb{E}[I_s(H_1, \tilde{H}_1)]$ and $\gamma''(0) = n_c \mathbb{E}[V_s(H_1, \tilde{H}_1)] + n_c^2 \mathbb{V}[I_s(H_1, \tilde{H}_1)]$, one can also evaluate the normal approximation in (40) from $p(\zeta)$ and its first and second derivatives evaluated at $\zeta = 0$.

### III. Massive MIMO Network

In this section, we consider a multiuser massive MIMO cellular network with $L$ cells, each served by a BS with $M$ antennas. We assume there are $K$ single-antenna users in each cell and focus on uplink transmission. As in Section II-A, we consider transmission over memoryless block-fading channels and use $n_c$ and $n_b$ to denote the number of symbols per block and the number of blocks spanned by each transmitted packet, respectively. We denote by $H_{\ell, i, k} \in \mathbb{C}^M$ the channel gain vector within the $\ell$th fading block between user $k$ in cell $i$ and the BS in cell $j$. We consider a spatially correlated Rayleigh fading model where $H^i_{\ell, i, k} \sim \mathcal{CN}(0, R^i_{\ell, i, k})$. The normalized trace $\beta^2_{i,k} = \text{tr}(R^i_{\ell, i, k})/M$ determines the average channel gain between user $k$ in cell $i$ and the BS in cell $j$, while the eigenstructure of $R^i_{\ell, i, k}$ describes its spatial channel correlation [9, Sec. 2.2].

#### A. Uplink pilot transmission

The $n_p$-dimensional pilot sequence transmitted by user $k$ in cell $j$ during fading block $\ell$ is denoted by the vector $x_{\ell, j, k} \in \mathbb{C}^{n_p}$. We assume that this vector satisfies $||x_{\ell, j, k}||^2 = n_p \rho$. Furthermore, we assume that the $LK$ users employ mutually orthogonal pilot sequences during each fading block. In particular, we set $n_p = LK$. During the pilot-transmission phase, the received signal $y_{\ell, j}^{(p)} \in \mathbb{C}^{M \times n_p}$ at the BS serving cell $j$ for fading block $\ell$ is given by
\[
y_{\ell, j}^{(p)} = \sum_{k=1}^{K} H_{\ell, i, k}^{i,j} (x_{\ell, j, k}^{(p)})^T + \mathbb{W}_{\ell, j}^{(p)}(k) \quad (55)
\]
where $\mathbb{W}_{\ell, j}^{(p)} \in \mathbb{C}^{M \times n_p}$ is the additive noise with i.i.d. $\mathcal{CN}(0, \sigma^2)$ entries.

We assume that the BS knows $R^i_{\ell, i, k}$, and that it can compute the MMSE channel estimates [9, Sec. 3.2]
\[
\hat{H}_{\ell, i, k} = R_{\ell, i, k}^{i,j} Q_{\ell, i, k}^{j} (y_{\ell, j}^{(p)} (x_{\ell, j, k}^{(p)})^T)
\] (56)
where
\[
Q_{\ell, i, k}^{j} = \left( \sum_{i' = 1}^{L} \sum_{k' = 1}^{K} R_{\ell, i', k'}^{i,j} (x_{\ell, i', k'}^{(p)})^T H_{\ell, i', k'}^{i,j} (x_{\ell, i', k'}^{(p)} + \sigma^2 I_M) \right)^{-1}
\] (57)

#### B. Uplink data transmission

To decode the signal transmitted from user $k$ in cell $j$ over the $\ell$th fading block, which we denote by $x_{\ell, j, k} \in \mathbb{C}^{n_c}$, where $n_c = n_c - n_p$, the BS serving cell $j$ uses the combining vector
\( V_{\ell,j} \in \mathbb{C}^M \) to compute the received vector \( Y_{\ell,j} \in \mathbb{C}^n \) as follows:

\[
Y_{\ell,j} = \left( V_{\ell,j}^H H_{\ell,j}^i \right) x_{\ell,i,j} + \sum_{k'=1}^K \left( V_{\ell,j,k}^H H_{\ell,j,k'}^i \right) x_{\ell,i,k'} + \sum_{i=1, i \neq j}^L \sum_{k'=1}^K \left( V_{\ell,j,k}^H H_{\ell,j,k'}^i \right) x_{\ell,i,k'} + V_{\ell,j,k} W_{\ell,j}.
\]  

(58)

Here, \( W_{\ell,j} \in \mathbb{C}^{M \times n} \) is the additive Gaussian noise on the \( \ell \)th fading block at the BS serving cell \( j \) with i.i.d. \( \mathcal{CN}(0, \sigma^2) \) entries.

We assume that the BS uses multicell-MMSE combiners, i.e.,

\[
V_{\ell,j} = \left( \sum_{i=1}^L \sum_{k=1}^K \hat{H}_{\ell,i,k}^j \left( \hat{H}_{\ell,i,k}^j \right)^H + Z_{\ell,j} \right)^{-1} \hat{H}_{\ell,j}^j.
\]  

(59)

where

\[
Z_{\ell,j} = \sum_{i=1}^L \sum_{k=1}^K \rho p R_{i,k}^j Q_{i,k}^j R_{i,k}^j + \frac{\sigma^2}{\rho} I_M.
\]  

(60)

Note that (58) has the same form as (2). Indeed, set \( Y_{\ell} = Y_{\ell,j,k} \), \( x_{\ell} = x_{\ell,j,k} \), \( H_{\ell} = V_{\ell,j,k} H_{\ell,j,k}^j \), \( \hat{H}_{\ell,j,k}^j \) and \( W_{\ell} = \sum_{k'=1}^K \sum_{k'=1}^K V_{\ell,j,k'}^H H_{\ell,j,k'}^j x_{\ell,j,k'} + \sum_{i=1, i \neq j}^L \sum_{k=1}^K V_{\ell,j,k}^H H_{\ell,i,k} x_{\ell,i,k} + V_{\ell,j,k} W_{\ell,j} \). Note also that, given \( \{ H_{\ell,i,k}^j, \tilde{H}_{\ell,i,k}^j \} \), the entries of the newly defined vector \( W_{\ell} \) are conditionally i.i.d. and follow a \( \mathcal{CN}(0, \sigma^2) \) distribution, with \( \sigma^2 = \sigma^2 \| V_{\ell,j,k} \|^2 + \rho \sum_{k'=1}^K \| V_{\ell,j,k}^H H_{\ell,j,k'}^j \|^2 \) + \rho \sum_{i=1, i \neq j}^L \sum_{k=1}^K \| V_{\ell,j,k}^H \tilde{H}_{\ell,i,k}^j \|^2 \).  

(61)

Hence, we can evaluate the uplink per-user error probability by using the information-theoretic bound in (4) and its normal and saddlepoint approximations discussed in Section II.

IV. NUMERICAL RESULTS AND DISCUSSION

In this section, we report numerical experiments to evaluate the accuracy and the numerical complexity of the introduced normal and saddlepoint approximations. Specifically, we will address the following three questions:

1) In typical scenarios, the number \( n_s \) of symbols per block is much larger than the number \( n_b \) of blocks spanned by a codeword. How large should \( n_b \) be for the approximations w.r.t. to \( n_s \) to be accurate?

2) Is the normal approximation (either w.r.t. \( n_s \) or w.r.t. \( n_b \)) sufficiently accurate in the URLLC regime, or should one use instead the saddlepoint approximations?

3) All the approximations presented in Section II require numerical methods such as Monte-Carlo averaging for the evaluation of terms that are not available in closed form. Which method has lower complexity for a given targeted accuracy?

In the following sections, we perform numerical experiments on SISO and massive MIMO setups to answer these questions.

A. SISO Setup

We start by considering a Rayleigh-fading scenario where the \( \{ H_{\ell} \}_{\ell=1}^{n_b} \) are generated independently from a \( \mathcal{CN}(0,1) \) distribution. Furthermore, we assume ML estimation of the channel at the receiver. Specifically, we set

\[
\hat{H}_{\ell} = \frac{1}{\rho p} \left( x_{\ell}^{(p)} \right)^H Y_{\ell}^{(p)}.
\]  

(62)

We assume that the noise variance \( \sigma^2 \) is equal to 1 for \( \ell \in \{1, \ldots, n_b\} \), and consider a blocklength \( n_p n_c \) of 288 channel uses. The results reported in this section are obtained after an optimization over the parameter \( s > 0 \) in (4) and over the number of pilots \( n_p \) within each block of \( n_c \) channel uses.

Accuracy: In Fig. 2, we report the transmit-power value \( \rho \) needed to achieve an error probability \( \epsilon = 10^{-5} \) for \( R = 0.104 \) bit per channel use; \( n_p \) and \( s \) are optimized.
Moving to the approximations w.r.t. to $n_b$, we note that the normal approximation does not provide accurate results even when $n_b = 48$. The saddlepoint approximation slightly underestimates the required transmit power $\rho$ for $n_b = 1$, but, perhaps surprisingly, returns accurate results already for $n_b$ as small as 2.

Note finally that for the scenario considered in the figure, the required $\rho$ is large for $n_b = 1$ and decreases rapidly until $n_b = 24$, after which it increases again. This behavior can be explained as follows. Increasing $n_b$ for a fixed product $n_b n_c$ yields an increase of the number of diversity branches, which is beneficial, but also of the total number of pilot symbols $n_p n_b$ which is detrimental because it increases the effective rate of the channel code one can use to protect the information bits. The first effect dominates for $n_b \leq 24$, whereas the second effect dominates when $n_b > 24$.

In Fig. 3, we report upper bounds on the packet-error probability $\epsilon$ as a function of $s$ for $R = 0.104$ bit per channel use, $n_b = 8$, $n_c n_b = 288$ and $\rho = 2.79$ dB. Here, we see that all approximations of the RCU bounds with the exception of the normal approximation w.r.t. $n_b$ are accurate for a large range of values of the parameter $s$. Moreover, the parameter $s$ can be optimized to obtain a tighter bound. This figure also shows that a moderate deviation from the optimal choice of $s$ does not affect tightness significantly.

In Fig. 4, we report upper bounds on the packet-error probability as a function of $n_p$ for $R = 0.104$ bit per channel use, $n_b = 8$, $n_c n_b = 288$ and $\rho = 2.79$ dB. Here, in parallel to the result in Fig. 3, all approximations except the normal approximation w.r.t. $n_b$, provide accurate results for a large range of values of $n_p$. This is crucial since $n_p$ is not necessarily optimized to minimize the packet error probability in every system. Thus, being able to use the approximations, regardless of the choice of $n_p$, is highly relevant.

**Complexity:** To address the third question, we assume again that all expectations that need to be evaluated numerically in the normal and saddlepoint approximations are computed via Monte-Carlo averaging. Since the channel and its estimate are jointly Gaussian random variables, as a proxy for numerical complexity, we count the minimum number of real Gaussian random variables that need to be generated to guarantee that the normalized mean squared difference between the error-probability value returned by the considered approximation and the actual error probability bound in (4) is less than a given threshold. Specifically, we compute each approximation $N_{\text{sim}}$ times for the $\rho$ value achieving $\epsilon_{\text{ub}} = 10^{-5}$ in (4), and let $\epsilon_{\text{app}}(N)$ be the error-probability estimate obtained in the $i$th trial, when evaluating the considered approximation for the case in which the Monte-Carlo averaging is performed using $N$ real Gaussian random variables. The normalized mean-squared difference is evaluated as follows:

$$
\epsilon(N) = \frac{1}{N_{\text{sim}}} \sum_{i=1}^{N_{\text{sim}}} \left( \frac{\epsilon_{\text{ub}} - \epsilon_{\text{app}}(N)}{\epsilon_{\text{ub}}} \right)^2.
$$

Clearly the smaller $\epsilon(N)$, the higher the accuracy.

In Fig. 5, we report the smallest value of $N$ necessary to guarantee that $\epsilon(N) \leq 0.5\%$ when $N_{\text{sim}} = 100$, as a function of $n_b$. In the figure, we assumed that $n_b n_c = 288$, $R = 0.104$ bit per channel use, and that a target error probability of $10^{-5}$ needs to be guaranteed for all values of $n_b$. The transmit power is set according to the RCU curve in Fig. 1. As shown in Fig. 5, the saddlepoint approximations w.r.t. $n_b$ requires around $2 \times 10^7$ real Gaussian samples. This is not surprising, since we want to evaluate accurately, via a Monte-Carlo procedure, an error probability of $10^{-5}$. Although the number of random variables that need to be generated increases with $n_b$, this increase translates in a larger value of $N$ only for small values of $n_b$.

On the contrary, the saddlepoint w.r.t. $n_b$ requires only around $10^5$ samples whenever $n_b \geq 4$. This is around two orders of magnitude fewer samples than the saddlepoint w.r.t. $n_s$. This suggests that the complexity of the Monte-Carlo procedure required to evaluate numerically the expectations in (44), (45), and (46), to the level of accuracy considered in this experiment, is much smaller than the complexity of the
Monte-Carlo procedure required to evaluate numerically the expectation over $\mathbf{H}$ and $\mathbf{H}$ in (8).

We do not report the complexity of the normal approximations since they do not achieve the targeted $\epsilon(N)$ because of their limited accuracy.

B. Massive MIMO Setup

Our simulation setup consists of $L$ square cells, each of size $75 \text{ m} \times 75 \text{ m}$, containing $K$ users each. The BSs, which are equipped with a uniform linear array with $M$ antenna elements separated by half a wavelength, are placed in the center of each cell. The antennas and the users are located in the same horizontal plane. Thus, the azimuth angle is sufficient to determine the directivity. We assume that the scatterers are uniformly distributed in the angular interval $[\phi_{i,k} - \Delta, \phi_{i,k} + \Delta]$, where $\phi_{i,k}$ is the nominal angle of arrival of user $k$ in cell $i$ and $\Delta$ is the angular spread, which we set to $\Delta = 25^\circ$. The $(m_1, m_2)$th entry of the matrix $R_{i,k}$ is then given by [9, Sec. 2.6]

$$R_{i,k}^{j} = \frac{\beta_{i,k}^{j}}{2\Delta} \int_{-\Delta}^{\Delta} e^{i\pi(m_1 - m_2)\sin(\phi_{i,k} + \varphi)} d\varphi. \quad (64)$$

Here, $\beta_{i,k}^{j}$ denotes the large-scale fading coefficient measured in dB

$$\beta_{i,k}^{j} = -35.3 - 37.6 \log_{10} \left( \frac{d_{i,k}^{j}}{1 \text{ m}} \right) \quad (65)$$

with $d_{i,k}^{j}$ being the distance between the BS in the cell $j$ and the user $k$ in cell $i$. The communication takes place over a 20 MHz bandwidth with a total receiver noise power of $\sigma^2 = -94$ dBm consisting of thermal noise and a noise figure of 7 dB in the receiver hardware.

In the next two subsections, we extend the accuracy and complexity analysis of the error-probability approximations performed for a SISO link in Section IV-A to the massive MIMO uplink. Since repeating the study carried out for SISO is unfeasible in a multi-cell multi-user MIMO setting, because of complexity constraints, we first focus in Section IV-B1 on a single-cell massive MIMO network with two users. We will then provide in Section IV-B2 an extension of this analysis to the multi-cell multiuser massive MIMO network for the special case in which the number of blocks $n_b$ is equal to 3.

1) Accuracy and Complexity Analysis for the Two-User Case: We consider the uplink of a single-cell massive MIMO network in which the BS serves two users ($L = 1$ and $K = 2$). The distance between the two users and the BS is $d_{1,1} = d_{1,2} = 36.4$ m. The nominal angle of user 1 w.r.t. the BS is $30^\circ$, and the nominal angle of user 2 w.r.t. the BS is $40^\circ$. We also assume that orthogonal pilot sequences are assigned to each user, that $n_p = 2$, and that MMSE spatial combining based on MMSE channel estimation is used at the BS. Finally, we set $n_b n_c = 144$ and $R = 2$ bit per channel use, which corresponds to 288 bits per packet. In Fig. 6 we report the smallest $\rho$ value needed to achieve an error probability of $10^{-5}$, as a function of $n_b$. All curves in the figures are obtained by performing a Monte-Carlo simulation involving the generation of $8 \times 10^{10}$ real Gaussian random variables. We observe that both saddlepoint approximations as well as the normal approximation over $n_b$ agree and are therefore assumed to be accurate for the $n_b$ values considered in the figure, including $n_b = 1$. On the contrary, the normal approximation w.r.t. $n_b$ does not appear to be accurate. Note that, because of the large spatial diversity available in this setup, increasing $n_b$ from 2 to 6 has only a limited benefit in terms of $\rho$, and increasing $n_b$ beyond 6 is actually deleterious, because of the reduction in the number of channel uses per block available for data transmission.

Focusing on both saddlepoint approximations and on the normal approximation w.r.t. $n_s$, we illustrate in Fig. 7, the minimum number of real Gaussian samples that need to be generated in the Monte-Carlo step required in all approxima-
Number of real Gaussian samples $N_e(N)$ are reported in Table I. These values of $n_p$ pilots over each block and set $n_{\text{BS}}$ of radius $L = 4$ massive MIMO network consisting of $K = 10$, $L = 4$, $N_{\text{sim}} = 100$, $R = 2$ bit per channel use, and $\epsilon = 10^{-5}$.

| Approximation          | $\rho$      |
|------------------------|-------------|
| Saddlepoint w.r.t. $n_b$| $-2.257$ dBm|
| Saddlepoint w.r.t. $n_s$| $-2.254$ dBm|
| Normal w.r.t. $n_b$    | $-2.500$ dBm|
| Normal w.r.t. $n_s$    | $-2.311$ dBm|

Table I: Required transmit powers

We finally consider a multi-cell, multi-user massive MIMO scenario. Here, $n_u = 30$, $n_p = 40$, $n_b = 3$, $K = 10$, $L = 4$, $N_{\text{sim}} = 100$, $R = 2$ bit per channel use, and $\epsilon = 10^{-5}$.

Fig. 8: Multi-cell, multi-user massive MIMO scenario. Here, $n_u = 30$, $n_p = 40$, $n_b = 3$, $K = 10$, $L = 4$, $N_{\text{sim}} = 100$, $R = 2$ bit per channel use, and $\epsilon = 10^{-5}$.

To assess the complexity of the two saddlepoint approximation, we take the arithmetic average of the transmit power evaluated with the two saddlepoint approximations as reference transmit power. In Fig. 8, we depict the normalized mean-square difference $e(N)$, defined in (63), as a function of the number of real Gaussian samples $N$ used in the Monte-Carlo step required for both saddlepoint approximations. We can observe that the saddlepoint approximation w.r.t. $n_s$ requires approximately 30 times fewer samples than the saddlepoint approximation w.r.t. $n_b$ to achieve $e(N) \leq 0.5\%$. This is again in accordance with the results reported in Fig. 5 for the SISO case, and Fig. 7 for the single-cell, two-user massive MIMO case.

V. CONCLUSION

We presented numerically efficient methods to evaluate an upper bound on the error probability achievable over SISO and massive MIMO memoryless block-fading channels when pilot-assisted transmission, scaled nearest-neighbor decoding, and i.i.d. Gaussian codebooks are used. Our methods include both normal and saddlepoint approximations w.r.t. to the number of data symbols per block $n_c$, as well as novel normal and saddlepoint approximations w.r.t. the number of fading blocks $n_b$ spanned by each codeword. All approximations involve the numerical evaluations of expectations that are not known in closed form and can be evaluated using Monte-Carlo methods. Our numerical experiments reveal that the saddlepoint approximation w.r.t. to $n_b$ yield accurate estimates of the error probability in URLLC scenarios of practical relevance. Furthermore, it involves a numerical complexity (measured in terms of

Fig. 7: Two-user, single-cell massive MIMO scenario: required number of real Gaussian samples over each block and set BS of radius $L = 4$ massive MIMO network consisting of $K = 10$, $L = 4$, $N_{\text{sim}} = 100$, $R = 2$ bit per channel use, and $\epsilon = 10^{-5}$.
total number of Monte-Carlo samples required to achieve a given accuracy) roughly two orders of magnitude lower than the complexity of the saddlepoint approximations in $n_s$. This holds for a variety of scenarios ranging from SISO to multi-cell, multiuser massive MIMO. Hence, this approximation should be preferred when evaluating error probabilities within URLLC optimization routines such as resource-allocation and scheduling algorithms. The normal approximations w.r.t. $n_b$ and $n_s$ are not viable alternatives as they often provide inaccurate results for the scenarios considered in this paper, at no advantage in terms of complexity compared to the saddlepoint approximation. An explicit characterization of the number of samples needed for the introduced approximations to be accurate is lacking. Indeed, obtaining such a characterization would be an interesting topic for future work.

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