## Message-Passing Decoding of Lattices Using Gaussian Mixtures

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Abstract— A lattice decoder which represents messages explicitly as a mixture of Gaussians functions is given. In order to prevent the number of functions in a mixture from growing as the decoder iterations progress, a method for replacing N Gaussian functions with M Gaussian functions, with  $M \leq N$ , is given. A squared distance metric is used to select functions for combining. A pair of selected Gaussians is replaced by a single Gaussian with the same first and second moments. The metric can be computed efficiently, and at the same time, the proposed algorithm empirically gives good results, for example, a dimension 100 lattice has a loss of 0.2 dB in signal-to-noise ratio at a probability of symbol error of  $10^{-5}$ .

Keywords— Lattice decoding, Gaussian mixtures

#### 1 Introduction

Lattices play a central role in many communication problems. While Shannon used a non-lattice, and non-constructive, Euclidean-space code to compute the capacity of the AWGN channel, recently Erez and Zamir showed that lattice encoding and decoding can also achieve the capacity of the AWGN channel [1]. Similarly, for the problem of communication with known noise, which has applications to multiuser communications and information hiding, lattice codes play an important role [2]. In source coding, lattices may be used for lossy compression of a real-valued source.

To approach theoretical capacities, it is necessary to let the dimension of the lattice or code become asymptotically large. However, for most lattices of interest, the decoder complexity is worse than linear in the dimension, and most studied lattices have small dimension. For example, a frequently cited reference on lattice decoding gives experimental results with a maximum dimension of 45 [3]. Other approaches use trellisbased lattices, which are exponentially complex in the number of states [4]. Historically, finite-field error correcting codes also suffered the same complexity limitation, however, with the advent of iteratively-decoded low-density parity check codes and turbo codes, the theoretical capacity of some binary-input communication channels can be achieved [5].

Recently, a new lattice construction and decoding algorithm, based upon the ideas of low-density parity check codes has been introduced. So-called low-density lattice codes (LDLC) are lattices defined by sparse inverse generator matrix with a pseudo-random construction. Decoding is performed iteratively using messagepassing, and complexity is linear in the block length. Sommer, Feder and Shalvi, who proposed this lattice and decoder, demonstrated decoding with dimension as high as  $10^6$ . However, the experiments considered decoding only for a special communications problem where the transmit power is unconstrained. Comments in their paper suggest that the algorithm did not converge when applied to the more important problem of general lattice decoding [6] [7].

When decoding on the AWGN channel, the LDLC decoder messages are continuous-valued functions, which can be exactly represented by a mixture of Gaussian functions. However, as iterations progress, the number of Gaussians in the mixture grows rapidly. A direct implementation of a decoder which exploits this property is infeasible, and so prior works quantize the messages, ignoring the Gaussian nature of the messages.

In this paper, the LDLC decoder messages are represented as Gaussian functions, and the growth in the number of Gaussians is reduced by a proposed Gaussian mixture reduction algorithm. This algorithm approximates a number of Gaussians N with a smaller number of Gaussians M. The algorithm combines Gaussians in a pair-wise fashion iteratively until a stopping condition is reached. A distance metric, which computes the squared difference between a pair of Gaussian functions, and the single Gaussian which has the same first and second moments, is used.

Section 2 gives a review of the construction and decoding algorithm for low-density lattice codes. If the channel noise is Gaussian, then messages in the decoding algorithm can be represented as a mixture of Gaussian functions. Section 3 gives a method for replacing a pair of Gaussians with a single Gaussian, which is applied to an algorithm which reduces a mixture of NGaussian functions to a mixture of M Gaussians. Section 4 applies this algorithm to the decoding of lowdensity lattice codes, and considers simulation results. Section 5 is the conclusion.

## 2 Low-density Lattice Codes

2.1 Lattices and Lattice Communication

A lattice is a regular infinite array of points in  $\mathbb{R}^n$ . *Definition* An *n*-dimensional lattice  $\Lambda$  is the set of points  $\mathbf{x} = (x_1, x_2, \dots, x_n)$  with

$$\mathbf{x} = G\mathbf{b},\tag{1}$$

where G is an n-by-n generator matrix and  $\mathbf{b} = (b_1, \ldots, b_n)$  is the set of all possible integer vectors,

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 $b_i \in \mathbb{Z}$ .

The following communications system is considered. Let the codeword  $\mathbf{x}$  be an arbitrary point of the lattice  $\Lambda$ . This codeword is transmitted over an AWGN channel with known noise variance  $\sigma^2$ , and received as the sequence  $\mathbf{y} = \{y_1, y_2, \dots, y_n\}$ :

$$y_i = x_i + z_i, i = 1, 2, \dots, n,$$
 (2)

where  $z_i$  is the AWGN. A maximum-likelihood decoder selects  $\hat{\mathbf{x}}$  as the estimated codeword:

$$\widehat{\mathbf{x}} = \arg \max_{\mathbf{x} \in \Lambda} \Pr(\mathbf{y} | \mathbf{x}) \tag{3}$$

The received codeword is correct if  $\mathbf{x} = \hat{\mathbf{x}}$  and incorrect otherwise. The power of the transmitted symbol,  $||\mathbf{x}||^2$  is unbounded. Instead, power is restricted by the volume of the Voronoi region, det(G).

For this system, Poltyrev [8] showed that for sufficiently large n, there exists a lattice for which the probability of error becomes arbitrarily small, if and only if,

$$\sigma^2 \quad < \quad \frac{|\det(G)|^{2/n}}{2\pi e}.\tag{4}$$

Poltyrev's result is in contrast to Shannon's theorem that the capacity of the Gaussian channel, subject to a transmission power constraint, is  $\frac{1}{2}\log(1 + \text{SNR})$ . To achieve capacity while observing the power constraint, the codepoints are on the surface of an *n*-sphere with high probability.

#### 2.2LDLC Definition

Definition A low-density lattice code is a lattice with a non-singular generator matrix G, for which  $H = G^{-1}$  is sparse.

Regular LDLC's have H matrices with constant row and column weight d. Although not necessary, it is convenient to assume that  $\det(H) = 1/\det(G) = 1$ . The non-zero entries are selected pseudo-randomly.

In a magic square LDLC, the absolute values of the d non-zero entries in each row and each column are drawn from the set  $\{h_1, h_2, \ldots, h_d\}$  with  $h_1 \ge h_2 \ge$  $\dots \geq h_d > 0$ . The signs of the entries of H are pseudorandomly changed to minus with probability 0.5. From here, (n, d) magic square LDLC's are considered with  $h_1 = 1$ , and  $h_i = 1/\sqrt{d}$  for  $i = 2, \ldots, d$ . Such codes resulted in only slightly worse performance than other weight sequences [7].

#### $\mathbf{2.3}$ LDLC Decoding

The LDLC decoding algorithm is based upon beliefpropagation, where messages are real functions corresponding to probability distributions on the symbols  $x_i$ . As with decoding low-density parity check codes, the decoding algorithm may be presented on a bipartite graph. There are nd variable-to-check messages  $q_k(z)$ , and *nd* check-to-variable messages  $r_k(z)$ ,  $k=1,2,\ldots,nd.$ 

With an AWGN channel, the initial message is:

$$q_k(z) = \frac{1}{\sqrt{2\pi\sigma}} e^{-\frac{(y_i-z)^2}{2\sigma^2}},$$
 (5)

for the edge k connected to variable node i.

#### **Check Node** 2.3.1

For the check node, note that (1) can be re-written as:

$$H\mathbf{x} = \mathbf{b}, \tag{6}$$

which defines a sparse system of equations:

$$h_{ij_1}x_{j_1} + h_{ij_2}x_{j_2} + \dots + h_{ij_1}x_{j_1} = b_i, \quad (7)$$

for  $i = 1, 2, \ldots, n$ , and  $j_k \in \mathcal{I}_i$ , where  $\mathcal{I}_i$  is the columns

of *H* which have a non-zero entry in position *i*. Let  $\tilde{x}_k = h_k x_k$ , so  $\sum_{i=1}^d \tilde{x}_i = b$ , where *b* is an integer. The input and output messages are  $q_k(z)$  and  $r_k(z)$ , respectively, for  $k = 1, 2, \ldots, d$ . From (7), for an arbitrary  $i, x_k =$ 

$$\frac{b - (h_1 x_1 + \dots + h_{k-1} x_{k-1} + h_{k+1} x_{k-1} + \dots + h_d x_d)}{h_k},$$

or,

$$x_k = \frac{1}{h_k} (b - \sum_{i=1}^{d \setminus k} \widetilde{x}_i).$$
(8)

The output message  $r_k(z)$  can be obtained from the input messages  $q_i(z), i = 1, \ldots, d, i \neq k$  in four steps, Unstretch, Convolution, Extension and Stretch.

Unstretch is multiplication by  $h_k$ . The message for  $\widetilde{x}_i$  is  $\widetilde{q}_k(z)$ ,

$$\widetilde{q}_k(z) = q_k(\frac{z}{h_k}). \tag{9}$$

Convolution The message for  $\sum_{i=1}^{d\setminus k} \tilde{x}_i$  is  $\tilde{r}_k(z)$ . The distribution of the sum of random variables is the convolution of distributions.

$$\widetilde{r}_k(z) = (\widetilde{q}_1 * \cdots * \widetilde{q}_{k-1} * \widetilde{q}_{k+1} * \cdots * \widetilde{q}_d)(z), (10)$$

where \* denotes real-number convolution.

*Extension* is a shift-and-repeat operation for the unknown integer b. Conditioned on a specific value of b, the distribution of  $b - \sum_{i=1}^{d \setminus k} \widetilde{x}_i$  is  $\widetilde{r}_k(b-z)$ . Assuming that b is an arbitrary integer with uniform a priori distribution.

$$\widetilde{r}'_k(z) = \sum_{b=-\infty}^{\infty} \widetilde{r}_k(b-z).$$
(11)

Stretching is multiplication by  $1/h_k$ . Finally the message  $r_k(z)$  which is the message for (8), is obtained as:

$$r_k(z) = \widetilde{r}'_k(h_k z) \tag{12}$$

Note that the above operations are linear and can be interchanged as is required for an implementation.

#### 2.3.2 Variable Node

At variable node i, take the product of incoming messages, and normalize.

Product:

$$\widehat{q}_k(z) = e^{-\frac{(y_i-z)^2}{2\sigma^2}} \prod_{i=1}^{d\setminus k} r_i(z).$$
(13)

Normalize:

$$q_k(z) = \frac{\widehat{q}_k(z)}{\int_{-\infty}^{\infty} \widehat{q}_k(z) dz}.$$
 (14)

#### 2.3.3 Estimated Codeword and Integer Sequence

The check node and variable node operations are repeated iteratively until a stopping condition is reached. Estimate the transmitted by codeword  $\hat{\mathbf{x}}$  by first computing the a posteriori message  $F_i(z)$  for the code symbol  $x_i$  as:

$$F_i(z) = e^{-\frac{(y_i-z)^2}{2\sigma^2}} \prod_{k=1}^d r_k(z).$$
 (15)

Find  $\hat{x}_i$  as:

$$\widehat{x}_i = \arg \max_{z \in \mathbb{R}} F_i(z).$$
(16)

The estimated integer sequence  $\widehat{\mathbf{b}}$  is:

$$\widehat{\mathbf{b}} = \langle H\widehat{\mathbf{x}} \rangle, \tag{17}$$

where  $\langle z \rangle$  denotes the integer closest to z.

#### 2.4 Gaussian Mixture Decoder

When the channel noise is Gaussian, all of the LDLC messages can be described as a mixture of Gaussian functions. From here, "Gaussians" will be used as shorthand for "Gaussian functions".

In this section, it is assumed that a message f(z) is a mixture of N Gaussians,

$$f(z) = \sum_{i=1}^{N} c_i \mathcal{N}(z; m_i, v_i), \qquad (18)$$

where  $c_i \ge 0$  are the mixing coefficients with  $\sum_{i=1}^{N} c_i = 1$ , and

$$\mathcal{N}(z;m,v) = \frac{1}{\sqrt{2\pi v}} e^{-\frac{(z-m)^2}{2v}}.$$
 (19)

In this way, the message f(z) can be described by a list of triples of means, variances and mixing coefficients,  $\{(m_1, v_1, c_1), \ldots, (m_N, v_N, c_N)\}$ 

In describing the Gaussian mixture decoder, initially assume that the input messages to a node consist of a single Gaussian, that is N = 1.

Check node Without loss of generality, consider check node inputs  $k = 1, 2, \ldots, d - 1$  and output d. Each input message  $q_k(z)$  is a single Gaussian  $\mathcal{N}(z; m_k, v_k)$ .

The message  $\tilde{q}_k(z)$  is obtained by multiplying by  $h_k$ , so  $\tilde{q}_k(z) = \mathcal{N}(z; h_k m_k, h_k^2 v_k)$ .

The message  $\tilde{r}_d(z)$  is the convolution of  $\tilde{q}_k(z), k = 1, \ldots, d-1$ . So:

$$\widetilde{r}_d(z) = \mathcal{N}\left(z; \sum_{k=1}^{d-1} h_k m_k, \sum_{i=1}^{d-1} h_k^2 v_k\right).$$
(20)

The message  $\tilde{r}'_d(z)$  is  $\tilde{r}_d(z)$  shifted over all possible integers:

$$\widetilde{r}'_d(z) = \sum_{b=-\infty}^{\infty} \mathcal{N}\left(z; \sum_{k=1}^{d-1} h_k m_k + b, \sum_{k=1}^{d-1} h_k^2 v_k\right).$$

The output message  $r_d(z)$  is obtained by scaling by  $-1/h_d$ , so:

$$r_d(z) = \sum_{b=-\infty}^{\infty} \mathcal{N}\left(z; -\frac{\sum_{k=1}^{d-1} h_k m_k + b}{h_d}, \frac{\sum_{k=1}^{d-1} h_k^2 v_k}{h_d^2}\right).$$

Variable Node. Let the check-to-variable node messages  $r_k(z), k = 1, \ldots, d-1$  be Gaussians  $\mathcal{N}(z; m_k, v_k)$ . For notational convenience, let  $m_0 = y_i$  be the symbol received from the channel at node *i* and let  $v_0 = \sigma^2$  be the channel variance, as in (5). The output message  $q_d(z)$ , the product of these input messages, will also be a Gaussian,

$$q_d(z) = k_d \mathcal{N}(z; m_d, v_d), \qquad (21)$$

where,

$$\frac{1}{v_d} = \sum_{k=0}^{d-1} \frac{1}{v_k},$$
(22)

$$\frac{m_d}{v_d} = \sum_{k=0}^{d-1} \frac{m_k}{v_k} \tag{23}$$

and,

$$k_d = \sqrt{\frac{v_d}{(2\pi)^{d-2} \prod_i v_i}} \exp\left(-\frac{v_d}{2} \sum_{i=0}^{d-2} \sum_{j=i+1}^{d-1} \frac{(m_i - m_j)^2}{v_i v_j}\right).$$

For the general case where the input consists of a mixture of Gaussians, at either the check node or the variable node, the output can be found by conditioning on one element from each input mixture and computing a single output Gaussian. The mixing coefficient for this Gaussian is the product of the input mixing coefficients. Then the output is the mixture of these single Gaussians created by conditioning all input combinations.

The number of Gaussians in each mixture grows rapidly as the iterations progress. At the variable node, if input k consists of a mixture of  $N_k$  Gaussian functions, then the output message will consist of  $N_1N_2\cdots N_{d-1}$  Gaussian functions. At the check node, even if the number of integer shifts is bounded, the number of Gaussian functions in the mixture also grows as  $O(N^{d-1})$ . A naive implementation of this Gaussian mixture decoder is prohibitively complex. The following section proposes a technique for approximating a large number of Gaussians.

### **3** Gaussian Mixture Reduction

This section describes an algorithm which approximates a mixture of Gaussian functions with a smaller number of Gaussian functions.

The algorithm input is a mixture of N Gaussians, f(z), as defined in (18), given as a list of triples. The algorithm output is a list of M triples of means, variances and mixing coefficients,  $\{(m_1^m, v_1^m, c_1^m), \ldots, (m_M^m, v_M^m, c_M^m)\}$  with  $\sum_{i=1}^M c_i^m = 1$ , that similarly forms a Gaussian mixture g(z). With  $M \leq N$ , the output mixture should be a good approximation of the input mixture:

$$f(z) \approx g(z) = \sum_{i=1}^{M} c_i^{\mathsf{m}} \mathcal{N}(z; m_i^{\mathsf{m}}, v_i^{\mathsf{m}}) \,.$$
(24)

First, a metric which describes the error due to replacing a two Gaussians with a single Gaussian is given. Then, this is incorporated into a greedy search algorithm which replaces N Gaussians with M Gaussians.

#### 3.1 Approximating a Mixture of Two Gaussians with a Single Gaussian

Definition The squared difference SD(p||q) between two distributions p(z) and q(z) with support  $\mathcal{Z}$  is defined as:

$$\mathrm{SD}(p||q) = \int_{z \in \mathcal{Z}} (p(z) - q(z))^2 dz \qquad (25)$$

Lemma The squared difference SD(p||q) has the following properties:

- $SD(p||q) \ge 0$  for any distributions p and q.
- SD(p||q) if and only if p = q.
- $\operatorname{SD}(p||q) = \operatorname{SD}(q||p).$

Lemma The squared difference between the Gaussian distributions  $\mathcal{N}(m_1, v_1)$  and  $\mathcal{N}(m_2, v_2)$  is given by  $SD(\mathcal{N}(m_1, v_1), \mathcal{N}(m_2, v_2)) =$ 

$$\frac{1}{2\sqrt{\pi v_1}} + \frac{1}{2\sqrt{\pi v_2}} - \frac{2}{\sqrt{2\pi(v_1 + v_2)}} e^{-\frac{(m_1 - m_2)^2}{2(s_1 + s_2)}}.$$
 (26)

Lemma The squared difference between a single Gaussian  $\mathcal{N}(m, v)$  and a mixture of two Gaussians  $c_1 \mathcal{N}(m_1, v_1) + c_2 \mathcal{N}(m_2, v_2)$ , with  $c_1 + c_2 = 1$ , is:

$$\frac{1}{2\sqrt{\pi v}} + \frac{c_1^2}{2\sqrt{\pi v_1}} + \frac{c_2^2}{2\sqrt{\pi v_2}} - \frac{2c_1}{\sqrt{2\pi(v+v_1)}}e^{-\frac{(m-m_1)^2}{2(v+v_1)}} - \frac{2c_2}{\sqrt{2\pi(v+v_2)}}e^{-\frac{(m-m_2)^2}{2(v+v_2)}} + \frac{2c_1c_2}{\sqrt{2\pi(v_1+v_2)}}e^{-\frac{(m_1-m_2)^2}{2(v_1+v_2)}}.$$
(27)

There is unfortunately no closed-form expression for the minimal squared difference in the previous lemma. However, minimizing the Kullback-Leibler divergence between the single Gaussian distribution and the mixture of two Gaussian distributions is tractable; it simply amounts to moment matching. Therefore, from now we will consider the moment-matched Gaussian approximation.

Lemma The mean m and variance v of a mixture of two Gaussian distributions  $c_1 \mathcal{N}(m_1, v_1) + c_2 \mathcal{N}(m_2, v_2)$ are given by:

$$m = c_1 m_1 + c_2 m_2 \tag{28}$$

$$s = c_1(m_1^2 + v_1) + c_2(m_2^2 + v_2) -c_1^2 m_1^2 - 2c_1 c_2 m_1 m_2 - c_2^2 m_2^2.$$
(29)

Let  $\overline{t}_i$ , i = 1, 2 denote the triple  $(m_i, v_i, \overline{c}_i)$ , where  $\overline{c}_1 + \overline{c}_2$  is not necessarily one, and let the normalized triple be  $t_i = (m_i, v_i, c_i/(c_1 + c_2))$ . The single Gaussian which satisfies the property of the Lemma is denoted as:

$$t = \mathrm{MM}(\bar{t}_1, \bar{t}_2), \tag{30}$$

where t = (m, v, 1), with m and v as given in (28) and (29).

Definition The Gaussian quadratic loss GQL(p) of a probability distribution p is defined as the squared difference between p and the Gaussian distribution with the same mean m and variance v as p:

$$GQL(p) = SD(p || \mathcal{N}(m, s)).$$
(31)

*Corollary* The Gaussian quadratic loss of a mixture of two Gaussian distributions,

$$GQL(t_1, t_2) = SD(c_1 \mathcal{N}(m_1, v_1) + c_2 \mathcal{N}(m_2, v_2) || \mathcal{N}(m, v)),$$

is obtained evaluating (27), with m and v as given in (28) and (29).

# **3.2** Approximating N Gaussians with M Gaussians

Here, we use the results from the previous subsection and propose an algorithm which approximates a mixture of N Gaussians with a mixture of N Gaussians.

Input: list  $\mathcal{L} = \{t_1, t_2, \ldots, t_N\}$  of N triples describing a Gaussian mixture, and two stopping parameters,  $\theta$  the allowable one-step error (measured by GQL) and M, the maximum number of allowable Gaussians in the output.

Algorithm

- 1. Initialize the current search list, C, with the input list:  $C \leftarrow L$ .
- 2. Initialize the current error,  $\theta^{c}$ , to the minimum GQL between all pairs of Gaussians:

$$\theta^{\mathsf{c}} = \min_{t_i, t_j \in \mathcal{C}, i \neq j} \operatorname{GQL}(t_i, t_j)$$

- 3. Initialize length of current list,  $M^{c} = N$ .
- 4. While  $\theta^{c} < \theta$  or  $M^{c} > M$ :
  - (a) Determine the pair of Gaussians  $(t_i, t_j)$  with the smallest GQL:

$$(t_i, t_j) = \arg \min_{t_i, t_j \in \mathcal{C}, i \neq j} \operatorname{GQL}(t_i, t_j)$$

(b) Add the single Gaussian with the same moment as  $t_i$  and  $t_j$  to the list:

$$\mathcal{C} \leftarrow \mathcal{C} \cup MM(t_i, t_j).$$

- (c) Delete  $t_i$  and  $t_j$  from list:  $\mathcal{C} \leftarrow \mathcal{C} \setminus \{t_i, t_j\}$ .
- (d) Recalcuate the minimum GQL:

$$\theta^{\mathsf{c}} = \min_{t_i, t_j \in \mathcal{C}, i \neq j} \operatorname{GQL}(t_i, t_j).$$

(e) Decrement the current list length:  $M^{c} \leftarrow M^{c} - 1$ .

5. Algorithm output: list of triples C

Note that two conditions must be satisfied for the algorithm to stop. That is, the one-step error may be greater than the threshold  $\theta$  if the minimum number of Gaussians is not yet met. On the other hand, the number of output Gaussians may be less than M, if the one-step error is sufficiently low.

# 4 Gaussian-Mixture Reduction Applied to LDLC Decoding

In this section, the Gaussian mixture reduction algorithm of Section 3 is applied to the LDLC decoding algorithm described in Section 2.4.

At the check node, observe that the message  $\tilde{r}_k(z)$ , as given in (10), can be computed recursively with  $a_k(z)$  and  $b_k(z)$  defined as:

$$a_1(z) = \widetilde{q}_1(z), \tag{32}$$

$$a_k(z) = a_{k-1}(z) * \widetilde{q}_k(z), k = 2..., d-1, (33)$$

and,

$$b_d(z) = \widetilde{q}_d(z), \tag{34}$$

$$b_k(z) = b_{k+1}(z) * \widetilde{q}_k(z), k = d - 1, \dots, 2.$$
 (35)

Then  $\tilde{r}_k(z)$  is found using a variation on the forwardbackward algorithm as:

$$\widetilde{r}_1(z) = b_2(z), \tag{36}$$

$$(z) = a_{k-1}(z) * b_{k+1}(z),$$

$$\widetilde{r}_d(z) = a_{d-1}(z).$$
 (37)  
(38)

(27)

The Gaussian mixture reduction algorithm is applied after the computation (33) and (35), for each k. For example if  $\overline{a}_k(z)$  is the mixture produced by applying the Gaussian mixture reduction algorithm to  $a_k(z)$ ,

$$\overline{a}_k(z) = \text{GMR}(a_k(z)), \tag{39}$$

then the forward recursion of the check node function may be stated as:

$$\overline{a}_1(z) = \widetilde{q}_1(z), \tag{40}$$

For  $k = 2, 3, \ldots, d - 1$ :

$$a_k(z) = \overline{a}_{k-1}(z) * \widetilde{q}_k(z), \tag{41}$$

$$\overline{a}_k(z) = \text{GMR}(a_k(z)), \tag{42}$$

and similarly for the backward recursion.

Similarly at the variable node, the product (13) can be decomposed into a forward and backward recursion. In this case as well, the Gaussian mixture reduction algorithm is applied after each step of the recursion.

In the Gaussian mixture reduction algorithm, it is desirable to repeat step 4 as long as the current reduced Gaussian function g(z) (represented by  $\mathcal{C}$ ) remains a good approximation of the input function f(z)(represented by  $\mathcal{L}$ ). In practice, it was found that using a "local" stopping condition of a threshold on the one-step error was sufficient to give a good "global" approximation  $f(z) \approx g(z)$ . In many cases, f(z) was wellapproximated by a single Gaussian, which was found by the proposed algorithm.

However, using an error threshold alone does not always restrict the number of output Gaussians, an important goal of the mixture reduction algorithm. Thus, a second stopping condition, which requires that the number of Gaussians be lower than some fixed threshold, is also enforced. Thus, the Gaussian combining may continue while  $M^{c} > M$ , even if the one-step error threshold has been exceeded. In practice, this did not appear to have a detrimental result for a wide range of symbol-error rates.

Simulation results comparing the proposed decoder with the quantized decoder [7] are shown in Fig. 1. A LDLC with n = 100, d = 5 was used. The symbol error rate of a cubic lattice used for transmission is labeled "Uncoded." The horizontal axis is the difference between the channel noise variance and the Poltyrev capacity,  $1/2\pi e$ , in dB.

For the parameter selection  $\theta = 0.5, 1.0$ , and  $M \leq 6$ , it was found that the proposed algorithm performed with a slight performance loss when the probability of symbol error was greater than  $10^{-5}$ . For example, with  $\theta = 0.5$  and M = 6, the loss at a symbol error rate of  $10^{-5}$  is less than 0.1 dB. For lower symbol error rates, an error floor appears. It may be helpful to consider this error floor as analogous to quantization error floors which appear in the decoding of low-density parity check codes when insufficient quantization levels are used.

Complexity In the Gaussian mixture reduction algorithm, the primary complexity is computing the initial error, which requires computing the GQL between N pairs, a complexity of  $O(N^2)$ . In the Gaussian mixture decoder, the primary complexity the pairwisecomputation of the outputs, which is  $O(M^2)$ . These



Figure 1: Symbol error rate of proposed Gaussian mixture (GM) decoder vs. quantized decoder for n = 100, d = 5.

numbers N and M are random variables which depend upon the nature of the messages, and the effectiveness of the Gaussian mixture reduction algorithm. In the simulations the maximum value of M was 6, and  $N \leq kM^2$ , where k is the constant number of integer shifts, k = 3 was used in the simulations. On the other hand, the complexity of the quantized algorithm is dominated by a discrete Fourier transform of size  $1/\Delta$  where  $\Delta$  is the quantization bin width,  $\Delta = 1/128$ was used in the simulations. It is difficult to directly make comparisons of the computational complexity of the two algorithms.

The memory required for the proposed algorithm, however, is significantly superior. The proposed algorithm requires storage of 3M (for the mean, variance and mixing coefficient), for each message, where  $M \leq 6$ . The quantized algorithm, however used 1024 quantization points for each message.

## 5 Conclusion

LDLC codes can be used for communication over unconstrained power channels. In this paper, we proposed a new LDLC decoding algorithm which exploits the Gaussian nature of the decoder messages. The core of the algorithm is a Gaussian mixture reduction method, which approximates a message by a smaller number of Gaussians. As a result, the LDLC algorithm which tracks the means, variances and mixing coefficients of the component Gaussians, rather than using quantized messages, was tractable. It was shown by computer simulation that this algorithm performs nearly as well as the quantized algorithm, when the dimension is n = 100, and the probability of symbol error is greater than  $10^{-5}$ .

#### References

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