Emerging Dynamics Arising From Coarse-Grained Quantum Systems

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The purpose of physics is to describe nature from elementary particles all the way up to cosmological objects like cluster of galaxies and black holes. Although a unified description for all this spectrum of events is desirable, an one-theory-fits-all would be highly impractical. To not get lost in unnecessary details, effective descriptions are mandatory. Here we analyze what are the dynamics that may emerge from a full quantum description when one does not have access to all the degrees of freedom of a system. More concretely, we describe the properties of the dynamics that arise from quantum mechanics if one has only access to a coarse grained description of the system. We obtain that the effective maps are not necessarily of Kraus form, due to correlations between accessible and non-accessible degrees of freedom, and that the distance between two effective states may increase under the action of the effective map. We expect our framework to be useful for addressing questions such as the thermalization of closed quantum systems, as well as the description of measurements in quantum mechanics.

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I. INTRODUCTION

It is widely accepted that quantum mechanics provides currently the best description we have of the physical world. However, the description of systems in our daily lives does not require the whole framework arising from quantum mechanics. In fact, our everyday life experiences heavily rely on effective (macroscopic) descriptions which are far less complex than their underlying intricate quantum characterization. For example, to describe the behavior of a macroscopic object, like the thermal expansions or compressions of a rail line, we do not need to specify the quantum states of all atoms composing such an object. In this situation we resort to the theory of thermodynamics [1], which is probably the clearest example of effective theories. Although the systems treated within this theory are composed by many quantum interacting particles, macroscopic variables – such as temperature, volume, and pressure– describe the systems well enough, allowing, for instance, for the design of thermal machines.

The idea of different scales is central in physics. But, how does the description in one scale emerge from the description in a deeper scale? Different ways of coarse graining the description of a system are often employed [2,6] in order to “zoom out” from one level and obtain an effective description. Coarse grainings frequently appear in statistical physics [6], and are arguably the central tool in the renormalization method developed by Kadanoff and Wilson [7,8]. Nevertheless, some of these early methods are sometimes based on not so well controlled approximations or on projections, leading thus to ill-defined and/or probabilistic effective dynamics when applied to quantum systems.

In the last decades, with the birth of the quantum information field, various tools were developed to deal with many-body quantum systems [9]. In particular, the theory of completely positive linear maps [10–12], which aims at describing the most general transformations that can be applied to a system (including the most general time evolution), became well established. This has been accompanied and supported by the formalization and development of a theory for quantum correlations [13], and by efficient descriptions of many-body quantum states [14,15]. The goal of the present contribution is to employ some of these tools in order to obtain effective descriptions of quantum systems and their dynamics. More concretely, see Fig. We give a system in the state represented by a density operator $\rho_0$ evolving under the unitary map $U_t$; what is the dynamics $\Gamma_t$ induced by a coarse graining $\Lambda_{CG}$? What types of dynamics might emerge when we departure from a full quantum
description of the systems?

In what follows we present a framework to address these questions. Its construction is closely related to that of open quantum systems \[16-19\]. In fact, concepts like the correlation between system and environment, and maps divisibility will play an important role here as well. Nevertheless, our framework encompasses and generalizes this previous formalism, as ours can be used in many other situations. It can, for instance, be used to describe closed systems from which just partial information is available, what might play a significant role in the thermalization of closed quantum systems \[20\] [21]. Our work is also related to recent articles by Kofler and Brukner \[22\] [23]. In these articles the authors analyze the effect of coarse-grained measurements in order to explain the emergence of the classical word. Their approach, however, is not dynamical, and that is exactly the gap we want to fill out.

Our article is organized as follows: In Section \[II\] we introduce two different characterizations of completely positive and trace preserving (CPTP) linear maps, which will allow us to describe generalized quantum dynamics and the coarse graining maps. In this contribution, a coarse graining map will simply be a CPTP linear map that reduces the dimension of the system. Such maps were recently used to obtain a sufficient criteria for the entanglement of high-dimensional bipartite states \[24\]. After that, in Section \[III\] we obtain the effective dynamics \(\Gamma_t\) induced by the coarse graining \(\Lambda_{CG}\), underlying evolution \(\Lambda_t\), and initial state \(\psi_0\). The properties of the effective map \(\Gamma_t\) are discussed in Section \[IV\]. In Section \[V\] we show that the distance between two effective states may increase under the action of the same effective map \(\Gamma_t\). This is in contrast with the usual contraction property of CPTP linear maps \[9\]. Finally, in Section \[VI\] we draw some final conclusions and hint to some possible applications of the developed formalism.

II. CPTP LINEAR MAPS: GENERAL DYNAMICS AND COARSE GRAINING

In order to define the coarse graining operations, which are the ones we are interested here, we will first briefly review some properties of CPTP linear maps. Comprehensive expositions can be found, for example, in \[9\] [12].

Let \(\mathcal{H}_D \simeq \mathcal{C}^D\) be the Hilbert space assigned to a \(D\)-dimensional quantum system. We define \(\mathcal{L}(\mathcal{H}_D)\) as the set of all linear operators acting on \(\mathcal{H}_D\), and \(\mathcal{D}(\mathcal{H}_D) = \{\psi \in \mathcal{L}(\mathcal{H}_D) | \psi \geq 0, \text{Tr}(\psi) = 1\}\) the convex set containing all the possible states of the system. Let \(\Lambda : \mathcal{L}(\mathcal{H}_D) \rightarrow \mathcal{L}(\mathcal{H}_C)\) be a linear map which abides by two constraints: i) it is trace preserving, meaning that \(\forall \psi \in \mathcal{L}(\mathcal{H}_D)\) we have \(\text{Tr}(\psi) = \text{Tr}(\Lambda(\psi))\); and ii) it is completely positive, i.e., for all positive operators \(\psi \in \mathcal{L}(\mathcal{H}_D \otimes \mathcal{H}_Z)\), with \(\mathcal{H}_Z\) an arbitrary finite dimensional Hilbert space, the linear map \(\Lambda \otimes \mathbb{1} : \mathcal{L}(\mathcal{H}_D \otimes \mathcal{H}_Z) \rightarrow \mathcal{L}(\mathcal{H}_C \otimes \mathcal{H}_Z)\) is such that \(\Lambda \otimes \mathbb{1}(\psi) \geq 0\) \[12\]. The first imposition guarantees that probabilities are conserved through the map action, while the completely positivity condition ensures that states are mapped into states even if the map acts only on a subsystem of the whole system. The following well-known theorem gives a very useful characterization of CPTP linear maps.

**Theorem 1** \([9\] [12]\). A linear map \(\Lambda : \mathcal{L}(\mathcal{H}_D) \rightarrow \mathcal{L}(\mathcal{H}_C)\) is completely positive and trace-preserving if and only if there exists a finite set of linear operators \(\{K_i\}_{i=1}^N\), with each \(K_i : \mathcal{H}_D \rightarrow \mathcal{H}_C\) known as a Kraus operator, such that \(\forall \psi \in \mathcal{L}(\mathcal{H}_D)\):

\[
\Lambda(\psi) = \sum_{i=1}^N K_i \psi K_i^\dagger \quad \text{with} \quad \sum_{i=1}^N K_i K_i^\dagger = \mathbb{1}_D.
\]

It is worth noticing that CPTP linear maps generalize the evolution of a quantum system, with the unitary evolution being a particular linear map \(U_t : \mathcal{L}(\mathcal{H}_D) \rightarrow \mathcal{L}(\mathcal{H}_D)\) with a single Kraus operator, namely the unitary \(U_t\) itself. In general, the number of Kraus operators is unlimited, but it is always possible to characterize a CPTP linear map \(\Lambda : \mathcal{L}(\mathcal{H}_D) \rightarrow \mathcal{L}(\mathcal{H}_C)\) with a set of Kraus operators with at most \(D.d\) elements \[9\], as this is the number of generators for the map. Moreover, the set of Kraus operators describing a given CPTP linear map is not unique. Given the two sets \(\{K_i\}_{i=1}^N\) and \(\{K_i'\}_{i=1}^M\), with \(N \geq M\), they represent the same CPTP linear map.
if, and only if, there exists a unitary $U \in \text{SU}(N)$ such that $K_i = \sum_j U_j K'_j$ (where, if necessary, we pad the smallest set with zeros) \cite{9, 12}.

This more general type of evolution allows for describing processes where there is a loss of information about the system, with pure states evolving to mixed ones. That is the case, for instance, when one is dealing with open quantum systems \cite{16}.

For the coarse graining operations we are going to employ below, the following (see Fig. 2) operational way to describe CPTP linear maps will turn handy.

\begin{figure}[h!]
\centering
\includegraphics[width=0.5\textwidth]{fig2.png}
\caption{Operational interpretation of a CPTP linear map $\Lambda$.}
\end{figure}

\textbf{Theorem 2} \cite{12}. Let $\Lambda : \mathcal{L}(\mathcal{H}_d) \to \mathcal{L}(\mathcal{H}_d)$ be a CPTP linear map. Then there exists an auxiliary Hilbert space $\mathcal{H}_r$, with dimension $r \leq d$, and a unitary $V$ acting on $\mathcal{H}_D \otimes \mathcal{H}_r \otimes \mathcal{H}_d$ such that $\forall \psi \in \mathcal{L}(\mathcal{H}_D)$

$$\Lambda(\psi) = \text{Tr}_{D'}[V(\psi \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|)V^\dagger].$$

Operationally, this theorem means that we can interpret CPTP linear maps $\Lambda$ as a unitary interaction among three systems, and further discarding of the first two parties. See Fig. 2. This interpretation is reminiscent of open quantum systems, where the system interacts unitarily with the environment, with the latter being discarded as we have no control about, or interest in, it. Here, however, the roles of system and environment are not so well delineated. As we want to allow for maps with different input-output dimensions, the partial trace is taken over the auxiliary system and also over the factor encoding the initial system state.

The theorems above provide equivalent characterization of CPTP linear maps, hence we will use them interchangeably. In fact, it is easy to relate them by setting $\forall |\psi\rangle \in \mathcal{H}_D, V(|\psi\rangle \otimes |0\rangle \otimes |0\rangle) = \sum_{i=1}^D \sum_{j=1}^r \chi_i \otimes |j\rangle \otimes K_{ij}(|\psi\rangle).$ This connection shows that the auxiliary system is necessary as to accommodate CPTP linear maps which require a number of Kraus operators bigger than $D$. We should stress that for a CPTP linear map with a set of Kraus operators $\{K_i\}_{i=1}^N$, we take the dimension of $\mathcal{H}_r$ as $r = \lceil N/D \rceil$, and must find an equivalent set of Kraus operators with $Dr$ elements, $\{K'_{ij}\}_{i=1}^{Dr}$. Hence, whenever $N > D$ the auxiliary dimension $r$ will be greater than one.

We are finally in position to establish the coarse graining operations. Roughly speaking, descriptions are named coarse-grained when some fine details of the underlying model are smoothed out, or replaced by average behaviors. In order to get valid descriptions of states after the coarse graining, we define it as a CPTP linear map that reduces the dimension of the system:

$$\Lambda_{CG} : \mathcal{L}(\mathcal{H}_D) \to \mathcal{L}(\mathcal{H}_d)$$

with $D > d$.

When one is not able to resolve the system in full detail, the coarse graining map gives an effective state for the system.

Resorting to the characterization of CPTP linear maps in Theorem 2 we know that there exists an auxiliary space $\mathcal{H}_r$ and a unitary $V : \mathcal{H}_D \otimes \mathcal{H}_r \otimes \mathcal{H}_d \to \mathcal{H}_D \otimes \mathcal{H}_r \otimes \mathcal{H}_d$, such that

$$\Lambda_{CG}(\psi) = \text{Tr}_{D'}[V(\psi \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|)V^\dagger].$$

Operationally, what the unitary $V$ accomplishes is to "write" the accessible degrees of freedom into the party in $\mathcal{H}_D$, while the unaccessible degrees of freedom are left in $\mathcal{H}_D \otimes \mathcal{H}_r$ to be later discarded. See Fig. 3. The intermediate states $\chi_0 = V(\psi \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|)V^\dagger$ and $\chi_t = V(\psi \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|)V^\dagger$, which live in $\mathcal{H}_D \otimes \mathcal{H}_r \otimes \mathcal{H}_d$, are virtual states, in the sense that they are mathematical abstractions. In this level the two contributions of degrees of freedom, accessible and non-accessible are split, but may be correlated.
A. Example: a blurred and saturated detector

In order to give a concrete example, let us consider a typical optical lattice experiment \[25\]–\[27\]. In these experiments a periodic oscillating potential is constructed by counter-propagating light beams, and individual atoms are trapped in each potential minimum. In the deep Mott insulator regime two hyperfine levels of each atom act as a qubit, and neighboring qubits interact with each other via a Heisenberg-like Hamiltonian. The measurement of each atom is made via a fluorescence technique: the atoms are shone with a laser in a way that if an atom is in the state, say, \(|1\rangle\), light is scattered by the atom, whereas if its state is \(|0\rangle\) no light is scattered. To resolve the light coming from each atom a powerful lens is necessary, and only recently a single-atom resolution was accomplished \[28\].

To simplify, consider the case with only two atoms. Suppose that the lens available is not good enough as to resolve the light coming from each individual atom. In this situation the states \(|01\rangle\) and \(|10\rangle\) cannot be distinguished. Moreover, imagine that the amount of light coming from a single atom is already sufficient to saturate the detector. Then, having two excitations, \(|11\rangle\), or one excitation, \(|01\rangle\) or \(|10\rangle\), leads to the same signal. In such conditions to describe the experiment with two atoms is superfluous, and an effective description becomes handy. These experimental conditions suggest the coarse graining presented in Table I.

| \(\Lambda_{CG}(|00\rangle)\) | \(|00\rangle\) | \(\Lambda_{CG}(|01\rangle)\) | \(|01\rangle\) | \(\Lambda_{CG}(|10\rangle)\) | \(|10\rangle\) | \(\Lambda_{CG}(|11\rangle)\) | \(|11\rangle\) |
|----------------|----------------|----------------|----------------|----------------|----------------|----------------|----------------|
| \(|00\rangle\) | \(|00\rangle\) | \(|01\rangle\) | \(|01\rangle\) | \(|10\rangle\) | \(|10\rangle\) | \(|11\rangle\) | \(|11\rangle\) |
| \(|01\rangle\) | \(|00\rangle\) | \(\sqrt{3}^{-1} |00\rangle\) | \(\sqrt{3}^{-1} |01\rangle\) | \(\sqrt{3}^{-1} |10\rangle\) | \(\sqrt{3}^{-1} |11\rangle\) |
| \(|10\rangle\) | \(|00\rangle\) | \(\sqrt{3}^{-1} |00\rangle\) | \(\sqrt{3}^{-1} |01\rangle\) | \(\sqrt{3}^{-1} |10\rangle\) | \(\sqrt{3}^{-1} |11\rangle\) |
| \(|11\rangle\) | \(|00\rangle\) | \(\sqrt{3}^{-1} |00\rangle\) | \(\sqrt{3}^{-1} |01\rangle\) | \(\sqrt{3}^{-1} |10\rangle\) | \(\sqrt{3}^{-1} |11\rangle\) |

TABLE I. Coarse graining for a blurred and saturated detector.

If a detector does not distinguish between the two systems, and does not differ between one or two excitations, this coarse graining gives the effective description of the system.

Note that as the detector does not distinguish between the states \(|01\rangle\), \(|10\rangle\), and \(|11\rangle\) there can be no coherence in this subspace. Furthermore, the \(1/\sqrt{3}\) factors are necessary to make \(\Lambda_{CG}\) a CPTP linear map. This signals that coherences in the effective description might decrease, but they do not necessarily vanish \[21\]. This can be readily seen by evaluating the action of \(\Lambda_{CG}\) over a general two-qubits pure state \(|\psi\rangle = \sum_{i,j=0}^{1} c_{ij} |ij\rangle\), with \(c_{ij} \in \mathbb{C}\), which gives:

\[
\Lambda_{CG}(|\psi\rangle) = \left( \begin{array}{cccc}
|c_{00}|^2 & |c_{01}|^2 & |c_{10}|^2 & |c_{11}|^2 \\
\sqrt{3} |c_{30}|^2 & \sqrt{3} |c_{31}|^2 & \sqrt{3} |c_{03}|^2 & \sqrt{3} |c_{13}|^2 \\
\sqrt{3} |c_{30}|^2 & \sqrt{3} |c_{31}|^2 & \sqrt{3} |c_{03}|^2 & \sqrt{3} |c_{13}|^2 \\
\sqrt{3} |c_{30}|^2 & \sqrt{3} |c_{31}|^2 & \sqrt{3} |c_{03}|^2 & \sqrt{3} |c_{13}|^2 \\
\end{array} \right).
\]

This effective state accounts for the statistics of all possible measurements that can be carried out by the detector here modeled. It is thus the description that really matters for this experimental condition, not carrying unaccessible information.

The Kraus operators for this map can be easily obtained by a quantum process tomography \[9\], and are given by:

\[
K_1 = \left( \begin{array}{cccc}
1 & 0 & 0 & 0 \\
0 & 1/\sqrt{3} & 1/\sqrt{3} & 1/\sqrt{3} \\
0 & 1/\sqrt{3} & 0 & -1/\sqrt{3} \\
0 & 1/\sqrt{3} & -1/\sqrt{3} & 0 \\
\end{array} \right);
\]

\[
K_2 = \left( \begin{array}{cccc}
0 & 0 & 0 & 0 \\
1/\sqrt{3} & 1/\sqrt{3} & 1/\sqrt{3} & 1/\sqrt{3} \\
0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 \\
\end{array} \right);
\]

\[
K_3 = \left( \begin{array}{cccc}
0 & 0 & 0 & 0 \\
0 & 1/\sqrt{3} & -1/\sqrt{3} & 0 \\
0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 \\
\end{array} \right);
\]

\[
K_4 = \left( \begin{array}{cccc}
0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 \\
1 & 1/\sqrt{3} & 1/\sqrt{3} & 0 \\
\end{array} \right).
\]

As we have four Kraus operators, \(N = 4\), and the dimension of the underlying system is also four, \(D = 4\), then the auxiliary system in \(\mathcal{H}_s\) can be taken as 1-dimensional and as such can be ignored. With the above Kraus operators, and neglecting the system in \(\mathcal{H}_s\), one can immediately obtain the corresponding unitary \(V\) for this example:

\[
V = \left( \begin{array}{cccc}
1 & 0 & 0 & 0 \\
0 & 0 & 1/\sqrt{3} & 1/\sqrt{3} \\
0 & 1 & 0 & 0 \\
0 & 0 & 1/\sqrt{3} & -1/\sqrt{3} \\
\end{array} \right).
\]

III. COARSE GRAINING INDUCED DYNAMICS

Now we address the central question of this contribution: what are the dynamics that might emerge from a fully quantum description if we are not able to resolve the system in all its details? More concretely, we look for an effective map \(\Gamma\) which makes the diagram in Fig. 1 consistent, i.e., in a way that \(\rho_0 = \Gamma(\rho_0) = \Lambda_{CG} \circ D_{\rho_0}(\psi_0)\), with \(\rho_0 = \Lambda_{CG}(\psi_0)\). The induced dynamics then emerges from a coarse grained description of the underlying dynamics.

To obtain the induced dynamics \(\Gamma\) acting on the effective state \(\rho_0\), we generalize the procedure suggested by
ployed the spectral decomposition $\omega$ on $V$. For the action of the unitary $V$, see Fig. 3 suggests the following decomposition:

$$\chi_0 = (\omega_0 \otimes \rho_0) + (\chi_0 - \omega_0 \otimes \rho_0),$$

where $\chi_0 = V(\psi_0 \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|) V^\dagger$, $\rho_0 = \Lambda_{CG}(\psi_0) = \text{Tr}_{Dr}(\chi_0)$, $\omega_0 = \text{Tr}_d(\chi_0)$. Note that $\omega_0$ is a state in $D(\mathcal{H}_D \otimes \mathcal{H}_d)$. Equation (1) is equivalent to Štelmachovič and Bužek decomposition in the abstract level $\mathcal{H}_D \otimes \mathcal{H}_C \otimes \mathcal{H}_d$, with the last term now representing the correlation between the degrees of freedom which can be accessed and those that cannot. As $V$ is unitary, we can equivalently write:

$$\psi_0 \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0| = V^\dagger (\omega_0 \otimes \rho_0) V +$$

$$+ V^\dagger (\chi_0 - \omega_0 \otimes \rho_0) V. \tag{2}$$

From the left hand side of Eq. (2) we get the evolved effective state by applying the underlying evolution map $\U_i$ onto the first tensor factor, followed by the application of $V$ and further partial trace of the two first tensor factors:

$$\rho_i = \Lambda_{CG} \otimes \U_i(\psi_0)$$

$$= \text{Tr}_{Dr}[V(\U_i(\psi_0) \otimes |0\rangle \langle 0| \otimes |0\rangle \langle 0|) V^\dagger]$$

$$= \Gamma_i(\rho_0).$$

The last equality comes from demanding consistence of the diagram in Fig. 1. Accordingly, assuming the underlying evolution map of the form $\U_i(\cdot) = U_i(\cdot)U_i^\dagger$, from the right hand side of Eq. (1) we get the effective evolution:

$$\Gamma_i(\rho_0) = \text{Tr}_{Dr}(W_i(\omega_0 \otimes \rho_0) W_i^\dagger) +$$

$$+ \text{Tr}_{Dr}(W_i(\chi_0 - \omega_0 \otimes \rho_0) W_i^\dagger), \tag{3}$$

where $W_i = V(U_i \otimes I \otimes I).V^\dagger$ is the unitary evolution operator in the level $\mathcal{H}_D \otimes \mathcal{H}_C \otimes \mathcal{H}_d$, i.e., $\chi_i = W_i \chi_0 W_i^\dagger$. See Fig. 3

The above evolution equation can be rewritten in a more meaningful way as

$$\Gamma_i(\rho_0) = \sum_{i,j} M_{ij} \rho_0 M_{ij}^\dagger + \text{Tr}_{Dr}(W_i(\chi_0 - \omega_0 \otimes \rho_0) W_i^\dagger), \tag{4}$$

with $M_{ij} = \sqrt{p_j}(|\phi_i \rangle \langle \phi_i| \otimes I)W_i(|\phi_j \rangle \langle \phi_j|)$, where we employed the spectral decomposition $\omega_0 = \sum p_j |\phi_j \rangle \langle \phi_j|$. This is the dynamics that emerges if one is not able, or does not wish, to resolve all the details of the underlying system.

The expression in Eq. (4) is composed by two contributions: the first one displays a Kraus form (see Theorem 1), with $\{M_{ij}\}$ the corresponding set of effective Kraus operators; the second one represents the evolution of the correlations between accessible and non-accessible degrees of freedom. This second term can be more clearly appreciated by evoking the Bloch representation of $\chi_0$:

$$\chi_0 = \frac{1}{Drd}(1 + (1d) \otimes \bar{\alpha} \bar{\sigma}_d +$$

$$+ \bar{\beta} \bar{\sigma}_d \otimes 1) + \sum_{i,j} \theta_{ij} \sigma^{(i)}_d \otimes \sigma^{(j)}_d W_i^\dagger.$$

where $\bar{\sigma}_d = (\sigma^{(1)}_d, \sigma^{(2)}_d, \ldots, \sigma^{(d^2-1)}_d)^T$ is a vector whose components are the $q \times q$ generalized Pauli matrices, $\bar{\alpha} \in \mathbb{R}^{d-1}$ is the Bloch vector of $\rho_0$, $\bar{\beta} \in \mathbb{R}^{(d^2-1)}$ is the Bloch vector of $\omega_0$, and the $(d^2-1)$ coefficients $\theta_{ij}$ in $\bar{\sigma}_d$ fix the correlation between accessible and non-accessible degrees of freedom. Defining the correlation matrix $|\Theta|_{ij} = (\theta_{ij} - \beta_{ij})/Drd$, the evolution of the coarse grained state can be written as:

$$\Gamma_i(\rho_0) = \sum_{i,j} M_{ij} \rho_0 M_{ij}^\dagger + \sum_{i,j} \Theta_{ij} \text{Tr}_{Dr}(W_i \sigma^{(i)}_d \otimes \sigma^{(j)}_d W_i^\dagger).$$

It can be easily verified that $\sum_{i,j} M_{ij}^\dagger M_{ij} = I_d$, and that

$$\text{Tr}_{d}(\text{Tr}_{Dr}(W_i \sigma^{(i)}_d \otimes \sigma^{(j)}_d W_i^\dagger)) = 0$$

as $W_i$ is unitary and the (generalized) Pauli matrices are traceless. These conditions guarantee that $\text{Tr}_{d}(\Gamma_i(\rho_0)) = 1$ for all times. The structure of this type of evolution is very similar to the one describing open quantum systems when system and environment are initially correlated [17, 18].

A. Example: effective dynamics as seen by a blurred and saturated detector

Consider again the situation described in subsection IA: two atoms in neighboring wells of an optical lattice being observed by a blurred and saturated detector. Suppose now that the atoms interact as specified by the Hamiltonian $H = JH i \sigma_i \otimes \sigma_j$, with $J$ a coupling constant in units of frequency. In such situation, an initial two-qubit pure state $|\psi_0\rangle = \sum_{i,j=0}^{1} c_{ij} |i j\rangle$ evolves to:

$$|\psi_t\rangle = (c_{00} |00\rangle + c_{11} |11\rangle)e^{-iJt} + (c_{01} |01\rangle + c_{10} |10\rangle)e^{iJt}.$$

The evolution of the effective state can then be easily evaluated via $\rho_t = \Lambda_{CG}(\psi_t)$, to give:

$$\rho_t = \begin{pmatrix}
|c_{00}|^2 & c_{00} c_{11} e^{-2iJt} & c_{10} e^{2iJt} & \sqrt{3} c_{01} \\
\sqrt{3} c_{00} e^{2iJt} & |c_{00}|^2 + |c_{11}|^2 & \sqrt{3} c_{01} & c_{00} e^{2iJt} \\
c_{01} e^{-2iJt} & \sqrt{3} c_{01} & |c_{01}|^2 + |c_{10}|^2 & c_{01} e^{-2iJt} \\
\sqrt{3} c_{10} & c_{01} e^{-2iJt} & c_{01} e^{2iJt} & \sqrt{3} c_{10} e^{-2iJt}
\end{pmatrix}.$$
As a concrete example, the effective evolution of a state $\psi_0$ with all coefficients $c_{ij}$ equal, $c_{ij} = 1/2$ for $i,j \in \{0,1\}$, is shown in the inset of Fig. 4. Figure 4 also shows how the purity, $\Tr(\rho_0^2)$, oscillates with time, exhibiting the alternation between pure and mixed state in the effective level. This is in clear contrast with the complete description of the system, where the state is pure for all times.

It is interesting to notice that the coefficients of $\psi_0$ define the state $\rho_0$, but also enter in the definition of the effective map $\Gamma_t$. In the above example this can be verified by evaluating $\alpha_i = \Tr_D \gamma_i 0$, which will also depend on the coefficients $c_{ij}$, That in turn, means that the effective Kraus operators $M_{ij}$ will also change with the $c_{ij}$ thus by changing $\rho_0$ the map may change. The same is true for the correlation matrix $\Theta_i j$. This interdependence of the parameters is treated in the next section, where the properties of $\Gamma_t$ are analyzed.

IV. PROPERTIES OF $\Gamma_t$

The effective map $\Gamma_t$ is generated by the underlying evolution $U_t$, the coarse graining map $\Lambda_{CG}$, and the state $\psi_0$. Equation (6), however, does not make explicit how the map depends on the elemental state $\psi_0$. For instance, how do we change the effective map $\Gamma_t$ for a fixed input state $\rho_0$? Or, how to change the effective input state keeping $\Gamma_t$ fixed? In what follows we address these and other questions making use of the Bloch representation for $\psi_0$:

$$\psi_0 = \frac{1}{D} (\mathbb{I}_D + \vec{\gamma}_0 \cdot \vec{\sigma}_D), \quad (7)$$

where $\vec{\gamma}_0 \in \mathbb{R}^{D^2-1}$ is the Bloch vector of $\psi_0$.

A. Fix $\rho_0$, change $\Gamma_t$

Fixed the coarse graining map, the Bloch vector $\vec{\alpha}$ of $\rho_0$ is obtained from $\vec{\gamma}_0$ by the linear relations:

$$\begin{aligned}
\alpha_1 &= \Tr(\Lambda_{CG}(\psi_0(\vec{\gamma}_0)) \sigma_i^{(1)}); \\
\alpha_2 &= \Tr(\Lambda_{CG}(\psi_0(\vec{\gamma}_0)) \sigma_i^{(2)}); \\
\vdots &\quad \vdots \\
\alpha_d^{(d-2)} &= \Tr(\Lambda_{CG}(\psi_0(\vec{\gamma}_0)) \sigma_i^{(d-2)}); \\
\alpha_d^{(d-1)} &= \Tr(\Lambda_{CG}(\psi_0(\vec{\gamma}_0)) \sigma_i^{(d-1)}).
\end{aligned} \quad (8)$$

In the $D^2 - 1$ dimensional space of Bloch vectors $\vec{\gamma}$ of $\psi_0$, these constraints represent hyperplanes whose intersection depicts the effective state $\rho_0$. It is important to notice that since $D > d$, the set of linear equations for the coefficients $\alpha_i$ is under-determined, meaning that various states $\psi_0$ lead to the same effective state $\rho_0$. Geometrically, in the “γ-space”, this many-to-one mapping is visualized as an hyper-surface of possible solutions.

Now, with this geometric perspective in mind, a fixed coarse graining and a fixed underlying evolution, it can be seen that changes in $\psi_0$, that move $\vec{\gamma}$ parallel to the hyper-planes within the solution hyper-surface will not affect the effective state $\rho_0$. Nevertheless, such change can induce modifications in $\vec{\alpha}$ or in $\Theta$, and as such $\Gamma_t$ will change. A simple example is presented in Fig. 5 and an abstract representation of the γ-space and the change in $\Gamma_t$ can be seen in Fig. 6.

B. Fix $\Gamma_t$, change $\rho_0$: the domain of $\Gamma_t$

To change the effective state $\rho_0$ we must change the value of the $\alpha_i$’s. Geometrically this is represented by moving the hyperplanes, defined in Eq. (8), in the γ-space. After the hyper-planes displacement, a new intersection is obtained representing now another effective state, say, $\rho_1$. As moving $\vec{\gamma}_0$ parallel to the hyper-planes changes the map, this time we must move $\vec{\gamma}_0$ perpendicular to the hyper-planes. This guarantees that only the effective state is changing. See Fig. 6).

It is important to notice that this change might in fact modify $\alpha_0$ or $\Theta$. This, however, comes only because of the change in the effective input state, as these quantities might be functions of $\vec{\alpha}$. Putting it differently, the effective Kraus operators might change, but this is only due to the change in the input of the effective map $\Gamma_t$. The dynamical equation (6) can be rewritten as to make this dependence explicit:

$$\Gamma_t(\rho_0(\alpha)) = \sum_{i,j} M_{ij}(\alpha)\rho_0(\alpha)M^\dagger_{ij}(\alpha) + \xi(\alpha), \quad (9)$$
where \( \zeta(\alpha) = \sum_{i,j} \Theta_{ij}(\alpha) \mathrm{Tr}_D(W_i \sigma_{D_0}^{(j)} \otimes \sigma_d^{(j)} W_i^*) \).

We can now determine the domain of a given \( \Gamma_t \). An effective map \( \Gamma_t \) is generated by an elemental state \( \psi_0 \), an underlying evolution map \( \Upsilon_t \), and a coarse graining map \( \Lambda_{CG} \). This information already gives the first element in the domain of \( \Gamma_t \), namely, \( \rho_0 = \Lambda_{CG}(\psi_0) \). The coarse graining \( \Lambda_{CG} \) fixes the hyper-planes in the \( \gamma \)-space through Eq. (8). Let \( \vec{v}_i \) be the normal vector for the \( i \)-th hyper-plane, and \( \vec{\gamma}_0 \) be the Bloch vector of \( \psi_0 \). The domain of \( \Gamma_t \) is then given by all \( \rho = \Lambda_{CG}(\psi) \) generated from \( \psi \) with Bloch vector \( \vec{\gamma} \) for which there exist coefficients \( c_i \in \mathbb{R} \) such that \( \vec{\gamma} = \vec{\gamma}_0 + \sum c_i \vec{v}_i \). The latter condition guarantees that the Bloch vector of all states in the domain of \( \Gamma_t \) can be reached from \( \vec{\gamma}_0 \) by moving it perpendicular to the hyper-planes in Eq. (8), and such not changing the effective map.

This immediately implies that the domain of \( \Gamma_t \) is convex: Let \( \rho_a = \Lambda_{CG}(\psi_a) \) and \( \rho_b = \Lambda_{CG}(\psi_b) \) be in the domain of \( \Gamma_t \). This means that there exist coefficients \( \{a_i\} \subset \mathbb{R} \) and \( \{b_i\} \subset \mathbb{R} \) such that the Bloch vectors of \( \psi_a \) and \( \psi_b \) can be written as \( \vec{\gamma}_a = \vec{\gamma}_0 + \sum a_i \vec{v}_i \) and \( \vec{\gamma}_b = \vec{\gamma}_0 + \sum b_i \vec{v}_i \), respectively. There are many states in \( \mathcal{L}(\mathcal{H}_d) \) which after the coarse graining lead to the convex combination \( \rho = p\rho_a + (1-p)\rho_b \), with \( p \in [0,1] \). In particular, the state \( \psi = p\psi_a + (1-p)\psi_b \) is such that \( \Lambda_{CG}(\psi) = \rho \) and it has Bloch vector \( \vec{\gamma}(\psi_0) + \sum (p a_i + (1-p) b_i) \vec{v}_i \). Therefore the convex combination \( \rho \) is also in the domain of \( \Gamma_t \).

C. Effective positivity and complete-positivity of \( \Gamma_t \)

Equation (9) clearly shows that in general \( \Gamma_t \) is not of “Kraus” form, like shown in Theorem 1. This means that if \( \Gamma_t \) is taken as a map between states from \( \mathcal{D}(\mathcal{H}_d) \) to itself, then \( \Gamma_t \) is not completely positive, possibly not even positive.

However, as we have just seen, the domain of a given \( \Gamma_t \) is not necessarily all the states in \( \mathcal{D}(\mathcal{H}_d) \). Restricting the action of \( \Gamma_t \) to its domain guarantees the positivity of the map. This can be immediately verified by the simple consistence of the diagram in Fig. 5, which demands

\[ \Gamma_t(\rho_0) = \Gamma_t \circ \Lambda_{CG}(\psi_0) = \Lambda_{CG} \circ \Upsilon_t(\psi_0). \]

As the right most part of this equation is a composition of positive maps, then the positivity of the first term is also guaranteed.
The same line of thought can be used as to argument for the complete positivity of \( \Gamma_t \). Indeed, as we are constructing our framework upon quantum mechanics, no contradiction with it can be obtained. However this argument should not go through without a caveat: not all extensions of effective states \( \rho \) in the domain of a given \( \Gamma_t \) into \( \omega \in \mathcal{D}(\mathcal{H}_A \otimes \mathcal{H}_B) \), with \( \mathcal{H}_A \) the Hilbert space of an auxiliary system, are possible. The possible extensions for \( \rho = \Lambda_{CG}(\psi) \) are those that can be obtained from states \( \Psi \in \mathcal{D}(\mathcal{H}_A \otimes \mathcal{H}_B) \) such that \( \text{Tr}_A(\Psi) = \psi \), which guarantees that \( \text{Tr}_A(\text{Id}_A \otimes \Lambda_{CG}(\psi)) = \rho \), and that \( \psi \) generates the map \( \Gamma_t \) (together with the underlying evolution \( \Lambda_{CG} \)). We call such a family of states \( \Psi \) as the set of effective complete positivity for \( \Gamma_t \). Physically this constraint comes from the fact that if one does not have control of all the degrees of freedom of a system, then not all the states can be generated. In another perspective, the entanglement that can be created in the fundamental level \( \mathcal{H}_A \otimes \mathcal{H}_B \) is, in general, decreased by the action of the coarse graining map \[24]\.

Strict complete positivity can be re-obtained if we demand that the composite map \( \mathcal{N} = \Lambda_{CG} \circ \text{Id}_A \) to be CP-divisible \[29\]. If that is the case, the definition of CP-divisible maps requires \( \mathcal{N} = \Gamma(t_j,t_{j+1}) \circ \mathcal{N}_j \) for all \( t_k \leq t_j \), with \( \Gamma(t_j,t_{j+1}) \) the effective evolution map for the time interval \( [t_j,t_{j+1}] \), completely positive (see Fig. 7). This shows a connection between the theory of coarse grained maps and the theory of non-Markovian maps \[19\] \[29\].

V. CONSEQUENCE: EFFECTIVE DISTANCE INCREASE BY \( \Gamma_t \)

A common property of CPTP linear maps \[1\], is that the distance between two input states cannot increase. Mathematically, let \( \Lambda : \mathcal{L}(\mathcal{H}_B) \to \mathcal{L}(\mathcal{H}_B) \) be a CPTP linear map and \( \psi \) and \( \psi' \) be states in \( \mathcal{L}(\mathcal{H}_B) \). Then \( ||\Lambda(\psi) - \Lambda(\psi')||_1 \leq ||\psi - \psi'||_1 \), where the 1-norm is defined as \( ||A||_1 : = \text{Tr}(\sqrt{A^\dagger A}) \). Physically, this means, for instance, that the discrimination between two unknown quantum states cannot be improved by any further processing of the states \[9\].

The effective map \( \Gamma_t \), as discussed in the previous section, is not in general of Kraus form. Can then the distance between two effective states increase? As argued before, no contradiction with quantum mechanics can arise. In fact, it is simple to check that the distance between two effective states is upper-bounded, for all times, by the distance between the underlying initial states. Let \( \rho_0 = \Lambda_{CG}(\psi_0) \) and \( \rho'_0 = \Lambda_{CG}(\psi'_0) \) be effective states in \( \mathcal{L}(\mathcal{H}_B) \) with respective evolved states \( \rho_t = \Gamma_t(\rho_0) \) and \( \rho'_t = \Gamma_t(\rho'_0) \). Then

\[
||\rho_t - \rho'_t||_1 = ||\Lambda_{CG}(\psi_t) - \Lambda_{CG}(\psi'_t)||_1;
\]

\[
\leq ||\psi_t - \psi'_t||_1 \tag{10};
\]

\[
= ||\text{Id}_A(\psi_0) - \text{Id}_A(\psi'_0)||_1;
\]

\[
\leq ||\psi_0 - \psi'_0||_1. \tag{11}
\]

The last inequality turns into an equality in the case of a unitary mapping \( U_t \), i.e., \( U_t(\cdot) = U_t(\cdot)U_t^\dagger \) for some unitary \( U_t \).

This, however, does not imply that a distance increasing between effective states is not allowed. In fact, it is possible to have an increase in distance between the effective states undergoing the same effective map. Take for example the coarse graining describing the blurred-saturated detector \([11\]) the underlying dynamics given by the Hamiltonian \( H = hJ\sigma_x \otimes \sigma_x \[31\] \) and select two states \( \psi_0 \) and \( \psi'_0 \) which generate the same effective map \([14\]). Figure 8A shows the distance evolution between the two effective states \( \rho_t = \Lambda_{CG}(\psi_t) \) and \( \rho'_t = \Lambda_{CG}(\psi'_t) \). A clear oscillation of the distance is observed. In this example, nevertheless, we have that \( ||\rho_0 - \rho'_0||_1 \geq ||\rho_t - \rho'_t||_1 \) for all times.

Now, switch on a transversal field, turning the Hamiltonian into \( H = hJ\sigma_x \otimes \sigma_x + 3(\sigma_x \otimes 1 + 1 \otimes \sigma_x) \), and take initial states \( \psi_0 \) and \( \psi'_0 \) in a way to have the same effective map. The evolution of the distance between the effective states is shown in Fig. 8B. In this case we see that the distance \( ||\rho_t - \rho'_t||_1 \) can go beyond \( ||\rho_0 - \rho'_0||_1 \) for some specific times.
we investigated what types of dynamics may emerge from a full quantum description when one does not have access to, or is not interested in, all the degrees of freedom of a given system. The presented formalism generalizes the theory of open quantum systems, as it works also for closed systems. Here the split between system and environment is substituted by the split between accessible and non-accessible degrees of freedom. The possibility of correlations between these two types of degrees of freedom may generate effective dynamics that are not of Kraus form – without violating any principle of quantum mechanics. This, in turn, allowed for the distance between two effective states to increase under the action of a fixed effective map – in contrast to what is achievable in the underlying quantum description.

Other aspects of this effective dynamics can be further explored. Most notably, the fact that the "Kraus operators" may depend on the state the map is acting on. This suggests a possible way to explain how non-linear dynamics may emerge from the quantum linear description: if one looks at the system only at time intervals for which the term quantifying correlations between accessible and non-accessible degrees of freedom, \( \zeta(\alpha) \) in Eq. (9), vanishes, i.e., for a coarse grained time \([30]\), then the non-linearity of the first term may become apparent. This (possible) non-linearity, together with the distance increase between effective states undergoing the same effective map may be the key to explain how chaotic systems arise from the underlying quantum mechanical description.

Lastly, we hope that the formalism here presented can shed some light on the quantum-to-classical transition: the higher the "zoom out" (stronger coarse graining, in the sense of larger difference between \( D \) and \( d \)), the more simplified becomes the description of the system and its dynamics, with quantum features fading away. We believe that these ideas can be of interest for areas as quantum thermodynamics – which tries to explain the thermalization of closed quantum systems \([20, 21]\) –, and even to address the measurement problem in quantum mechanics \([31]\).

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