# KEY RATE RELIABLE LOWER BOUND FOR BB84 PROTOCOL

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#### Abstract

Following George u. a. (2021) and Winick u. a. (2018) papers, a key rate reliable lower bound for the standard four states BB84 protocol is calculated in condition of:

- 1. depolarizing channel;
- 2. efficiency mismatch at receiver detectors.

The simulation includes the announcement of the bases between A and B, the sifting phase and a key map that stores the choice of the bases of A and B in a classic register. The goal of the simulation is to minimize the relative entropy using a semidefinite program (SDP) before and after the application of a pinching channel. The result is a lower bound for the key rate.

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# 1 Theory

Quantum key distribution (QKD) allow two distant parties, generally called Alice and Bob, to share a key with unconditional security. QKD protocols communicate upon a quantum channel and a classical channel. Quantum channel is untrusted, public and unauthenticated so an eavesdropper, typically called Eve, can manipulate the message. Classical channel is public and authenticated, so Alice and Bob recognize each other.

#### Source replacement scheme

In prepare and measure (P&M) schemes, Alice choose randomly the signal states  $\{|\phi_i\rangle\}_{i=1}^N$  with probability  $p_i$  in which encode the key bits and sends to Bob the sequence through a quantum channel; Bob randomly choose a detection basis. In the entangled based (EB) schemes, a bipartite state  $\psi_{AB}$  is sent to Alice and to Bob; Both Alice and Bob choose randomly the measurement basis.

Alice and Bob communicate via classical channel to perform error correction and privacy amplification in order to extract the key.

EB schemes are more natural to describe the joint system between Alice and Bob. P&M schemes can be transposed into EB schemes using the so-called source-replacement scheme.

In the source-replacement scheme Alice creates the bipartite entangled state(Ferenczi und Lütkenhaus (2012))

$$|\psi\rangle = \sum_{i=1}^{N} \sqrt{p_i} |i\rangle_A |\phi_i\rangle_{A'},$$

where  $\{|i\rangle\}_{i=1}^N$  is a basis of the Hilbert space  $\mathcal{H}^N$ . Alice keeps A and sends A' to Bob via quantum channel  $\mathcal{E}_A B$  which transforms  $A' \to B$ . For example, if the channel introduces depolarization with a probability p on a qubit  $\rho$ ,

$$\mathcal{E}_{AB}(\rho) = \sum_{i=1}^{4} E_i \rho E_i^{\dagger} \tag{1}$$

where, using the Pauli matrices  $\{\sigma_i\}$ ,  $E_1 = \sqrt{1 - \frac{3}{4}p}\mathbb{I}_2$ ,  $E_2 = \sqrt{p/4}\sigma_x$ ,  $E_3 = \sqrt{p/4}\sigma_y$  and  $E_4 = \sqrt{p/4}\sigma_z$ . The state shared between Alice and Bob is denoted by  $\rho_{AB}$ .

#### Post-procecessing

The post-processing phase can be divided into:

- 1. parameter estimation: Alice and Bob compute statistics on AN announced random data set and agree if proceed or aborting the communication;
- 2. public string announcement: Alice and Bob choose part of their bits, respectively  $\tilde{A}_i$  and  $\tilde{B}_i$  and communicate it over the classical channel. Alice and Bob also keep private part of the key in the sequences  $\bar{A}_i$  and  $\bar{B}_i$ ;
- 3. sifting phase: using the public string, Alice and Bob agree only on qubit in which they have used the same basis;
- 4. key map: Alice performs a key map function able to extract a key based on the public and private part of the keys of both Alice and Bob;
- 5. error correction: using classical methods, Alice and Bob agree on shorter but correlated key symbols. As proven by Shannon, this leakage is bounded by the mutual information  $I(A:B) = H(A) + H(B) H(AB) = 1 leak_{EC}$  where H is the Shannon entropy.

#### Protocol modelling

Eve can use whatever strategy to extract as much information as possible from  $\rho_{AB}$ :

- 1. individual attacks, Eve acts on a single qubit and performs her treatment on it before the classical post-processing phase;
- 2. collective attacks, Eve performs measurements after the post-processing phase on all the qubits she has gained.

The simulation works under the assumption of identically and independently distributed (i.i.d.) collective attack, so it is possible to simulate only one qubit reducing drastically the dimension of the system.

The aim is to estimate the best attack Eve can do on  $\rho_{AB}$ . Following George u. a. (2021) and Winick u. a. (2018), the constraints of the problem are given by Alice and Bob POVM  $\{P_i^A\}$  and  $\{P_i^B\}$ 

$$\operatorname{Tr}\left[\left(P_i^A \otimes P_j^B\right)\rho_{AB}\right] = p_{ij}.\tag{2}$$

For P&M schemes there are additional constraints, due to the fact that Eve cannot gain any information about the system *A*, so

$$\operatorname{Tr}\left[\left(\Theta_{i}^{A}\otimes\mathbb{I}^{B}\right)\rho_{AB}\right]=\theta_{i},\tag{3}$$

where  $\{\Theta_i\}_{i=1}^{\dim \mathcal{H}_A}$  is a complete set of hermitian operators of  $\mathcal{H}_A$ . For EB schemes the  $p_{ij}$ ,  $\theta_i$  constraints defines a set

$$S^{EB} = \left\{ \rho_{AB} | \operatorname{Tr} \left[ (P_i^A \otimes P_j^B) \rho_{AB} \right] = p_{ij} \right\}$$

and for P&M schemes

$$S^{P\&M} = \left\{ \rho_{AB} | \operatorname{Tr} \left[ (P_i^A \otimes P_j^B) \rho_{AB} \right] = p_{ij}, \operatorname{Tr} \left[ (\Theta_i^A \otimes \mathbb{I}^B) \rho_{AB} \right] = \theta_i \right\}.$$

To maintain a more general approach, I refer both to  $S^{EB}$  and  $S^{AB}$  with S.

The public string announcement is represented by the following Kraus representation for Alice

$$K_a^A = \sum_{\bar{a}} \sqrt{P_{a,\bar{a}}^A} \, |a\rangle_{\tilde{A}} \, |\bar{a}\rangle_{\bar{A}}$$

and for Bob

$$K_b^B = \sum_{\bar{t}_*} \sqrt{P_{b,\bar{b}}^B} \, |b\rangle_{\tilde{B}} \, |\bar{b}\rangle_{\bar{B}}$$

where (a,b) are the public parts of Alice and Bob key and  $\bar{a},\bar{b}$  are the private one. The two POVM acts on the system AB

$$\rho_{A\tilde{A}\tilde{A}B\tilde{B}\bar{B}}^{1} = \sum_{a,b} (K_{a}^{A} \otimes K_{b}^{B}) \rho_{AB} (K_{a}^{A} \otimes K_{b}^{B})^{\dagger} = \mathcal{A}(\rho_{AB}). \tag{4}$$

Then sifting phase is performed. Let **A** be the set of all announcements that are kept. Then the sifting phase can be seen as a projector

$$\Pi = \sum_{(a,b)\in\mathbf{A}} |a\rangle\langle a|_{\tilde{A}} \otimes |b\rangle\langle b|_{\tilde{B}} \to \rho_{A\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^2 = \frac{\Pi\rho_{A\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^1\Pi}{p_{pass}}$$
(5)

where  $p_{pass} = \text{Tr}[\rho_{AB}\Pi]$  gives the probability of passing the post selection.

Alice chooses a key map, *i.e.* a function that, depending on Alice public and private parts  $(a, \bar{a})$  and Bob announcement b, returns a value  $g(a, \bar{a}, b)$ . The value of the key map can be thought to be stored in a classical register called R, so the key map operator is represented by an isometry

$$V = \sum_{a,\bar{a},\bar{b}} |g(a,\bar{a},b)\rangle_R |a\rangle_{\bar{A}} |\bar{a}\rangle_{\bar{A}} |b\rangle_{\bar{B}} \to \rho_{RA\bar{A}\bar{A}B\bar{B}\bar{B}}^3 = V \rho_{A\bar{A}\bar{A}B\bar{B}\bar{B}}^2 V^{\dagger}.$$
 (6)

If *R* contains *J* symbols, the register turns out to be a classical register after applying a pinching channel on  $R \to Z^R$ 

$$\mathcal{Z}(\rho) = \sum_{j=1}^{J} (|j\rangle\langle j|_{R} \otimes \mathbb{I}) \rho(|j\rangle\langle j|_{R} \otimes \mathbb{I}) \to \rho_{Z^{R}A\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{4} = \mathcal{Z}(\rho_{RA\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{3})$$
 (7)

where the identity acts over all the other dimensions.

If the post-selection process involves only a sifting phase and base announcement, then the formula reduces only to system A and B without other registers. Suppose  $\{Z_i^A\}$  is the Alice key map POVM, then (9) reduces to

$$f(\rho_{AB}) = D\left(\rho_{AB} \bigg\| \sum_{i} (Z_i^A \otimes \mathbb{I}_B) \rho_{AB} (Z_i^A \otimes \mathbb{I}_B) \right). \tag{8}$$

#### Reliable lower bound

The function to be minimized is the relative entropy  $D(\rho \| \sigma)$  (also called Kullback–Leibler divergence), which is a measure of how one probability distribution is different from a second. In particular, the simulation consider the divergence

$$f(\rho_{AB}) = D\left(\rho_{RA\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{3} \middle\| \rho_{Z^{R}A\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{4}\right). \tag{9}$$

having defined the states as

$$\rho_{RA\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{3} = G(\rho_{AB}),$$

$$\rho_{Z^{R}A\tilde{A}\tilde{A}B\tilde{B}\tilde{B}}^{4} = \mathcal{Z}(G(\rho_{AB}))$$

where the *G* map is defined by

$$G(\rho_{AB}) = V \Pi \mathcal{A}(\rho_{AB}) \Pi V^{\dagger}$$

and  $\mathcal{Z}$  map is defined in (7).

The key rate is

$$K = p_{pass} \cdot \min_{\rho_{AB}} \left( H(Z^R | E\tilde{A}\tilde{B}) \right) - p_{pass} \cdot leak^{EC} = \min_{\rho_{AB}} D\left( G(\rho_{AB}) || \mathcal{Z}(G(\rho_{AB})) \right) - leak^{EC}$$

$$(10)$$

where  $Z^R$  is the classical register shared by Alice and Bob, and  $E\tilde{A}\tilde{B}$  is the portion of the system where Eve has complete access without any disturbance on A and B.

Quantum relative entropy is a convex function, thus  $f(\rho)$  is a convex function in  $\rho$  and can be minimized using a semidefinite program (SDP). Due to computer imprecision, the algorithm may exit before reaching the true lower bound. Thus, the problem is converted to its dual in order to find the upper bound. Then, a strategy can be split the algorithm into two steps: a primal problem and a dual problem. The primal problem with m variables is defined as

$$\min_{X \in S^m} : \langle C, X \rangle_{S^m}$$
subject to : $\langle A_i, X \rangle_{S^n} = b_i, \quad i = 1, ..., m$ 

$$X > 0. \tag{11}$$

The dual of this problem is

$$\max_{y \in \mathbb{R}^n} : \sum_{i} b_i y_i$$
subject to: 
$$\sum_{i} A_i y_i \le C.$$
(12)

These two steps have two different scopes:

step 1: estimation of the optimal attack Eve can apply to the qubits which gives an upper bound on the key rate; step 2: study the inverse problem finding the upper bound.

Let's see the step 1 algorithm:

#### Algorithm 1

Result: near optimal attack

**Parameters:**  $\epsilon > 0$ ,  $\rho_0 \in S$ ,  $maxIteration \in \mathbb{N}$  and i = 0

- 1. compute  $\Delta \rho = \arg \min_{\delta \rho} \operatorname{Tr}[\delta \rho \nabla f(\rho_i)]$ . subject to  $\Delta \rho + \rho_i \in S$ ,  $\operatorname{Tr}[\Delta \rho] = 0$  and hermiticity  $\Delta \rho = (\Delta \rho)^{\dagger}$ ;
- 2. **if**  $\text{Tr}[\Delta \rho \nabla f(\rho_i)] < \epsilon$  **then** go to step 2
- 3. find  $\lambda \in [0,1]$  that minimizes  $f(\rho_i + \lambda \Delta \rho_i)$ ;
- 4. set  $\rho_i \rightarrow \rho_i + \lambda \Delta \rho$ ;
- 5. set  $i \rightarrow i + 1$ ;
- 6. **if** i > maxIteration **then** go to step 2

I want to stress the fact that the estimation of the lower bound does not depend on the state shared by AB. Indeed, the initial density operator  $\rho_0$  can be written as

$$\rho_0 = \sum_{j=1}^{j} \tilde{p_l} \tilde{\Gamma}_l + \sum_{k=1}^{\dim \mathcal{H}_{AB}^2 - j} \omega_k \Omega_k$$

where  $\{\tilde{\Gamma}_l\}_{l=1}^j \cup \{\tilde{\Omega}_k\}_{k=1}^{\dim \mathcal{H}_{AB}^2 - j}$  is complete set of hermitian operators for the system AB. The first sector of this basis,  $\{\tilde{\Gamma}_l\}_{l=1}^j$ , is an orthonormal set of l < m hermitian operators which is created applying the Gram-Schmidt process to  $\{\Gamma_j\}_{j=1}^m$ , which are the m operators generating the constraints in (2) and (3);  $\{\Omega_k\}_{k=1}^{\dim \mathcal{H}_{AB}^2 - j}$  are used to complete the basis.

In step 2, the reliable lower bound is found as a maximization problem. Relative entropy is a highly non-linear function, so it may be difficult to find the dual problem; A solution can be computing the dual of the linearization of the original problem SDP, which reads

$$\max_{\vec{z} \in S^*(\rho)} \vec{z} \cdot \vec{\gamma}, \quad S^*(\rho) = \{z \in \mathbb{R} | \sum_{i=1}^m z_i \Gamma_i \le \nabla f(\rho) \}$$

where  $\vec{\gamma} = \{\gamma_1, \dots, \gamma_m\}$  are the *m*-constraints of the SDP. Finally, the reliable lower bound for the key rate is

$$\beta(\rho) = f(\rho) - \text{Tr}[\rho \nabla f(\rho)] + \max_{\vec{z} \in S^*(\rho)} \vec{z} \cdot \vec{\gamma},$$

in which is present the result  $f(\rho)$  of the step 1.

# 2 Code development

My simulation is performed on a BB84 protocol, which is a QKD protocol in which bits are encoded using two mutually unbiased bases (MUB)  $Z = \{|0\rangle, |1\rangle\}$  and  $X = \{|+\rangle, |-\rangle\}$  where  $|+\rangle = \frac{|0\rangle + |1\rangle}{\sqrt{2}}$  and  $|-\rangle = \frac{|0\rangle - |1\rangle}{\sqrt{2}}$ . Bit 0 can be encoded by the states  $\{|0\rangle, |+\rangle\}$ , while bit 1 by the states  $\{|1\rangle, |-\rangle\}$ .

To implement the theory into a PC simulation I have written a python class called QKD, which automatically elaborates the operators needed if not provided in the initialization of the object. Once the simulation object is created, it is possible to apply quantum channel passing the operators which represent the channel action on the qubit sent from A to B; Then, the class allow computing primal and dual SDP.

The convex optimization is performed in python by the library cvxpy using the solver MOSEK which can be obtained under license here. Otherwise, cvxpy allow the use of license-free solver called CVXOPT or SCS.

The following subsections will bring some examples, in particular:

- 1. EB standard BB84 protocol;
- 2. EB standard BB84 protocol with efficiency mismatch at detector.

#### EB standard BB84 protocol

Suppose both Alice and Bob perform measurements in the Z basis with probability  $p_z$  and in the X basis with probability  $p_x$ . Alice and Bob POVM are then symmetric

$$\begin{split} P_0^A = & p_z | 0 \rangle \langle 0 |_A, \\ P_1^A = & p_z | 1 \rangle \langle 1 |_A, \\ P_+^A = & (1-p_z) | + \rangle \langle + |_A, \\ P_-^A = & (1-p_z) | - \rangle \langle - |_A \end{split} \qquad \begin{split} P_0^B = & p_z | 0 \rangle \langle 0 |_B, \\ P_1^B = & p_z | 1 \rangle \langle 1 |_B, \\ P_+^B = & (1-p_z) | + \rangle \langle + |_B, \\ P_-^B = & (1-p_z) | - \rangle \langle - |_B. \end{split}$$

 $P_-^A=(1-p_z)|-\rangle\langle-|_A$  Then the public announcement is composed by the Kraus operators

$$\begin{split} K_1^A = & \sqrt{P_0^A} \, |0\rangle_{\tilde{A}} \, |0\rangle_{\tilde{A}} + \sqrt{P_1^A} \, |0\rangle_{\tilde{A}} \, |1\rangle_{\tilde{A}} \,, \quad K_2^A = \sqrt{P_+^A} \, |1\rangle_{\tilde{A}} \, |0\rangle_{\tilde{A}} + \sqrt{P_-^A} \, |1\rangle_{\tilde{A}} \, |1\rangle_{\tilde{A}} \,, \\ K_1^B = & \sqrt{P_0^B} \, |0\rangle_{\tilde{B}} \, |0\rangle_{\tilde{B}} + \sqrt{P_1^B} \, |0\rangle_{\tilde{B}} \, |1\rangle_{\tilde{B}} \,, \quad K_2^B = \sqrt{P_+^B} \, |1\rangle_{\tilde{B}} \, |0\rangle_{\tilde{B}} + \sqrt{P_-^B} \, |1\rangle_{\tilde{B}} \, |1\rangle_{\tilde{B}} \,, \end{split}$$

the sifting phase is defined by the projector

$$\Pi = |0\rangle\langle 0|_{\tilde{A}} \otimes |0\rangle\langle 0|_{\tilde{B}} + |1\rangle\langle 1|_{\tilde{A}} \otimes |1\rangle\langle 1|_{\tilde{B}}$$

and the key map is

$$V = |0\rangle_R \, |0\rangle\langle 0|_{\bar{A}} + |1\rangle_R \, |1\rangle\langle 1|_{\bar{A}}$$

with identity acting on the other subsystems. These operators define the *G* map and the problem is completely set. If the choice of these Kraus operators, sifting and isometry is standard, as in this case, calling the python class QKD will create all the structure described above

```
from src import qkd
import numpy as np
sim = qkd.QKD(dim_a=2, dim_b=2, n_of_signal_states=4, # of signal states
list_states_a=[qkd.zero, qkd.one, qkd.plus, qkd.minus], list_of_prob_a=[.25,.25,.25],
list_states_b=[qkd.zero, qkd.one, qkd.plus, qkd.minus], list_states_b=[.25,.25,.25])
```

Then, Eve presence in the communication is quantified by the quantum bit error rate Q (QBER) by means of the depolarization  $p = 2 \cdot Q$  using (1)

```
qber, key_primal, key_dual, key_th = np.linspace(0., .12, 15), [], [],
for ii in qber:
    # for theorical curve and leak^{EC}
hp = qkd.binary_entropy(ii)

# apply quantum channel
sim.apply_quantum_channel(qkd.depolarizing_channel(2*ii))

gamma = []
for jj in sim.povm: # get AB povm from QKD
    gamma.append(np.trace(jj @ sim.rho_ab)) # rho_AB is authomatically calculated by QKD

# set contraints
```

```
sim.set_constraints(gamma, sim.povm)

# compute
sim.compute_primal(epsilon=1e-11, maxit=50, finesse=100, "MOSEK") # finesse=finesse of the interval
for \lambda\in[0,1]
sim.compute_dual("MOSEK")

# getting result
key_th = 1 - hp - hp # hp is the leak from error correction
key_primal = sim.primal_sol - hp # hp is the leak from error correction
key_dual = sim.dual_sol - hp # hp is the leak from error correction
```

#### EB BB84 with efficiency mismatch on Bob's detectors

Consider the previous system with efficiency mismatch between the detection of bit 0 and 1 at Bob side. To treat the no-click event, the Hilbert space of system B must contain also the no-click event so  $dim\mathcal{H}_B = 3$ ; For this reason, if  $\eta$  is the efficiency of the detector, Bob POVM become

$$\begin{split} P_1^B = & p_z |0\rangle \langle 0|_B \oplus 0, \\ P_2^B = & p_z \eta \, |1\rangle \langle 1|_B \oplus 0, \\ P_3^B = & (1-p_z)|+\rangle \langle +|_B \oplus 0, \\ \end{split} \qquad \qquad P_5^B = & \mathbb{I}_B - \sum_{i=1}^4 P_i^B, \text{ where 0 indicates the no-click event,} \end{split}$$

while Alice POVM remains unchanged with respect to the previous case.

In the public string announcement, Bob's announcement Kraus representation is

$$K_1^B = \sqrt{P_1^B} \, |0\rangle_{\tilde{B}} \, |0\rangle_{\tilde{B}} + \sqrt{P_2^B} \, |0\rangle_{\tilde{B}} \, |1\rangle_{\tilde{B}} \,, \quad K_2^B = \sqrt{P_3^B} \, |1\rangle_{\tilde{B}} \, |0\rangle_{\tilde{B}} + \sqrt{P_4^B} \, |1\rangle_{\tilde{B}} \, |1\rangle_{\tilde{B}} \,, \quad K_3^B = \sqrt{P_5^B} \, |2\rangle_{\tilde{B}} \, |0\rangle_{\tilde{B}} \,.$$

With opportune adjustments, the sifting phase and the key map have the same structure as the standard BB84 simulation, so the *G* map can be defined inside QKD class. Depolarizing channel is also adapted to the 3-dimensional Hilbert space using the Kraus representation Ramzan und Khan (2011)

$$\begin{array}{lll} E_1 = \sqrt{1-p} \cdot \mathbb{I}_B & E_2 = \sqrt{p/8} \cdot Y & E_3 = \sqrt{p/8} \cdot Z \\ E_4 = \sqrt{p/8} \cdot Y^2 & E_5 = \sqrt{p/8} \cdot YZ & E_6 = \sqrt{p/8} \cdot Y^2Z \\ E_7 = \sqrt{p/8} \cdot YZ^2 & E_8 = \sqrt{p/8} \cdot Y^2Z^2 & E_9 = \sqrt{p/8} \cdot Z^2 \end{array}$$

where

$$Y = \begin{pmatrix} 0 & 1 & 0 \\ 0 & 0 & 1 \\ 1 & 0 & 0 \end{pmatrix}, \quad Z = \begin{pmatrix} 1 & 0 & 0 \\ 0 & \omega & 0 \\ 0 & 0 & \omega^2 \end{pmatrix}, \quad \omega = e^{2\pi i/3}.$$

The code is then

```
from src import qkd
2 # define states
states = [qkd.zero, qkd.one, qkd.plus, qkd.minus]
5 # with null clicks event
  states_b = [np.append(i, [0]) for i in states]
8 # ... iteration over efficiencies
  for depolarizing_prob in [0., .05, .1]:
       # iteration
10
       for efficiency in np.linspace(0.,1.,15):
           # ... definition of POVMB, KB, sifting and key_map
           # new simulation
           sim = qkd.QKD(dim_a=2, dim_b=3, n_of_singla_states=4,
14
               list_states_a=states, list_of_prob_a=[.25,.25,.25,.25]
15
               list\_states\_a = states\_b \;,\;\; list\_of\_prob\_a = [\;.25\,,.25\,,.25\,,.25\,] \;,
16
                             =POVMB,
               povm_b
               kraus_b
                             =KB,
18
19
               sifting_phase =sifting,
                             =V,
               key map
               announcement_register_b=[[1,0,0], [0,1,0], [0,0,1]])
           # apply quantum channel
           sim.apply_quantum_channel(qkd.depolarizing_channel(depolarizing_prob, 3))#dimension=3
2.4
25
           # compute primal and dual problem
           sim.compute_primal(1e-11, 1000, 20, solver_name="MOSEK")
26
           sim.compute_dual(solver_name="MOSEK")
```

### 3 Results

#### Standard BB84

In table 1 are reported the specific of time duration, primal problem result, dual problem result, precision (difference between primal and dual result) and tightness (difference between dual result and theoretical value). Looking at the table the precision and the tightness are well below the part per thousand.

In figure are represent the result of each step and the theoretical value of the secret key rate for the standard BB84.

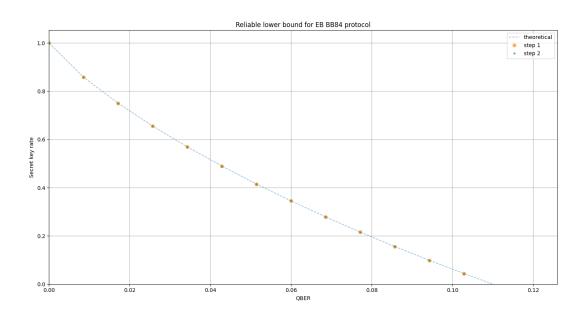


Figure 1: Standard BB84 secret key rate lower bound.

QBER	CPU time[s]	theoretical	primal	dual	precision	tightness
0.000	43.469	1.000	1.000	1.000	3.04e-09	3.04e-09
0.008	28.991	0,928	0.928	0.928	3.09e-07	3.10e-07
0.017	28.492	0,874	0.874	0.874	7.61e-04	7.58e-04
0.025	43.003	0,827	0.827	0.827	1.69e-04	1.69e-04
0.034	28.564	0,784	0.784	0.784	1.28e-06	1.28e-06
0.042	28.795	0,744	0.744	0.743	9.33e-04	9.28e-04
0.051	42.643	0,707	0.707	0.707	3.47e-04	3.47e-04
0.060	41.923	0,672	0.672	0.672	2.30e-06	2.30e-06
0.068	28.265	0,639	0.639	0.639	2.66e-06	2.66e-06
0.077	29.915	0,607	0.607	0.607	5.36e-04	5.34e-04
0.085	42.638	0,577	0.577	0.577	5.24e-06	5.22e-06
0.094	28.698	0,549	0.549	0.549	1.03e-04	1.03e-04
0.102	29.094	0,522	0.522	0.521	7.35e-04	7.32e-04
0.111	44.381	0,495	0.495	0.495	1.26e-04	1.26e-04
0.120	44.381	0,470	0.470	0.470	4.92e-06	4.92e-06

Table 1: exiting results from each simulation step

## Efficiency mismatch

In the tables 2, 3 and 4 are reported the results for each value of the efficiency  $\eta$  for depolarization probability of the quantum channel p = 0, 0.05, 0.1.

The figure represents the values of the dual problem solution which is the reliable lower bound for the secret key rate. The calculations of precision (difference between primal and dual problem solution) for all depolarization

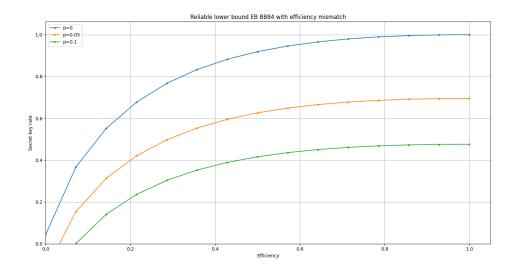


Figure 2: figure secret key rate changing the efficiency in [0,1] for three possible depolarization probabilities p = 0, 0.05, 0.1.

probabilities p = 0, 0.05, 0.1 are below the part per billion.

productives p = 0,0.00,0.1 are below the part per children									
Efficiency	CPU time [s]	primal	dual	precision	Efficiency	CPU time [s]	primal	dual	precision
0.000	73.824	0.045	0.045	8.55e-09	0.000	265.325	0.040	0.040	9.85e-07
0.071	272.240	0.368	0.368	4.37e-09	0.071	245.214	0.323	0.323	2.17e-07
0.142	136.864	0.552	0.552	3.96e-09	0.142	205.293	0.481	0.481	1.04e-07
0.214	71.155	0.677	0.677	1.91e-08	0.214	203.671	0.589	0.589	2.77e-07
0.285	167.272	0.767	0.767	9.30e-09	0.285	217.498	0.666	0.666	1.85e-07
0.357	223.856	0.833	0.833	1.12e-08	0.357	234.874	0.722	0.722	3.35e-07
0.428	288.619	0.882	0.882	3.49e-08	0.428	229.832	0.763	0.763	1.27e-06
0.500	238.178	0.918	0.918	7.22e-09	0.500	244.347	0.794	0.794	5.20e-07
0.571	244.388	0.945	0.945	9.03e-09	0.571	220.040	0.817	0.817	2.64e-07
0.642	170.951	0.965	0.965	1.28e-09	0.642	206.758	0.834	0.834	2.76e-07
0.714	175.240	0.979	0.979	5.97e-08	0.714	203.992	0.846	0.846	7.54e-07
0.785	596.072	0.989	0.989	7.65e-09	0.785	214.197	0.854	0.854	6.43e-07
0.857	187.316	0.995	0.995	1.71e-08	0.857	211.922	0.859	0.859	5.35e-07
0.928	192.420	0.998	0.998	4.32e-09	0.928	211.338	0.862	0.862	1.80e-06
1.000	457.271	0.999	0.999	8.38e-08	1.000	210.020	0.863	0.863	3.46e-07

Table 2: depolarization probability p=0

Table 3: depolarization probability p = 0.05

Efficiency	CPU time [s]	primal	dual	precision
0.000	212.277	0.036	0.036	3.87e-08
0.071	212.460	0.287	0.287	1.19e-06
0.142	209.758	0.427	0.427	3.61e-06
0.214	211.222	0.522	0.522	1.04e-06
0.285	210.727	0.589	0.589	1.30e-06
0.357	210.353	0.638	0.638	1.50e-06
0.428	208.802	0.675	0.675	1.68e-06
0.500	209.674	0.702	0.702	2.68e-06
0.571	208.272	0.722	0.722	4.37e-06
0.642	208.197	0.737	0.737	5.96e-06
0.714	209.123	0.747	0.747	3.79e-06
0.785	212.709	0.754	0.754	3.07e-06
0.857	212.752	0.759	0.759	6.16e-06
0.928	217.413	0.761	0.761	5.01e-06
1.000	280.391	0.762	0.762	1.73e-05

Table 4: depolarization probability p = 0.1

### 4 Self-evaluation

The standard BB84 simulation is in complete accordance with the theoretical curve and with over one part per thousand precision.

Efficiency mismatch simulation confirms the precision of the algorithm far over one part per billion.

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