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Master Thesis

Quantum Communication through finite dimensional lossy channels

Quantum capacity analysis of 4-dimensional
Multi-level Amplitude Damping channels

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Abstract

Quantum Information Theory (QIT) is a relatively new field of physics that is attracting a growing number of researchers. Although some elements of QIT can be found at the dawn of Quantum Mechanics (e.g. the EPR paradox), its first formal definition dates back to the 1960's[Gor62] (Bell published his paper on the EPR paradox in 1964[Bel64]).

The new framework provided by the QIT allowed for the study of open systems, i.e. systems subject to external noise; this is accomplished by defining any quantum process, or quantum channel, as a **Linear, Completely Positive and Trace-preserving** (LCPT) map, which interprets the set of initial density matrices as input and outputs new density matrices in a (possibly) different Hilbert space. This paved the way for an information theoretic approach to the study of quantum processes relevant to Quantum Computation and Quantum Communication, in the same way Shannon's noisy channel theorem did for classical communication and computation.

The question then becomes, in this context, what quantity can we interpret as "information"? And how well is this information preserved during the quantum process? In other words, how well can we store and reliably transmit information through a given quantum channel? It turns out that, in the quantum case, the amount of "stored information" of a system is given by the Von Neumann entropy, while there are a number of quantities called "capacities" that describe the amount of transmittable information through a given channel based on the nature of the information to be transmitted and the availability of additional resources (e.g. classical communication lines, entanglement...). This thesis work focuses on the study of some of these capacities, with a particular focus on quantum capacity and 2-way quantum capacity, for a specific family of channels, called Multilevel Amplitude Damping (MAD) channels.

MAD channels are the d -dimensional extension of the well understood 2-dimensional case of Amplitude Damping Channels. The foundations for how to treat these kind of channels were already laid out in an earlier work by V. Giovannetti and S. Chessa[CG21a], which explored the channels in the case of $d = 3$.

The quantum capacity is generally not computable, as its underlying quantity (the coherent information) is not subadditive when one considers multiple uses of a channel. The only cases in which the quantum capacity is computable correspond to degradable channels (i.e. there exists a LCPT map that connects the output state of the system to the output state of the environment, which takes on the role of the induced noise) and anti-degradable channels (i.e. there exists a LCPT map that connects the output state of the environment to the output state of the system; in this case the quantum capacity is 0). When neither of these conditions is satisfied, it is usually only possible to set upper and lower bounds on the quantum capacity.

In [CG21a] it has been shown that it is possible to find the value for the quantum

capacity of 3-dimensional MAD channels even in regions where those channels are neither degradable nor anti-degradable. This was accomplished by employing the pipeline inequalities, which led to the derivation of some monotonicity properties for the capacity that allowed to extend the values of the capacity found at the border of the degradability regions to the outside of these regions. Expanding upon this work, some general results for d -dimensional MAD channels were found (i.e. composition, monotonicity properties, inverse maps, degradability conditions), which enabled a generalization of the technique used to find the quantum capacity in non-degradability zones in [CG21a]. This paved the way for the numerical evaluation of the quantum capacity for various configurations of 4-dimensional MAD's, again, even where those channels are not degradable.

The 2-way quantum capacity of 3-dimensional MAD channels was also studied, as the literature was lacking in that regard, and lower bounds (which are always computable) were found.

The results of the thesis are to be added in the context of those related to the computation of the quantum capacity for qudit channels (which suggest that these channels could be more convenient with respect to their qubit counterparts) and to those related to the limits of repeater-less quantum communication. The techniques illustrated in the thesis can be ideally expanded to bigger dimension, the limit being the computational power of the machine used to arrive at the values of the capacities.

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Nomenclature

Abbreviations

Abbreviation	Definition
ADC	Amplitude Damping Channel
LCPT	Linear, Completely Positive, Trace-preserving
MAD	Multi-level Amplitude Damping
QCT	Quantum Communication Theory
QIT	Quantum Information Theory
ReMAD	Resonant MAD

1

Introduction

Theory of Quantum Information and Communication

In this chapter, the reader will be presented with all the tools necessary to understand this thesis work. The theories that encompass these tools are called Quantum Information Theory (QIT) and Quantum Communication Theory (QCT). The strength of these theories lies in the fact that they provide a framework that makes it possible to study the behavior of *open systems*, which will be defined in the following section. The reason behind the necessity of a coherent structure describing open system is a practical one: in any real world scenario, physicists will find themselves having to deal with *noise*, which is something that only an open system can take into account. The work of Shannon [Sha48] taught the scientific community how information is transmitted through a classical, noisy communication line. Quantum Information Theory allows for the study of *quantum noise*, which in turns paves the way for the development of, to name a few, Quantum Computers (whose components are very susceptible to external perturbations). Something something large number prime factoring something something quantum teleportation

2.1 Open Systems

Usually, in quantum mechanics, the time evolution of the system in a state can be summarized by defining a unitary operator, dependent on the Hamiltonian of the system. This is done through the Schrödinger equation:

$$i\hbar \frac{\partial}{\partial t} |\psi(t)\rangle = H |\psi(t)\rangle, \quad (2.1.1)$$

where $|\psi(t)\rangle$ is the state of the system at time t and H is the Hamiltonian operator. Solving (2.1.1) yields:

$$|\psi(t')\rangle = \exp [iH(t' - t)/\hbar] |\psi(t)\rangle \quad (2.1.2)$$

This picture only works for time-independent Hamiltonians¹ and for *closed systems*, i.e. systems whose interactions are confined within said systems. This is not generally the case in a real-world setting, as physicists often have to deal with noise, which can be represented as an interaction of the *laboratory system* (i.e. the system upon which one is able to make measurements) with an external system, usually called *environment*. Before defining a framework capable of describing this scenario, a precise definition of *quantum state* is needed. The most generic quantum state of a system \mathcal{H} takes the form of a *density matrix operator* ρ , which satisfies the following properties:

- ρ is a hermitian operator, $\rho^\dagger = \rho$.
- $\text{tr } \rho = 1$.
- $\text{tr } \rho^2 \leq 1$; if $\text{tr } \rho^2 = 1$ then the state is called a *pure state*, i.e. $\exists |\psi\rangle \in \mathcal{H} : \rho = |\psi\rangle\langle\psi|$. A state that is not pure is called *mixed state*.
- ρ is positive-semidefinite, $\rho \geq 0$.

The set of all the elements that satisfy the properties above is denoted as $\sigma(\mathcal{H})$. In order to characterize the interaction of a system \mathcal{H}_S with an environment \mathcal{H}_E , one needs to consider *joint states* $\rho_{SE} \in \sigma(\mathcal{H}_S \otimes \mathcal{H}_E)$, which represent the total states of the system laboratory-environment. In general, given a joint state $\rho_{AB} \in \sigma(\mathcal{H}_A \otimes \mathcal{H}_B)$, one could obtain the *reduced density matrices*, representing the states of the systems \mathcal{H}_A and \mathcal{H}_B through the use of the *partial trace* operation, defined as:

$$\begin{aligned}\rho_B &= \text{tr}_A \rho_{AB} \equiv \sum_i {}_A\langle\phi_i|\rho_{AB}|\phi_i\rangle_A \\ \rho_A &= \text{tr}_B \rho_{AB} \equiv \sum_i {}_B\langle\psi_i|\rho_{AB}|\psi_i\rangle_B\end{aligned}\tag{2.1.3}$$

where $\{|\phi_i\rangle_A\}_i$ and $\{|\psi_i\rangle_B\}_i$ are, respectively, bases of \mathcal{H}_A and \mathcal{H}_B . Then, let $\rho_S \in \sigma(\mathcal{H}_S)$ and let the state of the environment be a pure state² which will be denoted as $\tau_E = |0\rangle_E\langle 0| \in \sigma(\mathcal{H}_E)$, then the state of the composite system $\mathcal{H}_S \otimes \mathcal{H}_E$ is the tensor product³:

$$\rho_{SE} \equiv \rho_S \otimes |0\rangle_E\langle 0| \in \sigma(\mathcal{H}_S \otimes \mathcal{H}_E).\tag{2.1.4}$$

In section 2.3, the mechanism behind the evolution of ρ_S , in the setting illustrated above, will be explained.

¹A solution for time-dependent Hamiltonians (which might not commute between themselves at different times) exists and is called *Dyson series*, see [Dys49]

²This is because, since the environment represents anything that is external to the system A , even if τ_E was, in fact, a mixed state, it could always be *purified*, see 2.2

³In order to lighten the notation, when possible, the symbol \otimes of tensor product will usually be omitted, as indexes referring to specific quantum systems make it redundant.

2.2 Composite quantum states

Before moving on, it might be useful to label composite quantum states in a more precise way. Given a joint system $\mathcal{H}_S \otimes \mathcal{H}_R$, its quantum states ρ_{SR} can be classified in one of the three following categories:

- *Factorized states* present the form:

$$\rho_{SR} = \rho_S \tau_R, \quad (2.2.1)$$

- *Separable states* present the form:

$$\rho_{SR} = \sum_i p_i \rho_S^{(i)} \tau_R^{(i)}, \quad (2.2.2)$$

where $\{p_i\}_i$ forms a probability distribution, $\sum_i p_i = 1$. Separable states are convex combination of factorized states; furthermore, the convex combination of separable states is still a separable state. Note that (2.2.2) reduces to (2.2.1) if $p_i = \delta_{i,i_0}$

- *Non separable (entangled) states* are all the joint states that do not satisfy (2.2.2)

Purification

Let ρ_S be a state in \mathcal{H}_S , suppose that ρ_S is a mixed state and let \mathcal{H}_R be another system, called, in this case, *auxiliary system*. Through the *purification* process, it is **always** possible to build a composite state $\rho_{SR} \in \sigma(\mathcal{H}_S \otimes \mathcal{H}_R)$, $\text{tr}_R \rho_{SR} = \rho_S$ such that ρ_{SR} is a pure state. In fact, in its diagonal basis $\{|\psi_i\rangle_S\}_i$, ρ_S can be written as:

$$\rho_S = \sum_i p_i |\psi_i\rangle_S \langle \psi_i| \quad (2.2.3)$$

where p_i forms a probability distribution. Then, fixing $\dim \mathcal{H}_R = \dim \mathcal{H}_S$ and choosing a basis $\{|i\rangle_R\}_i$ for \mathcal{H}_R , one could build the pure state:

$$\begin{aligned} |\psi_i\rangle_{SR} &\equiv \sum_i \sqrt{p_i} |\psi_i\rangle_S |i\rangle_R \\ \rho_{SR} &= |\psi_i\rangle_{SR} \langle \psi_i| \end{aligned} \quad (2.2.4)$$

which clearly satisfies $\text{tr}_R \rho_{SR} = \rho_S$.

2.3 Stinespring Dilation

Since the composite system $\mathcal{H}_S \otimes \mathcal{H}_E$ is a closed system, its state in (2.1.4) will evolve following a unitary operator $U_{SE} : \mathcal{H}_S \otimes \mathcal{H}_E \mapsto \mathcal{H}_S \otimes \mathcal{H}_E$, which acts on density matrices as:

$$\rho_{SE} \mapsto U_{SE} (\rho_S |0\rangle_E \langle 0|) U_{SE}^\dagger. \quad (2.3.1)$$

Then, combining (2.3.1) with (2.1.3), one could find the evolved state of the system S by defining a *superoperator* (i.e. an object that maps linear operators into linear operators) Φ such that:

$$\rho_S \mapsto \Phi(\rho_S) = \text{tr}_E \left[U_{SE} (\rho_S |0\rangle_E \langle 0|) U_{SE}^\dagger \right]. \quad (2.3.2)$$

This equation defines a very specific family of maps: in fact any map Φ which can be cast in the form of (2.3.4) needs to satisfy the following properties:

- Φ is **Linear**,
- Φ is **Completely Positive**, meaning that its Choi matrix (see (2.4.12)) is positive [Cho75],
- Φ is **Trace preserving**, meaning that $\text{tr} \theta = \text{tr} \Phi(\theta)$ for all θ .

The maps satisfying these properties are named through the acronym LCPT maps, or, in the context of QIT, simply *quantum channels*.

One could also consider quantum channels whose input and output systems are different from each other. In this case, an input state $\rho_A \in \sigma(\mathcal{H}_A)$ is sent into an output state $\Phi(\rho_A) \in \sigma(\mathcal{H}_B)$. The channel

$$\Phi : \sigma(\mathcal{H}_A) \mapsto \sigma(\mathcal{H}_B) \quad (2.3.3)$$

can be defined by introducing an isometry $U_{BE \leftarrow AE}$, so that:

$$\rho_A \mapsto \Phi(\rho_A) = \text{tr}_E \left[U_{BE \leftarrow AE} (\rho_A |0\rangle_E \langle 0|) U_{BE \leftarrow AE}^\dagger \right]. \quad (2.3.4)$$

In general, the map $\rho_A \mapsto U_{BE \leftarrow AE} (\rho_A |0\rangle_E \langle 0|) U_{BE \leftarrow AE}^\dagger$ is called *Stinespring dilation* [Sti55], while the map in (2.3.4) is the *Stinespring representation* of the quantum channel Φ ; in fact, as will be explained in 2.4, a given LCPT map admits different but equivalent representations.

2.4 Channel representations

The Stinespring representation is not the only possible representation of quantum channels. In fact, it can be shown that there exist different representations equivalent to each other; the most important ones are reported below

2.4.a Stinespring representation

As seen in 2.3, the Stinespring dilation $\mathfrak{U}_{S \leftrightarrow SE}$:

$$\begin{aligned} \mathfrak{U}_{S \leftrightarrow SE}(\bullet) &\equiv U_{S \leftrightarrow SE} \bullet U_{S \leftrightarrow SE}^\dagger \\ U_{S \leftrightarrow SE} &\equiv U_{SE} |0\rangle_E \\ \rho_S &\xrightarrow{\mathfrak{U}_{S \leftrightarrow SE}} U_{SE} (\rho_S |0\rangle_E \langle 0|) U_{SE}^\dagger \end{aligned} \quad (2.4.1)$$

is an isometry that maps an input state in $\sigma(\mathcal{H}_S)$ onto a state in $\sigma(\mathcal{H}_S \otimes \mathcal{H}_E)$; tracing this *dilated* state over the environment E returns the output state $\Phi(\rho_S)$ (2.3.4). Note that the input and output system of the quantum channel do not need to coincide. This can be achieved by replacing the unitary U_{SE} with an isometry $U_{BE \leftarrow AE}$. This leads to a more general form of (2.4.1).

$$\begin{aligned} \mathfrak{U}_{A \rightarrow BE}(\bullet) &\equiv U_{BE \leftarrow A} \bullet U_{BE \leftarrow A}^\dagger \\ U_{BE \leftarrow A} &\equiv U_{BE \leftarrow AE} |0\rangle_E \\ \rho_A &\xrightarrow{\mathfrak{U}_{A \rightarrow BE}} U_{BE \leftarrow AE}(\rho_A |0\rangle_E \langle 0|) U_{BE \leftarrow AE}^\dagger \end{aligned} \quad (2.4.2)$$

2.4.b Kraus Representation

The Kraus representation is characterized by a set of operators $\mathcal{K} = \{K_i\}_i$ on \mathcal{H} such that:

$$\sum_i K_i^\dagger K_i = \mathbb{1} \quad (2.4.3)$$

where $\mathbb{1}$ is the identity operator on \mathcal{H} . The set \mathcal{K} describes a unique quantum channel Φ , defined by:

$$\Phi(\rho) = \sum_i K_i \rho K_i^\dagger \quad (2.4.4)$$

The set \mathcal{K} is called *Kraus set*, while the K_i 's are called *Kraus operators*. Note that unitary transformations are a particular class of quantum channels, as given the unitary operator U , one could build the trivial Kraus set $\mathcal{K} = \{U\}$ which satisfies (2.4.3) and the corresponding unitary transformation is given in the Kraus representation, $\rho \mapsto U\rho U^\dagger$.

It is possible, given a Stinespring representation of a channel, to derive (one of) its Kraus representations, in fact from (2.3.4):

$$\Phi(\rho_S) = \text{tr}_E \left[U_{SE} (\rho_S |0\rangle_E \langle 0|) U_{SE}^\dagger \right] = \sum_l {}_E \langle l | U_{SE} |0\rangle_E \rho_S |0\rangle_E \langle 0| U_{SE}^\dagger |l\rangle_E. \quad (2.4.5)$$

Then, one could define:

$$K_i \equiv {}_E \langle i | U_{SE} |0\rangle_E, \quad (2.4.6)$$

which is an operator acting on \mathcal{H}_S and that clearly satisfies (2.4.3), so that (2.4.4) is obtained. Conversely, one could obtain the Stinespring representation from a Kraus set by defining:

$$U_{SE} \equiv \sum_i K_i |i\rangle_E \langle 0| \quad (2.4.7)$$

Equivalence of Kraus representations

While a certain \mathcal{K} identifies a specific Φ , given a channel Φ one may find different Kraus sets associated with it. It is then reasonable to ask under which condition two Kraus

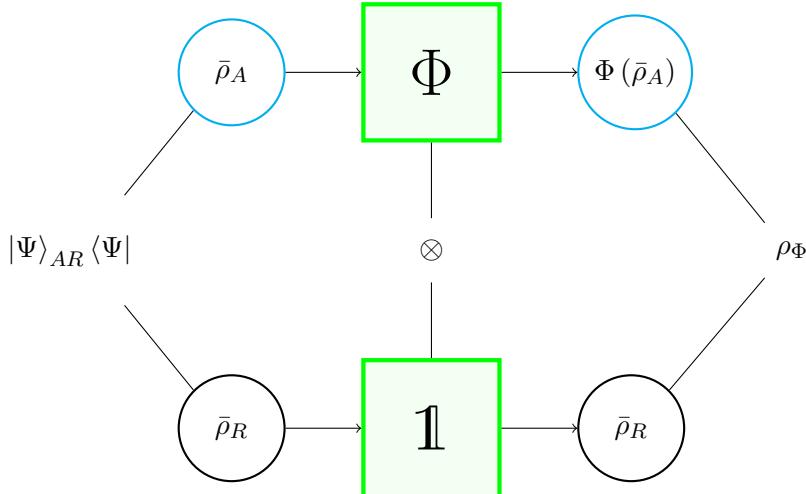


Figure 2.4.1: The Choi state ρ_Φ for a channel Φ can be built using the scheme described in this figure

sets represent the same quantum channel. The answer is that, given the Kraus sets $\mathcal{K} = \{K_i\}_i$ and $\mathcal{K}' = \{K'_i\}_i$, they represent the same channel if and only if:

$$K'_i = \sum_j \mathcal{U}_{ij} K_j, \quad (2.4.8)$$

where the \mathcal{U}_{ij} are the matrix elements of an isometry and they satisfy the conditions:

$$\sum_k \mathcal{U}_{ik} \mathcal{U}_{jk}^* = \sum_k \mathcal{U}_{ki} \mathcal{U}_{kj}^* = \delta_{ij}. \quad (2.4.9)$$

2.4.c Choi-Jamiołkowski representation

The Choi-Jamiołkowski isomorphism [Cho75], [Jam72] introduces a one-to-one mapping between a channel and a specific density matrix. Consider the channel Φ acting on the system \mathcal{H}_A of dimension d and a reference system \mathcal{H}_R ; given the completely chaotic state $\bar{\rho}_A$:

$$\bar{\rho}_A \equiv \frac{1}{d} \sum_i |i\rangle\langle i|, \quad (2.4.10)$$

it can be purified using the reference system R , leading to the maximally entangled state $|\Psi\rangle_{AR}$:

$$|\Psi\rangle_{AR} = \frac{1}{\sqrt{d}} \sum_i |i\rangle_A |i\rangle_R. \quad (2.4.11)$$

If the state in system A is sent through the channel Φ the corresponding state in AR is sent through $\Phi \otimes \mathbb{1}_R$; if the maximally entangled state in (2.4.11) is sent through $\Phi \otimes \mathbb{1}_R$, the output of this channel is called *Choi state* of the channel:

$$\begin{aligned} \rho_\Phi & \text{ Choi state} \\ \rho_\Phi & \equiv [\Phi \otimes \mathbb{1}_R] (|\Psi\rangle_{AR} \langle \Psi|) \end{aligned} \quad (2.4.12)$$

Figure 2.4.1 provides a visual depiction of Choi states. It is possible to prove (see [Cho75], [Jam72]) that:

$$\begin{aligned}\Phi(|\psi\rangle_A \langle\psi|) &= d_R \langle\psi^*|\rho_\Phi|\psi^*\rangle_R \\ &= d \operatorname{tr}_R \left(|\psi\rangle_R \langle\psi| (\rho_\Phi)^{T_A} \right)\end{aligned}\quad (2.4.13)$$

where

$$\begin{aligned}|\psi\rangle &= \sum_{i=0}^{d-1} \alpha_i |i\rangle \\ \Rightarrow |\psi^*\rangle &= \sum_{i=0}^{d-1} \alpha_i^* |i\rangle.\end{aligned}\quad (2.4.14)$$

and the operation of *partial transpose* has been introduced:

$$\begin{aligned}\theta_{AR} &= \sum_{ijmn} \theta_{ijmn} |i\rangle_A \langle j| |m\rangle_R \langle n| \\ \Rightarrow (\theta_{AR})^{T_A} &= \sum_{ijmn} \theta_{ijmn} |j\rangle_A \langle i| |m\rangle_R \langle n|.\end{aligned}\quad (2.4.15)$$

Due to the linearity of Φ , one could easily extend (2.4.13) to the mixed state case:

$$\Phi(\theta_A) = d \operatorname{tr}_R \left(\theta_R (\rho_\Phi)^{T_A} \right)\quad (2.4.16)$$

Choi theorem

The complete positiveness of a map can be verified by checking the positivity of the associated Choi matrix (2.4.12). This is a consequence of Choi's theorem illustrated in [Cho75], [Jam72]:

$$\Psi \text{ is completely positive} \Leftrightarrow \rho_\Psi \text{ is positive.}\quad (2.4.17)$$

2.5 Complementary channels

The Stinespring dilation 2.4.2 allows to define another type of quantum channels; in fact, one may wonder what happens when one traces the dilation of a state over the output system instead of the environment. This leads to the definition of the *complementary channels*, which act as the "output state of the environment".

$$\tilde{\Phi}(\rho_A) \equiv \operatorname{tr}_B \left(U_{BE \leftarrow AE} (\rho_A |0\rangle_E \langle 0|) U_{BE \leftarrow AE}^\dagger \right) \in \sigma(\mathcal{H}_E)\quad (2.5.1)$$

Complementary channels, just as their "standard" counterparts, are LCPT maps. In Figure 2.5.1 the reader can find a schematic depiction of quantum channels and complementary channels.

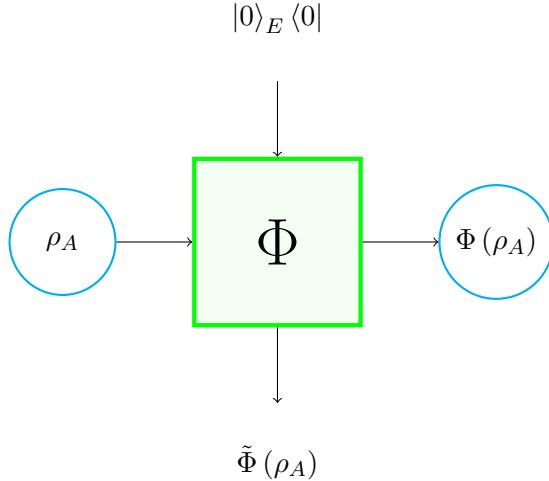


Figure 2.5.1: The quantum channel Φ depicted as a "black box" that takes as inputs ρ_S from the system S and $|0\rangle_E \langle 0|$ while it outputs $\Phi(\rho_S)$ in S and $\tilde{\Phi}(\rho_S)$ in E .

Complementary channel from Kraus representation

It is possible to find the complementary channel from a Kraus set by employing (2.4.6) into (2.5.1), yielding:

$$\tilde{\Phi}(\rho) = \sum_{ij} \text{tr} \left(K_i \rho K_j^\dagger \right) |i\rangle_E \langle j| \quad (2.5.2)$$

Different Kraus representations correspond to different complementary channel, which are unitarily equivalent; in fact, given the Kraus sets $\mathcal{K}, \mathcal{K}'$:

$$\begin{aligned} \tilde{\Phi}(\rho) &= \sum_{ij} \text{tr} \left(K_i \rho K_j^\dagger \right) |i\rangle_E \langle j| \\ \tilde{\Phi}'(\rho) &= \sum_{ij} \text{tr} \left(K'_i \rho K'^\dagger_j \right) |e_i\rangle_E \langle e_j|. \end{aligned} \quad (2.5.3)$$

Suppose $\mathcal{K}, \mathcal{K}'$, satisfy (2.4.8), then defining the isometry $U \equiv \sum_{i,j} \mathcal{U}_{ij} |e_i\rangle_E \langle j|$, one could verify that:

$$\tilde{\Phi}'(\rho) = U \tilde{\Phi}(\rho) U^\dagger \quad (2.5.4)$$

2.6 Degradability and antidegradability

A quantum channel Φ is said to be *degradable* if and only if, given its complementary channel $\tilde{\Phi}$, there exists a LCPT map Λ such that the composition $\Lambda \circ \Phi$ is equal to the complementary channel $\tilde{\Phi}$:

$$\Phi \text{ is degradable} \Leftrightarrow \exists \Lambda \text{ LCPT s.t. } \tilde{\Phi} = \Lambda \circ \Phi \quad (2.6.1)$$

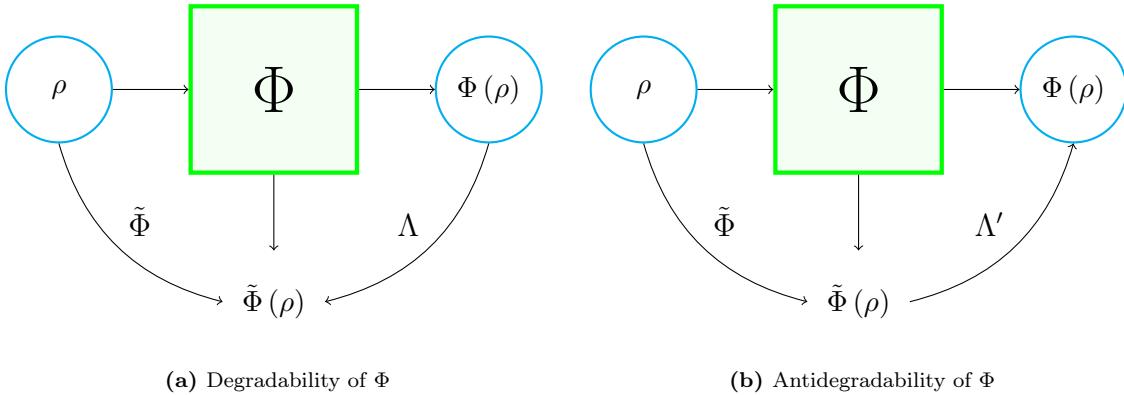


Figure 2.6.1: On the left, a visual representation for the degradability condition given in (2.6.1); on the right, a visual representation for the antidegradability condition given in (2.6.2);

Conversely Φ is said to be *antidegradable* if and only if there exists a LCPT map Λ' such that the composition $\Lambda' \circ \tilde{\Phi}$ is equal to the channel Φ :

$$\Phi \text{ is antidegradable} \Leftrightarrow \exists \Lambda' \text{ LCPT s.t. } \Phi = \Lambda' \circ \tilde{\Phi}, \quad (2.6.2)$$

Figure 2.6.1 provides a visual representation of the degradability and antidegradability conditions outlined in this section.

2.7 Covariant channels

Consider a quantum channel $\Phi : \sigma(\mathcal{H}_A) \mapsto \sigma(\mathcal{H}_B)$ and a group \mathfrak{G} which has unitary representations on \mathcal{H}_A , \mathcal{H}_B and the environment \mathcal{H}_E , so that $\forall g \in \mathfrak{G}, \exists U_g^X$ unitary operator on \mathcal{H}_X . The channel Φ is called *covariant* under the action of \mathfrak{G} if:

$$\Phi \left(U_g^A \rho U_g^{A\dagger} \right) = U_g^B \Phi(\rho) U_g^{B\dagger} \quad \forall g \in \mathfrak{G}, \forall \rho \in \sigma(\mathcal{H}_A) \quad (2.7.1)$$

Define $\tilde{\Phi}$ as the complementary channel of Φ ; if Φ satisfies (2.7.1), then [Hol07] also its complementary channel must be covariant under the same group \mathfrak{G} :

$$\tilde{\Phi} \left(U_g^A \rho U_g^{A\dagger} \right) = U_g^E \tilde{\Phi}(\rho) U_g^{E\dagger} \quad \forall g \in \mathfrak{G}, \forall \rho \in \sigma(\mathcal{H}_A) \quad (2.7.2)$$

$$\Phi \text{ covariant under } \mathfrak{G} \Rightarrow \tilde{\Phi} \text{ covariant under } \mathfrak{G} \quad (2.7.3)$$

2.8 Shannon entropy

In this section, the basic tool that allows for the study of (classical) information will be introduced. Given a probability distribution $X = \{p_x\}_x$, the *Shannon entropy* (or

classical entropy) is defined as:

$$H(X) \equiv - \sum_x p_x \log_2 p_x \quad (2.8.1)$$

From the Shannon entropy, a number of information theoretic quantities can be derived, as exposed in what follows. It is important to note that, if for some x_0 one finds $p_{x_0} = 0$, one needs to set a convention to treat the indefinite form of $0 \log_2 0$ found in (2.8.1). Therefore, in the context of entropies, the reader needs to assume:

$$0 \log_2 0 \equiv 0 \quad (2.8.2)$$

One can think of the Shannon entropy as the "classical information" content that can be gained by measuring the system, or the amount of *bits* needed to store the information contained in the system.

Classical relative entropy

The classical relative entropy provides a way to measure the closeness of two probability distributions $X = \{p_x\}_x$ and $Y = \{q_x\}_x$, which are defined over the same indexes. The relative entropy is defined as:

$$H(p_x || q_x) \equiv -H(X) - \sum_x p_x \log_2 q_x = \sum_x p_x \log_2 \frac{p_x}{q_x}. \quad (2.8.3)$$

An important property of the classical relative entropy is its non-negativity:

$$H(p_x || q_x) \geq 0 \quad (2.8.4)$$

Using this property, it is possible to find an upper bound for the Shannon entropy. In fact, using $q_x = 1/d \forall x$:

$$H(p_x || 1/d) = \log_2 d - H(X) \geq 0 \Rightarrow H(X) \leq \log_2 d \quad (2.8.5)$$

This upper bound is achievable by selecting $p_x = 1/d \forall x$.

Joint classical entropy

Given a pair of random variables x, y and a probability distribution over those variables $XY = \{p(x, y)\}_{x,y}$, then the *joint classical entropy* is a natural extension of (2.8.1):

$$H(XY) \equiv - \sum_{x,y} p(x, y) \log_2 p(x, y) \quad (2.8.6)$$

Classical conditional entropy

In the case of two random variables x, y , the knowledge of one of those variables, for example y , changes the information content that can be gained by measuring the system X ; this quantity is represented by the *conditional entropy*:

$$H(X|Y) \equiv H(XY) - H(Y) \quad (2.8.7)$$

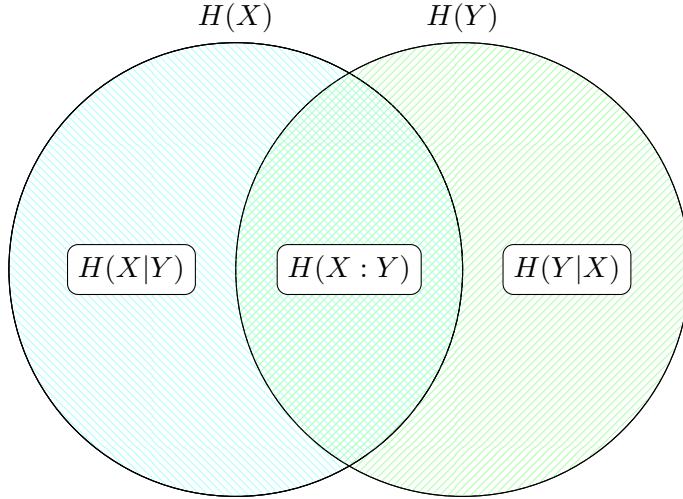


Figure 2.8.1: This pseudo-Venn diagram provides a powerful mnemonic tool that depicts the relationship between the classical entropic quantities.

Classical mutual information

The *mutual information* measures the information content in XY accessible to both the observers that only have access to either X or Y :

$$\begin{aligned} H(X : Y) &\equiv H(X) + H(Y) - H(XY) \\ H(X : Y) &\equiv H(X) - H(X|Y) \end{aligned} \tag{2.8.8}$$

The relationships between the classical entropic quantities can be (improperly) visually derived using the Venn diagram drawn in Figure 2.8.1.

2.9 Von Neumann entropy

The quantum equivalent of the Shannon entropy in (2.8.1) is the Von Neumann entropy; it describes the amount of *qubits* needed to store the information contained in a quantum system. The Von Neumann entropy is defined by:

$$S(\rho) \equiv -\text{tr}(\rho \log_2 \rho) \tag{2.9.1}$$

where ρ is a quantum state. The \log_2 in (2.9.1) is the logarithm base 2 of a matrix; formally speaking, it is the inverse operator of taking the power of that matrix base 2, while practically speaking, $\log_2 \theta$ returns a matrix θ' obtained by taking the element-wise \log_2 of θ in its diagonal basis and then transforming back to the original basis for θ . Since the trace operator is invariant under change of basis, this means that, denoting with $\{\lambda_x\}_x$ the eigenvalues of ρ , (2.9.1) can be reduced to the form of (2.8.1):

$$S(\rho) = -\sum_x \lambda_x \log_2 \lambda_x, \tag{2.9.2}$$

(2.8.2) still holds.

2.10 Quantum entropic quantities

In this Section, the most important quantum entropic quantities, built from the Von Neumann entropy, are listed.

Relative entropy

The quantum relative entropy is the extension to the quantum case of the quantity in (2.8.3). Its definition is given by:

$$S(\rho||\tau) \equiv \text{tr}(\rho \log_2 \rho) - \text{tr}(\rho \log_2 \tau) = -S(\rho) - \text{tr}(\rho \log_2 \tau) \quad (2.10.1)$$

The Klein's inequality guarantees that the relative entropy is non-negative:

$$S(\rho||\tau) \geq 0, \quad (2.10.2)$$

where the inequality is saturated if and only if $\rho = \tau$. The relation (2.10.2) offers a way of setting an upper bound on the Von Neumann entropy; in fact, assuming $\tau = \frac{1}{d}\mathbb{1}_{\mathcal{H}_d}$:

$$0 \leq S(\rho||\tau) = -S(\rho) + \log_2 d \quad (2.10.3)$$

$$S(\rho) \leq \log_2 d \quad (2.10.4)$$

This bound is achievable when $\rho = \frac{1}{d}\mathbb{1}_{\mathcal{H}_d}$; this state is often called the *completely chaotic state*.

Joint entropy

The joint quantum entropy is the extension of the Von Neumann entropy to a joint quantum system.

$$S(AB)_\rho \equiv \text{tr}_{AB}(\rho_{AB} \log_2 \rho_{AB}). \quad (2.10.5)$$

The index ρ signifies the state over which the entropy is to be calculated. This notation will also occur in later definitions.

Conditional entropy

When one part of a joint system AB is already known, the information content of the remaining system is given by the quantum conditional entropy:

$$S(A|B)_\rho \equiv S(AB)_\rho - S(B)_\rho, \quad (2.10.6)$$

where $\rho_{A/B} = \text{tr}_{B/A} \rho_{AB}$.

Mutual information

The quantum mutual information represents the amount of information stored in a joint state in the system AB accessible to both parties A, B :

$$I(A : B)_\rho \equiv S(A)_\rho + S(B)_\rho - S(AB)_\rho \quad (2.10.7)$$

where $\rho_{A/B} = \text{tr}_{B/A} \rho_{AB}$. If B is the output system of a channel $\Psi_{A \rightarrow B}$ whose input is the system A , then the mutual information can be written as:

$$\begin{aligned} I(A : B)_\rho &\equiv S(A)_\rho + S(B)_\rho - S(AB)_\rho \\ &\equiv I(\rho_A, \Psi_{A \rightarrow B}) = S(A)_\rho + S(B)_\rho - S(E)_\rho \\ &= S(\rho_A) + S(\Psi_{A \rightarrow B}(\rho_A)) - S(\tilde{\Psi}_{A \rightarrow E}(\rho_A)), \end{aligned} \quad (2.10.8)$$

refer to Figure 2.12.1 for a clearer understanding of the quantities in play. The main properties of the mutual information [Hol13] are listed below:

- $I(\rho, \Psi)$ is concave in ρ ,
- $I(\rho, \Psi)$ is concave in Ψ ,
- $I(\rho, \Psi)$ is subadditive,
- $I(\rho, \Psi)$ satisfies the following data-processing inequalities:

$$I(\rho, \Psi_2 \circ \Psi_1) \leq \min \{I(\rho, \Psi_1), I(\Psi_1(\rho), \Psi_2)\} \quad (2.10.9)$$

Coherent information

The coherent information is the basic building block for the quantum capacity, defined in 2.11; its definition is given by:

$$\begin{aligned} I_C(\Psi_{A \rightarrow B}, \rho_A) &\equiv S(B)_\rho - S(E)_\rho \\ &= S(\Psi_{A \rightarrow B}(\rho_A)) - S(\tilde{\Psi}_{A \rightarrow E}(\rho_A)) \end{aligned} \quad (2.10.10)$$

Uncomfortably, the coherent information does not satisfy many of the properties of the mutual information, such as the subadditivity property [DSS98]:

$$I_C(\rho_{12}, \Psi_1 \otimes \Psi_2) \not\leq I_C(\rho_1, \Psi_1) + I_C(\rho_2, \Psi_2), \quad (2.10.11)$$

where $\rho_{1/2} = \text{tr}_{2/1} \rho_{12}$, the concavity in the input states or one of the data-processing inequalities:

$$\begin{aligned} I_C(\rho, \Psi_2 \circ \Psi_1) &\leq I_C(\rho, \Psi_1), \\ I_C(\rho, \Psi_2 \circ \Psi_1) &\not\leq I_C(\Psi_1(\rho), \Psi_2). \end{aligned} \quad (2.10.12)$$

If a channel satisfies (2.6.1), then its coherent information is concave [YHD08]:

$$\Psi \text{ is degradable} \Rightarrow I_C \left(\sum_i p_i \rho_i, \Psi \right) \geq \sum_i p_i I_C(\rho_i, \Psi). \quad (2.10.13)$$

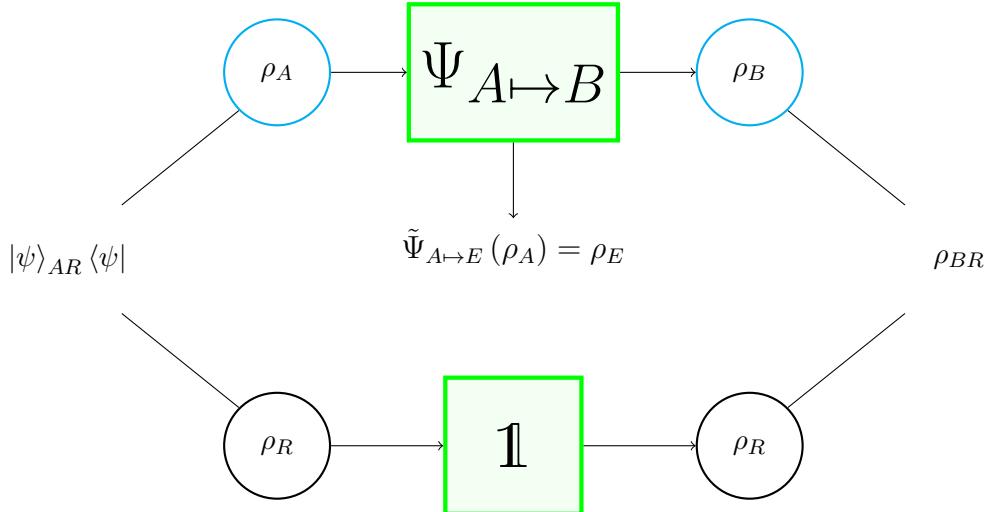


Figure 2.12.1: Figure 2.4.1 is a specific case of the general configuration illustrated in this scheme; Alice sends the message ρ_A to Bob, who receives the state ρ_B ; Alice also has access to a reference system in the state ρ_R , which is isospectral to ρ_A (due to the Schmidt decomposition, see e.g. [Hol13], page 37).

2.11 Quantum capacity

The quantum capacity of a quantum channel represents the rate for reliably transmitting quantum information, encoded in quantum states, through the channel. Its expression for a given noisy channel Ψ can be found in the paper by Devetak [Dev05]:

$$Q(\Psi) \equiv \lim_{n \rightarrow \infty} \frac{Q^{(n)}(\Psi)}{n} \equiv \lim_{n \rightarrow \infty} \frac{1}{n} \max_{\rho^{(n)} \in \sigma(\mathcal{H}^{\otimes n})} I_c(\rho^{(n)}, \Psi^{\otimes n}). \quad (2.11.1)$$

The quantity $Q^{(n)}(\Psi)$ represents the maximization of the coherent information for n uses of the channel over all possible inputs. Due to the property in (2.10.11), the limit for n uses of the channel can not be dropped, rendering the actual computation of the quantum capacity impossible. However [DS05], if a channel satisfies (2.6.1), then its coherent information is subadditive and the limit over n can be dropped, allowing to write:

$$\Psi \text{ is degradable} \Rightarrow Q(\Psi) = Q^{(1)}(\Psi) = \max_{\rho \in \sigma(\mathcal{H})} I_c(\rho, \Psi). \quad (2.11.2)$$

2.12 2-way capacity

The 2-way quantum capacity Q_2 represents the maximum rate for reliably transmitting quantum information under the assumption that the two parties involved are able to perform LOCC's (Local Operations and Classical Communication). At the moment

of writing, there is no known formula that outputs the Q_2 of a channel; however, it is always possible to find upper and lower bounds, as illustrated in [PLOB17]

2.12.a Lower bound

A natural lower bound of the 2-way capacity is the quantum capacity itself; another lower bound on the Q_2 of a quantum channel $\Psi_{A \rightarrow B}$ can be set by:

$$Q_2(\Psi_{A \rightarrow B}) \geq \max_{\rho_A} \{I_C(\Psi_{A \rightarrow B}, \rho_A), I_{RC}(\Psi_{A \rightarrow B}, \rho_A)\} \quad (2.12.1)$$

where I_C is the coherent information of the channel, while I_{RC} is called the reverse coherent information of the channel:

$$\begin{aligned} I_C(\Psi_{A \rightarrow B}, \rho_A) &= S(\text{tr}_R(\rho_{BR})) - S(\rho_{BR}) \\ I_{RC}(\Psi_{A \rightarrow B}, \rho_A) &\equiv S(\text{tr}_B(\rho_{BR})) - S(\rho_{BR}) \end{aligned} \quad (2.12.2)$$

Note that the expression for the coherent information I_C in (2.12.2) is equivalent to the definition (2.10.10). The system R is used to purify the input state ρ_A into $|\psi\rangle_{AR}\langle\psi|$. Refer to Figure 2.12.1 for a visual representation of the quantities in play. Since in (2.12.1) there is a maximization over all possible inputs, a fixed choice of ρ_A gives still rise to a (possibly non optimal) lower bound. In particular, a common choice of ρ_A is the completely chaotic state $\bar{\rho}_A$, so that ρ_{BR} is the Choi state [Cho75] of the channel ρ_Ψ . In this case, if the channel is finite dimensional and the output of the channel has the same dimensionality as its input (as it is the case for ReMAD's), then the reverse coherent information is always bigger than the coherent information, so that:

$$Q_2(\Psi_{A \rightarrow B}) \geq I_{RC}(\Psi_{A \rightarrow B}, \bar{\rho}_A) = \log_2 d - S(\rho_\Psi) \quad (2.12.3)$$

where $d = \dim \mathcal{H}_A$. The relation (2.12.3) was used to find lower bounds on the 2-way capacity of 3-dimensional MAD channels in 6.

2.12.b Upper Bound

The computation of the upper bound of the Q_2 of a channel is more difficult than that of the lower bound. One quantity that is always bigger than Q_2 under any condition is the squashed entanglement.

Squashed entanglement

Given a noisy channel \mathcal{N} , its squashed entanglement is defined by (see e.g. [KW20]):

$$\begin{aligned} E_{sq}(\mathcal{N}) &\equiv \sup_{|\psi\rangle_{AR}\langle\psi|} E_{sq}(R; B)_\psi \\ E_{sq}(R; B)_\psi &\equiv \frac{1}{2} \inf_{\Omega_{E \rightarrow E'}} \left(I(R; B|E')_\rho : \rho_{BRE'} = \Omega_{E \rightarrow E'}(|\psi\rangle_{BRE}\langle\psi|) \right) \end{aligned} \quad (2.12.4)$$

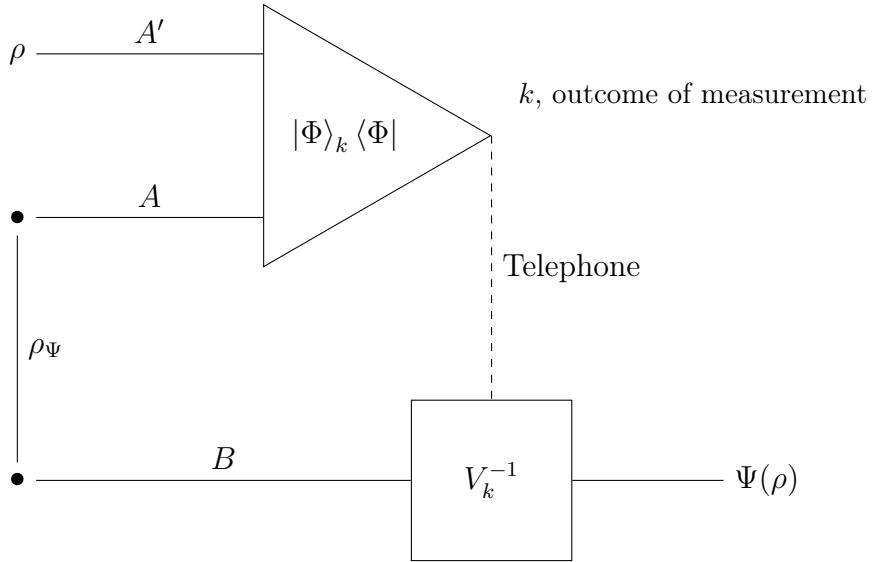


Figure 2.12.2: Schematic representation of the simulation of a telecovariant channel Ψ . Alice has access to the systems A , A' , while Bob has access to system B . They share two halves of a composite system AB , which is in the Choi state ρ_Ψ . Alice wants to send the state $\Psi(\rho)$ to Bob, while having the state ρ in her system A' . She makes a projective measurement of the Bell states in the system AA' , communicates the outcome to Bob who then performs an appropriate unitary transformation on his system B to obtain the desired state.

where $|\psi\rangle_{BRE} \langle \psi|$ is the output of the Stinespring dilation that corresponds to a pure state input $|\psi\rangle_{AR} \langle \psi|$, $(N_{A \rightarrow B} \otimes \mathbb{1}_R)(|\psi\rangle_{AR} \langle \psi|) = \text{tr}_E(|\psi\rangle_{BRE} \langle \psi|)$, while $I(R; B|E')_\rho$ is the quantum conditional mutual information:

$$\begin{aligned} I(R; B|E')_\rho &\equiv S(R|E')_\rho + S(B|E')_\rho - S(BR|E')_\rho \\ &= S(B|E')_\rho - S(B|RE')_\rho \end{aligned} \quad (2.12.5)$$

From [Chr06]:

$$Q_2(\mathcal{N}) \leq E_{sq}(\mathcal{N}) \quad (2.12.6)$$

The $\inf_{\Omega_{E \leftrightarrow E'}}$ in (2.12.4) can be dropped by choosing $\Omega_{E \leftrightarrow E'} = \mathbb{1}_{E \leftrightarrow E}$, the remaining quantity is still an upper bound, albeit possibly non optimal.

Teleportation covariance

One defines (see [PLOB17]) σ -stretchable channels as channels \mathcal{E} satisfying the condition:

$$\mathcal{E}(\rho) = \mathcal{T}(\sigma \otimes \rho) \quad \forall \rho \quad (2.12.7)$$

where σ is to be considered a shared resource between the two parties "Alice and Bob" while \mathcal{T} is a set of LOCC's. Accordingly, one can define "Choi-stretchable" channels that can be simulated over their Choi states $\rho_{\mathcal{E}}$ using LOCC's, i.e. $\exists \mathcal{T}$ LOCC:

$$\mathcal{E}(\rho) = \mathcal{T}(\rho_{\mathcal{E}} \otimes \rho) \quad \forall \rho \quad (2.12.8)$$

A channel \mathcal{E} is said to be teleportation covariant if, under a teleportation unitary U and another unitary (not necessarily a teleportation unitary) V one has:

$$\mathcal{E}(U\rho U^\dagger) = V\mathcal{E}(\rho)V^\dagger \quad (2.12.9)$$

A teleportation covariant channel is Choi-stretchable; the protocol with which these channels can be simulated is the same as the qudit teleportation protocol (see e.g. [PEW⁺15]), with the only difference being that the shared resource is not the maximally entangled state but the Choi matrix of the channel to be simulated. See [PLOB17] for details and Figure 2.12.2 for a visual representation of the protocol. A sufficient and necessary condition for tele-covariance is given by:

$$[\rho_{\mathcal{E}}, U_k \otimes V_k] = 0 \quad \forall k \quad (2.12.10)$$

where U_k 's are the teleportation unitaries and V_k 's are (possibly) different representations of the same group elements as the U_k 's. Amplitude damping channels are not tele-covariant; however, they can be mapped into bosonic lossy channels, which, since they are gaussian channels, are also Choi-stretchable.

Relative Entropy of Entanglement

The relative entropy of entanglement (REE) can be thought of a kind of distance from an input entangled state to its "closest" separable state (2.2.2) and is defined by:

$$E_R(\rho) \equiv \inf_{\sigma_s} S(\rho || \sigma_s) \quad \sigma_s \text{ separable} \quad (2.12.11)$$

If the channel \mathcal{E} satisfies (2.12.10), then an upper bound on $Q_2(\mathcal{E})$ can be found with relative ease:

$$Q_2(\mathcal{E}) \leq E_R(\rho_{\mathcal{E}}) \quad (2.12.12)$$

2.13 Pipeline inequalities

Given a generic capacity functional \mathfrak{C} , such as the quantum capacity Q or the 2-way capacity Q_2 , one may derive the pipeline (or data-processing) inequalities [HG12]: given two quantum channels Ψ_1, Ψ_2 and their composition $\Psi_2 \circ \Psi_1$ the capacity of the composed channel can not be bigger than those that of the single channels:

$$\mathfrak{C}[\Psi_2 \circ \Psi_1] \leq \min \{\mathfrak{C}[\Psi_1], \mathfrak{C}[\Psi_2]\}. \quad (2.13.1)$$

The relation (2.13.1) makes intuitive sense: the capacity represents the ability of the receiving party to decipher the information sent through the channel, therefore adding further "noise" (i.e. an additional channel) to a channel can not increase its capacity.

3

Previous research on finite dimensional lossy channels

As already stated in 1, the purpose of this thesis work is the computation of some capacity functionals for a specific family of quantum channels, called Multi-level Amplitude Damping (MAD) channels; the previous research on MAD channels is exposed in the present chapter.

3.1 Amplitude Damping Channels

Amplitude Damping Channels (see for example [NC10] page 380), or ADC's, are qubit-to-qubit channels defined by the Kraus set:

$$K_0 \equiv |0\rangle\langle 0| + \sqrt{1-\gamma} |1\rangle\langle 1| \quad K_1 \equiv \sqrt{\gamma} |0\rangle\langle 1|, \quad (3.1.1)$$

where γ is a real parameter satisfying $0 \leq \gamma \leq 1$; it completely identifies an ADC. Given an input density matrix ρ :

$$\rho = \begin{pmatrix} \rho_{00} & \rho_{01} \\ \rho_{01}^* & \rho_{11} \end{pmatrix}, \quad (3.1.2)$$

the corresponding output state of an ADC is:

$$\begin{aligned} \text{ADC}_\gamma(\rho) &\equiv K_0 \rho K_0^\dagger + K_1 \rho K_1^\dagger \\ \text{ADC}_\gamma(\rho) &= \begin{pmatrix} \rho_{00} + \gamma \rho_{11} & \sqrt{1-\gamma} \rho_{01} \\ \sqrt{1-\gamma} \rho_{01}^* & (1-\gamma) \rho_{11} \end{pmatrix} \end{aligned} \quad (3.1.3)$$

For a detailed review on the quantum capacity, classical and quantum entangled assisted capacity, and classical capacity of ADC's see [GF05]. In the present section, a more in-depth analysis will be given only for the quantum capacity (and, consequently, also for the classical private capacity). Upper and lower bounds for the two-way capacity of ADC's can be obtained following the results illustrated in [PLOB17].

3.1.a Composition of ADC's

Consider a composition $\text{ADC}_{\gamma''} \circ \text{ADC}_{\gamma'}$, its output channel is given by:

$$\text{ADC}_{\gamma''} \circ \text{ADC}_{\gamma'}(\rho) = \begin{pmatrix} \rho_{00} + (\gamma' + \gamma'' - \gamma'\gamma'')\rho_{11} & \sqrt{(1-\gamma')(1-\gamma'')}\rho_{01} \\ \sqrt{(1-\gamma')(1-\gamma'')}\rho_{01}^* & (1-\gamma')(1-\gamma'')\rho_{11} \end{pmatrix} \quad (3.1.4)$$

One could define a new parameter $\gamma = \gamma' + \gamma''(1 - \gamma')$, so that $\text{ADC}_{\gamma''} \circ \text{ADC}_{\gamma'}(\rho) = \text{ADC}_\gamma(\rho)$. This implies that ADC's are closed under composition, and the composition rules are:

$$\begin{aligned} \text{ADC}_\gamma &= \text{ADC}_{\gamma''} \circ \text{ADC}_{\gamma'} \\ \gamma &= \gamma' + \gamma''(1 - \gamma') \geq \gamma', \gamma'' \end{aligned} \quad (3.1.5)$$

Notice that these composition rules, coupled with the bottleneck inequalities (2.13.1), imply that any capacity functional \mathfrak{C} for ADC's must be monotonous non-increasing for increasing values of γ :

$$\mathfrak{C}(\text{ADC}_\gamma) \leq \mathfrak{C}(\text{ADC}_{\gamma'}) \quad \gamma \geq \gamma' \quad (3.1.6)$$

3.1.b Complementary channel of an ADC

Equation (2.5.2) can be used to find the output of the complementary channel of an ADC given an input matrix ρ of the form (3.1.2):

$$\widetilde{\text{ADC}}_\gamma(\rho) = \begin{pmatrix} \rho_{00} + (1-\gamma)\rho_{11} & \sqrt{\gamma}\rho_{01} \\ \sqrt{\gamma}\rho_{01}^* & \gamma\rho_{11} \end{pmatrix} \quad (3.1.7)$$

From (3.1.7) and (3.1.3), it is apparent that:

$$\text{ADC}_{1-\gamma} = \widetilde{\text{ADC}}_\gamma \quad (3.1.8)$$

3.1.c Degradability and antidegradability

The conditions (2.6.1) and (2.6.2) define what it means for a generic channel Φ to be, respectively, degradable or antidegradable. In the context of Amplitude Damping Channels, one finds, from (3.1.8), that the degradability condition reduces to:

$$\exists \Lambda : \text{ADC}_{1-\gamma} = \Lambda \circ \text{ADC}_\gamma, \quad (3.1.9)$$

while the antidegradability condition becomes:

$$\exists \Lambda' : \text{ADC}_\gamma = \Lambda' \circ \text{ADC}_{1-\gamma}. \quad (3.1.10)$$

Heuristically speaking, it would make sense if Λ and Λ' in (3.1.9) and (3.1.10) were both ADC's themselves. Following this hypothesis, one could rewrite (3.1.9) as:

$$\exists \lambda, 0 \leq \lambda \leq 1 : \text{ADC}_{1-\gamma} = \text{ADC}_\lambda \circ \text{ADC}_\gamma, \quad (3.1.11)$$

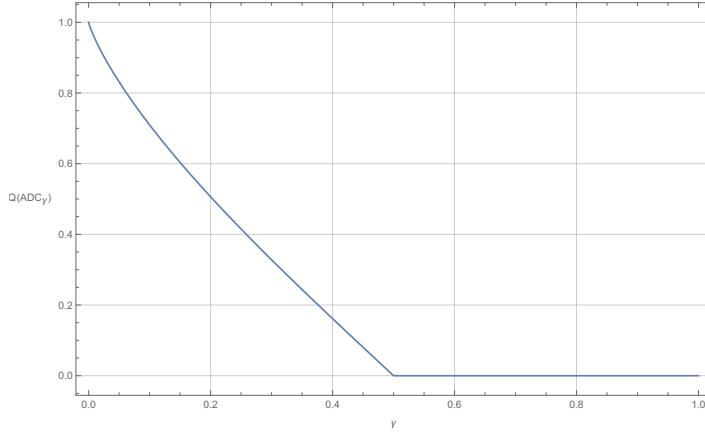


Figure 3.1.1: Quantum capacity for an Amplitude Damping Channel ADC_γ as a function of the parameter γ . Notice that (3.1.6) is satisfied.

Employing (3.1.5) into (3.1.11), one finds:

$$\lambda = \frac{1 - 2\gamma}{1 - \gamma}. \quad (3.1.12)$$

In order for λ to satisfy $0 \leq \lambda \leq 1$, one needs to set the condition $0 \leq \gamma \leq 1/2$; therefore, if $0 \leq \gamma \leq 1/2$ the channel ADC_γ is degradable. One could also follow this line of reasoning in order to find an antidegradability condition; in fact, assuming that the antidegrading channel is itself an ADC, one could rewrite (3.1.10) as:

$$\exists \lambda', 0 \leq \lambda' \leq 1 : \text{ADC}_\gamma = \text{ADC}_{\lambda'} \circ \text{ADC}_{1-\gamma}, \quad (3.1.13)$$

This leads to:

$$\lambda' = \frac{2\gamma - 1}{\gamma} \quad (3.1.14)$$

Therefore, ADC_γ is antidegradable if $1/2 \leq \gamma \leq 1$. This means that ADC's are either degradable or antidegradable:

$$\begin{cases} 0 \leq \gamma \leq 1/2 \Rightarrow \text{ADC}_\gamma \text{ is degradable} \\ 1/2 \leq \gamma \leq 1 \Rightarrow \text{ADC}_\gamma \text{ is antidegradable} \end{cases} \quad (3.1.15)$$

3.1.d Quantum Capacity of ADC's

Following the result (3.1.15), the quantum capacity of ADC's turns out to be relatively easy to compute. In fact, for $1/2 \leq \gamma \leq 1$, since the channel is antidegradable, the quantum capacity of ADC_γ is 0, while for $0 \leq \gamma \leq 1/2$, due to the property of degradable channels 2.11.2, the quantum capacity of ADC_γ corresponds to the maximum of the coherent information over all possible inputs for a single use of the channel:

$$Q(\text{ADC}_\gamma) = Q^{(1)}(\text{ADC}_\gamma) = \max_{\rho \in \sigma(\mathcal{H}_2)} I_c(\rho, \text{ADC}_\gamma) \quad \forall 0 \leq \gamma \leq 1/2 \quad (3.1.16)$$

Thanks to a property that MAD channels present for every dimension d (see (3.3.21)), the maximum in (3.1.16) can be computed over all diagonal density matrices $\rho^{(diag)}$:

$$\rho^{(diag)} \equiv p |0\rangle\langle 0| + (1-p) |1\rangle\langle 1| \quad (3.1.17)$$

$$\max_{\rho^{(diag)} \in \sigma(\mathcal{H}_2)} I_c(\rho, \text{ADC}_\gamma) = \max_p \left\{ \begin{array}{l} -(1-p)(1-\gamma) \log_2((1-p)(1-\gamma)) \\ -(\gamma+p(1-\gamma)) \log_2(\gamma+p(1-\gamma)) \\ +(1-\gamma(1-p)) \log_2(1-\gamma(1-p)) \\ +\gamma(1-p) \log_2(\gamma(1-p)) \end{array} \right\}. \quad (3.1.18)$$

The right-hand side of (3.1.18) can be computed using a numerical evaluation, obtaining the plot reported in Figure 3.1.1.

3.2 Partially Coherent Direct Sum channels

In [CG21b], the authors analyze channels of the form:

$$\Phi_{CC}(\Theta_{CC}) \equiv \left[\begin{array}{c|c} \Phi_{AA}(\Theta_{AA}) & \Phi_{AB}^{(off)}(\Theta_{AB}) \\ \hline \Phi_{BA}^{(off)}(\Theta_{BA}) & \Phi_{BB}(\Theta_{BB}) \end{array} \right] \quad (3.2.1)$$

$$\Theta_{CC} \equiv \left[\begin{array}{c|c} \Theta_{AA} & \Theta_{AB} \\ \hline \Theta_{BA} & \Theta_{BB} \end{array} \right];$$

these are called Partially Coherent Direct Sum (PCDS) channels. For this type of channels, it was shown that:

$$\Phi_{CC} \text{ is degradable} \Leftrightarrow \Phi_{AA}, \Phi_{BB} \text{ are both degradable} \quad (3.2.2)$$

If Φ_{CC} is degradable, its quantum capacity presents an upper bound:

$$Q(\Phi_{CC}) \leq \log_2 (2^{Q(\Phi_{AA})} + 2^{Q(\Phi_{BB})}) \quad (3.2.3)$$

Special case of $\Phi_{BB} = \mathbb{1}_{BB}$

Suppose Φ_{BB} is the identity channel on \mathcal{H}_B , $\mathbb{1}_{BB}$; the upper bound (3.2.3) becomes:

$$Q(\Phi_{CC}) \leq \log_2 (2^{Q(\Phi_{AA})} + d_B) \quad (3.2.4)$$

where $d_B = \dim \mathcal{H}_B$. In this case, a lower bound for $Q(\Phi_{CC})$ can also be found:

$$Q(\Phi_{CC}) \geq \log_2 (1 + d_B) \quad (3.2.5)$$

If the degradable channel Φ_{AA} has null quantum capacity, then the upper bound (3.2.4) and the lower bound (3.2.5) coincide:

$$Q(\Phi_{AA}) = 0 \Rightarrow Q(\Phi_{CC}) = \log_2 (1 + d_B) \quad (3.2.6)$$

3.2.a Direct Sum channels

PCDS channels provide a generalization of Direct Sum (DS) channels [FW07], which suppress the coherence terms between the Hilbert spaces \mathcal{H}_A and \mathcal{H}_B :

$$\begin{aligned}\Phi_{CC}^{DS}(\Theta_{CC}) &\equiv \left[\begin{array}{c|c} \Phi_{AA}(\Theta_{AA}) & 0 \\ \hline 0 & \Phi_{BB}(\Theta_{BB}) \end{array} \right] \\ \Theta_{CC} &\equiv \left[\begin{array}{c|c} \Theta_{AA} & \Theta_{AB} \\ \hline \Theta_{BA} & \Theta_{BB} \end{array} \right];\end{aligned}\tag{3.2.7}$$

The capacity of these channels is never greater than the capacity of PCDS channels:

$$Q(\Phi_{CC}^{DS}) = \max \{Q(\Phi_{AA}), Q(\Phi_{BB})\} \leq Q(\Phi_{CC})\tag{3.2.8}$$

3.3 3-dimensional Multi-level Amplitude Damping channels

Multi-level Amplitude Damping channels are the generalization to the qudit case of ADC channels; they were first explored in the article [CG21a], which is summarized in the present section.

3.3.a Settings for d -dimensional MAD channels

The paper [CG21a] outlines the settings for a d -dimensional MAD channel, which are also reported here for the sake of completeness. Given a d -dimensional Hilbert space \mathcal{H} , spanned by the basis $\{|i\rangle\}_i$, for $i = 0, \dots, d-1$, the MAD channel has a minimal Kraus representation given by the Kraus operators:

$$\begin{aligned}K_{ij} &\equiv \sqrt{\gamma_{ji}} |i\rangle\langle j| \quad 0 \leq i < j \leq d-1 \\ K_{00} &\equiv \sum_{j=0}^{d-1} \sqrt{\gamma_{jj}} |j\rangle\langle j|.\end{aligned}\tag{3.3.1}$$

The γ_{ji} 's in (3.3.1) describe the probabilities of decay from level $|j\rangle$ onto level $|i\rangle$, while γ_{jj} 's describe the probabilities that level $|j\rangle$ will not decay during the transformation; as such, these are real quantities satisfying:

$$\gamma_{jj} \equiv 1 - \sum_{i=0}^{j-1} \gamma_{ji}\tag{3.3.2}$$

$$\begin{cases} 0 \leq \gamma_{ji} \leq 1 & \forall 0 \leq i < j \leq d-1 \\ 0 \leq \gamma_{jj} \leq 1 & \forall 0 \leq j \leq d-1 \\ \gamma_{ji} = 0 & \forall i > j \end{cases}\tag{3.3.3}$$

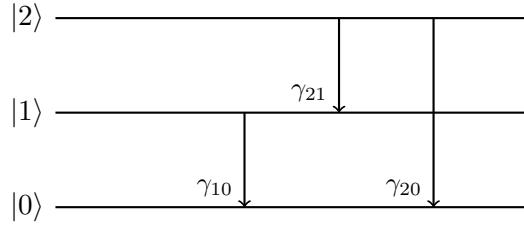


Figure 3.3.1: MAD channels represent decay processes, where each level of a system has a fixed probability of decaying onto a lower level. Here, a schematic depiction of a 3-dimensional MAD is reported

These quantities, which will be called *transition probabilities* in what follows, can be grouped into a matrix, which will be called *transition matrix*, defined by:

$$\Gamma \equiv \mathbb{1}_d + \sum_{j=1}^{d-1} \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle i| - \sum_{j=1}^{d-1} \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle j| \quad (3.3.4)$$

where $\mathbb{1}_d$ is the d -dimensional identity operator. There is a one-to-one relation between a specific MAD channel and its transition matrix, so that, given a transition matrix Γ , it is possible to identify the corresponding MAD channel by Φ_Γ :

$$\Phi_\Gamma \sim \Gamma \quad (3.3.5)$$

Finally, given an input $\rho \in \sigma(\mathcal{H})$, a MAD channel Φ_Γ outputs the state:

$$\Phi_\Gamma(\rho) = K_{00}\rho K_{00}^\dagger + \sum_{j=1}^{d-1} \sum_{i=0}^{j-1} K_{ij}\rho K_{ij}^\dagger. \quad (3.3.6)$$

Which, in terms of transition probabilities, translates to:

$$\Phi_\Gamma(\rho) = \sum_{m=0}^{d-1} \sum_{n=0}^{d-1} \sqrt{\gamma_{mm} \gamma_{nn}} \rho_{mn} |m\rangle\langle n| + \sum_{j=1}^{d-1} \sum_{i=0}^{j-1} \gamma_{ji} \rho_{ii} |i\rangle\langle i| \quad (3.3.7)$$

3.3.b Settings for 3-dimensional MAD channels

Set $d = 3$; a generic 3-dimensional MAD channel $\Phi_\Gamma^{(3)}$ is uniquely identified by its transition matrix Γ :

$$\Gamma = \begin{pmatrix} 1 & 0 & 0 \\ \gamma_{10} & 1 - \gamma_{10} & 0 \\ \gamma_{20} & \gamma_{21} & 1 - \gamma_{20} - \gamma_{21} \end{pmatrix}, \quad (3.3.8)$$

while the Kraus set for $\Phi_{\Gamma}^{(3)}$, taken from (3.3.1), is:

$$\begin{aligned} K_{00} &= \begin{pmatrix} 1 & 0 & 0 \\ 0 & \sqrt{1-\gamma_{10}} & 0 \\ 0 & 0 & \sqrt{1-\gamma_{20}-\gamma_{21}} \end{pmatrix} \\ K_{01} &= \begin{pmatrix} 0 & \sqrt{\gamma_{10}} & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} \quad K_{02} = \begin{pmatrix} 0 & 0 & \sqrt{\gamma_{20}} \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} \quad K_{12} = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & \sqrt{\gamma_{21}} \\ 0 & 0 & 0 \end{pmatrix} \end{aligned} \quad (3.3.9)$$

The transition probabilities $\gamma_{10}, \gamma_{20}, \gamma_{21}$ satisfy the conditions:

$$\begin{cases} 0 \leq \gamma_{10}, \gamma_{20}, \gamma_{21} \leq 1 \\ 0 \leq \gamma_{20} + \gamma_{21} \leq 1 \end{cases} \quad (3.3.10)$$

Given an input density matrix ρ :

$$\rho = \begin{pmatrix} \rho_{00} & \rho_{01} & \rho_{02} \\ \rho_{01}^* & \rho_{11} & \rho_{12} \\ \rho_{02}^* & \rho_{12}^* & \rho_{22} \end{pmatrix}, \quad (3.3.11)$$

the output state of $\Phi_{\Gamma}^{(3)}$ is:

$$\Phi_{\Gamma}^{(3)}(\rho) = \begin{pmatrix} \gamma_{10}\rho_{11} + \gamma_{20}\rho_{22} + \rho_{00} & \sqrt{1-\gamma_{10}}\rho_{01} & \sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{02} \\ \sqrt{1-\gamma_{10}}\rho_{01}^* & (1-\gamma_{10})\rho_{11} + \gamma_{21}\rho_{22} & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{12} \\ \sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{02}^* & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{12}^* & (1-\gamma_{20}-\gamma_{21})\rho_{22} \end{pmatrix}, \quad (3.3.12)$$

while the output state of the complementary channel $\tilde{\Phi}_{\Gamma}^{(3)}$ is

$$\tilde{\Phi}_{\Gamma}^{(3)}(\rho) = \begin{pmatrix} (1-\gamma_{10})\rho_{11} + (1-\gamma_{20}-\gamma_{21})\rho_{22} + \rho_{00} & \sqrt{\gamma_{10}}\rho_{01} & \sqrt{\gamma_{20}}\rho_{02} & \sqrt{1-\gamma_{10}}\sqrt{\gamma_{21}}\rho_{12} \\ \sqrt{\gamma_{10}}\rho_{01}^* & \gamma_{10}\rho_{11} & \sqrt{\gamma_{10}}\sqrt{\gamma_{20}}\rho_{12}^* & 0 \\ \sqrt{\gamma_{20}}\rho_{02}^* & \sqrt{\gamma_{10}}\sqrt{\gamma_{20}}\rho_{12}^* & \gamma_{20}\rho_{22} & 0 \\ \sqrt{1-\gamma_{10}}\sqrt{\gamma_{21}}\rho_{12}^* & 0 & 0 & \gamma_{21}\rho_{22} \end{pmatrix}. \quad (3.3.13)$$

Refer to Figure 3.3.1 for a schematic representation of 3-dimensional MAD channels.

3.3.c Composition rules

The paper [CG21a] shows that MAD channels in $d = 3$ are closed under channel composition, which means that the composition of two 3-dimensional MAD channels is itself a 3-dimensional MAD channel. This result can be generalized to an arbitrary dimension d , as shown in Section 4.2. The composition rules of 3-dimensional MAD channels are:

$$\begin{aligned} \Phi_{\Gamma}^{(3)} &= \Phi_{\Gamma''}^{(3)} \circ \Phi_{\Gamma'}^{(3)} \\ \begin{cases} \gamma_{10} = \gamma'_{10} + \gamma''_{10}(1-\gamma'_{10}) \\ \gamma_{20} = \gamma'_{20} + \gamma''_{20}(1-\gamma'_{20}) + \gamma'_{21}\gamma''_{10} \\ \gamma_{21} = \gamma'_{21}(1-\gamma''_{10}) + \gamma''_{21}(1-\gamma'_{20}-\gamma'_{21}) \end{cases} \end{aligned} \quad (3.3.14)$$

where $\gamma_{ji} = \langle j|\Gamma|i\rangle, \gamma'_{ji} = \langle j|\Gamma'|i\rangle, \gamma''_{ji} = \langle j|\Gamma''|i\rangle$. Intuitively, the composition rules (3.3.14) suggest that the transition probability γ_{ji} , resulting from a composition of two MAD channels, is the sum of the probabilities all the possible two-step decay "paths" from $|j\rangle$ to $|i\rangle$. This idea is corroborated in Section 4.2.

3.3.d Covariance

Consider a d -dimensional MAD channel $\Phi_\Gamma : \sigma(\mathcal{H}) \mapsto \sigma(\mathcal{H})$ and consider the d -dimensional unitary matrices U diagonal in the computational basis:

$$U \equiv \sum_{j=0}^{d-1} e^{i\varphi_j} |j\rangle\langle j| \quad \varphi_j \in \mathbb{R} \quad \forall j. \quad (3.3.15)$$

It is possible to verify that Φ_Γ is covariant under the action of the group of unitary operators whose elements in the representation in \mathcal{H} are in the form (3.3.15), i.e.:

$$\Phi_\Gamma(U\rho U^\dagger) = U\Phi_\Gamma(\rho)U^\dagger. \quad (3.3.16)$$

3.3.e Maximum of coherent information for degradable MAD channels

If a channel Φ is covariant under the action of a group \mathfrak{G} , by employing (2.7.3) and the invariance of the von Neumann entropy under unitary operations, one obtains:

$$\begin{aligned} \Phi : \sigma(\mathcal{H}_A) &\mapsto \sigma(\mathcal{H}_B) \text{ covariant under } \mathfrak{G} \\ \Rightarrow I_c(U_g^A \rho U_g^{A\dagger}, \Phi) &= I_c(\rho, \Phi) \end{aligned} \quad (3.3.17)$$

Furthermore, consider the state $\bar{\rho}_{\mathfrak{G}}$ obtained by taking the average over all the applications of elements $g \in \mathfrak{G}$ upon the input state ρ :

$$\bar{\rho}_{\mathfrak{G}} \equiv \int d\mu_g U_g^A \rho U_g^{A\dagger}, \quad (3.3.18)$$

where μ_g is a probability distribution over the group \mathfrak{G} ; if Φ is degradable, by the property (2.10.13) the coherent information is concave in the input state ρ , therefore:

$$I_c(\bar{\rho}_{\mathfrak{G}}, \Phi) \geq \int d\mu_g I_c(U_g^A \rho U_g^{A\dagger}, \Phi) \stackrel{(3.3.17)}{=} I_c(\rho, \Phi) \quad (3.3.19)$$

Regarding MAD channels Φ_Γ , as seen in (3.3.16), \mathfrak{G} could be replaced by the group of unitary operators diagonal in the computational basis; in this case, $\bar{\rho}_{\mathfrak{G}}$ corresponds to a diagonal density matrices $\rho^{(diag)}$, so that (3.3.19) becomes:

$$I_c(\rho^{(diag)}, \Phi_\Gamma) \geq I_c(\rho, \Phi_\Gamma) \quad (3.3.20)$$

Hence, for degradable MAD channels, the search for the maximum over input states in (2.11.2) can be restricted to diagonal input states:

$$Q(\Phi_\Gamma) = Q^{(1)}(\Phi_\Gamma) \equiv \max_{\rho} I_c(\rho, \Phi_\Gamma) \stackrel{(3.3.20)}{=} \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_\Gamma) \quad (3.3.21)$$

3.3.f Single decays

Consider the single decay channels $\Phi_{\Gamma_{10}(\gamma)}^{(3)}, \Phi_{\Gamma_{20}(\gamma)}^{(3)}, \Phi_{\Gamma_{21}(\gamma)}^{(3)}$, whose transition matrices present the same single transition probability:

$$\Gamma_{10}(\gamma) \equiv \begin{pmatrix} 1 & 0 & 0 \\ \gamma & 1-\gamma & 0 \\ 0 & 0 & 1 \end{pmatrix} \quad \Gamma_{20}(\gamma) \equiv \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ \gamma & 0 & 1-\gamma \end{pmatrix} \quad \Gamma_{21}(\gamma) \equiv \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & \gamma & 1-\gamma \end{pmatrix} \quad (3.3.22)$$

Define the swap matrices:

$$U_{10} \equiv \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{pmatrix}, \quad U_{21} \equiv \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix}; \quad (3.3.23)$$

the channels $\Phi_{\Gamma_{10}(\gamma)}^{(3)}, \Phi_{\Gamma_{20}(\gamma)}^{(3)}, \Phi_{\Gamma_{21}(\gamma)}^{(3)}$ can be mapped into each other via unitary channels corresponding to (3.3.23):

$$\mathcal{U}_{10}(\bullet) \equiv U_{10} \bullet U_{10} \quad \mathcal{U}_{21}(\bullet) \equiv U_{21} \bullet U_{21} \quad (3.3.24)$$

$$\begin{aligned} \mathcal{U}_{21} \circ \Phi_{\Gamma_{10}(\gamma)}^{(3)} \circ \mathcal{U}_{21} &= \Phi_{\Gamma_{20}(\gamma)}^{(3)} \\ \mathcal{U}_{10} \circ \Phi_{\Gamma_{20}(\gamma)}^{(3)} \circ \mathcal{U}_{10} &= \Phi_{\Gamma_{21}(\gamma)}^{(3)} \\ \mathcal{U}_{10} \circ \mathcal{U}_{21} \circ \Phi_{\Gamma_{10}(\gamma)}^{(3)} \circ \mathcal{U}_{21} \circ \mathcal{U}_{10} &= \Phi_{\Gamma_{21}(\gamma)}^{(3)} \end{aligned} \quad (3.3.25)$$

From the bottleneck inequality (2.13.1), (3.3.25) and the invertibility of unitary channels, one can infer that $\Phi_{\Gamma_{10}(\gamma)}^{(3)}, \Phi_{\Gamma_{20}(\gamma)}^{(3)}, \Phi_{\Gamma_{21}(\gamma)}^{(3)}$ have the same capacity functionals, which means that the capacity of a single decay 3-dimensional MAD channel only depends on the transition probability, not on the specific levels involved in the decay.

The composition rules for $\Phi_{\Gamma_{10}(\gamma)}^{(3)}, \Phi_{\Gamma_{20}(\gamma)}^{(3)}, \Phi_{\Gamma_{21}(\gamma)}^{(3)}$ are very similar to those found for ADC's (3.1.5):

$$\begin{aligned} \Phi_{\Gamma_{ji}(\gamma)}^{(3)} &= \Phi_{\Gamma_{ji}(\gamma'')}^{(3)} \circ \Phi_{\Gamma_{ji}(\gamma')}^{(3)} \\ \gamma &= \gamma' + \gamma''(1 - \gamma') \geq \gamma', \gamma'', \end{aligned} \quad (3.3.26)$$

where $0 \leq i < j \leq 1$.

3.3.g Monotonicity

Exploiting the composition rules derived in Subsection 3.3.c, one can infer monotonicity properties for the capacity functionals of 3-dimensional MAD channels. In fact, the relation (2.13.1), in the context of (3.3.14), becomes:

$$\mathfrak{C}\left(\Phi_{\Gamma}^{(3)}\right) \leq \min \left\{ \mathfrak{C}\left(\Phi_{\Gamma'}^{(3)}\right), \mathfrak{C}\left(\Phi_{\Gamma''}^{(3)}\right) \right\} \quad (3.3.27)$$

Define a generic 3-dimensional MAD channel:

$$\Phi_{\Gamma}^{(3)} \equiv \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} \quad (3.3.28)$$

where the dependency of Γ in (3.3.8) on the real parameters $\gamma_{10}, \gamma_{20}, \gamma_{21}$ has been made explicit; using (3.3.14), one can find decompositions of $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$, whose composing channels have higher capacity than $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$. Some of these decompositions are reported below:

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(0, \gamma_{20}, \gamma_{21})}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, 0, 0)}^{(3)}, \quad (3.3.29)$$

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(0, \bar{\gamma}_{20}, 0)}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)} \quad \bar{\gamma}_{20} \equiv \frac{\gamma_{20}}{1 - \gamma_{21}}, \quad (3.3.30)$$

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(0, 0, \bar{\gamma}_{21})}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)} \quad \bar{\gamma}_{21} \equiv \frac{\gamma_{21}}{1 - \gamma_{20}}, \quad (3.3.31)$$

Employing (3.3.26) into (3.3.29), (3.3.30) and (3.3.31), each single decay channel found in these decompositions is split into two new single decay channels with transition probabilities $\gamma'_{ji}, \gamma''_{ji} \leq \gamma_{ji}$:

$$\Phi_{\Gamma(\gamma_{10}, 0, 0)}^{(3)} = \Phi_{\Gamma(\gamma'_1 0, 0, 0)}^{(3)} \circ \Phi_{\Gamma(\gamma''_{10}, 0, 0)}^{(3)}, \quad (3.3.32)$$

$$\Phi_{\Gamma(0, \bar{\gamma}_{20}, 0)}^{(3)} = \Phi_{\Gamma(0, \bar{\gamma}'_{20}, 0)}^{(3)} \circ \Phi_{\Gamma(0, \bar{\gamma}''_{20}, 0)}^{(3)}, \quad (3.3.33)$$

$$\Phi_{\Gamma(0, 0, \bar{\gamma}_{21})}^{(3)} = \Phi_{\Gamma(0, 0, \bar{\gamma}'_{21})}^{(3)} \circ \Phi_{\Gamma(0, 0, \bar{\gamma}''_{21})}^{(3)}. \quad (3.3.34)$$

where:

$$\bar{\gamma}''_{21} \equiv \frac{\gamma''_{21}}{1 - \gamma_{20}} \quad \bar{\gamma}'_{21} \equiv \frac{\gamma'_{21}}{1 - \gamma_{20}} \quad (3.3.35)$$

$$\bar{\gamma}''_{20} \equiv \frac{\gamma''_{20}}{1 - \gamma_{21}} \quad \bar{\gamma}'_{20} \equiv \frac{\gamma'_{20}}{1 - \gamma_{21}}. \quad (3.3.36)$$

Therefore, (3.3.29), (3.3.30) and (3.3.31) become:

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(\gamma'_{10}, \gamma_{20}, \gamma_{21})}^{(3)} \circ \Phi_{\Gamma(\gamma''_{10}, 0, 0)}^{(3)} \quad \gamma'_{10} \leq \gamma_{10}, \quad (3.3.37)$$

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(0, \bar{\gamma}''_{20}, 0)}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, \gamma'_{20}, \gamma_{21})}^{(3)} \quad \gamma'_{20} \leq \gamma_{20}, \quad (3.3.38)$$

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} = \Phi_{\Gamma(0, 0, \bar{\gamma}''_{21})}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma'_{21})}^{(3)} \quad \gamma'_{21} \leq \gamma_{21}. \quad (3.3.39)$$

Combining (3.3.37), (3.3.38) and (3.3.39) with (3.3.27), the monotonicity rules w.r.t. the transition probabilities for the capacities of 3-dimensional MAD channels are found.

$$\begin{aligned} \mathfrak{C}\left(\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}\right) &\leq \mathfrak{C}\left(\Phi_{\Gamma(\gamma'_{10}, \gamma_{20}, \gamma_{21})}^{(3)}\right) \quad \forall \gamma_{10} \geq \gamma'_{10}, \\ \mathfrak{C}\left(\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}\right) &\leq \mathfrak{C}\left(\Phi_{\Gamma(\gamma_{10}, \gamma'_{20}, \gamma_{21})}^{(3)}\right) \quad \forall \gamma_{20} \geq \gamma'_{20}, \\ \mathfrak{C}\left(\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}\right) &\leq \mathfrak{C}\left(\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma'_{21})}^{(3)}\right) \quad \forall \gamma_{21} \geq \gamma'_{21}. \end{aligned} \quad (3.3.40)$$

This means that, when any transition probability of a 3-dimensional MAD channel increases, its capacity functionals decrease.

3.3.h Quantum capacity and private classical capacity

In Section 2.11 the quantum capacity of quantum channels was introduced; in most cases, it is not possible to compute the exact value of this quantity for a given channel, however, one notable exception to this rule stems from degradable channels, whose quantum capacity corresponds to the maximum over all input states of the coherent information of the channel, as seen in (2.11.2). Another important simplification for degradable MAD channels is provided by (3.3.21), which restricts the search for the maximum of the coherent information over diagonal input states. For degradable 3-dimensional MAD channels, this translates to:

$$\begin{aligned} Q\left(\Phi_{\Gamma}^{(3)}\right) &= \max_{\rho} I_c\left(\rho, \Phi_{\Gamma}^{(3)}\right) = \max_{p_0, p_1} I_c\left(\rho^{(diag)}, \Phi_{\Gamma}^{(3)}\right) \\ \rho^{(diag)} &\equiv \begin{pmatrix} p_0 & 0 & 0 \\ 0 & p_1 & 0 \\ 0 & 0 & 1 - p_0 - p_1 \end{pmatrix} \quad \left\{ \begin{array}{l} 0 \leq p_0, p_1 \leq 1 \\ 0 \leq 1 - p_0 - p_1 \leq 1 \end{array} \right. \end{aligned} \tag{3.3.41}$$

Equation (3.3.21) (and, particularly for $d = 3$, (3.3.41)) allows for the development of a *modus operandi* for the analysis of the quantum capacity of d -dimensional MAD channels Φ_{Γ} :

- Find degradability and antidegradability regions for Φ_{Γ} .
- Compute the quantum capacity in the degradability regions using (3.3.21).
- Try to extend the computation to non-degradable zones.

The last step in this process, in the context of 3, 4-dimensional MAD channels, is to be expanded individually in each degradable zone.

3.3.i Degradability regions

A generic algorithm for determining the degradability regions of a d -dimensional MAD channel is offered in 4.6; it consists on finding the right-inverse map for a d -dimensional MAD channel Φ_{Γ} , denoted by $\tilde{\Phi}_{\Gamma}^{(-1)}$ and defined in (4.5.7), and checking the positivity of the Choi matrix C_{Λ} corresponding to the map:

$$\Lambda \equiv \tilde{\Phi}_{\Gamma} \circ \Phi_{\Gamma}^{-1}. \tag{3.3.42}$$

By 2.4.17, $C_{\Lambda} \geq 0$ if and only if Λ is completely positive; since Λ is linear and trace preserving by construction¹, its complete positiveness would imply that it is a quantum channel, acting as the degrading channel of Φ_{Γ} , which, in light of this, would be a degradable channel. To simplify this computation-heavy task, a more heuristic approach

¹This is a consequence of the linearity and trace preservation of both $\Phi_{\Gamma}^{(-1)}$ (see (4.5.7)) and $\tilde{\Phi}_{\Gamma}$.

is needed and provided here for the specific case of $d = 3$. Given a 3-dimensional MAD channel $\Phi_{\Gamma}^{(3)}$, in order to find a connecting channel $\Lambda^{(3)}$ such that:

$$\Lambda^{(3)} \circ \Phi_{\Gamma}^{(3)} = \tilde{\Phi}_{\Gamma}^{(3)} \quad (3.3.43)$$

assume that the rank of the output density matrix is greater or equal than the rank of the density matrix of the environment:

$$\text{rank} \left(\Phi_{\Gamma}^{(3)}(\rho) \right) \geq \text{rank} \left(\tilde{\Phi}_{\Gamma}^{(3)}(\rho) \right) \quad \forall \rho \quad (3.3.44)$$

For a generic ρ , $\text{rank} \left(\Phi_{\Gamma}^{(3)}(\rho) \right) = 3$; given a minimal Kraus set \mathcal{K} of a channel Φ , from (2.5.2) it can be deduced that the rank of $\tilde{\Phi}(\rho)$ corresponds to the cardinality of \mathcal{K} , hence, $\text{rank} \left(\tilde{\Phi}_{\Gamma}^{(3)}(\rho) \right) = 4$; the rank of the environment can be reduced to be ≤ 3 by "turning off" at least one decay, i.e. by setting at least one $\gamma_{ji} = 0$ for some $i \leq j$. This breaks down the problem to three scenarios:

1. $\gamma_{10} = 0$, in which case $\Phi_{\Gamma}^{(3)}$ is degradable for $\gamma_{20} + \gamma_{21} \leq 1/2$, see [below](#).
2. $\gamma_{20} = 0$, in which case $\Phi_{\Gamma}^{(3)}$ is never degradable, see [below](#).
3. $\gamma_{21} = 0$, in which case $\Phi_{\Gamma}^{(3)}$ is degradable for $\gamma_{10} \leq 1/2 \wedge \gamma_{20} \leq 1/2$, see [below](#).

Degradability for $\gamma_{10} = 0$

The output state of $\Phi_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)}$ is:

$$\Phi_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)}(\rho) = \begin{pmatrix} \gamma_{20}\rho_{22} + \rho_{00} & \rho_{01} & \sqrt{1 - \gamma_{20} - \gamma_{21}}\rho_{02} \\ \rho_{01}^* & \gamma_{21}\rho_{22} + \rho_{11} & \sqrt{1 - \gamma_{20} - \gamma_{21}}\rho_{12} \\ \sqrt{1 - \gamma_{20} - \gamma_{21}}\rho_{02}^* & \sqrt{1 - \gamma_{20} - \gamma_{21}}\rho_{12}^* & (1 - \gamma_{20} - \gamma_{21})\rho_{22} \end{pmatrix} \quad (3.3.45)$$

While the state of the environment is:

$$\tilde{\Phi}_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)}(\rho) = \begin{pmatrix} (1 - \gamma_{20} - \gamma_{21})\rho_{22} + \rho_{00} + \rho_{11} & \sqrt{\gamma_{20}}\rho_{02} & \sqrt{\gamma_{21}}\rho_{12} \\ \sqrt{\gamma_{20}}\rho_{02}^* & \gamma_{20}\rho_{22} & 0 \\ \sqrt{\gamma_{21}}\rho_{12}^* & 0 & \gamma_{21}\rho_{22} \end{pmatrix} \quad (3.3.46)$$

Given the degrading map

$$\Lambda_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)} \equiv \tilde{\Phi}_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)} \circ \Phi_{\Gamma(0,\gamma_{20},\gamma_{21})}^{(3)-1}, \quad (3.3.47)$$

checking the positivity of the associated Choi matrix yields the degradability condition:

$$\gamma_{20} + \gamma_{21} \leq \frac{1}{2} \quad (3.3.48)$$

Degradability for $\gamma_{20} = 0$

The output state of $\Phi_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)}$ is:

$$\Phi_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)}(\rho) = \begin{pmatrix} \gamma_{10}\rho_{11} + \rho_{00} & \sqrt{1-\gamma_{10}}\rho_{01} & \sqrt{1-\gamma_{21}}\rho_{02} \\ \sqrt{1-\gamma_{10}}\rho_{01}^* & (1-\gamma_{10})\rho_{11} + \gamma_{21}\rho_{22} & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{21}}\rho_{12} \\ \sqrt{1-\gamma_{21}}\rho_{02}^* & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{21}}\rho_{12}^* & (1-\gamma_{21})\rho_{22} \end{pmatrix} \quad (3.3.49)$$

While the state of the environment is:

$$\tilde{\Phi}_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)}(\rho) = \begin{pmatrix} (1-\gamma_{10})\rho_{11} + (1-\gamma_{21})\rho_{22} + \rho_{00} & \sqrt{\gamma_{10}}\rho_{01} & \sqrt{1-\gamma_{10}}\sqrt{\gamma_{21}}\rho_{12} \\ \sqrt{\gamma_{10}}\rho_{01}^* & \gamma_{10}\rho_{11} & 0 \\ \sqrt{1-\gamma_{10}}\sqrt{\gamma_{21}}\rho_{12}^* & 0 & \gamma_{21}\rho_{22} \end{pmatrix} \quad (3.3.50)$$

Since the Choi matrix associated to

$$\Lambda_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)} \equiv \tilde{\Phi}_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, 0, \gamma_{21})}^{(3)-1}, \quad (3.3.51)$$

is positive semi-definite only if either one of γ_{10}, γ_{21} is set to 0 (and the other is $\leq 1/2$), the channel $\Phi_{\Gamma(\gamma_{10} \neq 0, 0, \gamma_{21} \neq 0)}^{(3)}$ is never degradable.

Degradability for $\gamma_{21} = 0$

The output state of $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)}$ is:

$$\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)}(\rho) = \begin{pmatrix} \gamma_{10}\rho_{11} + \gamma_{20}\rho_{22} + \rho_{00} & \sqrt{1-\gamma_{10}}\rho_{01} & \sqrt{1-\gamma_{20}}\rho_{02} \\ \sqrt{1-\gamma_{10}}\rho_{01}^* & (1-\gamma_{10})\rho_{11} & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}}\rho_{12} \\ \sqrt{1-\gamma_{20}}\rho_{02}^* & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}}\rho_{12}^* & (1-\gamma_{20})\rho_{22} \end{pmatrix} \quad (3.3.52)$$

While the state of the environment is:

$$\tilde{\Phi}_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)}(\rho) = \begin{pmatrix} (1-\gamma_{10})\rho_{11} + (1-\gamma_{20})\rho_{22} + \rho_{00} & \sqrt{\gamma_{10}}\rho_{01} & \sqrt{\gamma_{20}}\rho_{02} \\ \sqrt{\gamma_{10}}\rho_{01}^* & \gamma_{10}\rho_{11} & \sqrt{\gamma_{10}}\sqrt{\gamma_{20}}\rho_{12} \\ \sqrt{\gamma_{20}}\rho_{02}^* & \sqrt{\gamma_{10}}\sqrt{\gamma_{20}}\rho_{12}^* & \gamma_{20}\rho_{22} \end{pmatrix} \quad (3.3.53)$$

Given the degrading map

$$\Lambda_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)} \equiv \tilde{\Phi}_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)} \circ \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, 0)}^{(3)-1}, \quad (3.3.54)$$

checking the positivity of the associated Choi matrix yields the degradability condition:

$$\gamma_{10} \leq \frac{1}{2} \wedge \gamma_{20} \leq \frac{1}{2} \quad (3.3.55)$$

Degradability conditions

Compounding the results exposed in this Subsection, the degradability conditions for $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$ are found:

$$\begin{aligned} \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} \text{ is degradable} \\ \uparrow \\ \left(\gamma_{10} = 0 \wedge \gamma_{20} + \gamma_{21} \leq \frac{1}{2} \right) \vee \left(\gamma_{21} = 0 \wedge \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{20} \leq \frac{1}{2} \right) \end{aligned} \quad (3.3.56)$$

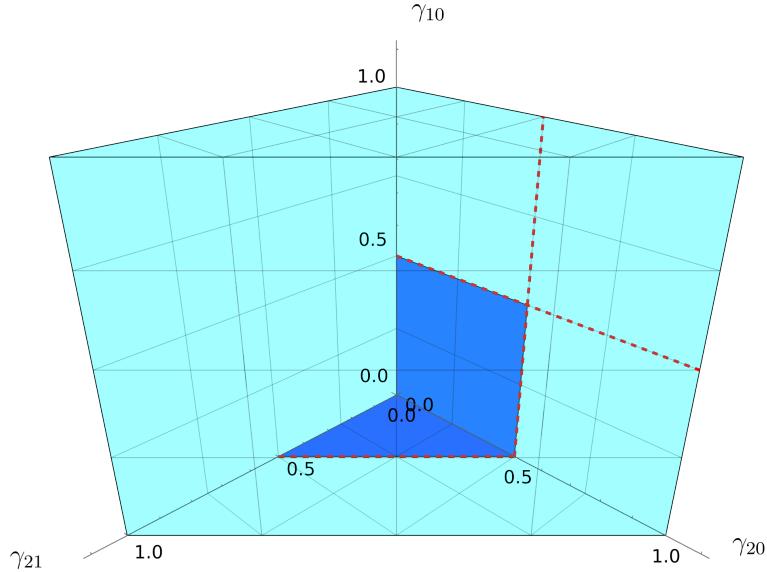


Figure 3.3.2: The plot is cast in the parameter space for a 3-dimensional MAD channel; all these channels satisfy the conditions (3.3.10), which are represented in cyan in the plot, while the blue areas correspond to the degradability conditions reported in (3.3.57)

The conditions in (3.3.56) are illustrated in Figure 3.3.2.

Additional degradable settings

One may wonder if the conditions in (3.3.56) encompass all the degradability regions for $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$; particularly, it is not clear whether there exists a configuration of $\{\gamma_{10}, \gamma_{20}, \gamma_{21}\}$ such that $\gamma_{10}, \gamma_{20}, \gamma_{21} \neq 0$ and $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$ is degradable. To tackle this question, it is possible to exploit the result in (4.4.2), which states that the channels forming an arbitrary decomposition of a degradable channel must be degradable themselves. The right-most channel in the decomposition (3.3.30) of a generic 3-dimensional MAD channel is only degradable if either one of γ_{10}, γ_{21} is set to 0, therefore it is not possible to find a configuration $\{\gamma_{10}, \gamma_{20}, \gamma_{21}\}$ such that $\gamma_{10}, \gamma_{20}, \gamma_{21} \neq 0$ and $\Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)}$ is degradable. Thus, (3.3.56) can be rewritten in a more strict form:

$$\begin{aligned} \Phi_{\Gamma(\gamma_{10}, \gamma_{20}, \gamma_{21})}^{(3)} \text{ is degradable} \\ \Updownarrow \\ \left(\gamma_{10} = 0 \wedge \gamma_{20} + \gamma_{21} \leq \frac{1}{2} \right) \vee \left(\gamma_{21} = 0 \wedge \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{20} \leq \frac{1}{2} \right) \end{aligned} \tag{3.3.57}$$

Note that this means that in the case of 3-dimensional MAD channels, the condition $\text{rank}(\Phi_{\Gamma}^{(3)}(\rho)) \geq \text{rank}(\tilde{\Phi}_{\Gamma}^{(3)}(\rho)) \forall \rho$ is a necessary condition for the degradability of

$\Phi_{\Gamma}^{(3)}$; this is not true in general, a counterexample for 4-dimensional MAD's is offered in (4.7.9).

3.3.j Quantum capacity of single decay 3-dimensional MAD channel

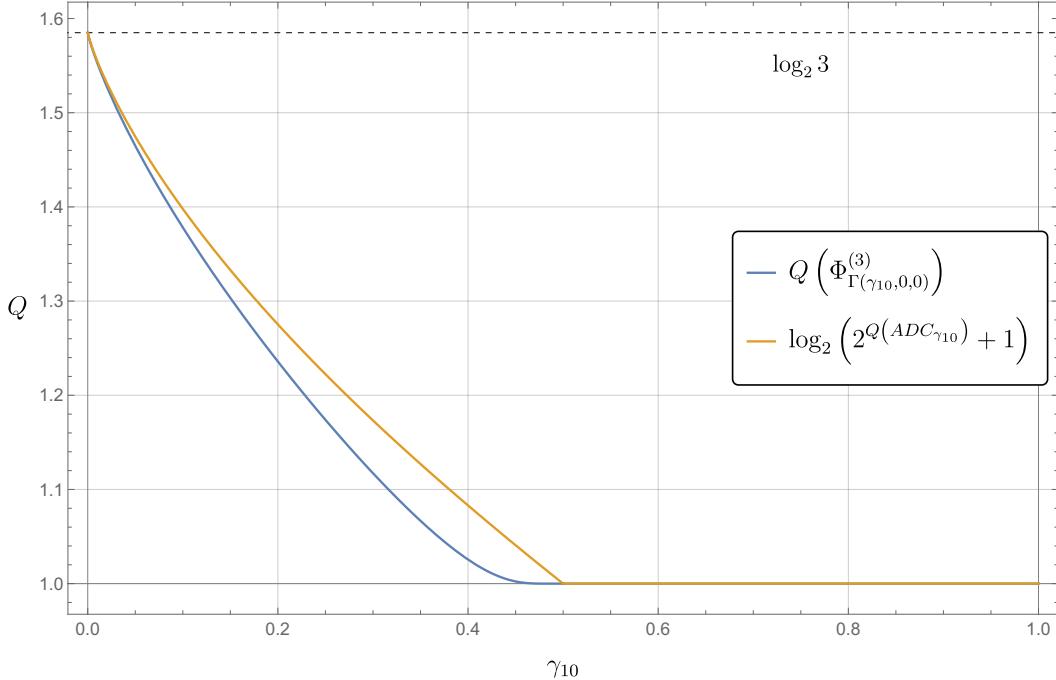


Figure 3.3.3: The quantum capacity of a single decay 3-dimensional MAD channel is plotted here in blue, while the upper bound corresponding to (3.2.4) is plotted in orange.

Consider the channel $\Phi_{\Gamma(\gamma_{10},0,0)}^{(3)}$; as a consequence of the findings in 3.3.f, the capacity functionals of this channel are the same as those of the other single decay 3-dimensional MAD's. Furthermore, according to (3.3.57), the channel is degradable if and only if $0 \leq \gamma_{10} \leq 1/2$. This channel exhibits the PCDS structure [CG21b] (3.2.1), where Φ_{AA} corresponds to the amplitude damping channel $ADC_{\gamma_{10}}$ and $\Phi_{BB} = \mathbb{1}_{BB}$, a property that will become useful in the analysis of its quantum capacity. For the degradable region $0 \leq \gamma_{10} \leq 1/2$, the quantum capacity of the channel $Q\left(\Phi_{\Gamma(\gamma_{10},0,0)}^{(3)}\right)$ can be computed using (3.3.41); at $\gamma_{10} = 1/2$, $ADC_{\gamma_{10}}$ is degradable and has null capacity, thus (3.2.6) holds:

$$Q\left(\Phi_{\Gamma(\gamma_{10}=1/2,0,0)}^{(3)}\right) = 1 \quad (3.3.58)$$

Additionally, since the channel presents a noiseless subspace spanned by $\{|0\rangle, |1\rangle\}$, 1 is also a lower bound for the quantum capacity of $\Phi_{\Gamma(\gamma_{10},0,0)}^{(3)}$:

$$Q\left(\Phi_{\Gamma(\gamma_{10},0,0)}^{(3)}\right) \geq 1. \quad (3.3.59)$$

Hence, combining (3.3.59), (3.3.58) and the monotonicity properties (3.3.40):

$$Q\left(\Phi_{\Gamma(\gamma_{10},0,0)}^{(3)}\right) = 1 \quad \forall \frac{1}{2} \leq \gamma_{10} \leq 1 \quad (3.3.60)$$

The computations for the quantum capacity of a single decay 3-dimensional MAD channel are plotted in Figure 3.3.3.

4

Properties of MAD channels and 4-dimensional MAD channels

This chapter will focus on the exposition of some general properties of Multi-Level Amplitude Damping channels that were found during this work of thesis. Most of these properties hold for any dimension d , potentially allowing future studies of MAD channels in higher dimensions to be built upon these results.

4.1 Settings for d -dimensional MAD channels

The settings for d -dimensional MAD channels have already been defined in Subsection 3.3.a. The relation (3.3.5), in particular, will turn out to be extremely useful in Section 4.2 for the derivation of the composition rules of MAD channels.

Settings for 4-dimensional MAD channels

Substituting $d = 4$ in the definitions given in 3.3.a, one obtains the settings that describe a 4-dimensional MAD channel $\Phi_{\Gamma}^{(4)}$, which is completely identified by the transition

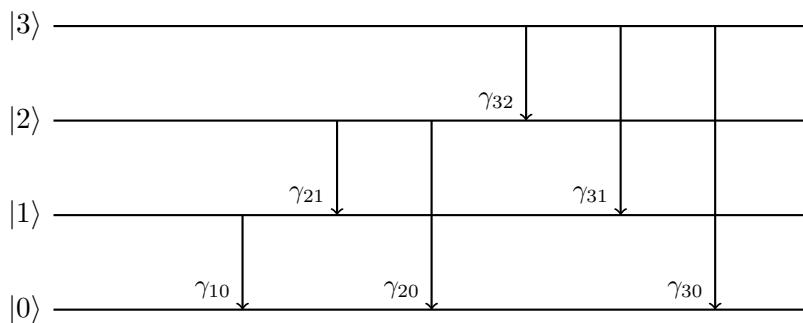


Figure 4.1.1: MAD channels represent decay processes, where each level of a system has a fixed probability of decaying onto a lower level. Here, a schematic depiction of a 4-dimensional MAD is reported

matrix Γ :

$$\Gamma \equiv \begin{pmatrix} 1 & 0 & 0 & 0 \\ \gamma_{10} & 1 - \gamma_{10} & 0 & 0 \\ \gamma_{20} & \gamma_{21} & 1 - \gamma_{20} - \gamma_{21} & 0 \\ \gamma_{30} & \gamma_{31} & \gamma_{32} & 1 - \gamma_{30} - \gamma_{31} - \gamma_{32} \end{pmatrix}, \quad (4.1.1)$$

whose elements satisfy are real, ≥ 0 and ≤ 1 , accordingly to (3.3.3). The Kraus set of $\Phi_{\Gamma}^{(4)}$ is composed by the Kraus operators:

$$\begin{aligned} K_{00} &= \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & \sqrt{1-\gamma_{10}} & 0 & 0 \\ 0 & 0 & \sqrt{1-\gamma_{20}-\gamma_{21}} & 0 \\ 0 & 0 & 0 & \sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}} \end{pmatrix} \\ K_{01} &= \begin{pmatrix} 0 & \sqrt{\gamma_{10}} & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} K_{02} = \begin{pmatrix} 0 & 0 & \sqrt{\gamma_{20}} & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} K_{12} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & \sqrt{\gamma_{21}} & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} \\ K_{03} &= \begin{pmatrix} 0 & 0 & 0 & \sqrt{\gamma_{30}} \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} K_{13} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & \sqrt{\gamma_{31}} \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} K_{23} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & \sqrt{\gamma_{32}} \\ 0 & 0 & 0 & 0 \end{pmatrix} \end{aligned} \quad (4.1.2)$$

Given a generic input $\rho \in \sigma(\mathcal{H}_4)$:

$$\rho \equiv \begin{pmatrix} \rho_{00} & \rho_{01} & \rho_{02} & \rho_{03} \\ \rho_{01}^* & \rho_{11} & \rho_{12} & \rho_{13} \\ \rho_{02}^* & \rho_{12}^* & \rho_{22} & \rho_{23} \\ \rho_{03}^* & \rho_{13}^* & \rho_{23}^* & \rho_{33} \end{pmatrix}, \quad (4.1.3)$$

the corresponding output state of the channel $\Phi_{\Gamma}^{(4)}$ is:

$$\Phi_{\Gamma}^{(4)}(\rho) = \begin{pmatrix} \rho_{00} + \gamma_{10}\rho_{11} + \gamma_{20}\rho_{22} + \gamma_{30}\rho_{33} & \sqrt{1-\gamma_{10}}\rho_{01} & \sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{02} & \sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}}\rho_{03} \\ \rho_{01}^*\sqrt{1-\gamma_{10}} & (1-\gamma_{10})\rho_{11} + \gamma_{21}\rho_{22} + \gamma_{31}\rho_{33} & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}-\gamma_{21}}\rho_{12} & \sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}}\rho_{13} \\ \rho_{02}^*\sqrt{1-\gamma_{20}-\gamma_{21}} & \rho_{12}^*\sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{20}-\gamma_{21}} & (1-\gamma_{20}-\gamma_{21})\rho_{22} + \gamma_{32}\rho_{33} & \sqrt{1-\gamma_{20}-\gamma_{21}}\sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}}\rho_{23} \\ \rho_{03}^*\sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}} & \rho_{13}^*\sqrt{1-\gamma_{10}}\sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}} & \rho_{23}^*\sqrt{1-\gamma_{20}-\gamma_{21}}\sqrt{1-\gamma_{30}-\gamma_{31}-\gamma_{32}} & (1-\gamma_{30}-\gamma_{31}-\gamma_{32})\rho_{33} \end{pmatrix} \quad (4.1.4)$$

while the output state of the complementary channel $\tilde{\Phi}_{\Gamma}^{(4)}$ is given in TODO (La matrice non ci entra bene, l'ho stampata in 7.2, in caso la tolgo). Refer to Figure 4.1.1 for a schematic representation of 4-dimensional MAD channels.

4.2 Composition of MAD channels

As seen in (3.3.14) and (3.1.5), MAD channels in $d = 2$ (ADC's) and $d = 3$ are closed under composition, meaning that the composition of two d -dimensional MAD channels for $d = 2, 3$ is still a d -dimensional MAD channel, whose transition probabilities can be found and depend on the transition probabilities of the starting channels. In this section,

a general rule for the composition of d -dimensional will be derived. The closeness of these channels under composition will be first assumed and then verified *a posteriori* using the composition rules that were found. First, one needs to define three transition matrices $\Gamma, \Gamma', \Gamma''$, whose transition probabilities are, respectively, $\gamma_{ji}, \gamma'_{ji}, \gamma''_{ji}$; then, one reminds that, substituting (3.3.1) into (2.4.4), MAD channels assume the form:

$$\Phi_\Gamma(\rho) = \sum_{i=0}^{j-1} \sum_{j=1}^{d-1} \gamma_{ji} \rho_{jj} |i\rangle\langle i| + \sum_{i=0}^{d-1} \sum_{j=1}^{d-1} \sqrt{\gamma_{ii}\gamma_{jj}} \rho_{ij} |i\rangle\langle j|. \quad (4.2.1)$$

One may try to write:

$$\Phi_\Gamma = \Phi_{\Gamma''} \circ \Phi_{\Gamma'} \quad (4.2.2)$$

and try to find a function f such that:

$$\Gamma = f(\Gamma', \Gamma''). \quad (4.2.3)$$

If one were to write $\Phi_{\Gamma''} \circ \Phi_{\Gamma'}(\rho)$ in the form (4.2.1), they would obtain:

$$\begin{aligned} \Phi_{\Gamma''} \circ \Phi_{\Gamma'}(\rho) &= \sum_{i=0}^{l-1} \sum_{l=1}^{j-1} \sum_{j=1}^{d-1} \gamma''_{li} \gamma'_{jl} \rho_{jj} |i\rangle\langle i| + \\ &\quad \sum_{i=0}^{j-1} \sum_{j=1}^{d-1} \gamma''_{ii} \gamma'_{ji} \rho_{jj} |i\rangle\langle i| + \\ &\quad \sum_{i=0}^{j-1} \sum_{j=1}^{d-1} \gamma''_{ji} \gamma'_{jj} \rho_{jj} |i\rangle\langle i| + \\ &\quad \sum_{i=0}^{d-1} \sum_{j=1}^{d-1} \sqrt{\gamma''_{ii} \gamma'_{ii} \gamma''_{jj} \gamma'_{jj}} \rho_{ij} |i\rangle\langle j|, \end{aligned} \quad (4.2.4)$$

which, employing the last property in (3.3.3), translates to

$$\Phi_{\Gamma''} \circ \Phi_{\Gamma'}(\rho) = \sum_{i=0}^{j-1} \sum_{j=1}^{d-1} \left(\sum_{l=0}^{d-1} \gamma''_{li} \gamma'_{jl} \right) \rho_{jj} |i\rangle\langle i| + \sum_{i=0}^{d-1} \sum_{j=1}^{d-1} \sqrt{\gamma''_{ii} \gamma'_{ii} \gamma''_{jj} \gamma'_{jj}} \rho_{ij} |i\rangle\langle j| \quad (4.2.5)$$

This would represent the right hand side of (4.2.2) for a generic input ρ . By direct computation, it is possible to verify that fixing the composition laws:

$$\gamma_{ji} = \sum_{l=0}^{d-1} \gamma''_{li} \gamma'_{jl} \quad \forall i \leq j \quad (4.2.6)$$

reduces (4.2.5) to (4.2.1). Therefore, it is possible to infer that the composition of two d -dimensional MAD channels is a d -dimensional MAD channel, whose transition probabilities depend on those of the initial channels as in (4.2.6). In terms of transition matrices, (4.2.6) can easily be obtained by setting $f(\Gamma', \Gamma'') \equiv \Gamma' \Gamma''$ in (4.2.3), so that the composition rules of MAD channels can be summarized by:

$$\begin{aligned} \Phi_\Gamma &= \Phi_{\Gamma''} \circ \Phi_{\Gamma'} \\ \Gamma &= \Gamma' \Gamma'' \end{aligned} \quad (4.2.7)$$

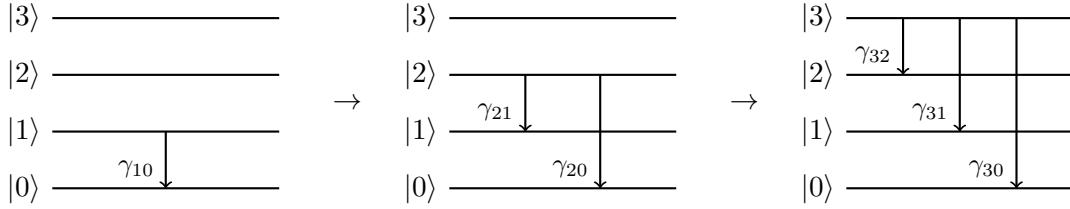


Figure 4.2.1: Visual representation of the decomposition of a 4-dimensional MAD channel using (4.2.11), read from left to right in "chronological" order.

4.2.a Useful decompositions of MAD channels

Following (4.2.7), one could find various decompositions of a generic MAD channels which may help to simplify the process of finding the capacity of the channel. In this section, the most intuitive decompositions will be listed.

Separated decays from increasing levels

This decomposition is the easiest to derive and also the most useful; it is found by defining the matrices:

$$\begin{aligned}\Gamma_k &\equiv \mathbb{1}_d + \sum_{i=0}^{k-1} \gamma_{ki} |k\rangle\langle i| - \sum_{i=0}^{k-1} \gamma_{ki} |k\rangle\langle k| \\ \Gamma^{(k)} &\equiv \mathbb{1}_d + \sum_{j=1}^{k-1} \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle i| - \sum_{j=1}^{k-1} \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle j|,\end{aligned}\tag{4.2.8}$$

where \$k < d\$ and \$\Gamma_k\$ represents a special kind of MAD channel where only the level \$|k\rangle\$ is allowed to decay, while \$\Gamma^{(k)}\$ represents a MAD channel where the decays from levels \$|k\rangle\$ and upwards are forbidden. In this setting, the most generic \$d\$-dimensional transition matrix \$\Gamma\$ in (3.3.4) is equal to \$\Gamma^{(d)}\$. By direct computation, setting \$k = d - 1\$ in (4.2.8) leads to:

$$\Gamma^{(d)} = \Gamma^{(d-1)} \Gamma_{d-1},\tag{4.2.9}$$

then, by iterating (4.2.9), one arrives at the decomposition:

$$\Gamma = \Gamma^{(d)} = \Gamma_1 \Gamma_2 \dots \Gamma_{d-1}\tag{4.2.10}$$

which, by employing (4.2.7), translates to:

$$\Phi_\Gamma = \Phi_{\Gamma_{d-1}} \circ \dots \circ \Phi_{\Gamma_1}.\tag{4.2.11}$$

Intuitively speaking, (4.2.11) means that in a MAD channel, the lower energy levels "have precedence" when decaying, as represented in Figure 4.2.1.

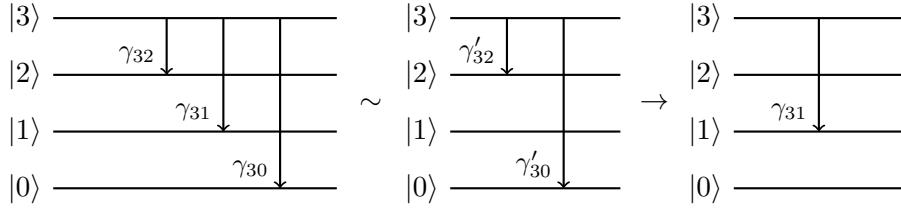


Figure 4.2.2: Example of the decomposition in (4.2.13), having fixed $k = 3, n = 1$ for 4 dimensional MAD channels.

The matrices Γ_k can be decomposed even further: one needs to define the matrices:

$$\begin{aligned}\Gamma_k^{(n)} &\equiv \mathbb{1}_d + \gamma'_{kn} |k\rangle\langle n| - \gamma'_{kn} |k\rangle\langle k| \\ T_k^{(n)} &\equiv \mathbb{1}_d + \sum_{\substack{i=0 \\ i \neq n}}^{k-1} \gamma_{ki} |k\rangle\langle i| - \sum_{\substack{i=0 \\ i \neq n}}^{k-1} \gamma_{ki} |k\rangle\langle k| \\ T_k^{(n)} &= \Gamma_k - \gamma_{kn} |k\rangle\langle n| + \gamma_{kn} |k\rangle\langle k|\end{aligned}\quad (4.2.12)$$

where $n < k$; these matrices compose in the following way:

$$\begin{aligned}\gamma'_{kn} &\equiv \gamma_{kn} \left(1 - \sum_{\substack{i=0 \\ i \neq n}}^{k-1} \gamma_{ki} \right)^{-1} \\ \Gamma_k &= T_k^{(n)} \Gamma_k^{(n)}\end{aligned}\quad (4.2.13)$$

The physical interpretation is clear: in this setting, given the decays from the level $|k\rangle$, the decay onto the level $|n\rangle$ needs to be performed at the end of the process. In order to achieve this, one needs to modify the amplitude γ_{kn} as in (4.2.13); a schematic graph of this decomposition is reported in Figure 4.2.2. Note that the denominator of γ'_{kn} in (4.2.13) does not depend on γ_{kn} .

MAD channel as composition of single decays

One may want to expand on the idea behind (4.2.13) in order to create a composition of single transition matrices. This can be achieved by a slight redefinition of the matrices in (4.2.12):

$$\begin{aligned}\Xi_k^{(n)} &\equiv \mathbb{1}_d + \gamma_{kn}^{(n+1)} |k\rangle\langle n| - \gamma_{kn}^{(n+1)} |k\rangle\langle k| \\ H_k^{(n)} &\equiv \mathbb{1}_d + \sum_{i=n}^{k-1} \gamma_{ki} |k\rangle\langle i| - \sum_{i=n}^{k-1} \gamma_{ki} |k\rangle\langle k|\end{aligned}\quad (4.2.14)$$

where $n < k$. The $\Xi_k^{(n)}$'s represent single decay MAD channels with modified amplitudes w.r.t. the original channel, while the $H_k^{(n)}$'s represent MAD channels where only the

$|k\rangle$ level can decay, with the same amplitudes as the original channel and having the transitions onto the lowest n levels forbidden. One can verify that:

$$H_k^{(0)} = \Gamma_k \quad (4.2.15)$$

which matches the descriptive definition given above. In order to find the values of $\gamma_{kn}^{(n+1)}$ in terms of the original amplitudes, one must try to solve the iterative equation:

$$H_k^{(n)} = H_k^{(n+1)} \Xi_k^{(n)} \quad (4.2.16)$$

The (4.2.16) fixes the values of $\gamma_{kn}^{(n+1)}$:

$$\gamma_{kn}^{(n+1)} \equiv \frac{\gamma_{kn}}{1 - \sum_{i=n+1}^{k-1} \gamma_{ki}} \quad (4.2.17)$$

which also implies that:

$$H_k^{(k-1)} = \Xi_k^{(k-1)} \quad (4.2.18)$$

Using the iterative (4.2.16) and employing (4.2.15), (4.2.18) allows for the decomposition of Γ_k in single level transition matrices:

$$\begin{aligned} \Gamma_k &= H_k^{(0)} \\ &= H_k^{(1)} \Xi_k^{(0)} \\ &= H_k^{(2)} \Xi_k^{(1)} \Xi_k^{(0)} \\ &\dots \\ &= \Xi_k^{(k-1)} \dots \Xi_k^{(0)} \end{aligned} \quad (4.2.19)$$

Combining (4.2.10) and (4.2.19), it is possible to consider a generic MAD channel as a composition of single-decay MAD channels, with appropriately modified amplitudes:

$$\Gamma = \prod_{\substack{k=1 \\ \rightarrow}}^{d-1} \left(\Xi_k^{(k-1)} \dots \Xi_k^{(0)} \right) \quad (4.2.20)$$

where \prod_{\rightarrow} indicates that the product is meant to be expanded from left to right for increasing k 's. The (4.2.20) translates to:

$$\Phi_\Gamma = \bigodot_{\substack{k=1 \\ \leftarrow}}^{d-1} \left(\Phi_{\Xi_k^{(0)}} \circ \dots \circ \Phi_{\Xi_k^{(k-1)}} \right) \quad (4.2.21)$$

where \bigodot_{\leftarrow} indicates a composition of channels that is meant to be expanded from right to left for increasing k 's.

4.3 Equivalence of single decay MAD channels

Expanding on the ideas introduced in Subsection 3.3.f, it is possible to prove that the capacity functionals of any single decay d -dimensional MAD channel only depend on the single transition probability, not on the levels involved in the decay. Consider the swap unitaries:

$$U_{mn} \equiv \mathbb{1}_{\mathcal{H}_d} - |m\rangle\langle m| - |n\rangle\langle n| + |m\rangle\langle n| + |n\rangle\langle m| \quad m, n \in \{0, \dots, d-1\}, \quad (4.3.1)$$

and the associated unitary quantum channels:

$$\mathcal{U}_{mn}(\bullet) \equiv U_{mn} \bullet U_{mn}; \quad (4.3.2)$$

note that, if $m = n$, then $\mathcal{U}_{mn} = \mathbb{1}_{\sigma(\mathcal{H}_d)}$. Let $\Gamma_{ji}(\gamma)$, $j > i$, be a single decay transition matrix of transition probability γ :

$$\Gamma_{ji}(\gamma) \equiv \mathbb{1}_{\mathcal{H}_d} + \gamma |j\rangle\langle i| - \gamma |j\rangle\langle j|, \quad (4.3.3)$$

It can be shown that:

$$\begin{aligned} \mathcal{U}_{kj} \circ \Phi_{\Gamma_{ji}(\gamma)} \circ \mathcal{U}_{kj} &= \Phi_{\Gamma_{ki}(\gamma)} \quad k > i \\ \mathcal{U}_{ki} \circ \Phi_{\Gamma_{ji}(\gamma)} \circ \mathcal{U}_{ki} &= \Phi_{\Gamma_{jk}(\gamma)} \quad k < j \end{aligned} \quad (4.3.4)$$

Utilizing the relations (4.3.4), it is possible to transform $\Phi_{\Gamma_{ji}(\gamma)}$ into another single decay MAD channel $\Phi_{\Gamma_{j'i'}(\gamma)}$, $j' > i'$. In order to achieve this, one needs to distinguish the two cases where $j' \geq j$ or $j' < j$:

$$\begin{aligned} j' \geq j \Rightarrow \Phi_{\Gamma_{j'i'}(\gamma)} &= \mathcal{U}_{i'i} \circ \mathcal{U}_{j'j} \circ \Phi_{\Gamma_{ji}(\gamma)} \circ \mathcal{U}_{j'j} \circ \mathcal{U}_{i'i} \\ j' < j \Rightarrow \Phi_{\Gamma_{j'i'}(\gamma)} &= \mathcal{U}_{j'j} \circ \mathcal{U}_{i'i} \circ \Phi_{\Gamma_{ji}(\gamma)} \circ \mathcal{U}_{i'i} \circ \mathcal{U}_{j'j} \end{aligned} \quad (4.3.5)$$

Employing (4.3.5) and the pipeline inequalities (2.13.1), one can conclude that the capacity functionals of single decay MAD channels depend only on their transition probability, not on the levels involved in the decay. These covariance relations were to be expected, as the unitaries involved represent a swap in the "labels" of the levels upon which they act. As long as changing the labels does not require modifying the nature of the channel (e.g. swapping the labels $|2\rangle$ and $|1\rangle$ for $\Phi_{\Gamma_{21}(\gamma)}$ implies that the channel is no longer a MAD channel, as the decay becomes a transition from a lower to a higher energy level), those operations are always permitted.

4.4 Composition of degradable channels

The topic of this section is not strictly related to MAD channels; however, its usefulness in the derivation of the degradability zone of MAD's will become apparent in section

[4.7.](#) Given two LCPT maps Ψ_1, Ψ_2 and their composition $\Psi \equiv \Psi_2 \circ \Psi_1$, it will be shown that:

$$\begin{aligned} \Psi &= \Psi_2 \circ \Psi_1 \\ \Psi \text{ degradable} &\Rightarrow \Psi_1, \Psi_2 \text{ degradable} \end{aligned} \quad (4.4.1)$$

Or, more eloquently:

$$\begin{aligned} \Psi &= \Psi_2 \circ \Psi_1 \\ \exists \Lambda \text{ LCPT} : \Lambda \circ \Psi = \tilde{\Psi} &\Rightarrow \begin{cases} \exists \Lambda_1 \text{ LCPT} : \Lambda_1 \circ \Psi_1 = \tilde{\Psi}_1 \\ \exists \Lambda_2 \text{ LCPT} : \Lambda_2 \circ \Psi_2 = \tilde{\Psi}_2 \end{cases} \end{aligned} \quad (4.4.2)$$

Complementary channels equivalence

The complementary channel $\tilde{\Psi}$ is unitarily equivalent to its counterpart written in terms of the channels $\Psi_{1,2}$:

$$\tilde{\Psi}'(\rho) = \tilde{\Psi}_1(\rho)_{E1} \otimes \tilde{\Psi}_2(\Psi_1(\rho))_{E2}, \quad (4.4.3)$$

i.e. the output of the complementary channel of Ψ , in this representation, corresponds to the tensor product of two environment states in the systems $E1, E2$. By [\(2.5.4\)](#), there must exist a unitary transformation \mathfrak{V} such that:

$$\tilde{\Psi}'(\rho) = \mathfrak{V} \left(\tilde{\Psi}(\rho) \right) = V \tilde{\Psi}(\rho) V^\dagger. \quad (4.4.4)$$

Ψ_2 is degradable

Define $\rho_1 \equiv \Psi_1(\rho)$; using this definition, the degradability hypothesis of Ψ reads:

$$\tilde{\Psi}(\rho) = \Lambda \circ \Psi_2(\rho_1). \quad (4.4.5)$$

One can then build the LCPT channel Ω_2 :

$$\Omega_2 \equiv \text{tr}_{E1} \circ \mathfrak{V}, \quad (4.4.6)$$

where tr_{E1} is the LCPT channel describing the partial trace over the $E1$ environment. Applying Ω_2 to [\(4.4.5\)](#) leads to:

$$\tilde{\Psi}_2(\rho_1) = \Omega_2 \circ \Lambda \circ \Psi_2(\rho_1) \quad (4.4.7)$$

Therefore, if one thought of Ψ_1 as a state preparation device, by defining the LCPT $\Lambda_2 \equiv \Omega_2 \circ \Lambda$ one would obtain:

$$\tilde{\Psi}_2 = \Lambda_2 \circ \Psi_2 \quad (4.4.8)$$

proving the degradability of Φ_2 .

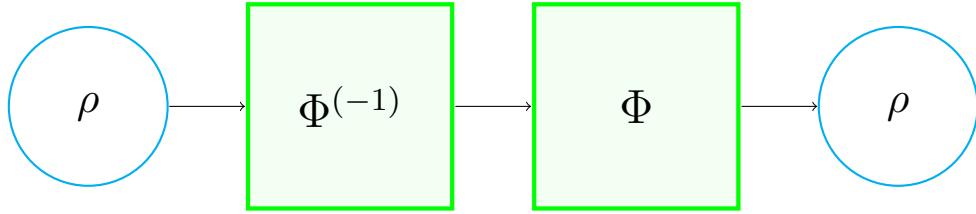


Figure 4.5.1: The map $\Phi^{(-1)}$ acts as the right inverse of the MAD channel Φ

Ψ_1 is degradable

Define the LCPT map:

$$\Omega_1 \equiv \text{tr}_{E2} \circ \mathfrak{V} \quad (4.4.9)$$

Starting from $\tilde{\Psi} = \Lambda \circ \Psi$, apply Ω_1 to both sides, which yields:

$$\tilde{\Psi}_1 = \Omega_1 \circ \Lambda \circ \Psi_2 \circ \Psi_1. \quad (4.4.10)$$

Then, one could define the LCPT

$$\Lambda_1 = \Omega_1 \circ \Lambda \circ \Psi_2 \quad (4.4.11)$$

which can be substituted into (4.4.10):

$$\tilde{\Psi}_1 = \Lambda_1 \circ \Psi_1. \quad (4.4.12)$$

(4.4.12) and (4.4.8) together prove (4.4.2).

4.5 Inverse maps of MAD channels

It is not always possible to find the (left or right) inverse map of a LCPT channel, a clear example of this fact are the channels which send a D -dimensional system into a d -dimensional system, where $d < D$. This is not the case for MAD channels; in fact it turns out that their inverse maps has a rather intuitive, albeit convoluted, definition. In the following, the right inverse map (represented in Figure 4.5.1) of a generic MAD channel will be derived.

4.5.a Inverse maps of ADC's

The inverse map of an ADC provides a valuable guide for the derivation of the inverse of a MAD channel. The reader can find a description of ADC's in 3.1. The scope of this subsection is to find the right inverse map of an ADC:

$$\begin{aligned} A_\gamma^{-1} : \sigma(\mathcal{H}_2) &\mapsto \sigma(\mathcal{H}_2) \\ A_\gamma \circ A_\gamma^{-1}(\rho) &= \rho \end{aligned} \quad (4.5.1)$$

The map A_γ^{-1} , while trace preserving and linear, is not expected to be completely positive, therefore it is not expected to be a quantum channel. In fact it could be cast in a "pseudo-Kraus" representation:

$$A_\gamma^{-1}(\theta) = \tilde{K}_0(\gamma)\theta\tilde{K}_0(\gamma)^\dagger - \tilde{K}_1(\gamma)\theta\tilde{K}_1(\gamma)^\dagger$$

$$\tilde{K}_0(\gamma) = \begin{pmatrix} 1 & 0 \\ 0 & \frac{1}{\sqrt{1-\gamma}} \end{pmatrix} \quad \tilde{K}_1(\gamma) = \begin{pmatrix} 0 & \sqrt{\frac{\gamma}{1-\gamma}} \\ 0 & 0 \end{pmatrix} \quad (4.5.2)$$

Notice that, while $\tilde{K}_0(\gamma)$ and $\tilde{K}_1(\gamma)$ form a Kraus set, the " $-$ " sign in (4.5.2) implies that A_γ^{-1} is **not** necessarily a quantum channel, as it is not written in the Kraus representation. One could also write (4.5.2) in matrix form:

$$A_\gamma^{-1}(\theta) = \begin{pmatrix} \theta_{00} - \frac{\gamma}{1-\gamma}\theta_{11} & \frac{1}{\sqrt{1-\gamma}}\theta_{01} \\ \frac{1}{\sqrt{1-\gamma}}\theta_{01}^* & \frac{1}{(1-\gamma)}\theta_{11} \end{pmatrix}. \quad (4.5.3)$$

The map defined in (4.5.1) is also the left-inverse of A_γ , which can be verified by direct computation.

4.5.b Inverse of single decay MAD channels

Consider the MAD channel $\Phi_{\Xi_k^{(n)}}$, where $\Xi_k^{(n)}$ is defined in (4.2.14) and the associated transition probability can be found in (4.2.17). This is a single decay MAD channel, which acts on the subspace spanned by $|k\rangle$ and $|n\rangle$ as an ADC with the same transition probability. Therefore, one might be tempted to define the right-inverse of $\Phi_{\Xi_k^{(n)}}$ as the embedding onto \mathcal{H}_d of the right-inverse of an ADC. Let $\Phi_{\Xi_k^{(n)}}^{(-1)}$ be the right inverse of $\Phi_{\Xi_k^{(n)}}$. Given $\theta \in \sigma(\text{span}\{|k\rangle, |n\rangle\})$, $\Phi_{\Xi_k^{(n)}}^{(-1)}$ needs to satisfy:

$$\Phi_{\Xi_k^{(n)}}^{(-1)}(\theta) = \begin{pmatrix} \theta_{nn} - \gamma_{kn}^{(n+1)} \left(1 - \gamma_{kn}^{(n+1)}\right)^{-1} \theta_{kk} & \left(1 - \gamma_{kn}^{(n+1)}\right)^{-1/2} \theta_{nk} \\ \left(1 - \gamma_{kn}^{(n+1)}\right)^{-1/2} \theta_{nk}^* & \left(1 - \gamma_{kn}^{(n+1)}\right)^{-1} \theta_{kk} \end{pmatrix} \quad (4.5.4)$$

This can be achieved by defining $\Phi_{\Xi_k^{(n)}}^{(-1)}$ as:

$$\Phi_{\Xi_k^{(n)}}^{(-1)}(\rho) \equiv \tilde{K}_0\left(\Xi_k^{(n)}\right)\rho\tilde{K}_0\left(\Xi_k^{(n)}\right)^\dagger - \tilde{K}_1\left(\Xi_k^{(n)}\right)\rho\tilde{K}_1\left(\Xi_k^{(n)}\right)^\dagger$$

$$\tilde{K}_0\left(\Xi_k^{(n)}\right) \equiv \mathbb{1}_d - \left(1 - \left(1 - \gamma_{kn}^{(n+1)}\right)^{-1/2}\right) |k\rangle\langle k| \quad (4.5.5)$$

$$\tilde{K}_1\left(\Xi_k^{(n)}\right) \equiv \left(\frac{\gamma_{kn}^{(n+1)}}{1 - \gamma_{kn}^{(n+1)}}\right)^{1/2} |n\rangle\langle k|$$

where $\rho \in \sigma(\mathcal{H}_d)$. It is possible to show, by direct computation, that:

$$\Phi_{\Xi_k^{(n)}} \circ \Phi_{\Xi_k^{(n)}}^{(-1)}(\rho) = \rho \quad \forall \rho \in \sigma(\mathcal{H}_d) \quad (4.5.6)$$

Notice that the map $\Phi_{\Xi_k^{(n)}}^{(-1)}$ is linear and trace preserving and is also the left-inverse of $\Phi_{\Xi_k^{(n)}}$.

4.5.c Inverse map as composition of inverse maps of single decays

The importance of (4.5.6) lies in the fact that it can be composed to generate the right-inverse map of a general MAD channel. In fact, recall that in (4.2.21) it was shown that a MAD channel can be seen as a composition of single decay MAD channels. Then, by "inverting" those single decay transitions one by one, the resulting channel must be the identity channel. This line of reasoning results in the definition:

$$\Phi_{\Gamma}^{-1} = \bigodot_{k=1}^{d-1} \left(\Phi_{\Xi_k^{(k-1)}}^{(-1)} \circ \dots \circ \Phi_{\Xi_k^{(0)}}^{(-1)} \right) \quad (4.5.7)$$

which, by construction, satisfies:

$$\Phi_{\Gamma} \circ \Phi_{\Gamma}^{-1} = \mathbb{1}. \quad (4.5.8)$$

Since each of the $\Phi_{\Xi_k^{(n)}}^{(-1)}$'s is linear and trace preserving, and a generic composition of such maps holds those same properties, then Φ_{Γ}^{-1} must also be linear and trace preserving.

Notice, finally, that Φ_{Γ}^{-1} , again, by construction, is also the left inverse of Φ_{Γ} .

4.6 Degradability of MAD channels

As seen in 2.11, the quantum capacity of degradable channels is computable and corresponds to the maximum (calculated over all possible inputs) of the coherent information 2.11.2. Therefore, it is important to find a complete characterization of the degradability conditions of MAD channels.

Since MAD channels admit a right inverse map, starting from (2.6.1), verifying the degradability of Φ_{Γ} reduces to verifying that Λ_{Γ} in:

$$\Lambda_{\Gamma} \equiv \tilde{\Phi}_{\Gamma} \circ \Phi_{\Gamma}^{(-1)} \quad (4.6.1)$$

is LCPT. The map Λ_{Γ} is linear and trace preserving by construction, while the complete positiveness is not guaranteed; using Choi's theorem (2.4.17), it is possible to verify whether or not Λ_{Γ} satisfies that last property:

$$\Lambda_{\Gamma} \text{ is completely positive} \Leftrightarrow C_{\Lambda_{\Gamma}} \text{ is positive} \quad (4.6.2)$$

where C_{Λ_Γ} is the Choi matrix of the map Λ_Γ , defined as in (2.4.12)¹.

$$\begin{aligned} C_{\Lambda_\Gamma} &\equiv \frac{1}{d} [\Lambda_\Gamma \otimes \mathbb{1}] (|\Omega\rangle\langle\Omega|) \\ |\Omega\rangle\langle\Omega| &\text{ maximally entangled state.} \end{aligned} \quad (4.6.3)$$

The computation of the eigenvalues of C_{Λ_Γ} for a generic 4-dimensional Γ turns out to be relatively demanding for a desktop computer (see 7.2 for details on computation methods), but all degradability regions can be found regardless by means of lateral thinking that simplifies the computation tasks, as demonstrated in 4.7.

4.7 Degradability of 4-dimensional MAD channels

In 3.3.i, the degradability regions for 3-dimensional MAD channels were determined by ensuring that the rank of the output state of the channel was not smaller than the rank of the state of the environment (3.3.44), which implied reducing the number of decays in the channel, and checking *a posteriori* whether or not less strict degradability conditions could be found by introducing additional decays to those regions. In the case of 4-dimensional MAD channels $\Phi_\Gamma^{(4)}$, (3.3.44) becomes:

$$\text{rank}(\Phi_\Gamma^{(4)}(\rho)) \geq \text{rank}(\tilde{\Phi}_\Gamma^{(4)}(\rho)) \quad \forall \rho, \quad (4.7.1)$$

which implies that the number of decays in the channel needs to not exceed 3. The number of 4-dimensional MAD channels with 3 decays can be found by counting the number of ways one can select 3 different couples (j, i) , where $j < i$; the number of such couples is $\binom{4}{2} = 6$, which means that the number of 3-decay configurations is $\binom{6}{3} = 20$; many of these configurations are unitarily equivalent.

In what follows, 4-dimensional MAD channels connected by unitary transformations of the form (4.3.2) will be grouped in *classes*; one can identify 9 classes of 3-decay channels, 4 classes of 2-decay channels and 1 class of single decay channels.

4.7.a Class 1A

This class identifies the single decay MAD channels; as seen in Section 4.3, if the transition probability is fixed, all possible configurations of single decay MAD's are unitarily equivalent. Let $\Phi_{\Gamma_{1A}(\gamma)}$ be a generic channel belonging to Class 1A, with transition probability γ . This class presents the PCDS structure (3.2.1), where $\Phi_{BB} =$

¹the factor $1/d$ is kept here for the sake of consistency, but is obviously not necessary when checking for positiveness

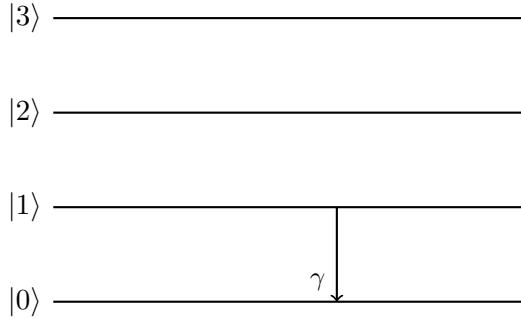


Figure 4.7.1: Class 1A consists of an ADC embedded into the higher dimensional \mathcal{H}_4

$\mathbb{1}_{\sigma(\mathcal{H}_2)}$ and $\Phi_{AA} = ADC_\gamma$ an amplitude damping channel; therefore, following (3.2.2), the degradability conditions of $\Phi_{\Gamma_{1A}(\gamma)}$ are the same as those given in (3.1.15):

$$\Phi_{\Gamma_{1A}(\gamma)} \text{ degradable} \Leftrightarrow \gamma \leq \frac{1}{2}. \quad (4.7.2)$$

The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{1A}(\gamma)}$ is never antidegradable.

4.7.b Class 2A

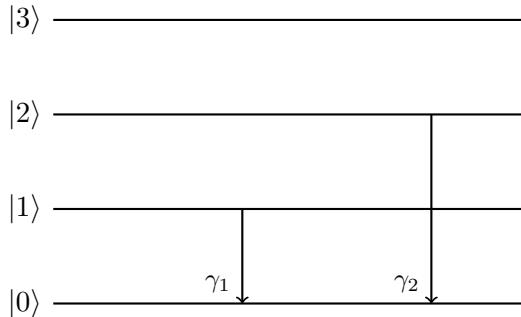


Figure 4.7.2: An example of channel belonging in Class 2A is provided by the embedding of $\Phi_{\Gamma(\gamma_1, \gamma_2, 0)}^{(3)}$ in \mathcal{H}_4 .

Channels belonging to this class are unitarily equivalent to the MAD channel identified by the transition matrix $\Gamma = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| - \gamma_1 |1\rangle\langle 1| + \gamma_2 |2\rangle\langle 0| - \gamma_2 |2\rangle\langle 2|$, depicted in Figure 4.7.2. Let a generic channel in this class be called $\Phi_{\Gamma_{2A}(\gamma_1, \gamma_2)}$; this channel presents the PCDS structure (3.2.1), where $\Phi_{BB} = \mathbb{1}_{\sigma(\mathcal{H}_1)}$ and $\Phi_{AA} = \Phi_{\Gamma(\gamma_1, \gamma_2, 0)}^{(3)}$, where $\Phi_{\Gamma(\gamma_1, \gamma_2, 0)}^{(3)}$ is the 3-dimensional MAD channel defined in (3.3.28). Following (3.2.2), the degradability conditions of $\Phi_{\Gamma_{2A}(\gamma_1, \gamma_2)}$ are the same as those given in (3.3.55):

$$\Phi_{\Gamma_{2A}(\gamma_1, \gamma_2)} \text{ degradable} \Leftrightarrow \gamma_1 \leq \frac{1}{2} \wedge \gamma_2 \leq \frac{1}{2} \quad (4.7.3)$$

The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{2A}(\gamma_1, \gamma_2)}$ is never antidegradable.

4.7.c Class 2B

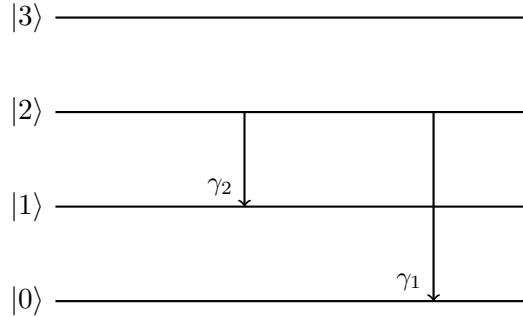


Figure 4.7.3: An example of channel belonging in **Class 2B** is provided by the embedding of $\Phi_{\Gamma(0,\gamma_1,\gamma_2)}^{(3)}$ in \mathcal{H}_4 .

Channels belonging to this class are unitarily equivalent to the MAD channel identified by the transition matrix $\Gamma = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |2\rangle\langle 0| + \gamma_2 |2\rangle\langle 1| - (\gamma_1 + \gamma_2) |2\rangle\langle 2|$, depicted in Figure 4.7.3. Let a generic channel in this class be called $\Phi_{\Gamma_{2B}(\gamma_1,\gamma_2)}$; this channel presents the PCDS structure (3.2.1), where $\Phi_{BB} = \mathbb{1}_{\sigma(\mathcal{H}_1)}$ and $\Phi_{AA} = \Phi_{\Gamma(0,\gamma_1,\gamma_2)}^{(3)}$. Following (3.2.2), the degradability conditions of $\Phi_{\Gamma_{2B}(\gamma_1,\gamma_2)}$ are the same as those given in (3.3.48):

$$\Phi_{\Gamma_{2B}(\gamma_1,\gamma_2)} \text{ degradable} \Leftrightarrow \gamma_1 + \gamma_2 \leq \frac{1}{2} \quad (4.7.4)$$

The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{2B}(\gamma_1,\gamma_2)}$ is never antidegradable.

4.7.d Class 2C

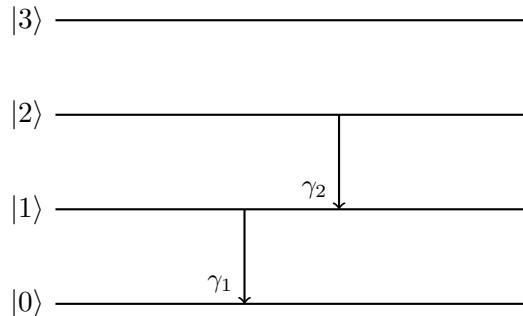


Figure 4.7.4: An example of channel belonging in **Class 2C** is provided by the embedding of $\Phi_{\Gamma(\gamma_1,0,\gamma_2)}^{(3)}$ in \mathcal{H}_4 .

Channels belonging to this class are unitarily equivalent to the MAD channel identified by the transition matrix $\Gamma = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| - \gamma_1 |1\rangle\langle 1| + \gamma_2 |2\rangle\langle 1| - \gamma_2 |2\rangle\langle 2|$, depicted in Figure 4.7.4. Let a generic channel in this class be called $\Phi_{\Gamma_{2C}(\gamma_1,\gamma_2)}$; this

channel presents the PCDS structure (3.2.1), where $\Phi_{BB} = \mathbb{1}_{\sigma(\mathcal{H}_1)}$ and $\Phi_{AA} = \Phi_{\Gamma(\gamma_1, 0, \gamma_2)}^{(3)}$. Following (3.2.2) and Subsection 3.3.i, it can be inferred that $\Phi_{\Gamma_{2C}(\gamma_1, \gamma_2)}$ is never degradable. The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{2C}(\gamma_1, \gamma_2)}$ is never antidegradable.

4.7.e Class 2D

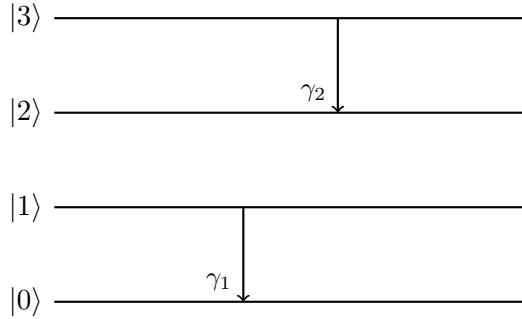


Figure 4.7.5: Class 2D is comprised of PCDS channels whose subchannels are ADC's

Channels belonging to this class are unitarily equivalent to the MAD channel identified by the transition matrix $\Gamma = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| - \gamma_1 |1\rangle\langle 1| + \gamma_2 |3\rangle\langle 2| - \gamma_2 |3\rangle\langle 3|$, depicted in Figure 4.7.4. Let a generic channel in this class be called $\Phi_{\Gamma_{2D}(\gamma_1, \gamma_2)}$; this channel presents the PCDS structure (3.2.1), where $\Phi_{AA} = ADC_{\gamma_1}$ and $\Phi_{BB} = ADC_{\gamma_2}$. Following (3.2.2) and (3.1.15), the degradability conditions of $\Phi_{\Gamma_{2D}(\gamma_1, \gamma_2)}$ are:

$$\Phi_{\Gamma_{2D}(\gamma_1, \gamma_2)} \text{ degradable} \Leftrightarrow \gamma_1 \leq \frac{1}{2} \wedge \gamma_2 \leq \frac{1}{2} \quad (4.7.5)$$

The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{2D}(\gamma_1, \gamma_2)}$ is never antidegradable.

4.7.f Class 3A

This class of channels consists PCDS channels whose subchannels are the identity channel $\Phi_{BB} = \mathbb{1}_{\sigma(\mathcal{H}_1)}$ and a 3-dimensional MAD channel $\Phi_{AA} = \Phi_{\Gamma(\gamma_1, \gamma_2, \gamma_3)}^{(3)}$. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow$ Class 3A \sim Class 2B
- $\gamma_2 = 0 \Rightarrow$ Class 3A \sim Class 2C
- $\gamma_3 = 0 \Rightarrow$ Class 3A \sim Class 2A

Let a generic channel belonging to this class be called $\Phi_{\Gamma_{3A}(\gamma_1, \gamma_2, \gamma_3)}$; following (3.2.2) and (3.3.57), $\Phi_{\Gamma(\gamma_1, \gamma_2, \gamma_3)}^{(3)}$ is never degradable if $\gamma_1, \gamma_2, \gamma_3 \neq 0$. This also implies, as a consequence of (4.4.2), that it is not possible to build a 4-decay degradable MAD channel starting from a MAD channel belonging to Class 3A. The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{3A}(\gamma_1, \gamma_2, \gamma_3)}$ is never antidegradable.

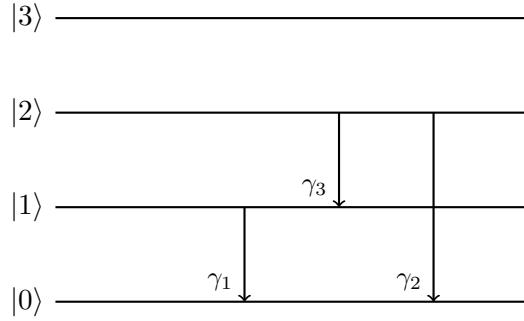


Figure 4.7.6: Class 3A consists of 3-dimensional MAD channels embedded into the higher dimensional \mathcal{H}_4

4.7.g Class 3B



Figure 4.7.7: Class 3B

Consider the transition matrix:

$$\Gamma_{3B}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |2\rangle\langle 0| + \gamma_2 |3\rangle\langle 0| + \gamma_3 |3\rangle\langle 1| - \gamma_1 |2\rangle\langle 2| - (\gamma_2 + \gamma_3) |3\rangle\langle 3|; \quad (4.7.6)$$

Class 3B is comprised of all the channels unitarily equivalent to $\Phi_{\Gamma_{3B}(\gamma_1, \gamma_2, \gamma_3)}$, depicted in Figure 4.7.7. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow$ Class 3A \sim Class 2B
- $\gamma_2 = 0 \Rightarrow$ Class 3A \sim Class 2D
- $\gamma_3 = 0 \Rightarrow$ Class 3A \sim Class 2A

The degradability conditions for this channel are:

$$\Phi_{\Gamma_{3B}(\gamma_1, \gamma_2, \gamma_3)} \text{ degradable} \Leftrightarrow \gamma_1 \leq \frac{1}{2} \wedge \gamma_2 + \gamma_3 \leq \frac{1}{2} \quad (4.7.7)$$

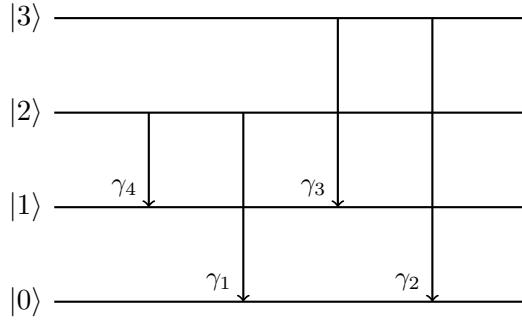


Figure 4.7.8: Class 4A

Starting from a channel in Class 3B, it is possible to build a degradable 4-decay MAD channel: consider the transition matrix:

$$\Gamma_{4A}(\gamma_1, \gamma_2, \gamma_3, \gamma_4) \equiv \Gamma_{3B}(\gamma_1, \gamma_2, \gamma_3) + \gamma_4 |2\rangle\langle 1| - \gamma_4 |2\rangle\langle 2| \quad (4.7.8)$$

The degradability conditions for $\Phi_{\Gamma_{4A}(\gamma_1, \gamma_2, \gamma_3, \gamma_4)}$ are:

$$\Phi_{\Gamma_{4A}(\gamma_1, \gamma_2, \gamma_3, \gamma_4)} \text{ degradable} \Leftrightarrow \gamma_1 + \gamma_4 \leq \frac{1}{2} \wedge \gamma_2 + \gamma_3 \leq \frac{1}{2} \quad (4.7.9)$$

This channel provides a counterexample to (4.7.1), proving that the relation is not a necessary condition for the degradability of a channel. The existence of a noiseless subspace for $\Phi_{\Gamma_{3B}(\gamma_1, \gamma_2, \gamma_3)}$ guarantees that the channel is never antidegradable.

4.7.h Class 3C

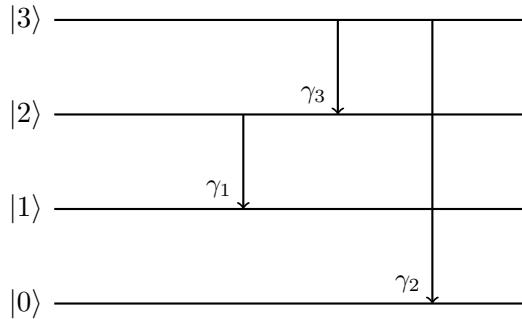


Figure 4.7.9: Class 3C

Consider the transition matrix:

$$\Gamma_{3C}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |2\rangle\langle 1| + \gamma_2 |3\rangle\langle 0| + \gamma_3 |3\rangle\langle 2| - \gamma_1 |2\rangle\langle 2| - (\gamma_2 + \gamma_3) |3\rangle\langle 3|; \quad (4.7.10)$$

Class 3C is comprised of all the channels unitarily equivalent to $\Phi_{\Gamma_{3C}(\gamma_1, \gamma_2, \gamma_3)}$, depicted in Figure 4.7.9. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow \text{Class 3C} \sim \text{Class 2B}$
- $\gamma_2 = 0 \Rightarrow \text{Class 3C} \sim \text{Class 2C}$
- $\gamma_3 = 0 \Rightarrow \text{Class 3C} \sim \text{Class 2D}$

$\Phi_{\Gamma_{3C}(\gamma_1, \gamma_2, \gamma_3)}$ is never degradable when $\gamma_1, \gamma_2, \gamma_3 \neq 0$. The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{3C}(\gamma_1, \gamma_2, \gamma_3)}$ is never antidegradable.

4.7.i Class 3D

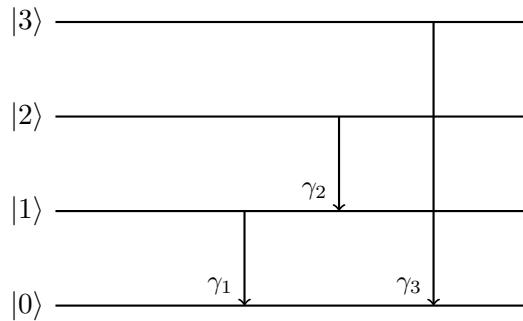


Figure 4.7.10: Class 3D

Consider the transition matrix:

$$\Gamma_{3D}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| + \gamma_2 |2\rangle\langle 1| + \gamma_3 |3\rangle\langle 0| - \gamma_1 |1\rangle\langle 1| - \gamma_2 |2\rangle\langle 2| - \gamma_3 |3\rangle\langle 3|; \quad (4.7.11)$$

Class 3D is comprised of all the channels unitarily equivalent to $\Phi_{\Gamma_{3D}(\gamma_1, \gamma_2, \gamma_3)}$, depicted in Figure 4.7.10. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow \text{Class 3D} \sim \text{Class 2D}$
- $\gamma_2 = 0 \Rightarrow \text{Class 3D} \sim \text{Class 2A}$
- $\gamma_3 = 0 \Rightarrow \text{Class 3D} \sim \text{Class 2C}$

$\Phi_{\Gamma_{3D}(\gamma_1, \gamma_2, \gamma_3)}$ is never degradable when $\gamma_1, \gamma_2, \gamma_3 \neq 0$.

4.7.j Class 3E

Consider the transition matrix:

$$\Gamma_{3E}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| + \gamma_2 |2\rangle\langle 0| + \gamma_3 |3\rangle\langle 0| - \gamma_1 |1\rangle\langle 1| - \gamma_2 |2\rangle\langle 2| - \gamma_3 |3\rangle\langle 3|; \quad (4.7.12)$$

$\Phi_{\Gamma_{3E}(\gamma_1, \gamma_2, \gamma_3)}$ is the only channel in Class 3E, depicted in Figure 4.7.11. Turning off one of the decays reduces this class to one of the 2-decay classes:

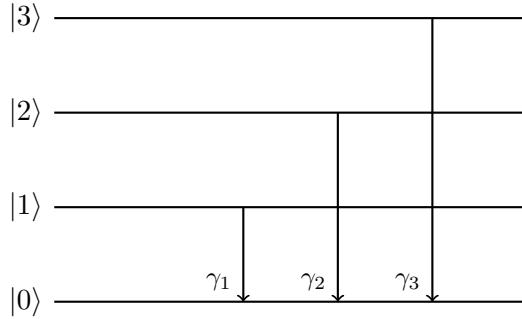


Figure 4.7.11: Class 3E

- $\gamma_1 = 0 \Rightarrow \text{Class 3E} \sim \text{Class 2A}$
- $\gamma_2 = 0 \Rightarrow \text{Class 3E} \sim \text{Class 2A}$
- $\gamma_3 = 0 \Rightarrow \text{Class 3E} \sim \text{Class 2A}$

The degradability conditions for $\Phi_{\Gamma_{3E}(\gamma_1, \gamma_2, \gamma_3)}$ are:

$$\Phi_{\Gamma_{3E}(\gamma_1, \gamma_2, \gamma_3)} \text{ degradable} \Leftrightarrow \gamma_1 \leq \frac{1}{2} \wedge \gamma_2 \leq \frac{1}{2} \wedge \gamma_3 \leq \frac{1}{2} \quad (4.7.13)$$

4.7.k Class 3F

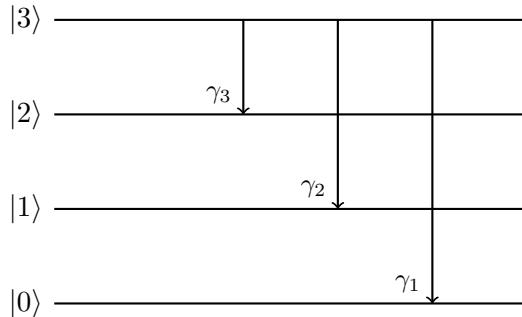


Figure 4.7.12: Class 3F

Consider the transition matrix:

$$\Gamma_{3E}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |3\rangle\langle 0| + \gamma_2 |3\rangle\langle 1| + \gamma_3 |3\rangle\langle 2| - (\gamma_1 + \gamma_2 + \gamma_3) |3\rangle\langle 3|; \quad (4.7.14)$$

$\Phi_{\Gamma_{3F}(\gamma_1, \gamma_2, \gamma_3)}$ is the only channel in Class 3F, depicted in Figure 4.7.12. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow \text{Class 3F} \sim \text{Class 2B}$
- $\gamma_2 = 0 \Rightarrow \text{Class 3F} \sim \text{Class 2B}$

- $\gamma_3 = 0 \Rightarrow \text{Class 3F} \sim \text{Class 2B}$

The degradability conditions for $\Phi_{\Gamma_{3F}(\gamma_1, \gamma_2, \gamma_3)}$ are:

$$\Phi_{\Gamma_{3F}(\gamma_1, \gamma_2, \gamma_3)} \text{ degradable} \Leftrightarrow \gamma_1 \leq \frac{1}{2} \wedge \gamma_2 \leq \frac{1}{2} \wedge \gamma_3 \leq \frac{1}{2} \quad (4.7.15)$$

The existence of a noiseless subspace guarantees that $\Phi_{\Gamma_{3F}(\gamma_1, \gamma_2, \gamma_3)}$ is never antidegradable.

4.7.1 Class 3G

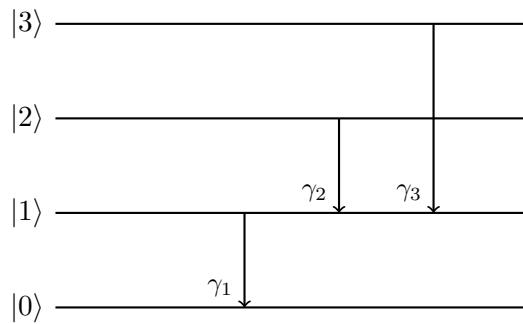


Figure 4.7.13: Class 3G

Consider the transition matrix:

$$\Gamma_{3G}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| + \gamma_2 |2\rangle\langle 1| + \gamma_3 |3\rangle\langle 1| - \gamma_1 |1\rangle\langle 1| - \gamma_2 |2\rangle\langle 2| - \gamma_3 |3\rangle\langle 3|; \quad (4.7.16)$$

$\Phi_{\Gamma_{3G}(\gamma_1, \gamma_2, \gamma_3)}$ is the only channel in Class 3G, depicted in Figure 4.7.13. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow \text{Class 3G} \sim \text{Class 2A}$
- $\gamma_2 = 0 \Rightarrow \text{Class 3G} \sim \text{Class 2C}$
- $\gamma_3 = 0 \Rightarrow \text{Class 3G} \sim \text{Class 2C}$

$\Phi_{\Gamma_{3G}(\gamma_1, \gamma_2, \gamma_3)}$ is never degradable when $\gamma_1, \gamma_2, \gamma_3 \neq 0$.

4.7.m Class 3H

Consider the transition matrix:

$$\Gamma_{3H}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |2\rangle\langle 0| + \gamma_2 |2\rangle\langle 1| + \gamma_3 |3\rangle\langle 2| - (\gamma_1 + \gamma_2) |2\rangle\langle 2| - \gamma_3 |3\rangle\langle 3|; \quad (4.7.17)$$

$\Phi_{\Gamma_{3H}(\gamma_1, \gamma_2, \gamma_3)}$ is the only channel in Class 3H, depicted in Figure 4.7.14. Turning off one of the decays reduces this class to one of the 2-decay classes:

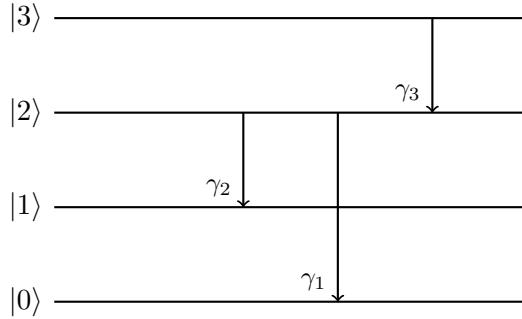


Figure 4.7.14: Class 3H

- $\gamma_1 = 0 \Rightarrow$ Class 3H \sim Class 2C
- $\gamma_2 = 0 \Rightarrow$ Class 3H \sim Class 2C
- $\gamma_3 = 0 \Rightarrow$ Class 3H \sim Class 2B

$\Phi_{\Gamma_{3H}(\gamma_1, \gamma_2, \gamma_3)}$ is never degradable when $\gamma_1, \gamma_2, \gamma_3 \neq 0$.

4.7.n Class 3I

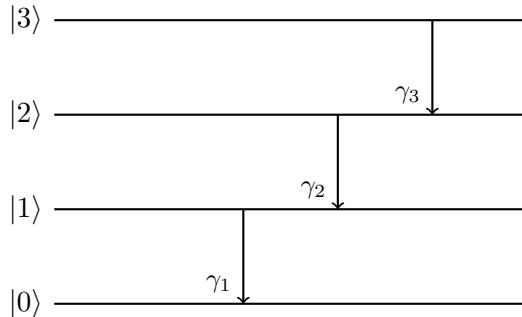


Figure 4.7.15: Class 3I

Consider the transition matrix:

$$\Gamma_{3I}(\gamma_1, \gamma_2, \gamma_3) = \mathbb{1}_{\mathcal{H}_4} + \gamma_1 |1\rangle\langle 0| + \gamma_2 |2\rangle\langle 1| + \gamma_3 |3\rangle\langle 2| - \gamma_1 |1\rangle\langle 1| - \gamma_2 |2\rangle\langle 2| - \gamma_3 |3\rangle\langle 3| ; \quad (4.7.18)$$

$\Phi_{\Gamma_{3I}(\gamma_1, \gamma_2, \gamma_3)}$ is the only channel in Class 3I, depicted in Figure 4.7.15. Turning off one of the decays reduces this class to one of the 2-decay classes:

- $\gamma_1 = 0 \Rightarrow$ Class 3I \sim Class 2C
- $\gamma_2 = 0 \Rightarrow$ Class 3I \sim Class 2D
- $\gamma_3 = 0 \Rightarrow$ Class 3I \sim Class 2C

$\Phi_{\Gamma_{3I}(\gamma_1, \gamma_2, \gamma_3)}$ is never degradable when $\gamma_1, \gamma_2, \gamma_3 \neq 0$.

4.8 Degradability of a MAD channel embedded in a higher dimensional system

The structure defined in (3.2.1) lends itself very well to the case of MAD channels; in fact, given a d -dimensional MAD channel $\Phi_\Gamma : \sigma(\mathcal{H}_d) \mapsto \sigma(\mathcal{H}_d)$, one may consider another MAD channel $\Phi_{\bar{\Gamma}} : \sigma(\mathcal{H}_{d+1}) \mapsto \sigma(\mathcal{H}_{d+1})$ whose transition probabilities are:

$$\bar{\gamma}_{ji} = \begin{cases} \gamma_{ji} & \text{if } 0 \leq i < j \leq d-1 \\ 0 & \text{if } 0 \leq i < j = d \end{cases} \quad (4.8.1)$$

In the context of (3.2.1), the Φ_{AA} and Φ_{BB} channels of $\Phi_{\bar{\Gamma}}$ are Φ_Γ and $\mathbb{1}_{\mathcal{H}_1}$. Then, (3.2.2) implies that $\Phi_{\bar{\Gamma}}$ is degradable if and only if Φ_Γ is degradable.

4.9 Monotonicity properties

In [CG21a], using the pipeline inequalities (2.13.1) and the monotonicity of the 3-dimensional MAD channels in their parameters, proven in 3.3.g, the authors were able to compute some capacity functionals for certain configurations of the channels that were not degradable nor anti-degradable. In order to generalize the property of monotonicity of the capacity functionals to the d -dimensional case, one needs to consider two transition matrices, Γ, Γ' , whose elements $\gamma_{ji}, \gamma'_{ji}$ differ only for a single pair of indices (j_0, i_0) , so that:

$$\gamma'_{j_0, i_0} \geq \gamma_{j_0, i_0}. \quad (4.9.1)$$

Then, on account of the pipeline inequalities (2.13.1), one could conclude that the capacities MAD channels are non increasing in the parameter $\gamma'_{j_0, i_0} \geq \gamma_{j_0, i_0}$ if one were able to find two LCPT maps Λ_L, Λ_R such that:

$$\Lambda_L \circ \Phi_\Gamma \circ \Lambda_R = \Phi_{\Gamma'} \quad (4.9.2)$$

Finding Λ_L, Λ_R is not an easy task and generally one has to resort to heuristic methods to simplify the problem. Recalling the property of closure under composition for MAD channels, one might assume that Λ_L, Λ_R are MAD's themselves; however, for $d > 3$, this line of reasoning only allows to derive monotonicity properties under the transition probabilities $\gamma_{d-1, i_0}, \forall 0 \leq i_0 < d-1$ and γ_{10} :

- **Monotonicity under γ_{d-1, i_0} :** Consider $\Lambda_R = \mathbb{1}_{\mathcal{H}_d}$ the identity channel and $\Lambda_L = \Phi_{\Gamma_{\lambda_{d-1, i_0}}}$ a single decay MAD channel from $|d-1\rangle$ to $|i_0\rangle$ with transition probability λ_{d-1, i_0} . From (4.2.7), the resulting channel in (4.9.2), $\Phi_{\Gamma'} = \Phi_{\Gamma_{\lambda_{d-1, i_0}}} \circ \Phi_\Gamma$ has the same transition probabilities γ'_{ji} of Φ_Γ aside from $\gamma'_{d-1, i_0} \geq \gamma_{d-1, i_0}$
- **Monotonicity under γ_{10} :** Consider $\Lambda_L = \mathbb{1}_{\mathcal{H}_d}$ the identity channel and $\Lambda_R = \Phi_{\Gamma_{\lambda_{10}}}$ a single decay MAD channel from $|1\rangle$ to $|0\rangle$ with transition probability λ_{10} . From (4.2.7), the resulting channel in (4.9.2), $\Phi_{\Gamma'} = \Phi_\Gamma \circ \Phi_{\Gamma_{\lambda_{10}}}$ has the same transition probabilities γ'_{ji} of Φ_Γ aside from $\gamma'_{10} \geq \gamma_{10}$

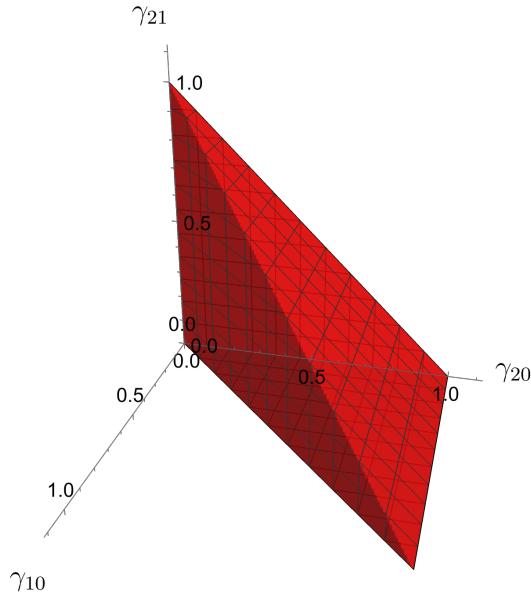


Figure 4.9.1: Monotonicity region under γ_{21} (4.9.7). Note that γ_{21}, γ_{20} are still bounded by (3.3.3), meaning that $0 \leq \gamma_{20} + \gamma_{21} \leq 1$.

Monotonicity properties in $d = 4$

It is unclear whether a 4-dimensional MAD channel presents monotonous capacities under the transition probabilities γ_{20}, γ_{21} . As stated above, resorting to (4.9.2) is only useful if additional assumptions are made. Assume that either one of $\Lambda_L, \Lambda_R = \mathbb{1}_{\mathcal{H}_d}$, then, utilizing the inverse map (4.5.7), from (4.9.2):

$$\Lambda_R = \mathbb{1}_{\mathcal{H}_d} \Rightarrow \Lambda_L = \Phi_{\Gamma'} \circ \Phi_{\Gamma}^{(-1)} \quad (4.9.3)$$

$$\Lambda_L = \mathbb{1}_{\mathcal{H}_d} \Rightarrow \Lambda_R = \Phi_{\Gamma}^{(-1)} \circ \Phi_{\Gamma'} \quad (4.9.4)$$

Note that, by construction, both of these maps are linear and trace preserving; verifying their complete positiveness would imply that they are quantum channels, which can be done by verifying the positive semi-definiteness of their Choi matrices, due to (2.4.17).

$$\gamma'_{21} \geq \gamma_{21}$$

Assume that Γ' and Γ only differ in the $|2\rangle\langle 1|$ element, so that:

$$\begin{aligned} \gamma'_{21} &\equiv \gamma_{21} + \varepsilon_{21} \geq \gamma_{21} \\ 0 &\leq \varepsilon_{21} \leq \gamma_{22} \end{aligned} \quad (4.9.5)$$

where the last condition is needed in order for $\gamma_{22}' = \gamma_{22} - \varepsilon_{21} \geq 0$ to be satisfied. Computing the eigenvalues of the Choi matrices of Λ_L, Λ_R in (4.9.3) and (4.9.4), one finds that the former is only positive semi-definite if:

$$\gamma_{32} = 0, \quad (4.9.6)$$

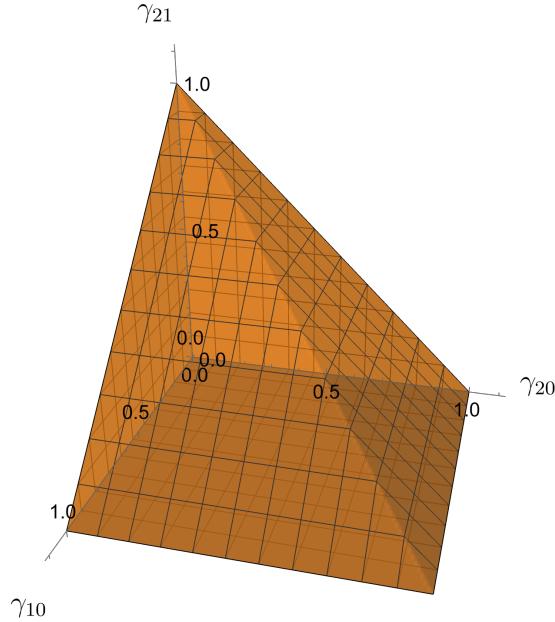


Figure 4.9.2: Monotonicity region under γ_{20} (4.9.10). Note that γ_{21}, γ_{20} are still bounded by (3.3.3), meaning that $0 \leq \gamma_{20} + \gamma_{21} \leq 1$.

while the latter is positive semi-definite under the following condition:

$$\gamma_{20} - \gamma_{10} \geq 0 \quad (4.9.7)$$

which means that the capacity functionals of a 4-dimensional MAD channel are monotonous under γ_{21} if $\gamma_{32} = 0$ or $\gamma_{20} \geq \gamma_{10}$. The region described by (4.9.7) is illustrated in Figure 4.9.1.

$$\gamma'_{20} \geq \gamma_{20}$$

Assume that Γ' and Γ only differ in the $|2\rangle\langle 0|$ element, so that:

$$\begin{aligned} \gamma'_{20} &\equiv \gamma_{20} + \varepsilon_{20} \geq \gamma_{20} \\ 0 &\leq \varepsilon_{20} \leq \gamma_{22} \end{aligned} \quad (4.9.8)$$

where the last condition is needed in order for $\gamma_{22}' = \gamma_{22} - \varepsilon_{20} \geq 0$ to be satisfied. Computing the eigenvalues of the Choi matrices of Λ_L, Λ_R in (4.9.3) and (4.9.4), one finds that the former is only positive semi-definite if:

$$\gamma_{32} = 0, \quad (4.9.9)$$

while the latter is positive semi-definite under the following condition:

$$1 - \gamma_{21} - \gamma_{10} \geq 0 \quad (4.9.10)$$

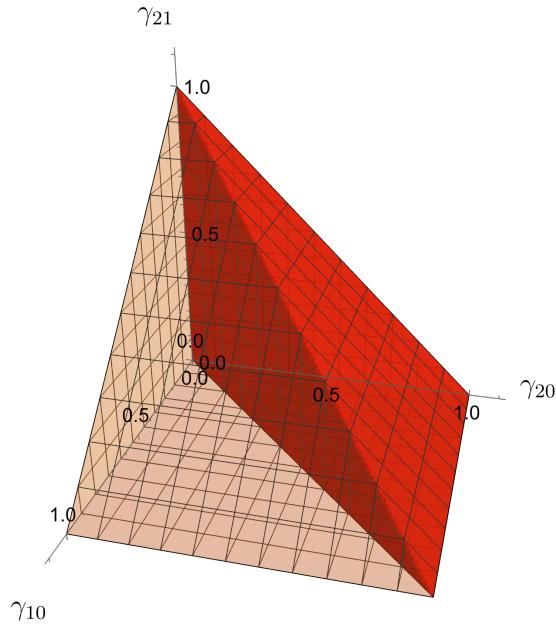


Figure 4.9.3: The monotonicity region under γ_{21} (4.9.7) is completely contained within the monotonicity region under γ_{20} (4.9.10).

which means that the capacity functionals of a 4-dimensional MAD channel are monotonous under γ_{20} if $\gamma_{32} = 0$ or $\gamma_{21} \leq 1 - \gamma_{10}$. The region described by (4.9.10) is illustrated in Figure 4.9.2. Note that the region described by (4.9.7) is completely contained within the region described by (4.9.10), as can be seen in Figure 4.9.3.

4.10 Complete damping of level $|3\rangle$ in 4-dimensional MAD's

Consider the 4-dimensional MAD channel $\Phi_{\Gamma_{CD3}}$ identified by the transition matrix:

$$\Gamma_{CD3} = \mathbb{1}_{\mathcal{H}_4} + \sum_{j=1}^3 \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle i| - \sum_{j=1}^3 \sum_{i=0}^{j-1} \gamma_{ji} |i\rangle\langle j| \quad (4.10.1)$$

$$\text{where } \gamma_{33} = 1 - \gamma_{30} - \gamma_{31} - \gamma_{32} = 0$$

The quantum capacity of this channel can be shown to be equivalent to the quantum capacity of a 3-dimensional MAD channel $\Phi_{\Gamma_{MAD3}}^{(3)}$, whose transition matrix is:

$$\Gamma_{MAD3} = \mathbb{1}_{\mathcal{H}_3} + \sum_{j=1}^2 \sum_{i=0}^{j-1} \gamma_{ji} |j\rangle\langle i| - \sum_{j=1}^2 \sum_{i=0}^{j-1} \gamma_{ji} |i\rangle\langle j| \quad (4.10.2)$$

In order to reach this result, one needs to find coinciding upper and lower bounds on $Q(\Phi_{\Gamma_{CD3}})$.

Lower bound

The lower bound on $Q(\Phi_{\Gamma_{CD3}})$ is trivial: if one were to encode information only on the subspace spanned by $\{|0\rangle, |1\rangle, |2\rangle\}$, the channel $\Phi_{\Gamma_{CD3}}$ would be equivalent to $\Phi_{\Gamma_{MAD3}}^{(3)}$; of course, this choice of encoding is not guaranteed to be optimal, therefore:

$$Q(\Phi_{\Gamma_{CD3}}) \geq Q\left(\Phi_{\Gamma_{MAD3}}^{(3)}\right) \quad (4.10.3)$$

Upper bound

The upper bound on $Q(\Phi_{\Gamma_{CD3}})$ can be obtained using the pipeline inequalities (2.13.1) and the Direct Sum channels' property (3.2.8); consider the DS channel:

$$\begin{aligned} \Phi_{\Gamma_{MAD3}}^{DS}(\theta) &\equiv \left[\begin{array}{c|c} \Phi_{\Gamma_{MAD3}}^{(3)}(p\theta^{(3)}) & 0 \\ \hline 0 & 1-p \end{array} \right] \\ \theta &\equiv \left[\begin{array}{c|c} p\theta^{(3)} & \theta_{j3} \\ \hline \theta_{j3}^* & 1-p \end{array} \right] \\ \theta &\in \sigma(\mathcal{H}_4), \theta^{(3)} \in \sigma(\mathcal{H}_3) \\ 0 &\leq p \leq 1 \end{aligned} \quad (4.10.4)$$

and the Level Erasure channel:

$$\begin{aligned} LE_{\Gamma_{CD3}}(\rho) &\equiv \begin{pmatrix} \rho_{00} + \gamma_{30}\rho_{33} & \rho_{01} & \rho_{02} & 0 \\ \rho_{01}^* & \rho_{11} + \gamma_{31}\rho_{33} & \rho_{12} & 0 \\ \rho_{02}^* & \rho_{12}^* & \rho_{22} + \gamma_{32}\rho_{33} & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix} \\ \rho &\equiv \begin{pmatrix} \rho_{00} & \rho_{01} & \rho_{02} & \rho_{03} \\ \rho_{01}^* & \rho_{11} & \rho_{12} & \rho_{13} \\ \rho_{02}^* & \rho_{12}^* & \rho_{22} & \rho_{23} \\ \rho_{03}^* & \rho_{13}^* & \rho_{23}^* & \rho_{33} \end{pmatrix}. \end{aligned} \quad (4.10.5)$$

The composition of these two channel is equal to $\Phi_{\Gamma_{CD3}}$:

$$\Phi_{\Gamma_{CD3}} = LE_{\Gamma_{CD3}} \circ \Phi_{\Gamma_{MAD3}}^{DS}. \quad (4.10.6)$$

The quantum capacity of $\Phi_{\Gamma_{MAD3}}^{DS}$ is given in (3.2.8), which in this case translates to:

$$Q(\Phi_{\Gamma_{MAD3}}^{DS}) = Q\left(\Phi_{\Gamma_{MAD3}}^{(3)}\right) \quad (4.10.7)$$

The pipeline inequalities (2.13.1) ensure that (4.10.7) is an upper bound for $Q(\Phi_{\Gamma_{CD3}})$:

$$Q(\Phi_{\Gamma_{CD3}}) \leq Q\left(\Phi_{\Gamma_{MAD3}}^{(3)}\right) \quad (4.10.8)$$

Combining (4.10.3) and (4.10.8) leads to the conclusion:

$$Q(\Phi_{\Gamma_{CD3}}) = Q\left(\Phi_{\Gamma_{MAD3}}^{(3)}\right) \quad (4.10.9)$$

Capacity computations for MAD channels in $d = 4$

Building upon the results in Chapters 2, 3 and 4, applying the techniques described in 7.2, an analysis of the quantum, private classical and two-way capacities 4-dimensional MAD channel becomes feasible. This analysis was performed and the consequent results are reported in this chapter.

5.1 Single decay 4-dimensional MAD channel

All single decay 4-dimensional MAD channel are unitarily equivalent, all belonging to Class 1A defined in 4.7.a.

5.1.a Class 1A

In order to compute the capacities of single decay 4-dimensional MAD channels, consider the sample channel $\Phi_{\Gamma_{1A}(\gamma_{10})}$ identified by the transition matrix:

$$\Gamma_{1A}(\gamma_{10}) \equiv \mathbf{1}_{\mathcal{H}_4} + \gamma_{10} |1\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1| \quad (5.1.1)$$

As stated in 4.7.a, $\Phi_{\Gamma_{1A}(\gamma_{10})}$ obeys the PCDS structure (3.2.1), with $\Phi_{BB} = \mathbf{1}_{\sigma(\mathcal{H}_2)}$ and $\Phi_{AA} = ADC_{\gamma_{10}}$, and is degradable if and only if $\gamma_{10} \leq 1/2$; in this region, the quantum capacity can be computed directly using (3.3.21):

$$Q(\Phi_{\Gamma_{1A}(\gamma_{10})}) = \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{1A}(\gamma_{10})}) \quad \forall \gamma_{10} \leq \frac{1}{2} \quad (5.1.2)$$

For $\gamma_{10} = 1/2$, $ADC_{\gamma_{10}}$ has null quantum capacity, therefore the capacity of $\Phi_{\Gamma_{1A}(\gamma_{10})}$ corresponds to its lower bound provided by (3.2.5), which in this case is $\log_2 3$. Using the monotonicity properties exposed in 4.9, it can be inferred that:

$$Q(\Phi_{\Gamma_{1A}(\gamma_{10})}) = C_P(\Phi_{\Gamma_{1A}(\gamma_{10})}) = \log_2 3 \quad \forall \frac{1}{2} \leq \gamma_{10} \leq 1 \quad (5.1.3)$$

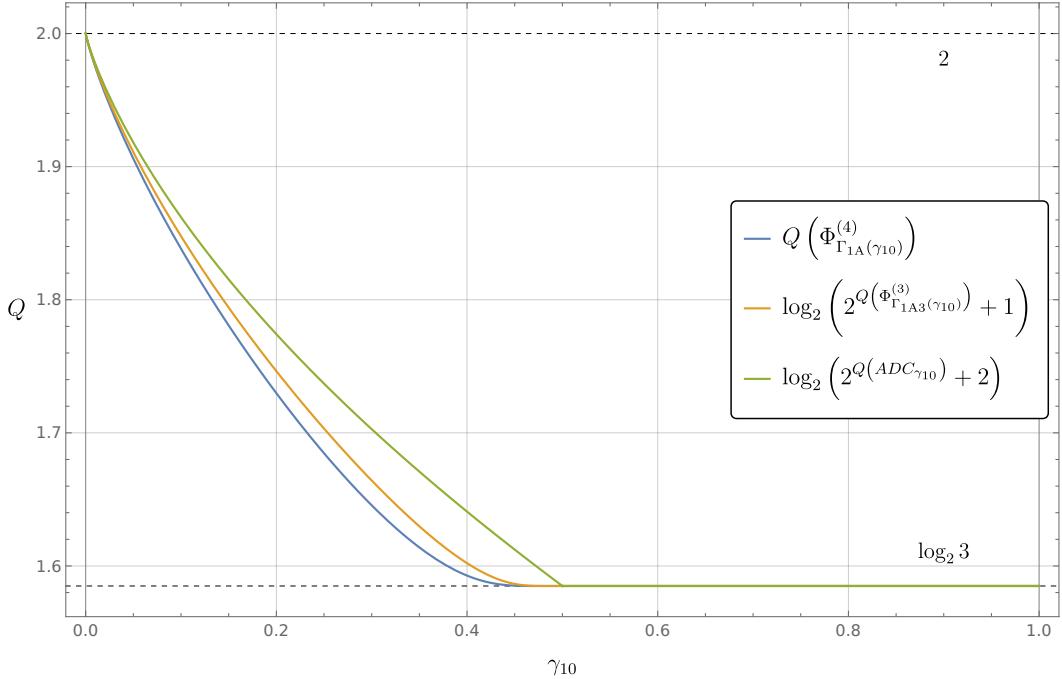


Figure 5.1.1: Quantum capacity of $\Phi_{\Gamma_{1A}(\gamma_{10})}$, compared with its upper bounds (3.2.4) given by the PCDS structure of the channel.

These results are plotted in Figure 5.1.1, alongside the upper bounds on the quantum capacity of PCDS channels (3.2.4), which are computed starting from the quantum capacity of an ADC and the quantum capacity of a 3-dimensional MAD channel $\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}$, identified by the transition matrix:

$$\Gamma_{1A3}(\gamma_{10}) \equiv \mathbb{1}_{\mathcal{H}_3} + \gamma_{10} |1\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1|. \quad (5.1.4)$$

5.2 2-decay 4-dimensional MAD channels

There are 4 classes of 2-decay 4-dimensional MAD channels, whose capacities are reported in this section.

5.2.a Class 2A

Consider the sample channel $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ belonging in Class 2A, identified by the transition matrix:

$$\Gamma_{2A}(\gamma_{10}, \gamma_{20}) = \mathbb{1}_{\mathcal{H}_4} + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |2\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1| - \gamma_{20} |2\rangle\langle 2| \quad (5.2.1)$$

As stated in 4.7.a, $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ obeys the PCDS structure (3.2.1), with $\Phi_{BB} = \mathbb{1}_{\sigma(\mathcal{H}_1)}$ and $\Phi_{AA} = \Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}$ is a 3-dimensional MAD channel identified by the transition

matrix:

$$\Gamma_{2A3}(\gamma_{10}, \gamma_{20}) = \mathbb{1}_{\mathcal{H}_3} + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |2\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1| - \gamma_{20} |2\rangle\langle 2| \quad (5.2.2)$$

The degradability conditions for $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ are given in (4.7.3), which in this case translate to:

$$\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})} \text{ degradable} \Leftrightarrow \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{20} \leq \frac{1}{2} \quad (5.2.3)$$

In this region, the quantum and classical private capacities can be computed using (3.3.21):

$$\begin{aligned} Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}) &= \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}) \\ \forall \gamma_{10}, \gamma_{20} : \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{20} \leq \frac{1}{2} \end{aligned} \quad (5.2.4)$$

$Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})})$ presents a symmetry under swap of $(\gamma_{10}, \gamma_{20})$, which is a consequence of the relation:

$$\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})} = \mathcal{U}_{21} \circ \Phi_{\Gamma_{2A}(\gamma_{20}, \gamma_{10})} \circ \mathcal{U}_{21}, \quad (5.2.5)$$

where \mathcal{U}_{21} is defined in (4.3.2).

It is possible to extend the computation of the quantum capacity of $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ to the whole region $0 \leq \gamma_{10} \leq 1 \wedge 0 \leq \gamma_{20} \leq 1$.

Extension to $\frac{1}{2} \leq \gamma_{20} \leq 1$

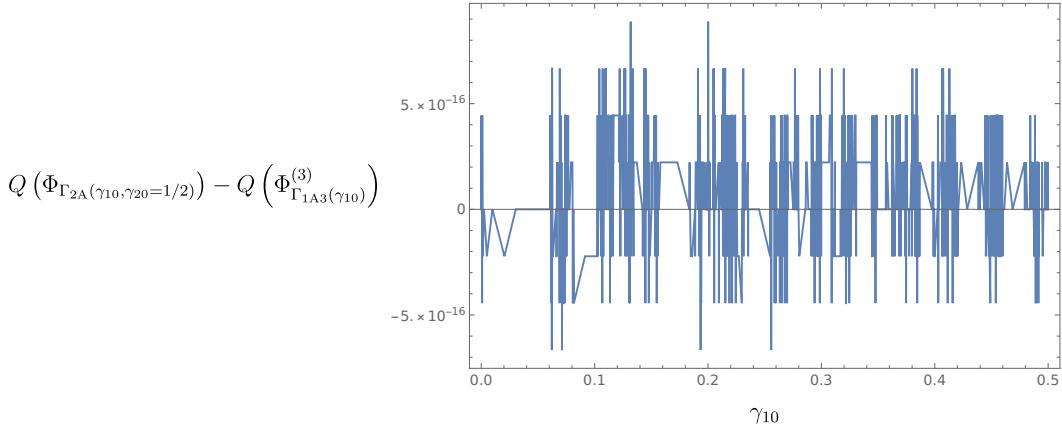


Figure 5.2.1: Difference between $Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20}=1/2)})$ and $Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20}=1/2)})$ for $0 \leq \gamma_{10} \leq 1/2$; it is reasonable to assume, due to their very low values, that the discrepancies from 0 in the plot are due to the computational resolution of the machine used for the evaluation.

The channel $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ is covariant to $\Phi_{\Gamma'_{2A}(\gamma_{10}, \gamma_{20})}$:

$$\begin{aligned} \Gamma'_{2A}(\gamma_{10}, \gamma_{20}) &\equiv \mathbb{1}_{\mathcal{H}_4} + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |3\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1| - \gamma_{20} |3\rangle\langle 3| \\ \Phi_{\Gamma'_{2A}(\gamma_{10}, \gamma_{20})} &= \mathcal{U}_{23} \circ \Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})} \circ \mathcal{U}_{23} \end{aligned} \quad (5.2.6)$$

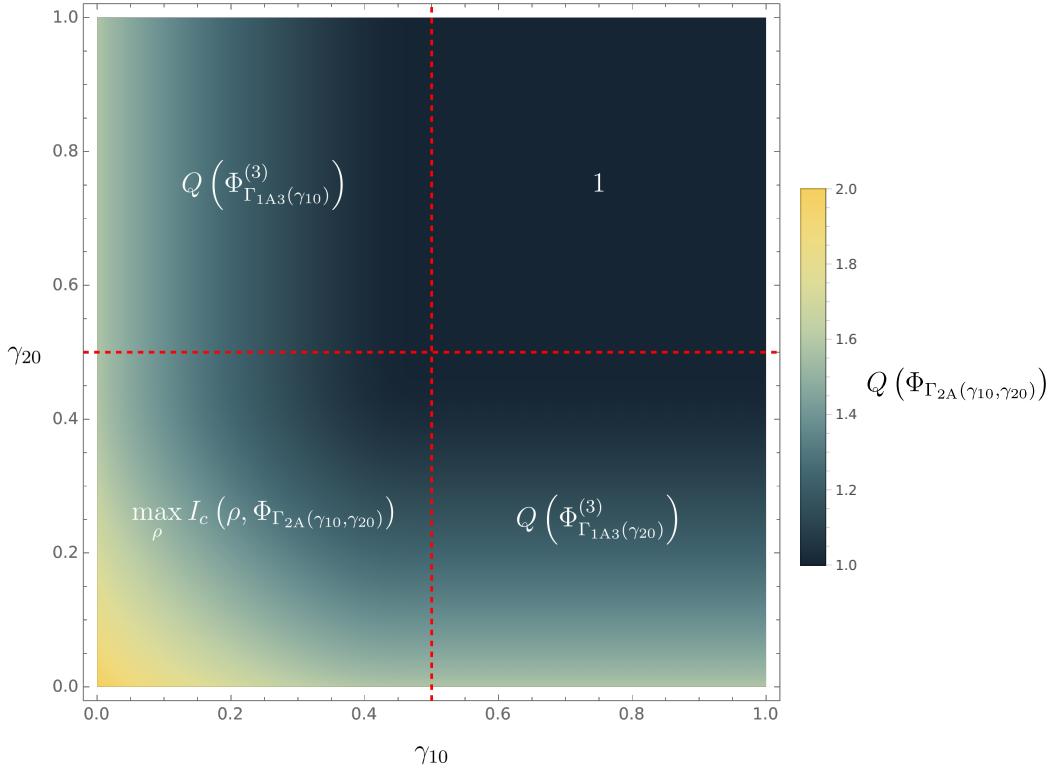


Figure 5.2.2: Density plot representing the quantum capacity of $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$.

When $\gamma_{20} = 1$, applying (5.2.6) and (4.10.9), one finds that:

$$\begin{aligned} Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20}=1)}) &= Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \\ \Gamma_{1A3}(\gamma_{10}) &\equiv \mathbb{1}_{\mathcal{H}_3} + \gamma_{10}|1\rangle\langle 0| - \gamma_{10}|1\rangle\langle 1| \end{aligned} \quad (5.2.7)$$

Where $\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}$ is a 3-dimensional single decay MAD channel. Furthermore, computing the difference between $Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20}=1)})$ and $Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20}=1/2)})$ yields the plot reported in Figure 5.2.1. The discrepancies from 0 in the plot are likely due to the computational resolution of the machine used for the calculation. This result, coupled with the monotonicity of the capacities in γ_{10} for MAD channels, as seen in 4.9, implies that:

$$Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}) = Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \quad \forall 0 \leq \gamma_{10} \leq \frac{1}{2}, \frac{1}{2} \leq \gamma_{20} \leq 1 \quad (5.2.8)$$

$Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)})$ itself is computable in $0 \leq \gamma_{10} \leq 1$ following the results of [CG21a].

Quantum capacity in $\frac{1}{2} \leq \gamma_{20}, \gamma_{10} \leq 1$

When $\gamma_{20}, \gamma_{10} = 1/2$ the computation of the quantum capacity of $\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}$ is equal to the \log_2 of the dimension of the noiseless subspace of the channel:

$$\Phi_{\Gamma_{2A}(\gamma_{10}=1/2, \gamma_{20}=1/2)} = 1 \quad (5.2.9)$$

This value is a lower bound on the capacity of the channel; hence, given the monotonicity of the channel w.r.t. γ_{10}, γ_{20} implies that:

$$Q(\Phi_{\Gamma_{2A}(\gamma_{10}, \gamma_{20})}) = 1 \quad \forall \frac{1}{2} \leq \gamma_{10}, \gamma_{20} \leq 1 \quad (5.2.10)$$

Quantum capacity plot

The results in (5.2.10), (5.2.8), (5.2.5) and (5.2.4) combine to shape the plot of the quantum capacity for channels belonging in Class 2A reported in Figure 5.2.2.

5.2.b Class 2B

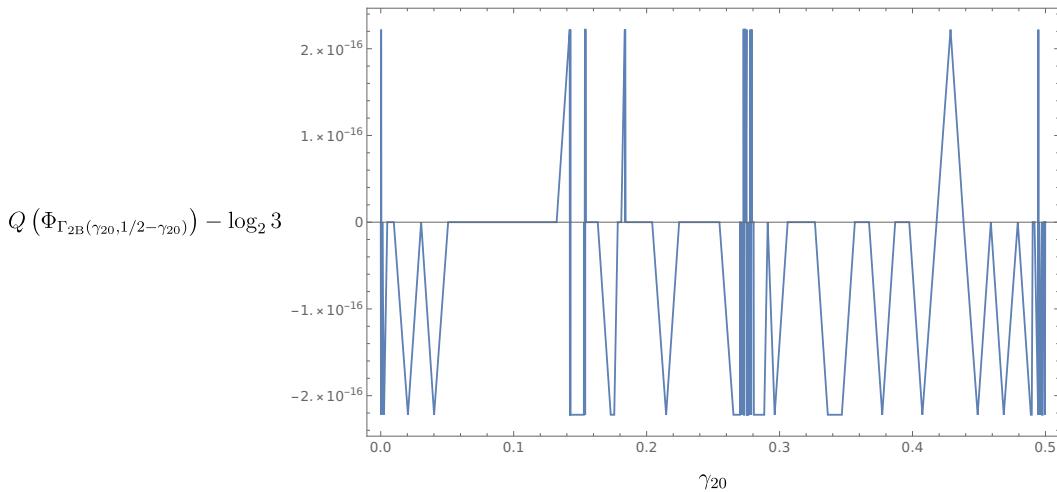


Figure 5.2.3: Difference between $Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21}=1-\gamma_{20})}) = \log_2 3$ and $Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21}=1/2-\gamma_{20})})$ for $0 \leq \gamma_{20} \leq 1/2$; it is reasonable to assume, due to their very low values, that the discrepancies from 0 in the plot are due to the computational resolution of the machine used for the evaluation.

Consider the sample channel $\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}$ belonging in Class 2B, identified by the transition matrix:

$$\Gamma_{2A}(\gamma_{20}, \gamma_{21}) = \mathbb{1}_{\mathcal{H}_4} + \gamma_{20} |2\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| - (\gamma_{20} + \gamma_{21}) |2\rangle\langle 2| \quad (5.2.11)$$

The channel $\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}$ obeys the PCDS structure (3.2.1); its degradability conditions are reported in (4.7.4), which in this case translate to:

$$\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})} \text{ degradable} \Leftrightarrow \gamma_{20} + \gamma_{21} \leq \frac{1}{2}. \quad (5.2.12)$$

In this region, the quantum capacity of $\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}$ can be computed using (3.3.21):

$$\begin{aligned} Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}) &= \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}) \\ \forall \gamma_{20}, \gamma_{21} : \gamma_{20} + \gamma_{21} &\leq \frac{1}{2}. \end{aligned} \quad (5.2.13)$$

$Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})})$ is symmetric in the pair $(\gamma_{20}, \gamma_{21})$ due to the covariance of the channel under the unitary transformation \mathcal{U}_{10} . It is possible to extend the computation of the quantum capacity performed in $0 \leq \gamma_{20} + \gamma_{21} \leq \frac{1}{2}$ to the region $\frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1$.

Extension to $\frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1$

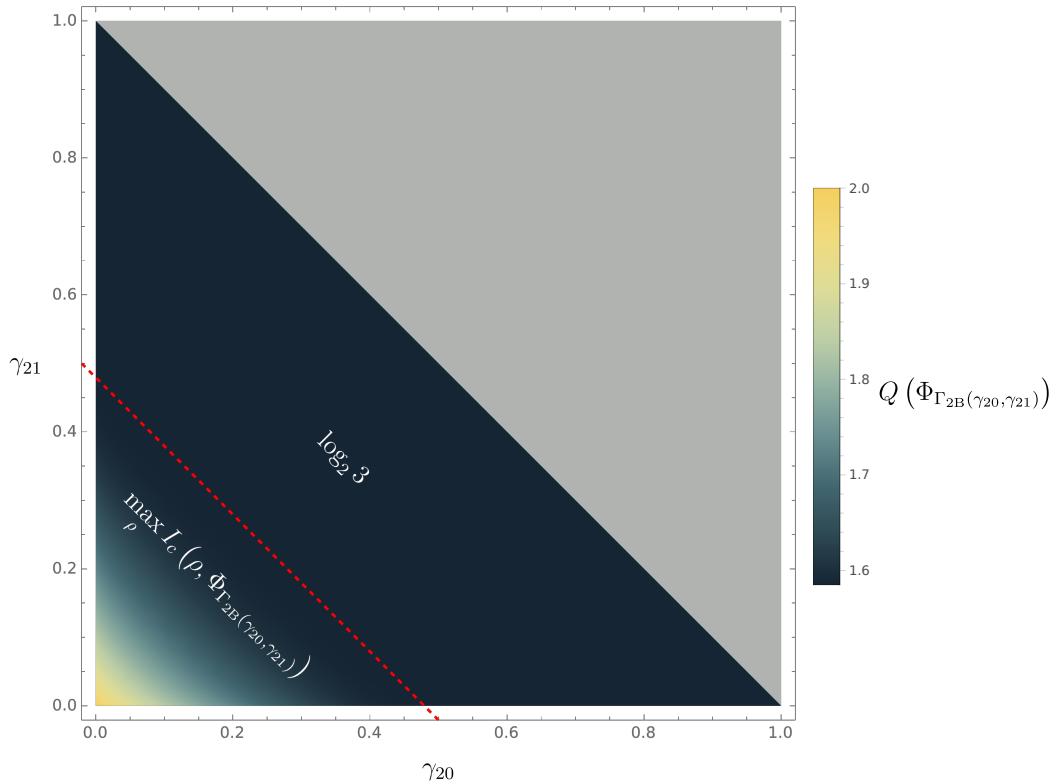


Figure 5.2.4: Density plot representing the quantum capacity of $\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}$.

The channel $\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}$ is covariant to $\Phi_{\Gamma'_{2B}(\gamma_{20}, \gamma_{21})}$:

$$\begin{aligned} \Gamma'_{2B}(\gamma_{20}, \gamma_{21}) &\equiv \mathbb{1}_{\mathcal{H}_4} + \gamma_{20} |3\rangle\langle 0| + \gamma_{21} |3\rangle\langle 1| - (\gamma_{20} + \gamma_{21}) |3\rangle\langle 3| \\ \Phi_{\Gamma'_{2B}(\gamma_{20}, \gamma_{21})} &= \mathcal{U}_{23} \circ \Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})} \circ \mathcal{U}_{23}. \end{aligned} \quad (5.2.14)$$

When $\gamma_{20} + \gamma_{21} = 1$ applying (5.2.14) and (4.10.9), one finds that:

$$Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21}=1-\gamma_{20})}) = Q(\mathbb{1}_{\sigma(\mathcal{H}_3)}) = \log_2 3 \quad (5.2.15)$$

Furthermore, computing the difference between $Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21}=1-\gamma_{20})})$ and $Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{20}=1/2-\gamma_{20})})$ yields the plot reported in Figure 5.2.3, from which one could infer that:

$$Q(\Phi_{\Gamma_{2B}(\gamma_{20}, \gamma_{21})}) = \log_2 3 \quad \forall \frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1 \quad (5.2.16)$$

Quantum capacity plot

The results in (5.2.16) and (5.2.13) combine to shape the plot of the quantum capacity for channels belonging in Class 2B, reported in Figure 5.2.4.

5.2.c Class 2C

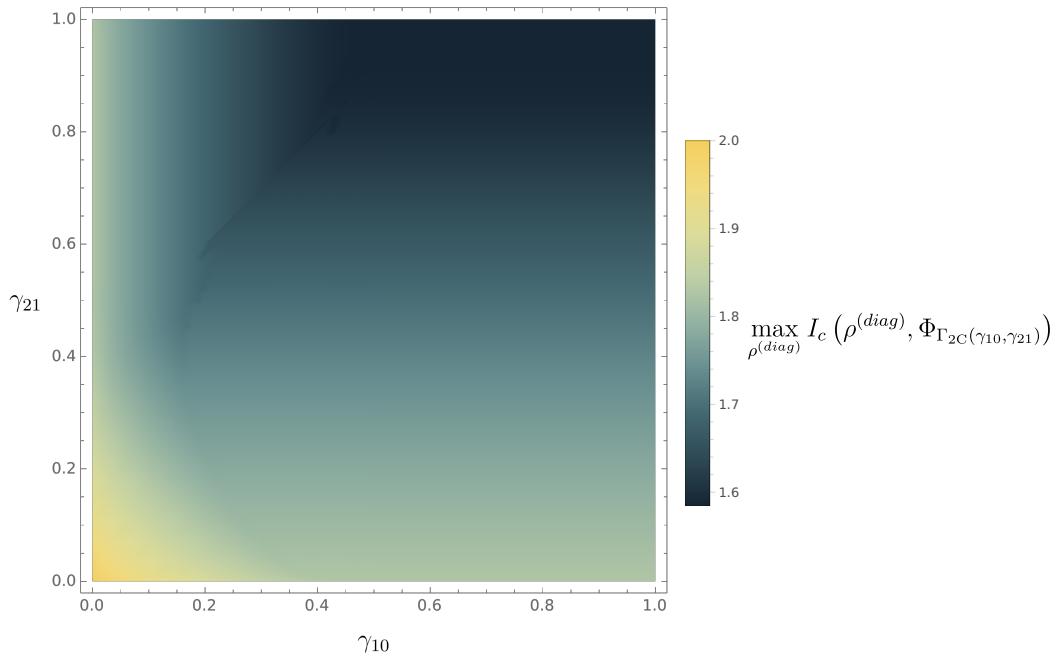


Figure 5.2.5: Density plot of the lower bound on the quantum capacity of $\Phi_{\Gamma_{2C}(\gamma_{10}, \gamma_{21})}$, obtained from the maximization of the coherent information for a single use on the channel over all possible diagonal input matrices.

Consider the sample channel $\Phi_{\Gamma_{2C}(\gamma_{10}, \gamma_{21})}$ belonging in Class 2C, identified by the transition matrix:

$$\Gamma_{2C}(\gamma_{10}, \gamma_{21}) = \mathbb{1}_{\mathcal{H}_4} + \gamma_{10} |1\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| - \gamma_{10} |1\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2| \quad (5.2.17)$$

This channel is never degradable. One needs to settle for computing a lower bound on the quantum capacity, provided by the maximization over all possible diagonal input matrices of the coherent information:

$$\max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{2C}(\gamma_{10}, \gamma_{21})}) \leq Q^{(1)}(\Phi_{\Gamma_{2C}(\gamma_{10}, \gamma_{21})}) \leq Q(\Phi_{\Gamma_{2C}(\gamma_{10}, \gamma_{21})}). \quad (5.2.18)$$

The lower bound in (5.2.18) is plotted in Figure 5.2.5.

5.2.d Class 2D

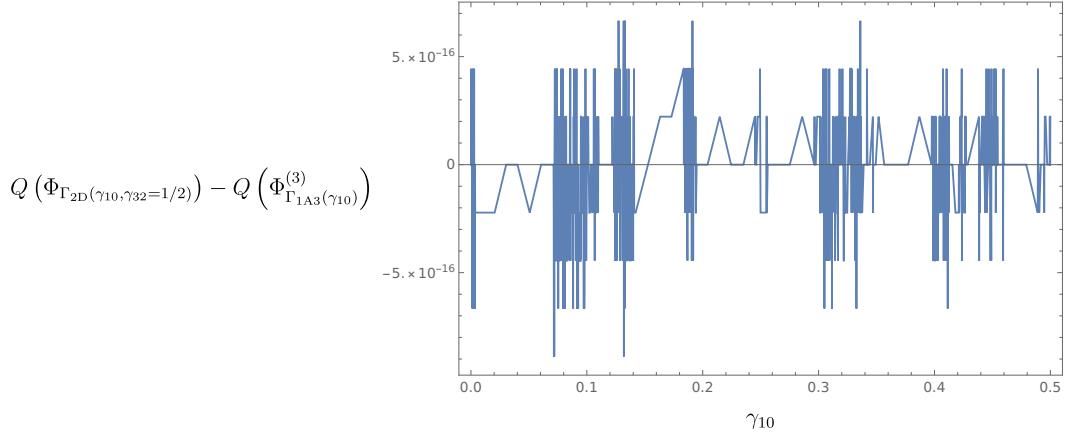


Figure 5.2.6: Difference between $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32}=1/2)})$ and $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32}=1)})$ for $0 \leq \gamma_{10} \leq 1/2$; it is reasonable to assume, due to their very low values, that the discrepancies from 0 in the plot are due to the computational resolution of the machine used for the evaluation.

Consider the sample channel $\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}$ belonging in Class 2D, identified by the transition matrix:

$$\Gamma_{2D}(\gamma_{10}, \gamma_{32}) = \mathbb{1}_{\mathcal{H}_4} + \gamma_{10}|1\rangle\langle 0| + \gamma_{32}|3\rangle\langle 2| - \gamma_{10}|1\rangle\langle 1| - \gamma_{32}|3\rangle\langle 3| \quad (5.2.19)$$

Following (4.7.5), the channel $\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}$ has degradability conditions:

$$\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})} \text{ degradable} \Leftrightarrow \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{32} \leq \frac{1}{2} \quad (5.2.20)$$

In this region, the quantum capacity can be computed using (3.3.21):

$$\begin{aligned} Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}) &= \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}) \\ &\forall \gamma_{10}, \gamma_{32} : \gamma_{10} \leq \frac{1}{2} \wedge \gamma_{32} \leq \frac{1}{2}. \end{aligned} \quad (5.2.21)$$

The quantum capacity $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$ is invariant under swap of $(\gamma_{10}, \gamma_{32})$ as a consequence of the covariance of the channel under the unitary transformation $\mathcal{U}_{21} \circ \mathcal{U}_{10} \circ \mathcal{U}_{32} \circ \mathcal{U}_{21}$. It is possible to extend the computation of the quantum capacity performed in the degradability region to rest of the surface $0 \leq \gamma_{10} \leq 1 \wedge 0 \leq \gamma_{32} \leq 1$.

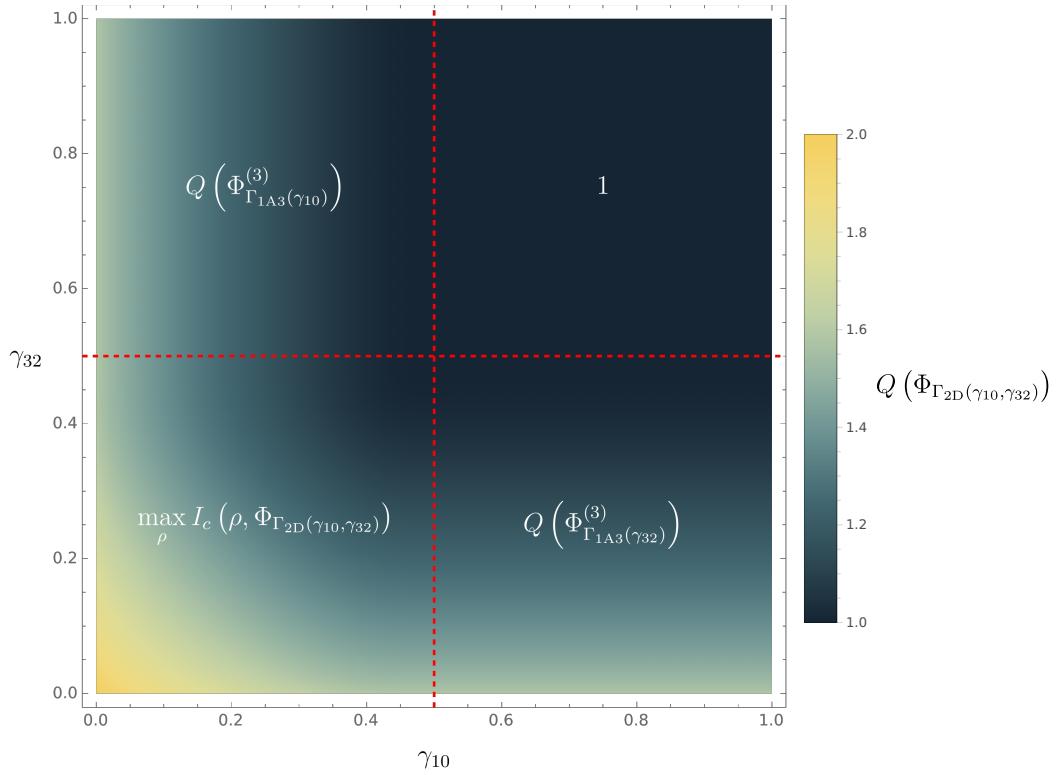


Figure 5.2.7: Density plot representing the quantum capacity of $\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}$.

Extension to $\frac{1}{2} \leq \gamma_{32} \leq 1$

When $\gamma_{32} = 1$, applying (4.10.9), one finds that:

$$Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32}=1)}) = Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}), \quad (5.2.22)$$

where $\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}$ is the 3-dimensional MAD channel identified by the transition matrix defined in (5.2.7). One can verify that $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32}=1/2)}) = Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32}=1)})$ by computing their difference; the results are reported in Figure 5.2.6. By employing the monotonicity of $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$ in its parameters, it is possible to infer that:

$$Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}) = Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \quad \forall 0 \leq \gamma_{10} \leq \frac{1}{2}, \frac{1}{2} \leq \gamma_{32} \leq 1 \quad (5.2.23)$$

Quantum capacity in $\frac{1}{2} \leq \gamma_{10}, \gamma_{32} \leq 1$

From the fact that $Q(\Phi_{\Gamma_{2D}(\gamma_{10}=1/2, \gamma_{32}=1/2)}) = 1$, 1 being a natural lower bound of $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$, as it is the \log_2 of the noiseless subspace of the channel, and the non-increasing behavior of $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$ in the transition probabilities, one can infer that:

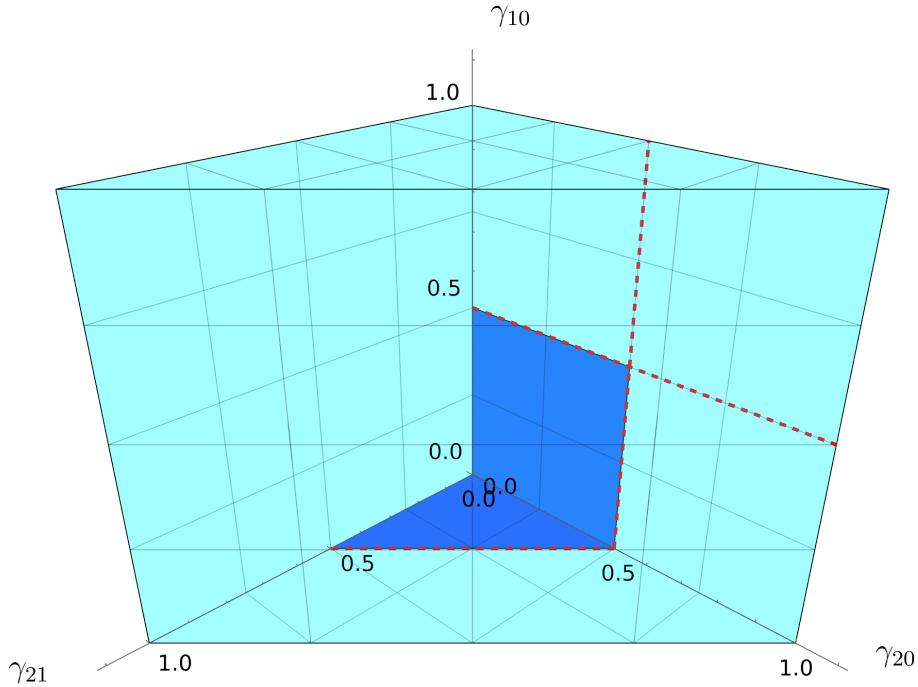


Figure 5.3.1: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3A}}(\gamma_{10}, \gamma_{20}, \gamma_{21})$ and its degradability regions (blue).

$$Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}) = 1 \quad \forall \frac{1}{2} \leq \gamma_{10}, \gamma_{32} \leq 1. \quad (5.2.24)$$

Quantum capacity plot

Combining (5.2.24), (5.2.23) and (5.2.21) and the symmetry of $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$ in the transition probabilities, the quantum capacity of $\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})}$ can be found on the whole surface $0 \leq \gamma_{10}, \gamma_{32} \leq 1$. The plot of $Q(\Phi_{\Gamma_{2D}(\gamma_{10}, \gamma_{32})})$ is reported in Figure 5.2.7.

5.3 3-decay 4-dimensional MAD channels

There are 9 classes of 3-decay 4-dimensional MAD channels, whose capacities are reported in this section.

5.3.a Class 3A

MAD channels belonging to Class 3A present the PCDS structure (3.2.1), with one of the channels forming the structure being a 3-dimensional MAD channel. Consider the sample channel $\Phi_{\Gamma_{3A}}(\gamma_{10}, \gamma_{20}, \gamma_{21})$ in class 3A, where the identifying transition matrix is:

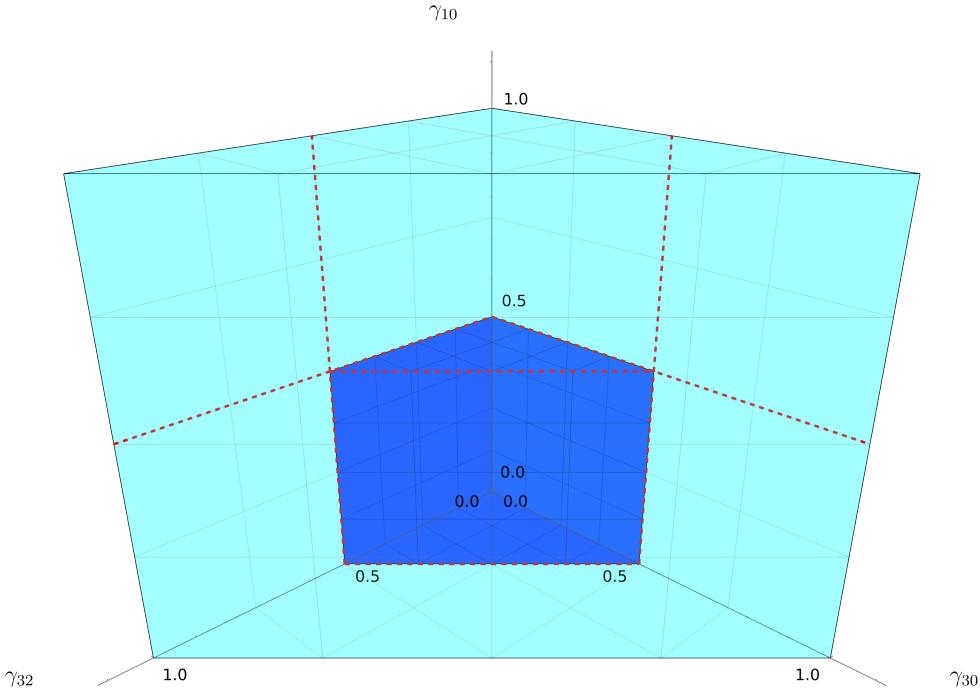


Figure 5.3.2: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3B}(\gamma_{20}, \gamma_{30}, \gamma_{31})}$ and its degradability regions (blue).

$$\begin{aligned} \Gamma_{3A}(\gamma_{10}, \gamma_{20}, \gamma_{21}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |2\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| \\ & - \gamma_{10} |1\rangle\langle 1| - (\gamma_{20} + \gamma_{21}) |2\rangle\langle 2| \end{aligned} \quad (5.3.1)$$

The results in 4.7.f indicate that a necessary condition for the degradability of $\Phi_{\Gamma_{3A}(\gamma_{10}, \gamma_{20}, \gamma_{21})}$ is $\gamma_{10} = 0 \vee \gamma_{21} = 0$, at which point the channel reduces to Class 2B and Class 2A respectively. A graph representing the parameter space of $\Phi_{\Gamma_{3A}(\gamma_{10}, \gamma_{20}, \gamma_{21})}$, together with its degradability regions, is reported in Figure 5.3.1.

5.3.b Class 3B

Consider the sample channel $\Phi_{\Gamma_{3B}(\gamma_{20}, \gamma_{30}, \gamma_{31})}$ belonging in class 3B; this channel is identified by the transition matrix:

$$\begin{aligned} \Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{30} |3\rangle\langle 0| + \gamma_{32} |3\rangle\langle 2| \\ & - \gamma_{10} |1\rangle\langle 1| - (\gamma_{30} + \gamma_{32}) |3\rangle\langle 3| \end{aligned} \quad (5.3.2)$$

The degradability conditions for class 3B, reported in (4.7.7), translate to :

$$\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})} \text{ degradable} \Leftrightarrow 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{32} + \gamma_{30} \leq \frac{1}{2}; \quad (5.3.3)$$

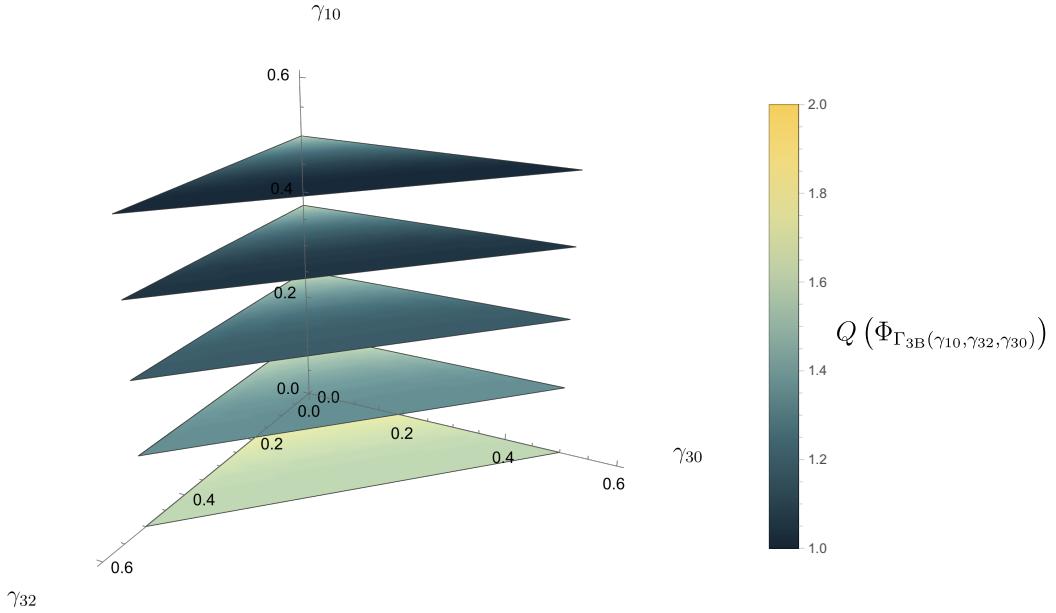


Figure 5.3.3: Slices of the quantum capacity in the 3-dimensional degradable region of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$.

this conditions are rendered in Figure 5.3.2. It is possible to extend the computation performed in the degradability zones to the remainder of the parameter region.

Quantum capacity in degradability region

One can directly compute the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ in the blue region rendered in Figure 5.3.2 using the result (3.3.21).

$$\begin{aligned} Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}) &= \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}) \\ &\forall \gamma_{10}, \gamma_{30}, \gamma_{32} : 0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2} \wedge 0 \leq \gamma_{10} \leq \frac{1}{2}. \end{aligned} \quad (5.3.4)$$

This computation was performed for 5 choices of γ_{10} in that region and plotted in Figure 5.3.3.

Quantum capacity on borders of the parameter region

Consider the 2-dimensional surfaces of Figure 5.3.2; it is possible to compute the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ on any of these surfaces.

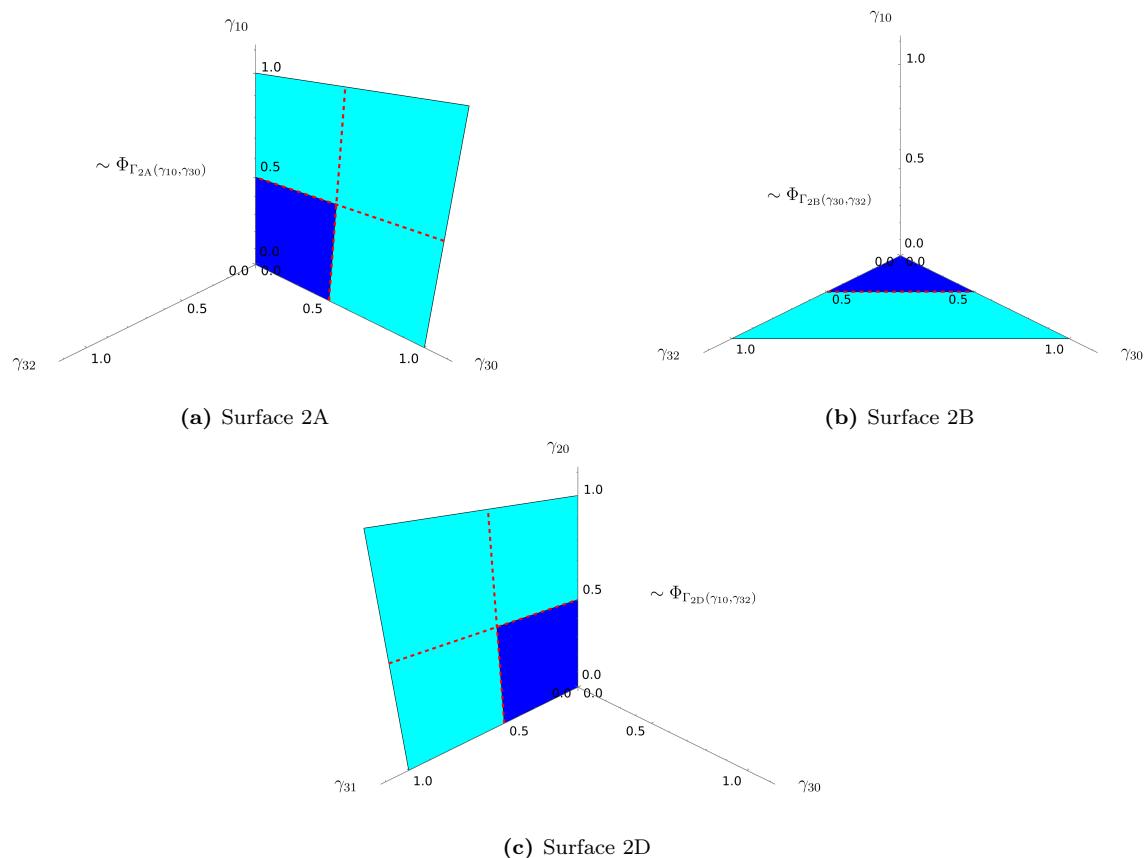


Figure 5.3.4: Surfaces in the parameter space of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ corresponding to setting one of the transition probabilities $\gamma_{10}, \gamma_{30}, \gamma_{32}$ to 0.

When either $\gamma_{10}, \gamma_{30}, \gamma_{32}$ is 0 (these surfaces are represented in Figure 5.3.4) the channel reduces to channels belonging to, respectively, Class 2B (Figure 5.3.4b), Class 2D (Figure 5.3.4c) and Class 2A (Figure 5.3.4a). The results in 5.2.a, 5.2.b, 5.2.d allow for the computation of the quantum capacity on those surfaces.

When $\gamma_{30} + \gamma_{32} = 1$, the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ is computable thanks to the result in (4.10.9), which implies that:

$$Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1 - \gamma_{30})}) = Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \quad (5.3.5)$$

where $\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}$ is defined in (5.2.7).

When $\gamma_{10} = 1$, it can be useful to work with another channel belonging in Class 3B, identified by the transition matrix:

$$\begin{aligned} \Gamma'_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32}) = & \mathbb{1}_4 + \gamma_{10}|3\rangle\langle 0| + \gamma_{30}|2\rangle\langle 0| + \gamma_{32}|2\rangle\langle 1| \\ & - \gamma_{10}|3\rangle\langle 3| - (\gamma_{30} + \gamma_{32})|2\rangle\langle 2| \end{aligned} \quad (5.3.6)$$

For a fixed choice of the transition probabilities, $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ and $\Phi_{\Gamma'_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ are unitarily covariant. The quantum capacity of $\Phi_{\Gamma'_{3B}(\gamma_{10}=1, \gamma_{30}, \gamma_{32})}$ can be found using (4.10.9):

$$Q(\Phi_{\Gamma_{3B}(\gamma_{10}=1, \gamma_{30}, \gamma_{32})}) = Q(\Phi_{\Gamma'_{3B}(\gamma_{10}=1, \gamma_{30}, \gamma_{32})}) = Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)}), \quad (5.3.7)$$

where $\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)}$ is a 3-dimensional MAD channel identified by the transition matrix:

$$\Gamma_{2B3}(\gamma_{30}, \gamma_{32}) = \mathbb{1}_3 + \gamma_{30}|2\rangle\langle 0| + \gamma_{32}|2\rangle\langle 1| - (\gamma_{30} + \gamma_{32})|2\rangle\langle 2|. \quad (5.3.8)$$

The quantum capacity of $\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)}$ can be obtained using the results in [CG21a].

Extension to $\frac{1}{2} \leq \gamma_{10} \leq 1$ and $0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2}$

$Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)})$ is a lower bound on $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})})$, as it is the emergent quantum capacity from a fixed choice of encoding. Furthermore, numerically computing the difference between these values at $\gamma_{10} = \frac{1}{2}$ and $0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2}$, one finds that (refer to Figure 5.3.5):

$$\begin{aligned} Q(\Phi_{\Gamma_{3B}(\gamma_{10}=1/2, \gamma_{30}, \gamma_{32})}) = & Q(\Phi_{\Gamma_{3B}(\gamma_{10}=1, \gamma_{30}, \gamma_{32})}) \stackrel{(5.3.7)}{=} Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)}) \\ \forall \gamma_{30}, \gamma_{32} : & 0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2} \end{aligned} \quad (5.3.9)$$

The Equation (5.3.9), together with the non-increasing monotonous property of $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})})$ in its parameters, allow to infer the value of the latter quantity in the region $0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2}, \frac{1}{2} \leq \gamma_{10} \leq 1$, plotted in Figure 5.3.6.

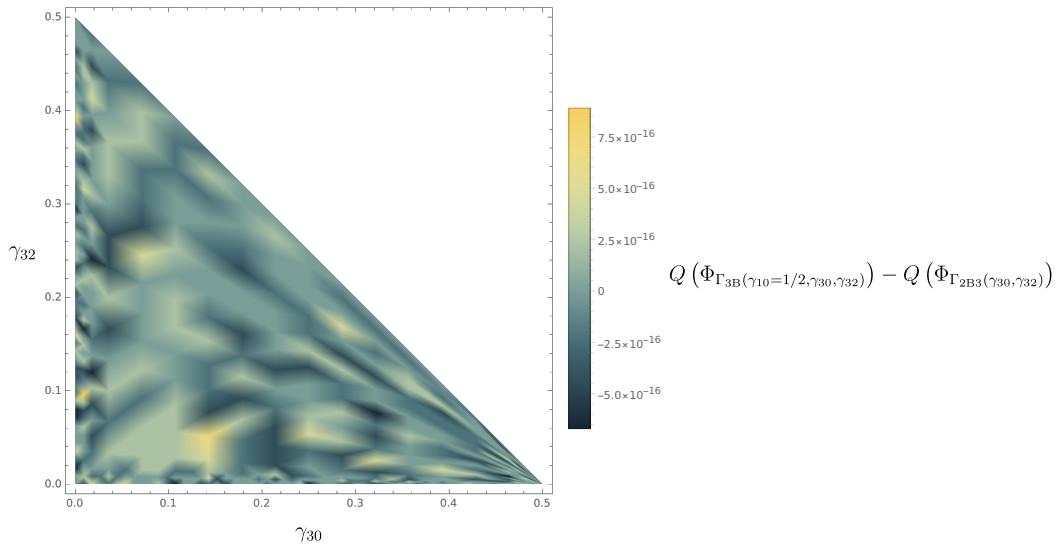


Figure 5.3.5: Density plot of the difference between $Q(\Phi_{\Gamma_{3B}(1/2, \gamma_{30}, \gamma_{32})})$ and $Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)})$. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

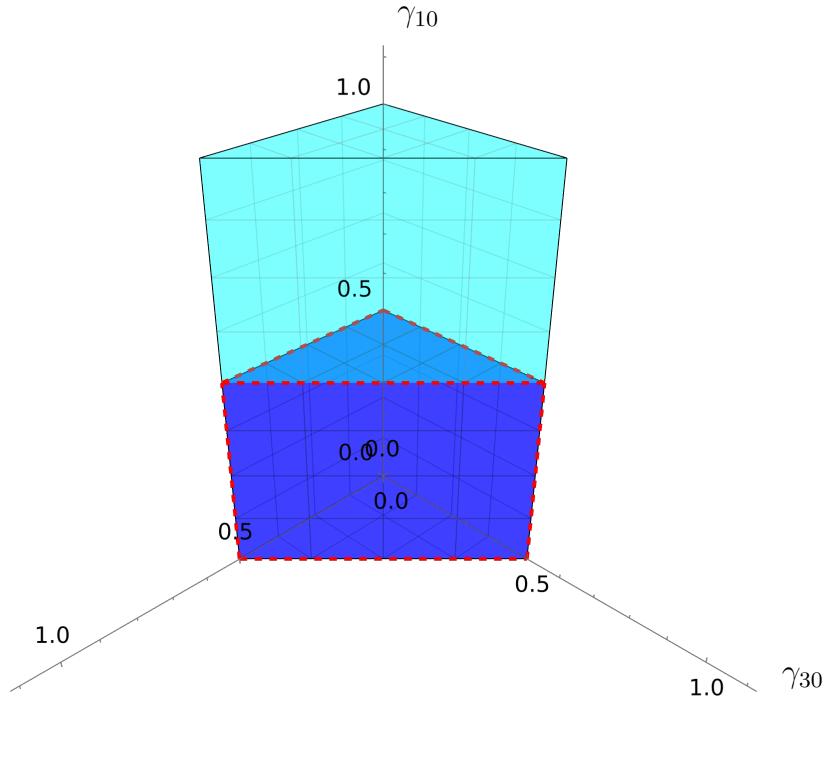


Figure 5.3.6: (5.3.10) permits the calculation of the quantum capacity in the cyan region, extending the computation performed in the degradable (blue) region

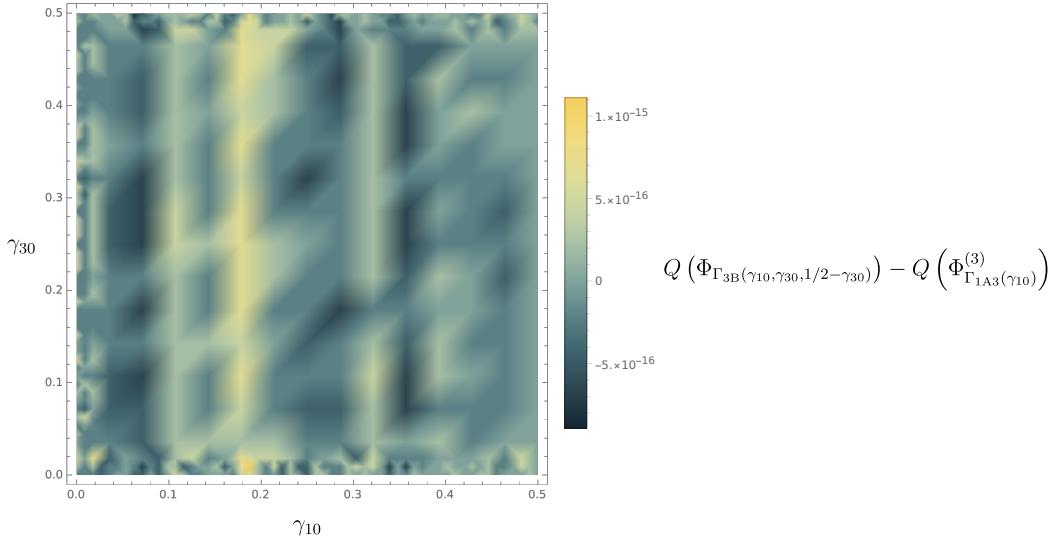


Figure 5.3.7: Density plot of the difference between $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1/2 - \gamma_{30})})$ and $Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)})$. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

$$\begin{aligned} Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}) &= Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{32})}^{(3)}) \\ \forall \gamma_{10}, \gamma_{30}, \gamma_{32} : 0 \leq \gamma_{30} + \gamma_{32} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{10} \leq 1. \end{aligned} \quad (5.3.10)$$

Extension to $\frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1, 0 \leq \gamma_{10} \leq \frac{1}{2}$

Following along the previous line of reasoning, one could calculate the difference between $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1 - \gamma_{30})}) \stackrel{(5.3.5)}{=} Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)})$ and $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1/2 - \gamma_{30})})$, which is plotted in Figure 5.3.7.

$$Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1/2 - \gamma_{30})}) = Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, 1 - \gamma_{30})}) \stackrel{(5.3.5)}{=} Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \quad (5.3.11)$$

The monotonicity properties of $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})})$, together with (5.3.11) allow to obtain the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ in the region $\frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1, 0 \leq \gamma_{10} \leq \frac{1}{2}$, represented in Figure 5.3.8.

$$\begin{aligned} Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}) &= Q(\Phi_{\Gamma_{1A3}(\gamma_{10})}^{(3)}) \\ \forall \gamma_{10}, \gamma_{30}, \gamma_{32} : \frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1 \wedge 0 \leq \gamma_{10} \leq \frac{1}{2} \end{aligned} \quad (5.3.12)$$

Extension to $\frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1, \frac{1}{2} \leq \gamma_{10} \leq 1$

The remaining region in which the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ is yet to be found is $\frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1, \frac{1}{2} \leq \gamma_{10} \leq 1$. One could compute this quantum capacity on

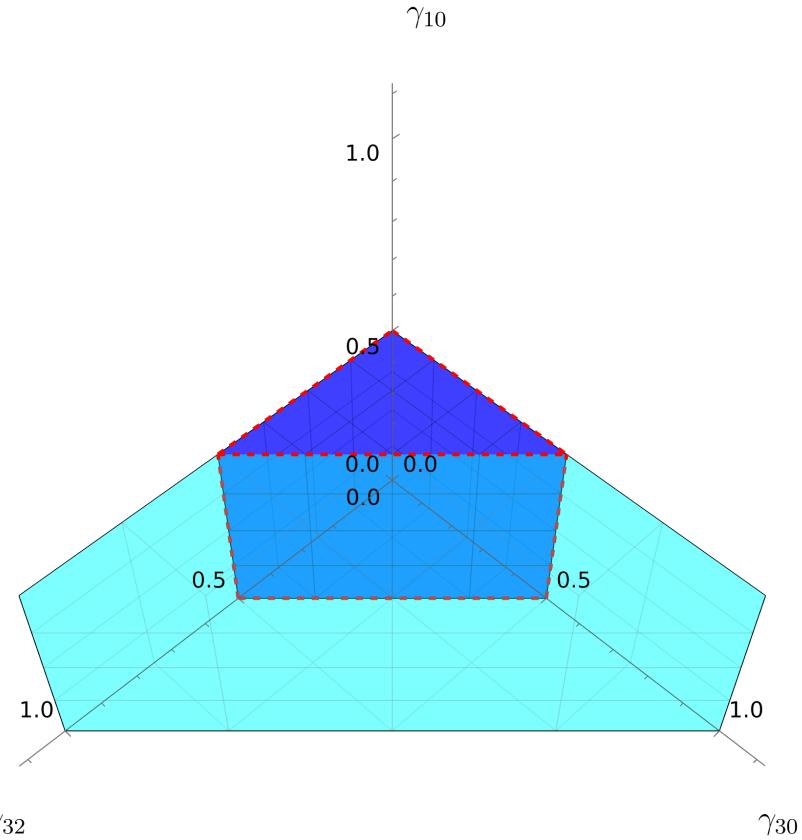


Figure 5.3.8: (5.3.10) permits the calculation of the quantum capacity in the cyan region, extending the computation performed in the degradable (blue) region

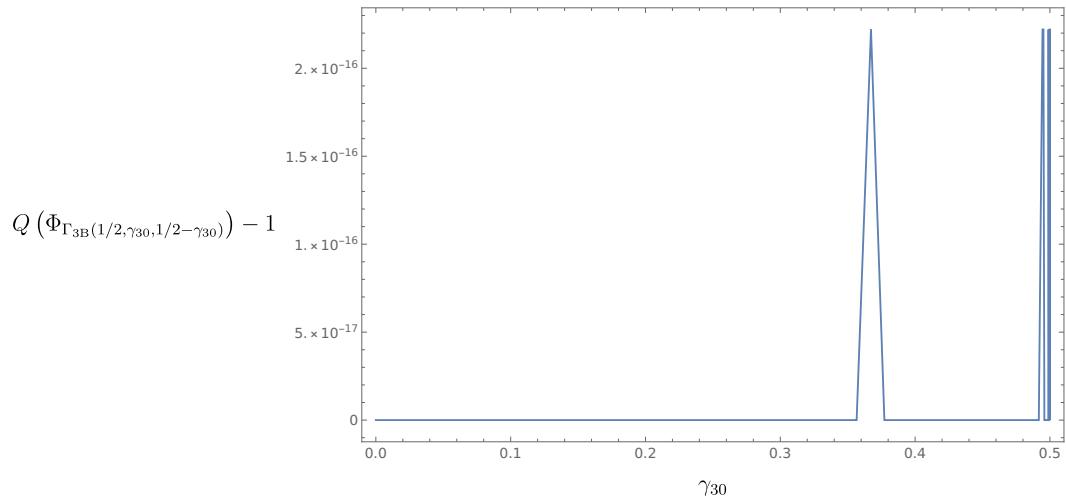


Figure 5.3.9: Plot of the difference between $Q(\Phi_{\Gamma_{3B}(1/2, \gamma_{30}, 1/2 - \gamma_{30})})$ and 1. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

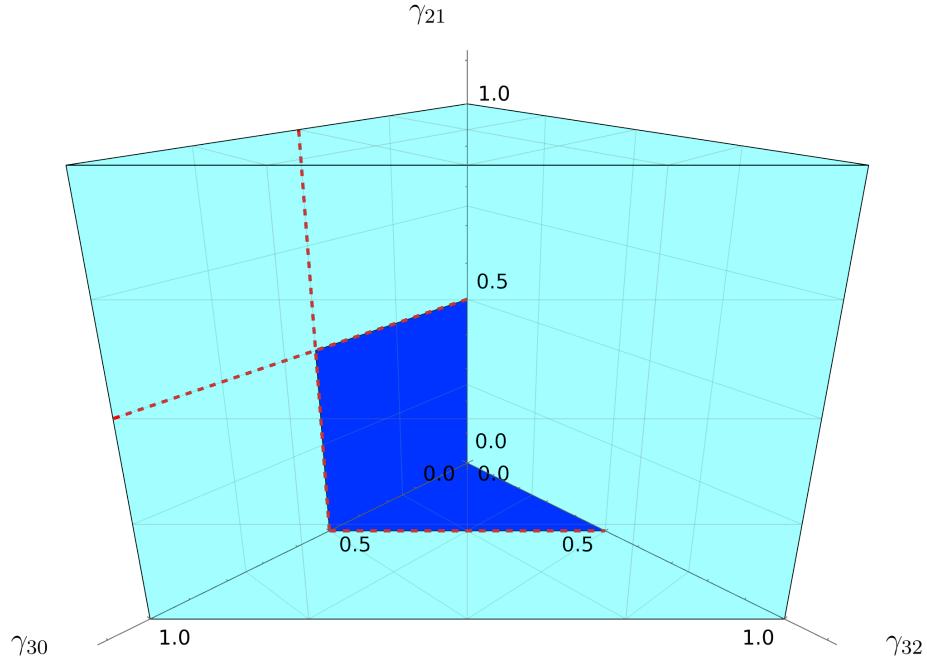


Figure 5.3.10: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32})}$ and its degradability regions (blue).

$\gamma_{10} = \frac{1}{2}$, $\gamma_{30} + \gamma_{32} = \frac{1}{2}$ and verify that it is equal to 1 (see Figure 5.3.9), with 1 being a lower bound on the quantum capacity as its the \log_2 of the noiseless subspace of the channel. Since $Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})})$ is non-increasing in its arguments, it is equal to 1 in the whole region $\frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1$, $\frac{1}{2} \leq \gamma_{10} \leq 1$:

$$\begin{aligned} Q(\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}) &= 1 \\ \forall \gamma_{10}, \gamma_{30}, \gamma_{32} : \frac{1}{2} \leq \gamma_{30} + \gamma_{32} \leq 1 \wedge \frac{1}{2} \leq \gamma_{10} \leq 1 \end{aligned} \tag{5.3.13}$$

Combining (5.3.13), (5.3.12), (5.3.10), and (5.3.4) one is able to find the quantum capacity of $\Phi_{\Gamma_{3B}(\gamma_{10}, \gamma_{30}, \gamma_{32})}$ in its whole parameter space.

5.3.c Class 3C

Consider the sample channel $\Phi_{\Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32})}$ belonging in Class 3C, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32}) &= \mathbb{1}_4 + \gamma_{21} |2\rangle\langle 1| + \gamma_{30} |3\rangle\langle 0| + \gamma_{32} |3\rangle\langle 2| \\ &\quad - \gamma_{21} |2\rangle\langle 2| - (\gamma_{30} + \gamma_{32}) |3\rangle\langle 3| \end{aligned} \tag{5.3.14}$$

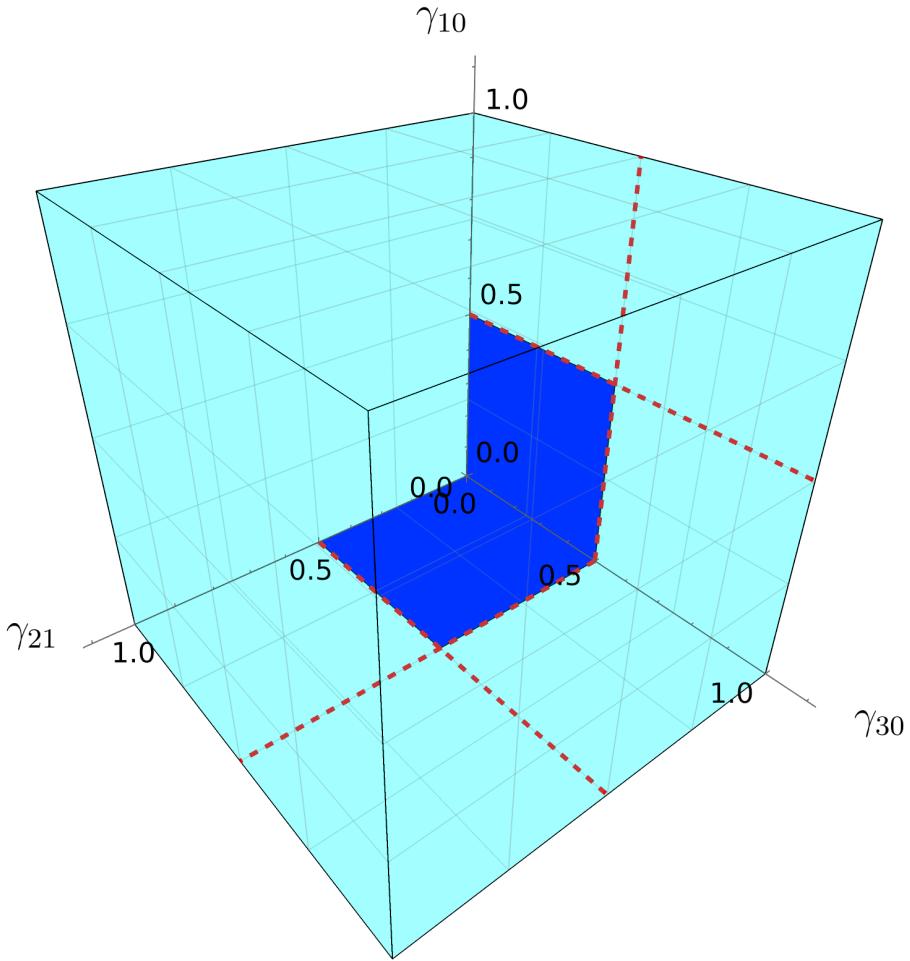


Figure 5.3.11: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32})}$ and its degradability regions (blue).

As a consequence of the results of 4.7.h, $\Phi_{\Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32})}$ is never degradable unless either one of γ_{21} , γ_{32} is 0, at which point the channel reduces, respectively, to a channel belonging in Class 2B and Class 2D. The degradability regions in the parameter space of $\Phi_{\Gamma_{3C}(\gamma_{21}, \gamma_{30}, \gamma_{32})}$ are plotted in Figure 5.3.10.

5.3.d Class 3D

Consider the sample channel $\Phi_{\Gamma_{3D}(\gamma_{10}, \gamma_{21}, \gamma_{30})}$ belonging in Class 3D, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3D}(\gamma_{10}, \gamma_{21}, \gamma_{30}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| + \gamma_{30} |3\rangle\langle 0| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2| - \gamma_{30} |3\rangle\langle 3| \end{aligned} \quad (5.3.15)$$

As a consequence of the results of 4.7.i, $\Phi_{\Gamma_{3D}(\gamma_{10}, \gamma_{21}, \gamma_{30})}$ is never degradable unless

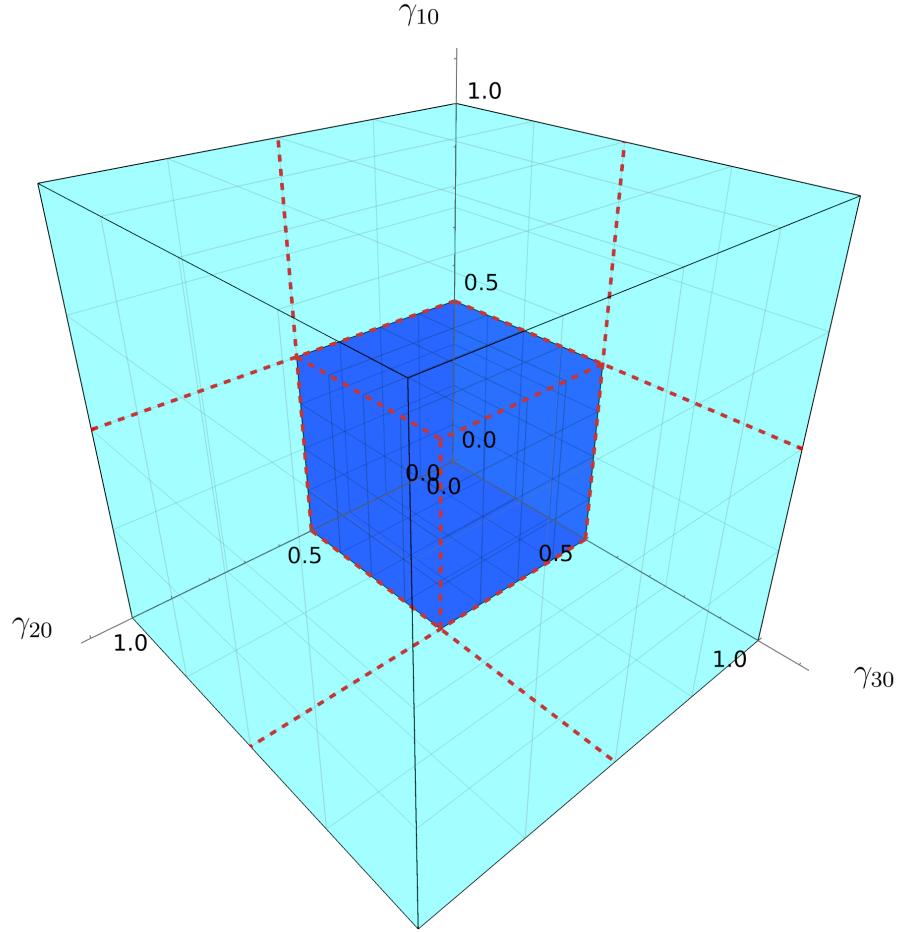


Figure 5.3.12: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3D}(\gamma_{10}, \gamma_{21}, \gamma_{30})}$ and its degradability regions (blue).

either one of γ_{10} , γ_{21} is 0, at which point the channel reduces, respectively, to a channel belonging in Class 2D and Class 2A. The degradability regions in the parameter space of $\Phi_{\Gamma_{3D}(\gamma_{10}, \gamma_{21}, \gamma_{30})}$ are plotted in Figure 5.3.11.

5.3.e Class 3E

Consider the sample channel $\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}$ belonging in Class 3E, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |2\rangle\langle 0| + \gamma_{30} |3\rangle\langle 0| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{20} |2\rangle\langle 2| - \gamma_{30} |3\rangle\langle 3| \end{aligned} \quad (5.3.16)$$

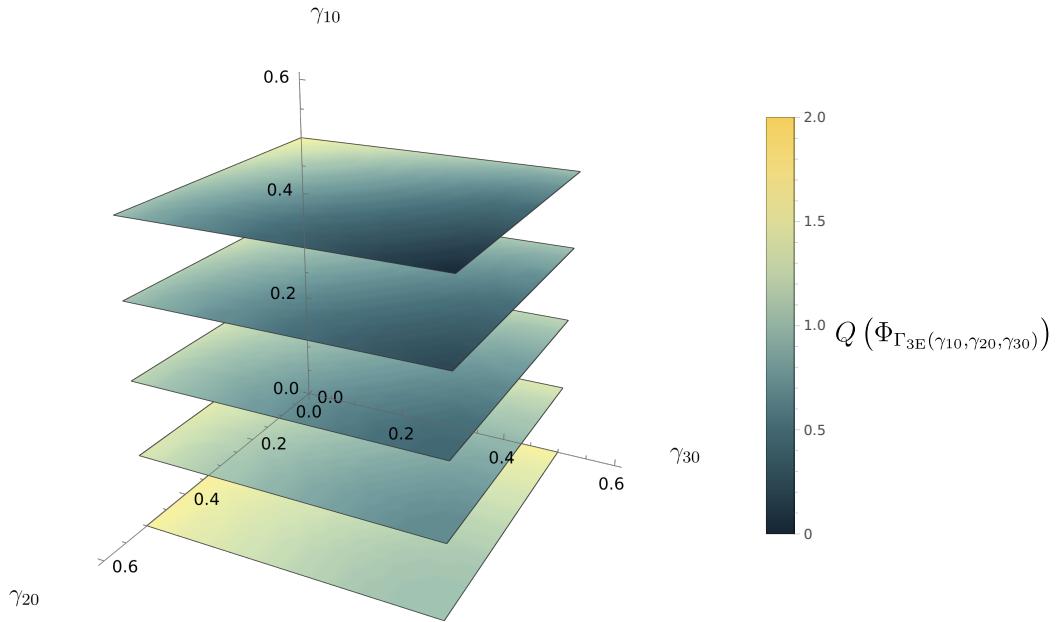


Figure 5.3.13: Slices of the quantum capacity in the 3-dimensional degradable region of $\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}$.

The degradability conditions for class 3E, reported in (4.7.13), translate to :

$$\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})} \text{ degradable} \Leftrightarrow 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{20} \leq \frac{1}{2} \wedge 0 \leq \gamma_{30} \leq \frac{1}{2}; \quad (5.3.17)$$

these conditions are rendered in Figure 5.3.12. Note that the quantum capacity of $\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}$ is left invariant under any permutation of $(\gamma_{10}, \gamma_{20}, \gamma_{30})$. It is possible to extend the computation performed in the degradability zones to the remainder of the parameter region.

Quantum capacity in degradability region

One can directly compute the quantum capacity of $\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}$ in the blue region rendered in Figure 5.3.12 using the result (3.3.21).

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}) &= \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{20} \leq \frac{1}{2} \wedge 0 \leq \gamma_{30} \leq \frac{1}{2}. \end{aligned} \quad (5.3.18)$$

This computation was performed for 5 choices of γ_{10} in the degradability region and plotted in Figure 5.3.13.

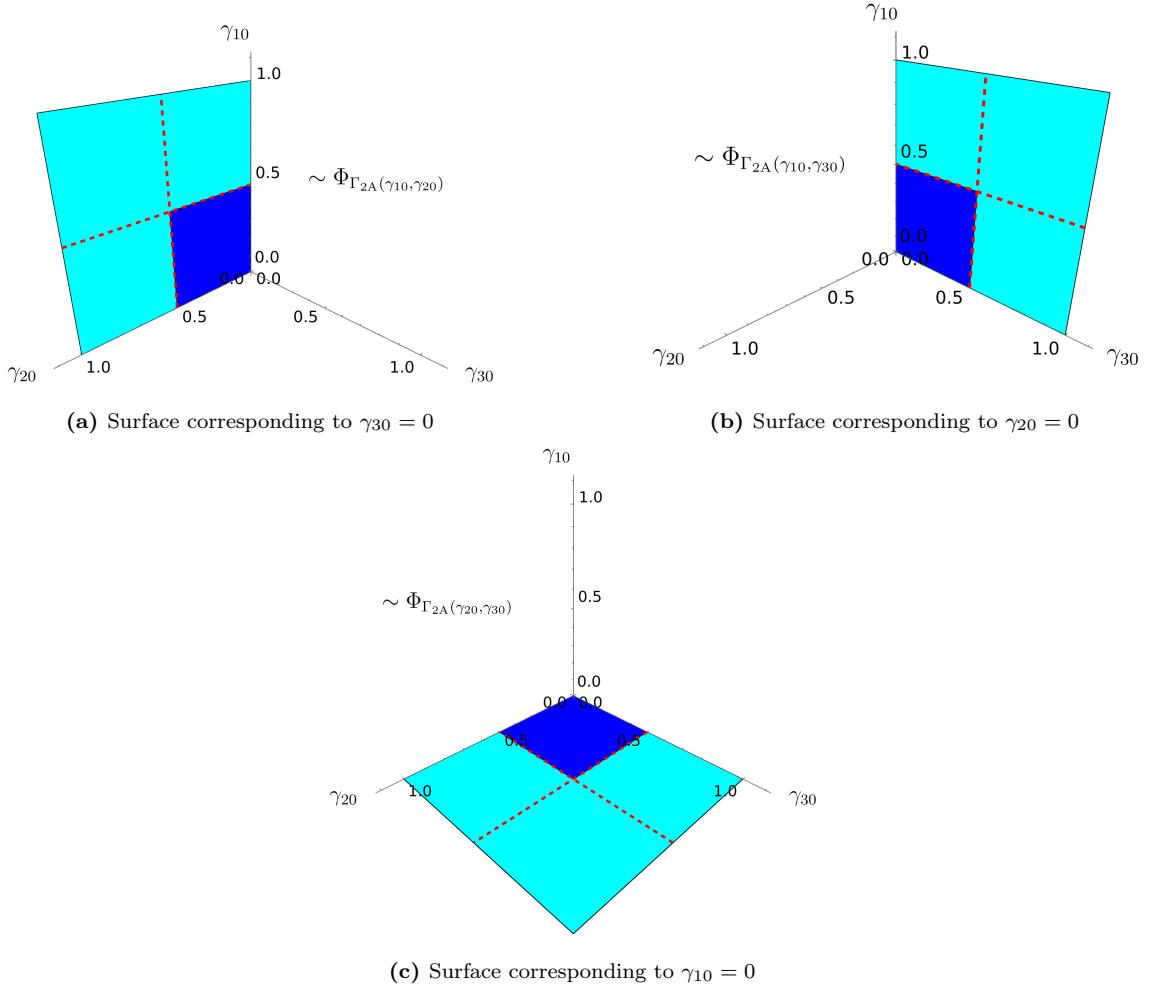


Figure 5.3.14: Surfaces in the parameter space of $\Phi_{\Gamma_{3E}}(\gamma_{10}, \gamma_{20}, \gamma_{30})$ corresponding to setting one of the transition probabilities $\gamma_{10}, \gamma_{20}, \gamma_{30}$ to 0.

Quantum capacity on borders of the parameter region

Consider the 2-dimensional borders of the cyan region in Figure 5.3.12; it is possible to compute the quantum capacity of $\Phi_{\Gamma_{3E}}(\gamma_{10}, \gamma_{20}, \gamma_{30})$ on any of those surfaces.

When either $\gamma_{10}, \gamma_{20}, \gamma_{30}$ is 0 (these surfaces are represented in Figure 5.3.14), $\Phi_{\Gamma_{3E}}(\gamma_{10}, \gamma_{20}, \gamma_{30})$ reduces to a channel belonging in Class 2A. The results in 5.2.a allow for the computation of the quantum capacity on those surfaces. Assume complete damping of a single level of the qudit, i.e. either one of $\gamma_{10}, \gamma_{20}, \gamma_{30}$ is 1; given the invariance of $Q(\Phi_{\Gamma_{3E}}(\gamma_{10}, \gamma_{20}, \gamma_{30}))$ under swap of the parameters, one could study, without loss of generality, the case $\gamma_{30} = 1$. Using (4.10.9), the quantum capacity of $\Phi_{\Gamma_{3E}}(\gamma_{10}, \gamma_{20}, \gamma_{30})$ in

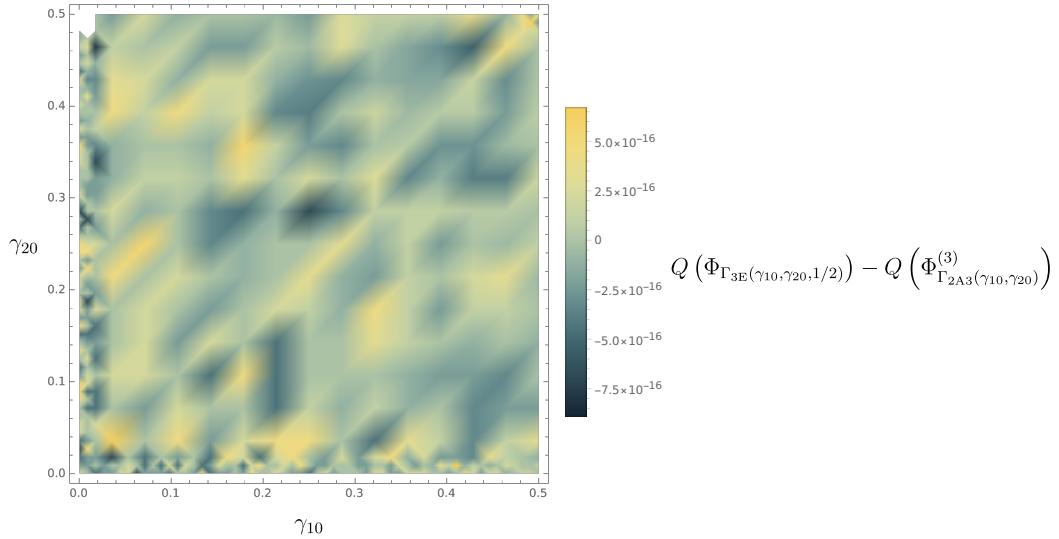


Figure 5.3.15: Density plot of the difference between $Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30}=1/2)})$ and $Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)})$. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

the case of complete damping for a level of the qudit is:

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30}=1)}) &= Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}), \\ Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}=1, \gamma_{30})}) &= Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{30})}^{(3)}), \\ Q(\Phi_{\Gamma_{3E}(\gamma_{10}=1, \gamma_{20}, \gamma_{30})}) &= Q(\Phi_{\Gamma_{2A3}(\gamma_{20}, \gamma_{30})}^{(3)}), \end{aligned} \quad (5.3.19)$$

where $\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}$ is a 3-dimensional MAD channel whose identifying transition matrix is:

$$\Gamma_{2A3}(\gamma_{10}, \gamma_{20}) = \mathbb{1}_3 + \gamma_{10}|1\rangle\langle 0| + \gamma_{20}|2\rangle\langle 0| - \gamma_{10}|1\rangle\langle 1| - \gamma_{20}|2\rangle\langle 2| \quad (5.3.20)$$

The quantum capacity of $\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}$ can be obtained exploiting the results from [CG21a].

Extension to $\frac{1}{2} \leq \gamma_{30} \leq 1$, $0 \leq \gamma_{20} \leq \frac{1}{2}$, $0 \leq \gamma_{10} \leq \frac{1}{2}$

$Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)})$ is a lower bound on $Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})})$, as it is the emergent quantum capacity from a fixed choice of encoding. Furthermore, numerically computing the difference between these values at $\gamma_{30} = \frac{1}{2}$ and $0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{20} \leq \frac{1}{2}$, one finds that (refer to Figure 5.3.15):

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30}=1/2)}) &= Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30}=1)}) \stackrel{(5.3.19)}{=} Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}) \\ \forall \gamma_{10}, \gamma_{20} : 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{20} \leq \frac{1}{2}. \end{aligned} \quad (5.3.21)$$

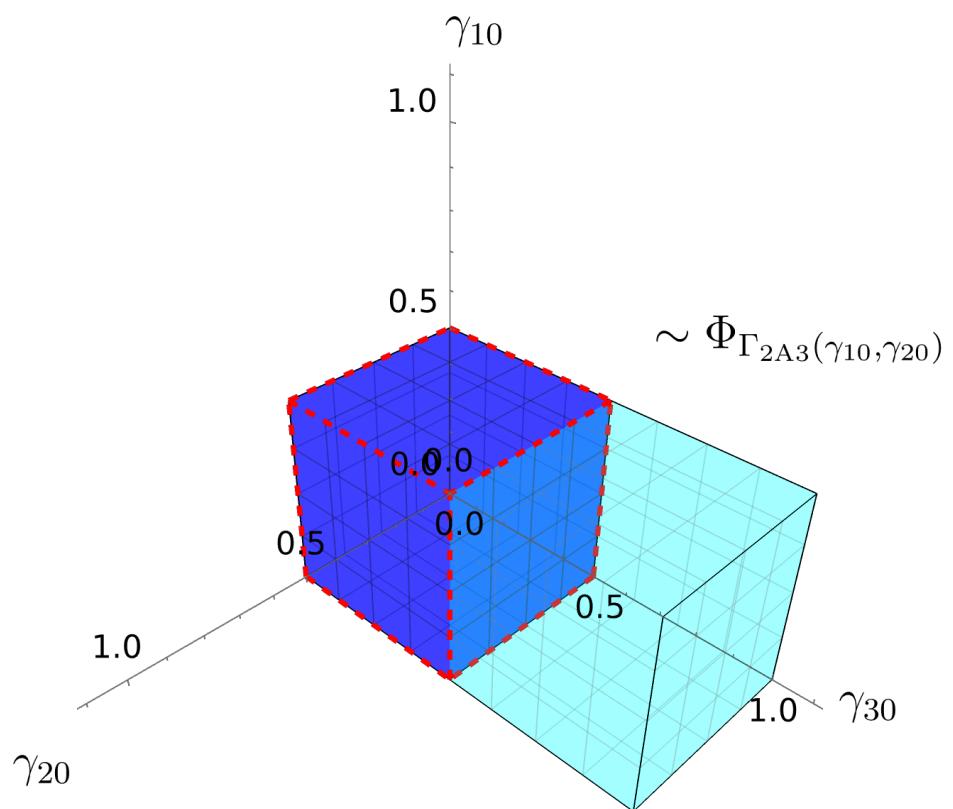


Figure 5.3.16: (5.3.22) permits the calculation of the quantum capacity in the cyan region, extending the computation performed in the degradable (blue) region

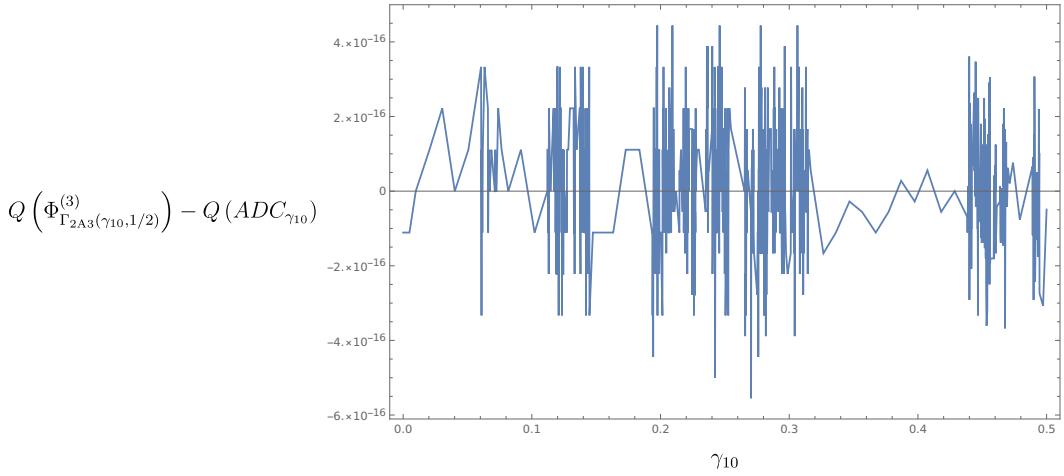


Figure 5.3.17: plot of the difference between $Q\left(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20}=1/2)}^{(3)}\right)$ and $Q(ADC_{\gamma_{10}})$. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

The Equation (5.3.21), together with the non-increasing monotonous property of $Q\left(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}\right)$ in its parameters and its invariance under swap of the parameters, allow to infer the value of the latter quantity in the regions , the first of which is plotted in Figure 5.3.16.

$$\begin{aligned} Q\left(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}\right) &= Q\left(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20})}^{(3)}\right) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{20} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{30} \leq 1. \end{aligned} \quad (5.3.22)$$

$$\begin{aligned} Q\left(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}\right) &= Q\left(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{30})}^{(3)}\right) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge 0 \leq \gamma_{30} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{20} \leq 1. \end{aligned} \quad (5.3.23)$$

$$\begin{aligned} Q\left(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}\right) &= Q\left(\Phi_{\Gamma_{2A3}(\gamma_{20}, \gamma_{30})}^{(3)}\right) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{20} \leq \frac{1}{2} \wedge 0 \leq \gamma_{30} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{10} \leq 1. \end{aligned} \quad (5.3.24)$$

Extension to $\frac{1}{2} \leq \gamma_{30} \leq 1$, $\frac{1}{2} \leq \gamma_{20} \leq 1$, $0 \leq \gamma_{10} \leq \frac{1}{2}$

Following along the previous line of reasoning, one could prove that:

$$\begin{aligned} Q\left(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}\right) &= Q(ADC_{\gamma_{10}}) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{10} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{20} \leq 1 \wedge \frac{1}{2} \leq \gamma_{30} \leq 1, \end{aligned} \quad (5.3.25)$$

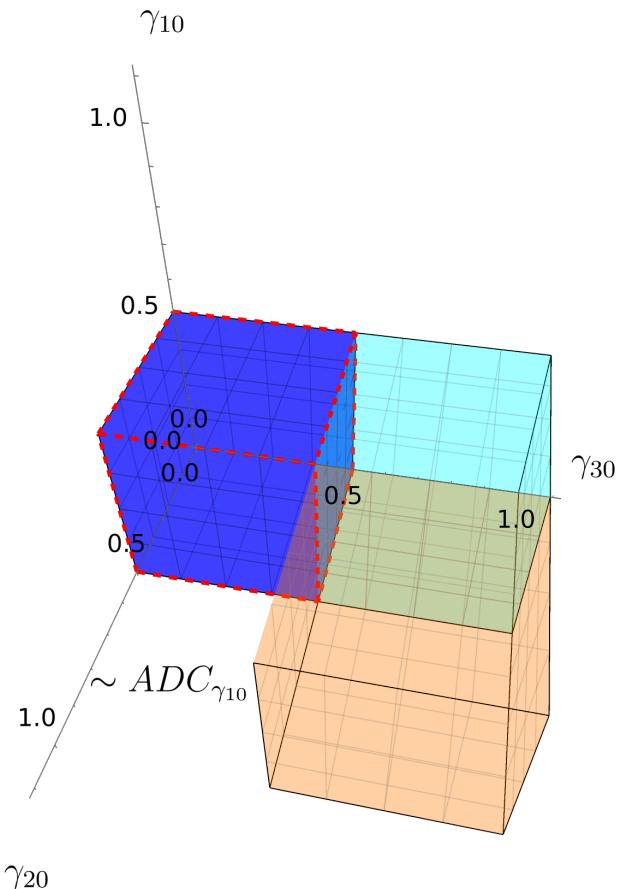


Figure 5.3.18: (5.3.25) permits the calculation of the quantum capacity in the orange region, extending the computation performed in the cyan region, possible thanks to (5.3.22).

where $ADC_{\gamma_{10}}$ was defined in (3.1.3), by proving that

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}=1/2, 1/2 \leq \gamma_{30} \leq 1)}) &\stackrel{(5.3.22)}{=} Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20}=1/2)}^{(3)}) \\ &= Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20}=1)}^{(3)}) \\ &\stackrel{[CG21a]}{=} Q(ADC_{\gamma_{10}}) \\ \forall \gamma_{10} : 0 \leq \gamma_{10} \leq \frac{1}{2}. \end{aligned} \quad (5.3.26)$$

A numerical computation of the difference between $Q(\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20}=1/2)}^{(3)})$ and $Q(ADC_{\gamma_{10}})$ was performed and plotted in Figure 5.3.18. Note that, as a consequence of the invariance of $Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})})$ under swap of its parameters, from (5.3.25) one could also infer that:

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}) &= Q(ADC_{\gamma_{20}}) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{20} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{10} \leq 1 \wedge \frac{1}{2} \leq \gamma_{30} \leq 1, \end{aligned} \quad (5.3.27)$$

$$\begin{aligned} Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}) &= Q(ADC_{\gamma_{30}}) \\ \forall \gamma_{10}, \gamma_{20}, \gamma_{30} : 0 \leq \gamma_{30} \leq \frac{1}{2} \wedge \frac{1}{2} \leq \gamma_{10} \leq 1 \wedge \frac{1}{2} \leq \gamma_{20} \leq 1. \end{aligned} \quad (5.3.28)$$

Combining the symmetry under swap of the parameters of $Q(\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})})$, (5.3.25), (5.3.22) and (5.3.18), the quantum capacity of $\Phi_{\Gamma_{3E}(\gamma_{10}, \gamma_{20}, \gamma_{30})}$ is found in the whole of its parameter space.

5.3.f Class 3F

Consider the sample channel $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ belonging in Class 3F, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32}) &= \mathbb{1}_4 + \gamma_{30}|3\rangle\langle 0| + \gamma_{31}|3\rangle\langle 1| + \gamma_{32}|3\rangle\langle 2| \\ &\quad - (\gamma_{30} + \gamma_{31} + \gamma_{32})|3\rangle\langle 3| \end{aligned} \quad (5.3.29)$$

The degradability conditions for class 3E, reported in (4.7.15), translate to :

$$\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})} \text{ degradable} \Leftrightarrow 0 \leq \gamma_{30} + \gamma_{31} + \gamma_{32} \leq \frac{1}{2} \quad (5.3.30)$$

this conditions are rendered in Figure 5.3.19. Note that the quantum capacity of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ is left invariant under any permutation of $(\gamma_{30}, \gamma_{31}, \gamma_{32})$. It is possible to extend the computation performed in the degradability zones to the remainder of the parameter region.

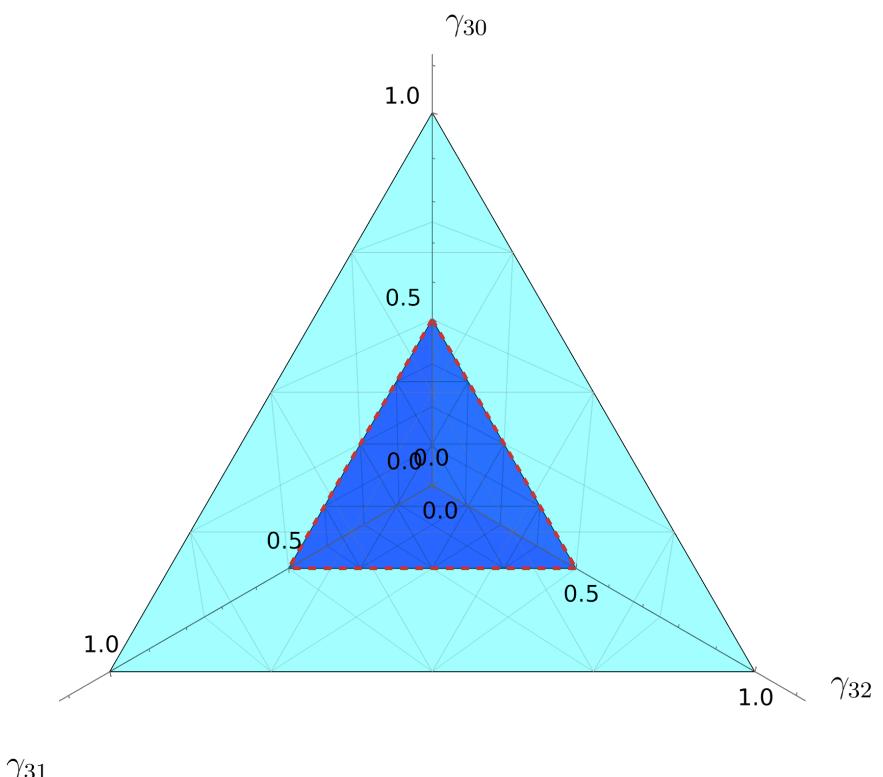


Figure 5.3.19: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ and its degradability region (blue).

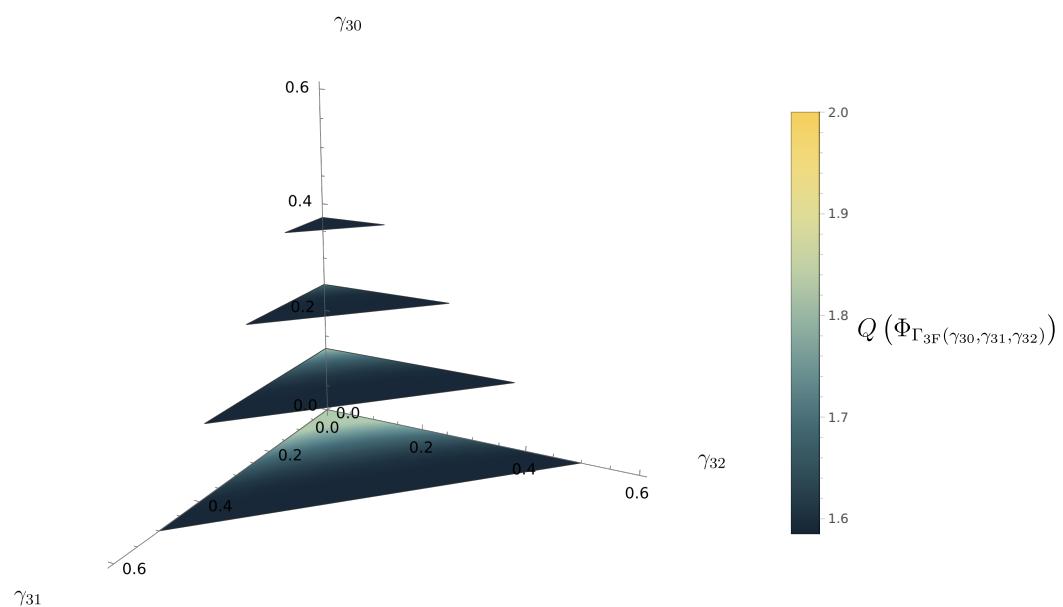


Figure 5.3.20: Slices of the quantum capacity in the 3-dimensional degradable region of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$.

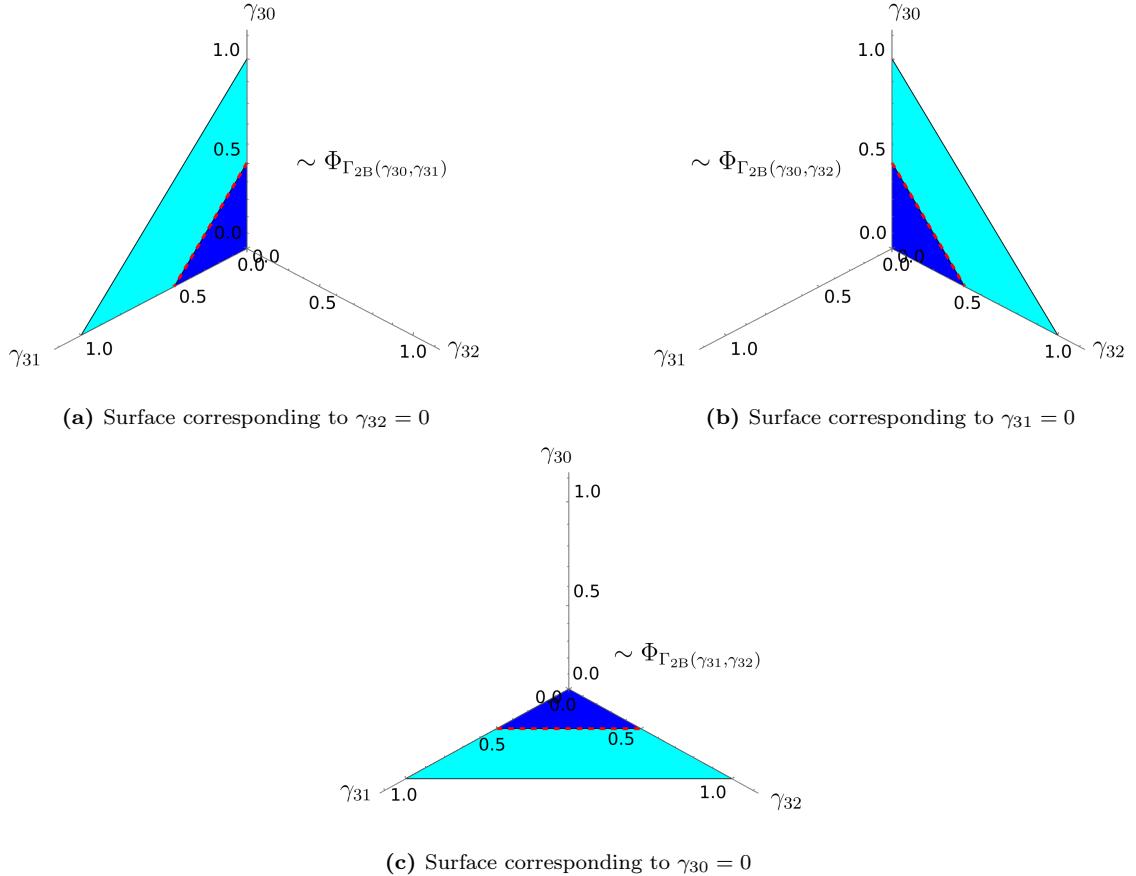


Figure 5.3.21: Surfaces in the parameter space of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ corresponding to setting one of the transition probabilities $\gamma_{30}, \gamma_{31}, \gamma_{32}$ to 0.

Quantum capacity in degradability region

One can directly compute the quantum capacity of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ in the blue region rendered in Figure 5.3.19 using the result (3.3.21).

$$Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}) = \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}) \quad (5.3.31)$$

$$\forall \gamma_{30}, \gamma_{31}, \gamma_{32} : 0 \leq \gamma_{30} + \gamma_{31} + \gamma_{32} \leq \frac{1}{2}$$

This computation was performed for 4 choices of γ_{30} in the degradability region and plotted in Figure 5.3.20.

Quantum capacity on borders of the parameter region

Consider the 2-dimensional borders of the cyan region in Figure 5.3.19; it is possible to compute the quantum capacity of $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ on any of those surfaces. When either

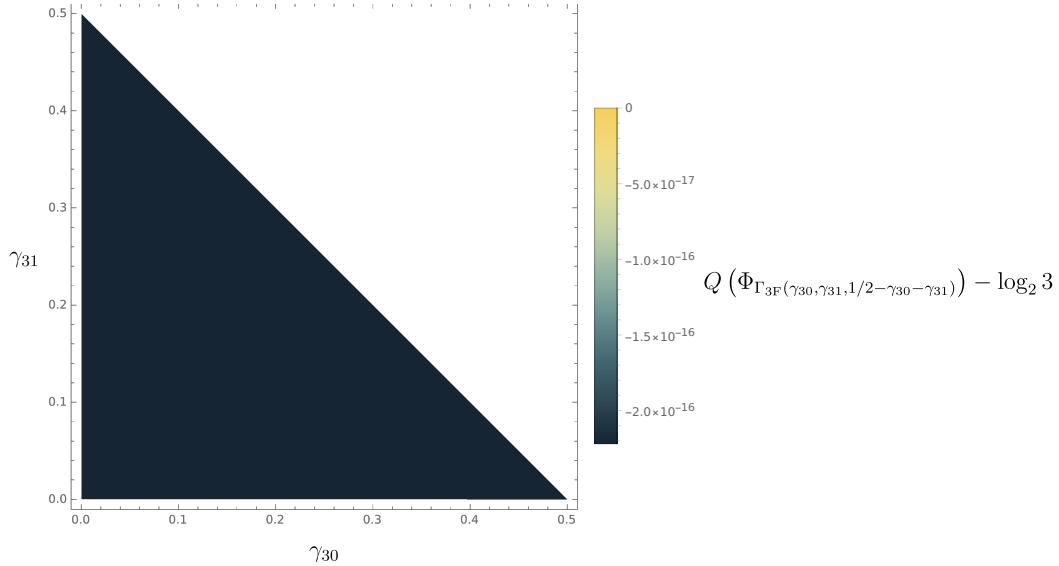


Figure 5.3.22: Density plot of the difference between $Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, 1/2 - \gamma_{30} - \gamma_{31})})$ and $\log_2 3$. The discrepancies from 0 are to be attributed to the precision of the machine used for the computation.

$\gamma_{30}, \gamma_{31}, \gamma_{32}$ is 0 (these surfaces are represented in Figure 5.3.21), $\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}$ reduces to a channel belonging in Class 2B. The results in 5.2.b allow for the computation of the quantum capacity on those surfaces. In the case of complete damping of level $|3\rangle$, i.e. when $\gamma_{30} + \gamma_{31} + \gamma_{32} = 1$, from the result (4.10.9) one can infer that:

$$Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, 1 - \gamma_{30} - \gamma_{31})}) = \log_2 3 \quad (5.3.32)$$

Extension to $\frac{1}{2} \leq \gamma_{30} + \gamma_{31} + \gamma_{32} \leq 1$

A natural lower bound on $Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})})$ is $\log_2 3$, as it is the emergent quantum capacity from a fixed choice of encoding. Furthermore, numerically computing the difference between these values at $\gamma_{32} = \frac{1}{2} - \gamma_{31} - \gamma_{30}$ and $0 \leq \gamma_{30} + \gamma_{31} \leq \frac{1}{2}$, one finds that (refer to Figure 5.3.22):

$$Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, 1/2 - \gamma_{30} - \gamma_{31})}) = Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, 1 - \gamma_{30} - \gamma_{31})}) \stackrel{(5.3.32)}{=} \log_2 3 \quad (5.3.33)$$

$$\forall \gamma_{30}, \gamma_{31} : 0 \leq \gamma_{30} + \gamma_{31} \leq \frac{1}{2}.$$

The Equation (5.3.33), together with the non-increasing monotonous property of $Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})})$ in its parameters, allow to infer the value of the latter quantity in the cyan region of Figure 5.3.19.

$$Q(\Phi_{\Gamma_{3F}(\gamma_{30}, \gamma_{31}, \gamma_{32})}) = \log_2 3 \quad (5.3.34)$$

$$\forall \gamma_{30}, \gamma_{31}, \gamma_{32} : \frac{1}{2} \leq \gamma_{30} + \gamma_{31} + \gamma_{32} \leq 1.$$

The equations (5.3.34) and (5.3.31) combined offer the quantum capacity of channels in Class 3F in their entire parameter space.

5.3.g Class 3G

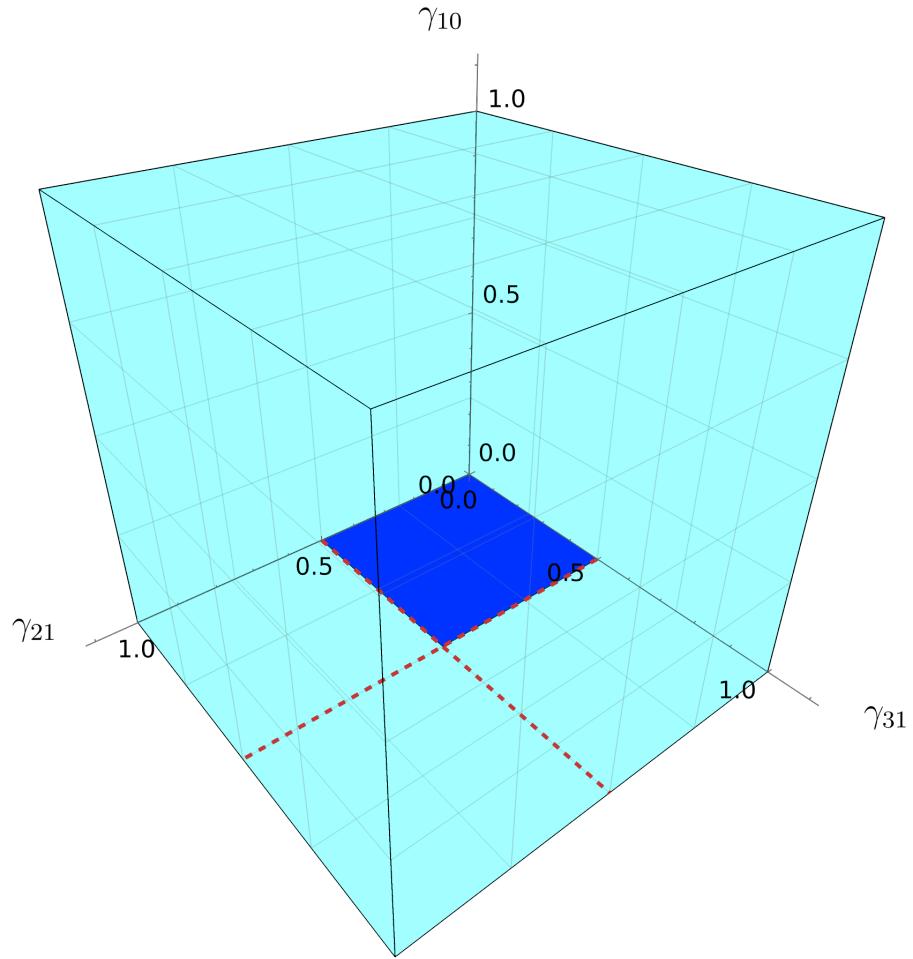


Figure 5.3.23: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3G}(\gamma_{10}, \gamma_{21}, \gamma_{31})}$ and its degradability regions (blue).

Consider the sample channel $\Phi_{\Gamma_{3G}(\gamma_{10}, \gamma_{21}, \gamma_{31})}$ belonging in Class 3G, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3G}(\gamma_{10}, \gamma_{21}, \gamma_{31}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| + \gamma_{31} |3\rangle\langle 1| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2| - \gamma_{31} |3\rangle\langle 3| \end{aligned} \quad (5.3.35)$$

As a consequence of the results of 4.7.1, $\Phi_{\Gamma_{3G}(\gamma_{10}, \gamma_{21}, \gamma_{31})}$ is never degradable unless γ_{10} is 0, at which point the channel reduces to a channel belonging in Class 2A. The degradability regions in the parameter space of $\Phi_{\Gamma_{3G}(\gamma_{10}, \gamma_{21}, \gamma_{31})}$ are plotted in Figure 5.3.23.

5.3.h Class 3H

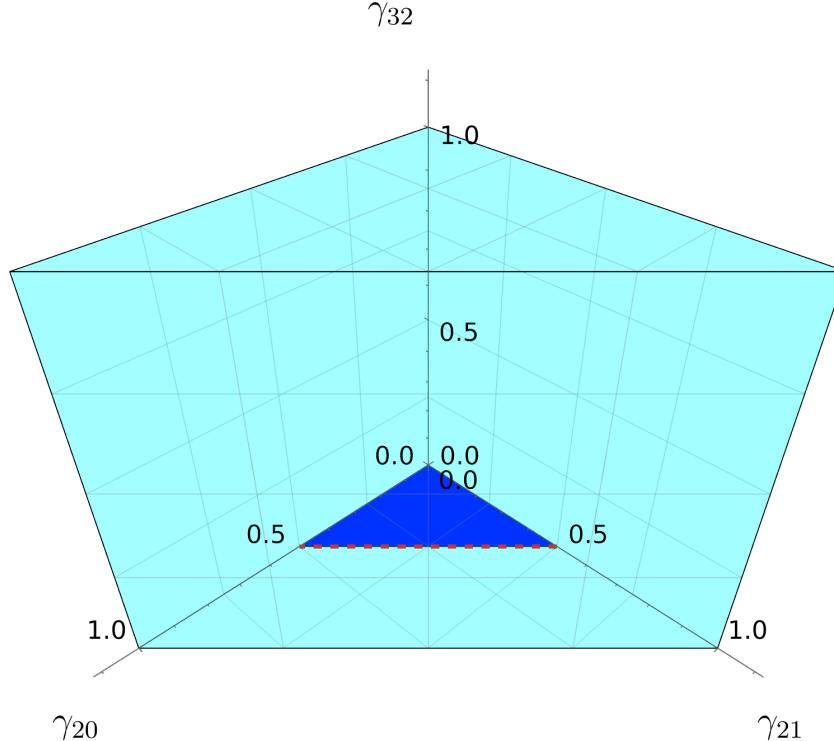


Figure 5.3.24: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3H}(\gamma_{20}, \gamma_{21}, \gamma_{32})}$ and its degradability regions (blue).

Consider the sample channel $\Phi_{\Gamma_{3H}(\gamma_{20}, \gamma_{21}, \gamma_{32})}$ belonging in Class 3H, identified by the transition matrix:

$$\begin{aligned} \Gamma_{3H}(\gamma_{20}, \gamma_{21}, \gamma_{32}) = & \mathbb{1}_4 + \gamma_{20} |2\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| + \gamma_{32} |3\rangle\langle 2| \\ & - (\gamma_{20} + \gamma_{21}) |2\rangle\langle 2| - \gamma_{32} |3\rangle\langle 3| \end{aligned} \quad (5.3.36)$$

As a consequence of the results of 4.7.m, $\Phi_{\Gamma_{3H}(\gamma_{20}, \gamma_{21}, \gamma_{32})}$ is never degradable unless γ_{32} is 0, at which point the channel reduces to a channel belonging in Class 2B. The degradability regions in the parameter space of $\Phi_{\Gamma_{3H}(\gamma_{20}, \gamma_{21}, \gamma_{32})}$ are plotted in Figure 5.3.24.

5.3.i Class 3I

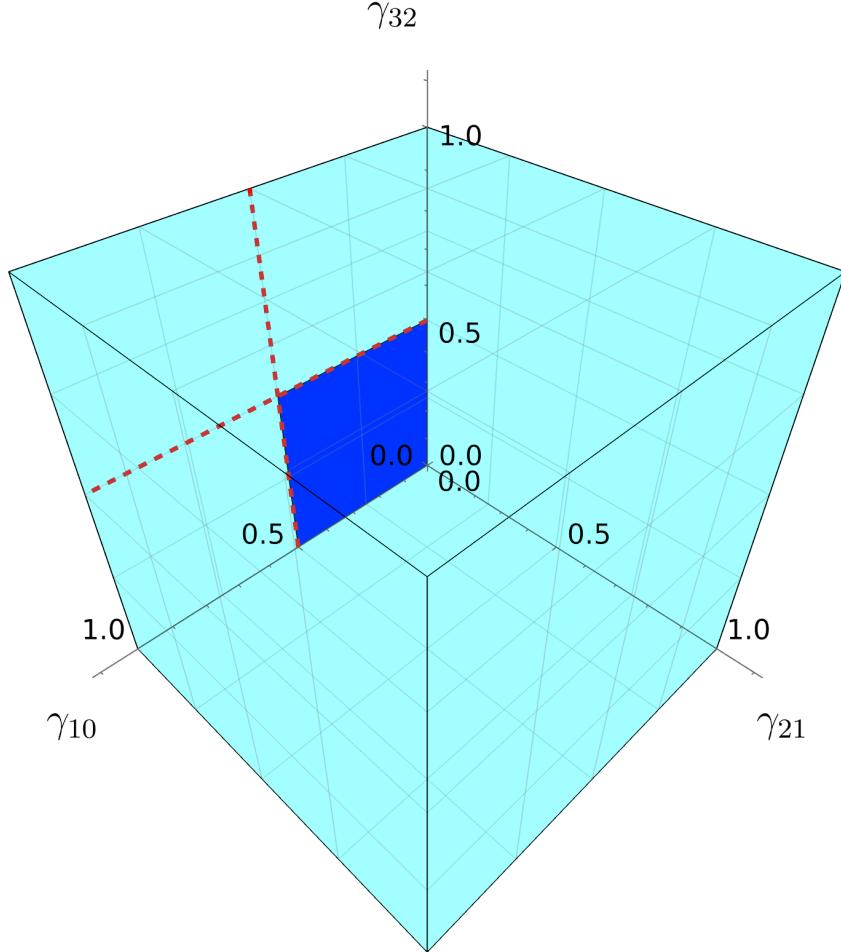


Figure 5.3.25: The graph represents the parameter region (cyan) of $\Phi_{\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32})}$ and its degradability regions (blue).

Consider the sample channel $\Phi_{\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32})}$ belonging in Class 3I, identified by the transition matrix:

$$\begin{aligned}\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32}) = & \mathbb{1}_4 + \gamma_{10} |1\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| + \gamma_{32} |3\rangle\langle 2| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2| - \gamma_{32} |3\rangle\langle 3|\end{aligned}\tag{5.3.37}$$

As a consequence of the results of 4.7.n, $\Phi_{\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32})}$ is never degradable unless γ_{21} is 0, at which point the channel reduces to a channel belonging in Class 2D. The degradability regions in the parameter space of $\Phi_{\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32})}$ are plotted in Figure

5.3.25. Note that Class 3I is the only Class of 3-decay 4-dimensional MAD channels whose capacities were not shown to be monotonous in all their parameters; in fact, $Q(\Phi_{\Gamma_{3I}(\gamma_{10}, \gamma_{21}, \gamma_{32})})$ is generally not monotonous in γ_{21} under (4.9.7) in every parameter configurations.

5.4 4-decay 4-dimensional MAD channels

As mentioned in 4.7.g, there exists a Class of 4-decay 4-dimensional MAD channels which present 4-dimensional degradability conditions. This Class is unique for 4-dimensional MAD's and can be obtained by "turning on" an additional decay starting from a channel belonging in Class 3B.

5.4.a Class 4A

Consider a the channel belonging in [Class 3B](#) identified by the transition matrix:

$$\begin{aligned}\Gamma'_{3B}(\gamma_{20}, \gamma_{30}, \gamma_{31}) = & \mathbb{1}_4 + \gamma_{20} |2\rangle\langle 0| + \gamma_{30} |3\rangle\langle 0| + \gamma_{31} |3\rangle\langle 1| \\ & - \gamma_{20} |2\rangle\langle 2| - (\gamma_{30} + \gamma_{31}) |3\rangle\langle 3|;\end{aligned}\quad (5.4.1)$$

by allowing for an additional decay from level $|2\rangle$ to level $|1\rangle$, one obtains a channel belonging in Class 4A $\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}$, identified by the transition matrix:

$$\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31}) = \Gamma'_{3B}(\gamma_{20}, \gamma_{30}, \gamma_{31}) + \gamma_{21} |2\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2| \quad (5.4.2)$$

The degradability conditions for class 4A, reported in (4.7.9), translate to :

$$\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})} \text{ degradable} \Leftrightarrow 0 \leq \gamma_{20} + \gamma_{21} \leq \frac{1}{2} \wedge 0 \leq \gamma_{31} + \gamma_{30} \leq \frac{1}{2}; \quad (5.4.3)$$

For obvious reasons, a visual rendering of these conditions in a 4-dimensional parameter space is not possible in this context. It is possible to extend the computation of the quantum capacity $\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}$ in its entire parameter space. Note that the quantum capacity $Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})})$ presents the following symmetries:

$$Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) = Q(\Phi_{\Gamma_{4A}(\gamma_{30}, \gamma_{31}, \gamma_{20}, \gamma_{21})}), \quad (5.4.4)$$

$$Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) = Q(\Phi_{\Gamma_{4A}(\gamma_{21}, \gamma_{20}, \gamma_{31}, \gamma_{30})}). \quad (5.4.5)$$

Quantum capacity in degradability region

The quantum capacity in the region described in (5.4.3) can be computed directly using (3.3.21):

$$\begin{aligned}Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) = & \max_{\rho^{(diag)}} I_c(\rho^{(diag)}, \Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) \\ \forall \gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31} : & 0 \leq \gamma_{20} + \gamma_{21} \leq \frac{1}{2} \wedge 0 \leq \gamma_{31} + \gamma_{30} \leq \frac{1}{2}.\end{aligned}\quad (5.4.6)$$

Quantum capacity on borders of the parameter region

Aside from the 3-dimensional borders obtained by setting one of the decay probabilities of $\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}$ to 0, which reduce it to a channel belonging in Class 3B, one may find the quantum capacities on the borders corresponding to $\gamma_{20} + \gamma_{21} = 1$ and, by (5.4.4), to $\gamma_{30} + \gamma_{31} = 1$. Suppose then $\gamma_{30} + \gamma_{31} = 1$; by (4.10.9), the quantum capacity on this border is:

$$Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, 1-\gamma_{30})}) = Q(\Phi_{\Gamma_{2B3}(\gamma_{20}, \gamma_{21})}^{(3)}), \quad (5.4.7)$$

where $\Phi_{\Gamma_{2B3}(\gamma_{20}, \gamma_{21})}^{(3)}$ is defined in (5.3.8) and whose capacity can be found using the results in [CG21a].

Equation (5.4.4) implies that:

$$Q(\Phi_{\Gamma_{4A}(1-\gamma_{20}, \gamma_{30}, \gamma_{31})}) = Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{31})}^{(3)}). \quad (5.4.8)$$

Extension to $\frac{1}{2} \leq \gamma_{30} + \gamma_{31} \leq 1 \wedge 0 \leq \gamma_{20} + \gamma_{21} \leq \frac{1}{2}$

$Q(\Phi_{\Gamma_{2B3}(\gamma_{20}, \gamma_{21})}^{(3)})$ is a lower bound on $Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})})$ as it is the emergent quantum capacity from a fixed choice of encoding. Furthermore, a numerical evaluation of the difference between $Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, 1/2-\gamma_{30})})$ and $Q(\Phi_{\Gamma_{2B3}(\gamma_{20}, \gamma_{21})}^{(3)})$ yields discrepancies of the order $\lesssim 10^{(-15)}$, likely due to the machine's precision, which, together with the monotonicity rules 4.9, implies that:

$$\begin{aligned} Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) &= Q(\Phi_{\Gamma_{2B3}(\gamma_{20}, \gamma_{21})}^{(3)}) \\ \forall \gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31} : \frac{1}{2} \leq \gamma_{30} + \gamma_{31} \leq 1 \wedge 0 \leq \gamma_{20} + \gamma_{21} \leq \frac{1}{2}. \end{aligned} \quad (5.4.9)$$

The (5.4.4) also implies that:

$$\begin{aligned} Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) &= Q(\Phi_{\Gamma_{2B3}(\gamma_{30}, \gamma_{31})}^{(3)}) \\ \forall \gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31} : \frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1 \wedge 0 \leq \gamma_{30} + \gamma_{31} \leq \frac{1}{2}. \end{aligned} \quad (5.4.10)$$

Extension to $\frac{1}{2} \leq \gamma_{30} + \gamma_{31} \leq 1 \wedge \frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1$

A natural lower bound on $Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})})$ is offered by the \log_2 of the dimension of the noiseless subspace of the channel, which corresponds to 1; a numerical evaluation yields $Q(\Phi_{\Gamma_{4A}(1/2-\gamma_{20}, \gamma_{30}, 1/2-\gamma_{30})}) = Q(\Phi_{\Gamma_{2B3}(1/2, 1/2-\gamma_{20})}^{(3)})$ [CG21a] $= 1$. By the non-increasing properties of $Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})})$ in its arguments, it is possible to conclude that:

$$\begin{aligned} Q(\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}) &= 1 \\ \forall \gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31} : \frac{1}{2} \leq \gamma_{20} + \gamma_{21} \leq 1 \wedge \frac{1}{2} \leq \gamma_{30} + \gamma_{31} \leq 1. \end{aligned} \quad (5.4.11)$$

Combining (5.4.9), (5.4.10), (5.4.11) and (5.4.6), one is able to obtain the quantum capacity of $\Phi_{\Gamma_{4A}(\gamma_{20}, \gamma_{21}, \gamma_{30}, \gamma_{31})}$ in the entirety of its parameter space. TODO graph if you have time.

6

2-way capacity for 3-dimensional MAD channels

The relation (2.12.3) provides an easier to compute lower bound for the 2-way capacity of a quantum channel; in the case of 3-dimensional MAD channels (2.12.3) offers a lower bound on Q_2 for any given choice of the decay probabilities $(\gamma_{10}, \gamma_{20}, \gamma_{21})$. Thanks to the results in [CG21a], it is possible to compare this bound to the natural lower bound given by the quantum capacity Q where this quantity was found. This comparison was performed and reported in the following sections.

6.1 2-way capacity for Class 1A3 MAD's - lower bound

Consider a single decay 3-dimensional MAD channel $\Phi_{\Gamma_{1A3}(\gamma_{10},0,0)}^{(3)}$, identified by the transition matrix:

$$\Gamma_{1A3}(\gamma_{10}, 0, 0) \equiv \mathbb{1}_{\mathcal{H}_3} + \gamma_{10} |1\rangle\langle 0| - \gamma_{10} |1\rangle\langle 1|. \quad (6.1.1)$$

Comparing (2.12.3) with the results of [CG21a] one obtains the plot reported in Figure 6.1.1

6.2 2-way capacity for Class 2A3 MAD's - lower bound

Consider a double decay 3-dimensional MAD channel $\Phi_{\Gamma_{2A3}(\gamma_{10},\gamma_{20},0)}^{(3)}$, identified by the transition matrix:

$$\begin{aligned} \Gamma_{2A3}(\gamma_{10}, 0, 0) \equiv & \mathbb{1}_{\mathcal{H}_3} + \gamma_{10} |1\rangle\langle 0| + \gamma_{20} |2\rangle\langle 0| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{20} |2\rangle\langle 2|. \end{aligned} \quad (6.2.1)$$

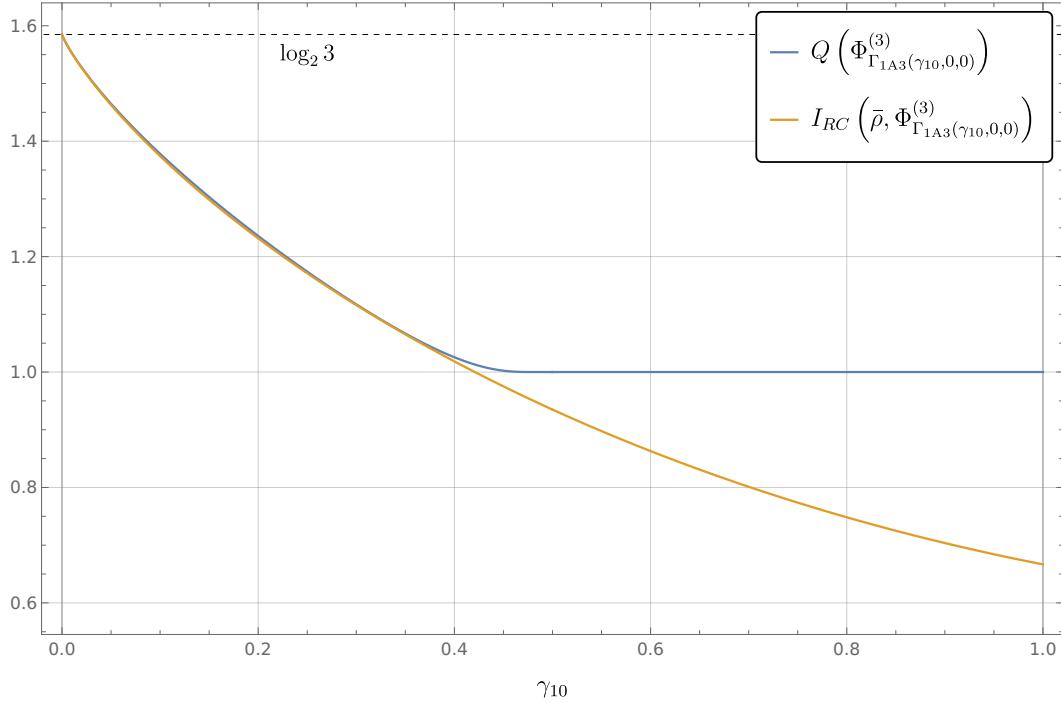


Figure 6.1.1: Comparison of lower bounds for the 2-way capacity of $\Phi_{\Gamma_{1A3}(\gamma_{10}, 0, 0)}^{(3)}$.

Comparing (2.12.3) with the results of [CG21a] one obtains the plot reported in Figure 6.2.1

6.3 2-way capacity for Class 2B3 MAD's - lower bound

Consider a double decay 3-dimensional MAD channel $\Phi_{\Gamma_{2B3}(0, \gamma_{20}, \gamma_{21})}^{(3)}$, identified by the transition matrix:

$$\begin{aligned} \Gamma_{2B3}(\gamma_{10}, 0, 0) \equiv & \mathbb{1}_{\mathcal{H}_3} + \gamma_{20} |2\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| \\ & - (\gamma_{20} + \gamma_{21}) |2\rangle\langle 2|. \end{aligned} \tag{6.3.1}$$

Comparing (2.12.3) with the results of [CG21a] one obtains the plot reported in Figure 6.3.1

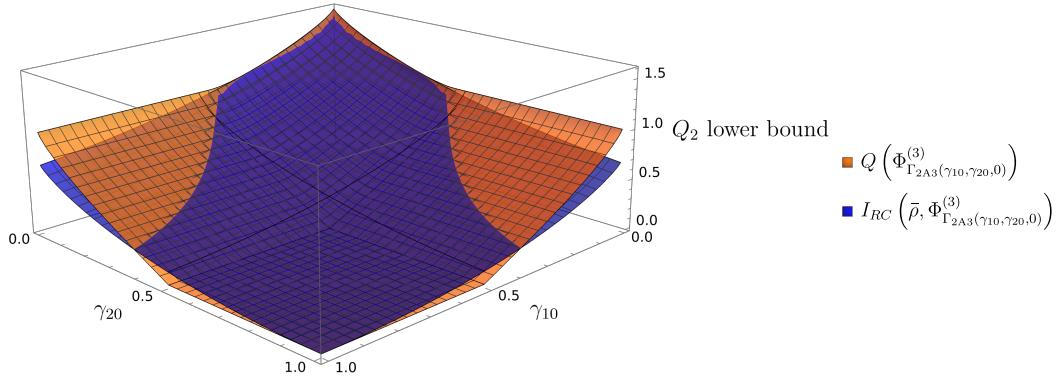


Figure 6.2.1: Comparison of lower bounds for the 2-way capacity of $\Phi_{\Gamma_{2A3}(\gamma_{10}, \gamma_{20}, 0)}^{(3)}$.

6.4 2-way capacity for Class 2C3 MAD's - lower bound

Consider a double decay 3-dimensional MAD channel $\Phi_{\Gamma_{2C3}(\gamma_{10}, 0, \gamma_{21})}^{(3)}$, identified by the transition matrix:

$$\begin{aligned} \Gamma_{2C3}(\gamma_{10}, 0, 0) \equiv & \mathbb{1}_{\mathcal{H}_3} + \gamma_{10} |1\rangle\langle 0| + \gamma_{21} |2\rangle\langle 1| \\ & - \gamma_{10} |1\rangle\langle 1| - \gamma_{21} |2\rangle\langle 2|. \end{aligned} \quad (6.4.1)$$

Comparing (2.12.3) with the results of [CG21a] one obtains the plot reported in Figure 6.4.1

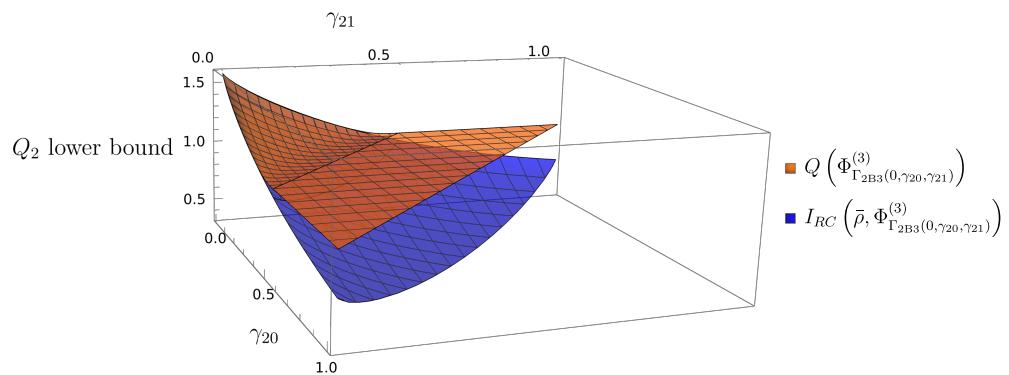


Figure 6.3.1: Comparison of lower bounds for the 2-way capacity of $\Phi_{\Gamma_{2B3}(0,\gamma_{20},\gamma_{21})}^{(3)}$.

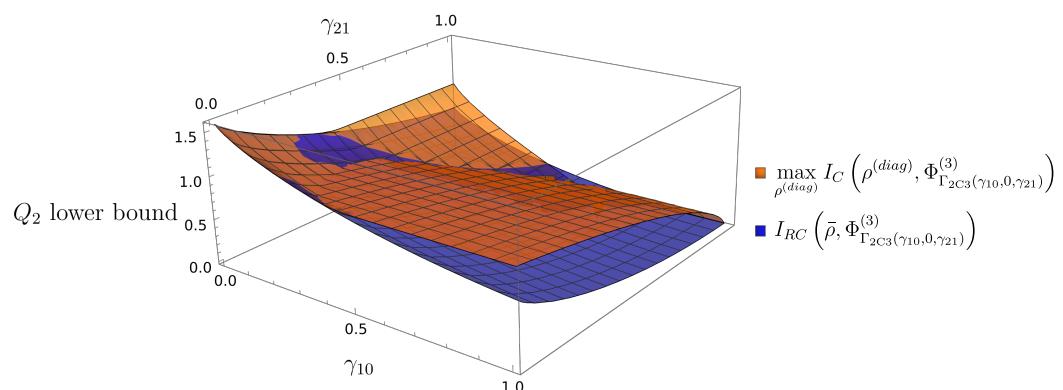


Figure 6.4.1: Comparison of lower bounds for the 2-way capacity of $\Phi_{\Gamma_{2C3}(\gamma_{10},0,\gamma_{21})}^{(3)}$.

Conclusion

What has been exposed in this Thesis work is the continuation of an effort of the scientific community to better understand the long distance communication capabilities of qudit-based systems. In fact, MAD channels describe an energy decay process when transmitting messages over long distances, and the capacities of those channels quantify how much the induced noise degrades the signal. Given that energy decays in these processes are very likely to happen (although they are not the only possible kind of noise), the study of their capacities would serve as an important headlight for the development of quantum communication technologies based upon qudits.

7.1 Summary

The study mentioned above, for 4-dimensional MAD channels, was achieved for the quantum capacity during this Thesis work in 4 main steps:

- Explore general properties of MAD channels, such as composition rules and monotonicity of capacity functionals.
- Find degradability regions in the parameter space for MAD channels.
- Compute the quantum capacity in the degradability regions.
- Extend the computations performed in the degradability regions to some non-degradable regions.

During the first step of this process, a number of results were found which were reported in Chapter 4. Particularly, the results in Section 4.9 on the monotonicity of the capacities of MAD channels meant that the technique for extending the computation of the capacities first developed in [CG21a] could also be applied to the wide range of Classes of 4-dimensional MAD channels, listed in Section 4.7, which admit degradability regions. These degradability regions were found thanks to the relation (4.6.1) using the methods outlined in Section 7.2. The computation of the quantum capacity, thanks to

the result (3.3.21), was performed as illustrated in Section 7.2. The resulting plots are reported in 5, along with the explanations of the extension to non-degradable regions.

Finally, a lower bound on the 2-way capacity was computed using (2.12.3)

7.2 Future research

The next step in the research of MAD channels consists in setting upper bounds on the 2-way capacity, as to this date there exists no literature in this regard. MAD channels do not seem to be telecovariant (see (2.12.9)), therefore the bound (2.12.12) would not work for this family of channels. The upper bound given by the squashed entanglement (2.12.6) is valid for any channel, however it is hard to compute; a simplification may be possible on MAD channel: in [PLOB17], it was shown that the search of the sup in (2.12.4), for the case of ADC's, can be restricted to purification of diagonal density matrices. This may also hold for higher dimensional settings. Furthermore, one could also study a kind of channels closely related to MAD channels, Resonant Multi-level Amplitude Damping (ReMAD) channels. These channels describe an energy decay process for a qudit under the assumption that the qudit is encoded in a harmonic oscillator, meaning that the energy gaps between successive energy levels of the qudit are fixed. Just as MAD's, they can be identified by a transition matrix (3.3.4), but their Kraus set is different:

$$\begin{aligned} K_{\Gamma}^{(i)} &\equiv \sum_{l=0}^{d-i-1} \sqrt{\gamma_{i+l,l}} |l\rangle\langle i+l| \quad \forall 0 \leq i \leq d-1 \\ \Phi_{\Gamma}^{(\text{ReMAD})}(\rho) &= \sum_{i=0}^{d-1} K_{\Gamma}^{(i)} \rho K_{\Gamma}^{(i)\dagger} \end{aligned} \tag{7.2.1}$$

The analysis of the quantum capacity of ReMAD channels for $d = 3$ has been performed in [CG23] and is based upon the same concept as the quantum capacity analysis of MAD channels. The literature for the 2-way capacity of ReMAD channel, however, is lacking.

The lower bound on the 2-way capacity of ReMAD can always be found using (2.12.1). The squashed entanglement is always an upper bound on the 2-way capacity, and the simplification on its computation given in [PLOB17] for ADC's may also hold for ReMAD's. However, for a specific kind of ReMAD's, it might be possible to set an upper bound on Q_2 using (2.12.12). By setting the transition probabilities so that:

$$\gamma_{jk}^{(\eta)} \equiv \binom{j}{k} \eta^{j-k} (1-\eta)^k, \tag{7.2.2}$$

the corresponding ReMAD channel resembles a beam-splitter of transmittance η which gets passed through by a harmonic oscillator. In this setting, the ReMAD channel can be mapped into a gaussian channel (see e.g. [CEGH08]); gaussian channels are

always telecovariant [PLOB17], therefore their 2-way capacity is upper-bounded by the relative entropy of entanglement and, as a consequence of (2.13.1) and of the mapping between beam-splitter-type ReMAD's and gaussian channels, this bound is also an upper bound for that type of ReMAD's. As a final note, the capacity analysis of MAD's and ReMAD's seems to be highly scalable: in fact, the quantum capacities in some non-degradable regions found in this Thesis work for MAD channels reduced to quantum capacities in lower dimensional systems found in [CG21a]. The "new" computations corresponded to those of the maxima of the coherent information of the degradable regions. This suggests that, when considering higher dimensional systems, much of the work previously performed could be easily exported to future researches on this type of lossy finite-dimensional channels.

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Appendix

Big Matrices

$$\begin{pmatrix}
\rho_{00} + (1 - \gamma_{10})\rho_{11} + (1 - \gamma_{20} - \gamma_{21})\rho_{22} + (1 - \gamma_{30} - \gamma_{31} - \gamma_{32})\rho_{33} & \sqrt{\gamma_{10}}\rho_{01} & \sqrt{1 - \gamma_{10}}\sqrt{\gamma_{20}}\rho_{12} & \sqrt{\gamma_{30}}\rho_{03} & \sqrt{1 - \gamma_{10}}\sqrt{\gamma_{31}}\rho_{13} \\
(\rho_{01})^* \frac{\sqrt{\gamma_{10}}}{\sqrt{\gamma_{20}}} & (\rho_{12})^* \frac{\sqrt{\gamma_{10}}\sqrt{\gamma_{20}}}{\sqrt{\gamma_{20}}} & \frac{\sqrt{\gamma_{10}}\sqrt{\gamma_{20}}\rho_{12}}{\gamma_{20}\rho_{22}} & \frac{\sqrt{\gamma_{10}}\sqrt{\gamma_{30}}\rho_{13}}{\sqrt{\gamma_{20}}\sqrt{\gamma_{30}}\rho_{23}} & 0 \\
(\rho_{02})^* \frac{\sqrt{\gamma_{20}}}{\sqrt{\gamma_{10}}} & 0 & 0 & 0 & 0 \\
(\rho_{12})^* \frac{\sqrt{-\gamma_{10}}}{\sqrt{\gamma_{20}}} & 0 & \gamma_{21}\rho_{22} & 0 & \sqrt{\gamma_{21}}\sqrt{\gamma_{31}}\rho_{23} \\
(\rho_{03})^* \frac{\sqrt{\gamma_{10}}}{\sqrt{\gamma_{30}}} & (\rho_{13})^* \frac{\sqrt{\gamma_{10}}\sqrt{\gamma_{30}}}{\sqrt{\gamma_{30}}\sqrt{\gamma_{31}}} & (\rho_{23})^* \frac{\sqrt{\gamma_{20}}\sqrt{\gamma_{30}}}{0} & \gamma_{31}\rho_{33} & 0 \\
(\rho_{13})^* \frac{\sqrt{1 - \gamma_{10}}}{\sqrt{\gamma_{31}}} & 0 & (\rho_{23})^* \frac{\sqrt{\gamma_{21}}\sqrt{\gamma_{31}}}{0} & 0 & \gamma_{31}\rho_{33} \\
(\rho_{23})^* \frac{\sqrt{1 - \gamma_{20} - \gamma_{21}}}{\sqrt{\gamma_{32}}} & 0 & 0 & 0 & \gamma_{32}\rho_{33}
\end{pmatrix}$$

Computation methods

The values for the quantum capacities and the lower bounds on the 2-way capacity reported in this document were obtained thanks to the `Mathematica` framework. The version that was used for the computations is `Mathematica 13.0`. The most basic function necessary for quantum channels, such as the creation of a channel from its Kraus set, were provided by the [QuantumUtils](#) package.

```
"      Needs["QuantumUtils`"]
```

However, most of the functions needed were created specifically for this project. The notebooks can be found at [my GitHub page](#). The problem of finding the degradability regions of MAD channels 4.7 reduces to finding the eigenvalues of the Choi matrix of the map built in (4.6.1). The Choi matrix of this map was built using the following function:

```
" ChoiMatrixConnecting[mat_]:=(
nkr=nK[mat];
CM1=ConstantArray[0,{d nkr,d nkr}];
dm=ConstantArray[0,{d,d}];
Do[
dm[[ik,jk]]=1;
dm1=InverseMAD[dm,mat];
dm2=Comp1MAD[dm1,Kraus[KrausList[mat]]];
dm[[ik,jk]]=0;
Do[
CM1[[((ik-1) nkr+iik,(jk-1) nkr+jjk)]] = FullSimplify[dm2[[iik,jjk]]],
{iik,1,nkr},{jjk,1,nkr}
],
{ik,1,d},{jk,1,d}
];
Simplify[CM1]
);
```

where `nkr` is the number of Kraus operators and `mat` is the transition matrix of the channel. Once the Choi matrix is created, it is possible to use the built-in `Mathematica` function `Eigenvalues` to compute the eigenvalues of the matrix; if those eigenvalues are ≥ 0 , the channel is degradable.

The computation of the maximum of the coherent information in the degradable regions, corresponding to the quantum capacity in those regions, was a bit trickier. In order to help `Mathematica` understand the bounds where the search of the maximum was to be performed, a trigonometric formulation of the input density matrices was used:

```
"      BuildDensity[par_]:=(
ρSys=ConstantArray[0,{d,d}];
Do[
ρSys[[j1,j1]]=(  
    Product[Sin[ToExpression[ToString[par]<>ToString[j2-1]]]^2,
{j2=1,j1-1}];  
    Cos[ToExpression[ToString[par]<>ToString[j1-1]]]^2,
{j1,1,d-1}]);
ρSys[[d,d]]=Product[Sin[ToExpression[ToString[par]<>ToString[j2-1]]]^2,
{j2=1,d-1}];  
Simplify[ρSys]
)
```

Then, the numerical maxima over the parameters (par_j) were reported in the plots of 5.