

Linear response theory and entropic fluctuations in repeated interaction quantum systems

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Abstract

We study linear response theory and entropic fluctuations of finite dimensional non-equilibrium Repeated Interaction Systems (RIS). More precisely, in a situation where the temperatures of the probes can take a finite number of different values, we prove analogs of the Green-Kubo fluctuation-dissipation formula and Onsager reciprocity relations on energy flux observables. Then we prove a Large Deviation Principle, or Fluctuation Theorem, and a Central Limit Theorem on the full counting statistics of entropy fluxes. We consider two types of non-equilibrium RIS: either the temperatures of the probes are deterministic and arrive in a cyclic way, or the temperatures of the probes are described by a sequence of i.i.d. random variables with uniform distribution over a finite set.

1 Introduction

In this paper we are interested in the linear response theory and entropic fluctuation for a particular class of open quantum systems called Repeated Interaction Systems (RIS), see Section 3 for a precise description. Our study fits in the wider framework of non-equilibrium quantum statistical mechanics. In this context, linear response theory and entropic fluctuation have attracted lot of attention in the last decades, see e.g. [LS, DS, JOP1, JOP2, JOP3, JOP4, JPP, dRM, dR, JOPP, BPR] and references therein.

Repeated interaction systems consist of a small system \mathcal{S} coupled to an environment made of a chain of independent probes \mathcal{E}^n with which \mathcal{S} will interact in a sequential way, i.e. \mathcal{S} interacts with \mathcal{E}^1 during the time interval $[0, \tau_1[$, then with \mathcal{E}^2 during the interval $[\tau_1, \tau_1 + \tau_2[$, etc. While \mathcal{S} interacts with a given probe \mathcal{E}^n the other ones evolve freely according to their intrinsic (uncoupled) dynamics. Formally, if $H_{\mathcal{S}}$ and $H_{\mathcal{E}^n}$ denote the non-interacting hamiltonians of \mathcal{S} and the \mathcal{E}^n 's and V_n denotes the coupling operator between \mathcal{S} and \mathcal{E}^n then the hamiltonian of the full system is the time-dependent, piecewise constant, operator

$$H(t) := H_{\mathcal{S}} + H_{\mathcal{E}^n} + V_n + \sum_{p \neq n} H_{\mathcal{E}^p}, \quad t \in [\tau_1 + \dots + \tau_{n-1}, \tau_1 + \dots + \tau_n[. \quad (1.1)$$

In the simplest case all the probes are identical (each \mathcal{E}^n is a copy of the same \mathcal{E} with the same initial state $\rho_{\mathcal{E}}$, e.g. a thermal state) and interact with \mathcal{S} by means of the same coupling operator V on $\mathcal{S} + \mathcal{E}$ and for the same duration τ . The dynamics restricted to the small system is shown to be determined by a map \mathcal{L} , see (2.2), which assigns $\rho(\tau) = \mathcal{L}(\rho)$ to ρ as the result of the interaction of \mathcal{S} with one subsystem \mathcal{E} for the duration τ . Heuristically, from the point of view of the small system, all subsystems interacting in sequence with \mathcal{S} are equivalent, so that the result of $n \in \mathbb{N}$ repeated interactions amounts to iterating n times the map \mathcal{L} on the initial state $\rho_{\mathcal{S}}$. This expresses the markovian character of repeated interactions in discrete time, see e.g. [BJM3] for an introduction to these RIS.

The typical physical situation of repeated interaction models is that of the one atom maser, see e.g. [FJM, MWM, RBH, WBKM]. Here \mathcal{S} is the quantized electromagnetic field in a cavity through which a beam of atoms, the \mathcal{E}^n 's, is shot in such a way that, at least with very high probability [HBR], no more than one atom is present in the cavity at any time. Such systems play a fundamental role in the experimental and theoretical investigations of basic matter-radiation processes. On the mathematical side various aspects of RIS, or quantum Markov chains, have been studied in the literature, see e.g. [KM, AP, AJ, BJM1, BJM2, Pe, PeP, NPe, BBB, Wo, vHG, GK, BFFS, CP, HJPR] and references therein. We mention also [BP, vHG2, Bru] for the analysis of a specific model related to the one-atom maser and [HMO] for related results in the framework of correlated quantum spin chains. We refer the reader to the lecture notes [Wo] and references therein for related results in the context of quantum information theory.

In order to consider a non-equilibrium situation we shall naturally consider here the more interesting situation where the probes are not always identical but allowed to vary over a finite set, especially via the temperatures of their initial states $\rho_{\mathcal{E}^n}$. The picture one should have in mind is that the system \mathcal{S} is coupled to finitely many reservoirs $\mathcal{R}_1, \dots, \mathcal{R}_M$ ($M \geq 2$) which are initially in thermal equilibrium at possibly different temperatures. Each probe is then associated to one of these M reservoirs, all the probes of a given reservoir being identical. After n interactions the state of \mathcal{S} is thus given by

$$\rho_n = \mathcal{L}_{j_n} \circ \dots \circ \mathcal{L}_{j_1}(\rho) \quad (1.2)$$

where $j_1, \dots, j_n \in \{1, \dots, M\}$ describes the *ordered* sequence of the probes.

Our first results deal with the linear response theory for energy fluxes out of the reservoirs. Linear response theory describes thermodynamics when the system under consideration is close to equilibrium. To first order the non-equilibrium energy fluxes depend linearly on the driving forces and we are interested in the corresponding response (or kinetic) coefficients, see (4.13). The main results in linear response theory are concerned with the Green-Kubo fluctuation-dissipation formula, Onsager reciprocity relations and a central limit theorem. Green-Kubo formula expresses the kinetic coefficients in terms of dynamical flux-flux correlations *at equilibrium*, Onsager relations expresses a symmetry in these coefficients and the central limit theorem relates the kinetic coefficients to fluctuations of the energy fluxes *at equilibrium*. We address all these questions in the framework of repeated interaction systems. Our main results are Theorem 4.4 and 5.9.

The second series of results deals with the fluctuation of entropy production and the so-called fluctuation relations. By fluctuation of entropy production we mean a large deviation principle for the statistics of entropy production, see Theorem 5.8. To do so we adopt the two-time measurement protocol introduced independently by Kurchan [Ku] and Tasaki [Ta].

A very similar large deviation principle as well as a central limit theorem similar to the one we have mentioned are proven in [vHG] for output statistics of quantum Markov chains. The only difference is that their statistics concern a single, post interaction, measurement while we consider a double, pre and post interaction, measurement. However from a mathematical perspective this causes only minor changes. Our main new result here concerns two symmetries satisfied by the large deviation rate function, or in equivalent way by the moment generating function. The first symmetry is what is usually called the fluctuation relation and can be understood as a generalization of the linear response results for systems far from equilibrium [Ga]. The second symmetry is a translation symmetry which is a refinement of energy conservation. To the best of our knowledge this translation symmetry goes back to [AGMT]. These two symmetries are fundamental results with deep consequences on the second and first law of thermodynamics respectively. We mention that in [vHG] no such fluctuation relations were proven.

In the framework of continuous-time quantum dynamical semigroups $\mathcal{L}_t := e^{tL}$, $L = L_1 + \dots + L_M$ and where L_j is the Lindblad generator describing the interaction with the j -th reservoir, those questions have been initiated in [LS] and then more recently studied in [DdrM, dRM, JPW]. Eq. (1.2) suggests that the situation should be very similar, if not identical, for RIS except that we have a discrete-time dynamics. As we shall see this is only partly correct. If the global strategy of the proofs largely follows those in [LS, JPW, vHG], RIS however have several specificities that have to be taken into account.

The first one can easily be seen from (1.2). As mentioned above there is a specific order in which \mathcal{S} interacts with the various probes. In order to make sure that \mathcal{S} interacts as much with each reservoir the first, maybe naive, idea is to make the order of interactions cyclic: \mathcal{S} interacts first with a probe associated to \mathcal{R}_1 , then to \mathcal{R}_2 and so on up to \mathcal{R}_M and then \mathcal{R}_1 again etc. In this case, if it is possible to derive a Green-Kubo type formula we will see that the usual Onsager reciprocity relations will fail, and similarly for the fluctuation relation of entropy production. The reason is simply that the cyclic order of interactions breaks time-reversal invariance even if we suppose that each interaction is time-reversal invariant. Due to the cyclic order of the interactions it is for example not surprising that a change of temperature in \mathcal{R}_1 will have a greater influence on the energy flux out of \mathcal{R}_2 than a change of temperature in \mathcal{R}_2 will have on the flux out of \mathcal{R}_1 . What one gets actually resembles the Onsager-Casimir relations one can find in the presence of a magnetic field. Namely, one has to compare the kinetic coefficients of the cyclic model with those of the *reversed cyclic* model in which the order of interactions is exactly the opposite, see (4.16) and (5.10). The study of this cyclic model provides a simple example which shows how much time-reversal invariance is fundamental, in particulier in the derivation of the usual fluctuation relations.

To remedy this lack of global time-reversal invariance we shall therefore also consider the situation where the probes associated to the various reservoirs interact with \mathcal{S} in a random order. To make it simple we shall consider here the case where the probes are chosen independently at each time and with a uniform distribution (so that on average \mathcal{S} interacts as much with each reservoir). Our results can be easily generalized to the case of an arbitrary i.i.d. distribution. The case where the distribution of the probes is given by a more general Markov process will be considered in [BJP].

The second specificity is related to the fact that the RIS hamiltonian (1.1) is time-dependent. As a consequence, even in the ideal case where all the probes are identical, this may lead to a non-vanishing of entropy production which is usually considered as the

signature of a non-equilibrium situation. In this paper we are interested in the response of the system to the presence of thermal forces. This forces us to impose some extra assumption called Assumption **(NE)**, see Section 3.4, which guarantees that the case where all the temperatures of the various reservoirs are equal is indeed an equilibrium situation. A general linear response theory for RIS should also take into account a departure from this assumption. It is however not clear at the moment how to quantify this or, said differently, what is the natural quantity one can associate to a generic RIS and the vanishing of which would correspond to the fulfilment of Assumption **(NE)**. We therefore postpone this question to future work. Regarding this point we finally mention that the question of linear response theory for time-dependent quantum hamiltonians have been considered in [DS], in a weak-coupling regime, but there only the perturbation was time-dependent contrary to what happens in RIS.

Finally, to avoid technical issues we shall stick here to the case where all the subsystem's Hilbert spaces, for \mathcal{S} and the \mathcal{E}^n 's, are finite dimensional. Most results can easily be extended to infinite dimension provided the various assumptions are adapted in an *ad hoc* way, in particular the ergodic Assumption **(ER \sharp)** of Section 3.3.2 has to require the existence of a spectral gap. It is however difficult to find physically relevant models to which these assumptions apply. The model for the one-atom maser studied in [BP, Bru] does not have a spectral gap for example. Nevertheless it is still possible to prove the Green-Kubo formula and Onsager relations for this model, see [Bo]. We also mention [BDBP] which considers a RIS type model for the motion of a tight-binding electron and where an analog of the fluctuation relation is proven for the position increments of the electron.

The paper is organized as follows. In Section 2 we briefly recall some basic concepts of open systems. In Section 3 we describe the non-equilibrium RIS model and state the various assumptions which we will use. We will in particular describe in more detail the above mentioned Assumption **(NE)** and discuss its origin and some of its consequences. Section 4 is devoted to the linear response theory and in particular the derivation of Green-Kubo formula and of Onsager reciprocity relations. These are stated in Theorem 4.4. In Section 5 we consider the fluctuation of entropy production and prove a fluctuation relation for RIS. We also complement linear response theory with a Central Limit Theorem. Our main results in this section are Theorems 5.4, 5.5, 5.8 and 5.9. Finally the proofs are given in Section 6.

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2 Preliminaries

2.1 Observables, states and their evolution

All the Hilbert spaces considered in the paper are finite dimensional. We denote by $\mathbb{1}$ the identity operator (we shall sometimes indicate $\mathbb{1}_{\mathcal{H}}$ to specify the underlying Hilbert space \mathcal{H} if any confusion is possible). Observables on \mathcal{H} are described by self-adjoint elements in $\mathcal{B}(\mathcal{H})$.

States are given by density matrices, i.e. positive elements with unit trace, and we shall use the same symbol for a density matrix ρ and for the associated positive unital linear functional $A \mapsto \text{Tr}(\rho A)$, i.e. $\rho(A)$ will stand for $\text{Tr}(\rho A)$ the expectation value of the observable A in the state ρ . A state ρ is said to be faithful if it is positive definite.

A bounded linear map Φ acting on $\mathcal{B}(\mathcal{H})$ is called positive (or positivity preserving) if $\Phi(A)$ is non-negative for any non-negative $A \in \mathcal{B}(\mathcal{H})$. Φ is said to be completely positive if, for any $d > 0$, $\Phi \otimes \mathbb{1}_{\mathcal{B}(\mathbb{C}^d)}$ is a positive map. It is called unital if $\Phi(\mathbb{1}) = \mathbb{1}$ and trace-preserving if $\text{Tr}(\Phi(\rho)) = \text{Tr}(\rho)$ for any ρ . Note that Φ is trace-preserving iff its dual map Φ^* is unital, and where Φ^* is such that $\text{Tr}(\Phi(\rho)A) = \text{Tr}(\rho\Phi^*(A))$ for any ρ and A .

In the markovian description of open quantum systems the evolution of a state is described by a completely positive and trace-preserving (CPTP) map (Schrödinger picture) while the one of an observable is described by a completely positive and unital map (Heisenberg picture), see Section 2.2. If \mathcal{L} is a CPTP map then $(\mathcal{L}^n)_{n \in \mathbb{N}}$ is called a quantum dynamical semigroup. Obviously the large time ($n \rightarrow \infty$) limit of quantum dynamical semigroups is closely related to the spectral properties of the map \mathcal{L} . It is known that positive maps satisfy Perron-Frobenius type results, see e.g. [EHK, Wo]. In particular their spectral radius is always an eigenvalue and there is a non-negative corresponding eigenvector. If moreover \mathcal{L} is CPTP then its spectral radius is 1 so that \mathcal{L} admits an invariant state. We recall that any completely positive map can be written in the form

$$\mathcal{L}(\cdot) = \sum_{i=1}^m V_i \cdot V_i^*, \quad (2.1)$$

where the V_i 's are in $\mathcal{B}(\mathcal{H})$ and $m \in \mathbb{N}$. Such a form is called a Kraus representation of \mathcal{L} [Kr], it is however not unique. The following notion of primitive CPTP map will play an important role in the paper.

Definition 2.1. *Let \mathcal{L} be a completely positive map on $\mathcal{B}(\mathcal{H})$ given by (2.1). Then \mathcal{L} is primitive if there exists $n \in \mathbb{N}$ such that $\text{Span}\{V_{i_1} \cdots V_{i_n} \mid i_1, \dots, i_n \in I\} = \mathcal{B}(\mathcal{H})$.*

Remark 2.1. There are actually several equivalent definitions of primitive positive maps. The one given here is the simplest for our purpose. We refer the reader to e.g. [Wo] for a more detailed discussion on the subject.

The importance of the notion of primitive map for our purpose is due to the following Proposition, see e.g. [Wo],

Proposition 2.1. *Let \mathcal{L} be a completely positive map and denote by r its spectral radius. \mathcal{L} is primitive iff its spectral radius is a simple dominant eigenvalue, i.e. all other eigenvalues λ satisfy $|\lambda| < r$, with positive definite left and right eigenvectors.*

When \mathcal{L} is CPTP its spectral radius is 1 and its dual map \mathcal{L}^* is unital, so that $\mathbb{1}$ is a left eigenvector. Hence the above proposition can simply be rephrased as “a CPTP map \mathcal{L} is primitive if and only if 1 is a simple dominant eigenvalue and \mathcal{L} admits a (unique) faithful invariant state”. Using spectral decomposition the notion of primitive CPTP map is thus immediately related to strong ergodic properties of the corresponding quantum dynamical semigroup.

Proposition 2.2. *Let $(\mathcal{L}^n)_{n \in \mathbb{N}}$ be a quantum dynamical semigroup. \mathcal{L} is primitive if and only if for any state ρ one has $\lim_{n \rightarrow +\infty} \mathcal{L}^n(\rho) = \rho_+$, where ρ_+ is the unique faithful invariant state of \mathcal{L} . In other words ρ_+ is mixing for the semigroup generated by \mathcal{L} .*

A slightly stronger notion is sometimes useful. A map \mathcal{L} is called positivity improving if for any non-negative and non-zero A its image $\mathcal{L}(A)$ is positive definite. It is easy to see that if a CP map is positivity improving then it is primitive.

Finally, if $\mathcal{H}_1, \mathcal{H}_2$ are two Hilbert spaces and ρ is a density matrix on $\mathcal{H}_1 \otimes \mathcal{H}_2$, we denote by $\rho_1 := \text{Tr}_{\mathcal{H}_2}(\rho)$ the partial trace of ρ with respect to \mathcal{H}_2 . It is the unique density matrix on \mathcal{H}_1 such that for all $A \in \mathcal{B}(\mathcal{H}_1)$ one has $\text{Tr}(\rho \times A \otimes \mathbb{1}_{\mathcal{H}_2}) = \text{Tr}(\rho_1 A)$. Similarly given a density matrix ρ_2 on \mathcal{H}_2 and an observable $A \in \mathcal{B}(\mathcal{H}_1 \otimes \mathcal{H}_2)$. Then $A_1 := \text{Tr}_{\mathcal{H}_2}(\mathbb{1}_{\mathcal{H}_1} \otimes \rho_2 \times A)$ is the unique element in $\mathcal{B}(\mathcal{H}_1)$ such that for any state ρ_1 on \mathcal{H}_1 one has $\text{Tr}(\rho_1 \otimes \rho_2 \times A) = \text{Tr}(\rho_1 A_1)$. A_1 is called the partial trace of A w.r.t. the state ρ_2 , and will be denoted $A_1 = \text{Tr}_{\rho_2}(A)$.

2.2 Open quantum systems

A quantum system \mathcal{S} is said to be open when it interacts with another quantum system \mathcal{E} . \mathcal{S} is sometimes called the small system, \mathcal{E} the environment (which can further be made of several components), and $\mathcal{S} + \mathcal{E}$ is the global system. If $\mathcal{H}_{\mathcal{S}}$ and $\mathcal{H}_{\mathcal{E}}$ are the Hilbert spaces describing \mathcal{S} and \mathcal{E} respectively then the Hilbert space of the global system is $\mathcal{H}_{\mathcal{S}+\mathcal{E}} = \mathcal{H}_{\mathcal{S}} \otimes \mathcal{H}_{\mathcal{E}}$, and if the joint system $\mathcal{S} + \mathcal{E}$ is in the state $\rho_{\mathcal{S}+\mathcal{E}}$ then \mathcal{S} is in the reduced state $\rho_{\mathcal{S}} = \text{Tr}_{\mathcal{H}_{\mathcal{E}}}(\rho_{\mathcal{S}+\mathcal{E}})$. Similarly, if the system is initially decoupled, i.e. $\mathcal{S} + \mathcal{E}$ is in a state $\rho_{\mathcal{S}+\mathcal{E}} = \rho_{\mathcal{S}} \otimes \rho_{\mathcal{E}}$ where $\rho_{\mathcal{S}}$ and $\rho_{\mathcal{E}}$ are the states of \mathcal{S} and \mathcal{E} respectively, then for an observable $A \in \mathcal{B}(\mathcal{H}_{\mathcal{S}} \otimes \mathcal{H}_{\mathcal{E}})$ of the global system, the observable “seen by \mathcal{S} ” is $\text{Tr}_{\rho_{\mathcal{E}}}(A) = \text{Tr}_{\mathcal{H}_{\mathcal{E}}}(\mathbb{1} \otimes \rho_{\mathcal{E}} \times A)$.

The non-interacting dynamics of \mathcal{S} and \mathcal{E} are described by hamiltonians $H_{\mathcal{S}}$ and $H_{\mathcal{E}}$ respectively and the interaction between \mathcal{S} and \mathcal{E} by some interaction operator V acting on $\mathcal{H}_{\mathcal{S}} \otimes \mathcal{H}_{\mathcal{E}}$. Then $H := H_{\mathcal{S}} \otimes \mathbb{1} + \mathbb{1} \otimes H_{\mathcal{E}} + V$ is the hamiltonian for the joint evolution. In the sequel we will often omit the inessential factors $\mathbb{1}$ in the tensor products.

Fix now some initial (or reference) state $\rho_{\mathcal{E}}$ of the environment and suppose the system is in some initially decoupled state $\rho \otimes \rho_{\mathcal{E}}$. Then the state of the global system $\mathcal{S} + \mathcal{E}$ after some time $\tau > 0$ is $U \times \rho \otimes \rho_{\mathcal{E}} \times U^*$ where $U := e^{-i\tau H}$. Hence the small system \mathcal{S} is in the state

$$\mathcal{L}(\rho) = \text{Tr}_{\mathcal{H}_{\mathcal{E}}}(U \times \rho \otimes \rho_{\mathcal{E}} \times U^*). \quad (2.2)$$

One easily checks that \mathcal{L} defines a CPTP map. It is called the reduced dynamics map of \mathcal{S} associated to the open quantum system $\mathcal{S} + \mathcal{E}$ and for the duration τ . Its dual map describes the evolution of observables $A \in \mathcal{B}(\mathcal{H}_{\mathcal{S}})$ and is given by

$$\mathcal{L}^*(A) = \text{Tr}_{\rho_{\mathcal{E}}}(U^* \times A \otimes \mathbb{1} \times U).$$

2.3 Entropy production in an open system

The entropy observable of a quantum system in the faithful state ρ is defined as $S(\rho) := -\log \rho$ where \log denotes the natural logarithm. The von Neumann entropy of ρ is then the expectation value of $S(\rho)$ in the state ρ , i.e. $\text{Ent}(\rho) := -\text{Tr}(\rho \log \rho) = \rho(S(\rho))$. The relative entropy of ρ relatively to the faithful state ν is then

$$\text{Ent}(\rho | \nu) := \text{Tr}(\rho(\log \rho - \log \nu)) = \rho(S(\nu) - S(\rho)).$$

It is well-known that $\text{Ent}(\rho, \nu) \geq 0$ and $\text{Ent}(\rho, \nu) = 0$ if and only if $\rho = \nu$.

Consider an open system $\mathcal{S} + \mathcal{E}$ as in Section 2.2, interacting for a duration $\tau > 0$. Denote by U the unitary evolution of the joint system and \mathcal{L} the reduced dynamics on \mathcal{S} as

given by (2.2). Assume moreover that the system \mathcal{E} is initially in thermal equilibrium, i.e. $\rho_{\mathcal{E}} := \frac{e^{-\beta H_{\mathcal{E}}}}{\text{Tr}(e^{-\beta H_{\mathcal{E}}})}$ is a Gibbs state at some inverse temperature β . Then a simple calculation shows that the following entropy balance equation holds, see also [HJPR],

$$\text{Ent}(\mathcal{L}(\rho)) - \text{Ent}(\rho) = \text{Ent}(U \times \rho \otimes \rho_{\mathcal{E}} \times U^* | \mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}) - \beta \text{Tr}((\rho \otimes \rho_{\mathcal{E}})(U^* H_{\mathcal{E}} U - H_{\mathcal{E}})). \quad (2.3)$$

The second term of the right-hand side can be interpreted as the entropy flux coming from \mathcal{E} (it is β times the energy variation in \mathcal{E}). This motivates the following

Definition 2.2. We define the entropy production of \mathcal{S} during its interaction with \mathcal{E} by

$$\sigma := \text{Ent}(U \times \rho \otimes \rho_{\mathcal{E}} \times U^* | \mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}). \quad (2.4)$$

The entropy production is meant to be always nonnegative and to be zero iff there is no energy flux between \mathcal{S} and \mathcal{E} . As a matter of fact one easily infers that

$$\sigma = 0 \iff \text{Tr}(\rho \otimes \rho_{\mathcal{E}}(U^* H_{\mathcal{E}} U - H_{\mathcal{E}})) = \text{Ent}(\mathcal{L}(\rho)) - \text{Ent}(\rho) = 0.$$

Indeed, see also Section 2 in [JP], $\sigma = 0$ iff $U \times \rho \otimes \rho_{\mathcal{E}} \times U^* = \mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}$ so that if $\sigma = 0$ then on one hand

$$\text{Tr}(\rho \otimes \rho_{\mathcal{E}}(U^* H_{\mathcal{E}} U - H_{\mathcal{E}})) = \text{Tr}(\mathcal{L}(\rho) \otimes \rho_{\mathcal{E}} H_{\mathcal{E}} - \rho \otimes \rho_{\mathcal{E}} H_{\mathcal{E}}) = \text{Tr}(\mathcal{L}(\rho) - \rho) \times \text{Tr}(\rho_{\mathcal{E}} H_{\mathcal{E}}) = 0,$$

because \mathcal{L} is trace preserving, and on the other hand

$$\text{Ent}(\rho) + \text{Ent}(\rho_{\mathcal{E}}) = \text{Ent}(\rho \otimes \rho_{\mathcal{E}}) = \text{Ent}(U \times \rho \otimes \rho_{\mathcal{E}} \times U^*) = \text{Ent}(\mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}) = \text{Ent}(\mathcal{L}(\rho)) + \text{Ent}(\rho_{\mathcal{E}}).$$

The other implication follows directly from (2.3).

The fact that $\sigma = 0$ iff $U \times \rho \otimes \rho_{\mathcal{E}} \times U^* = \mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}$ can also be rephrased as follows: the entropy production vanishes iff the joint system $\mathcal{S} + \mathcal{E}$ is left non-entangled by the dynamics with moreover the state of \mathcal{E} being invariant. This leads to the following definition.

Definition 2.3. Let U be a unitary operator on $\mathcal{H}_{\mathcal{S}} \otimes \mathcal{H}_{\mathcal{E}}$, ρ a state on $\mathcal{H}_{\mathcal{S}}$ and $\rho_{\mathcal{E}}$ a state on $\mathcal{H}_{\mathcal{E}}$. We say that the triple $(U, \rho, \rho_{\mathcal{E}})$ satisfies the Non-Entanglement condition if

$$U \times \rho \otimes \rho_{\mathcal{E}} \times U^* = \mathcal{L}(\rho) \otimes \rho_{\mathcal{E}}, \quad (2.5)$$

in other words if the entropy production of the associated open system vanishes.

The importance of this Non-Entanglement condition for RIS will be made more transparent in Section 3.4. We mention that it also appears in [HJPR] in the context of adiabatic RIS. At this point we simply note that if ρ is an invariant state of \mathcal{L} then the triple $(U, \rho, \rho_{\mathcal{E}})$ satisfies the Non-Entanglement condition if and only if $\rho \otimes \rho_{\mathcal{E}}$ is an invariant state of the interacting dynamics, i.e.

$$U \times \rho \otimes \rho_{\mathcal{E}} \times U^* = \rho \otimes \rho_{\mathcal{E}}. \quad (2.6)$$

As a consequence the entropy observable

$$S(\rho \otimes \rho_{\mathcal{E}}) = -\log(\rho) \otimes \mathbb{1} - \mathbb{1} \otimes \log(\rho_{\mathcal{E}}) \quad (2.7)$$

is a constant of motion.

3 Non-equilibrium Repeated Interaction Systems

3.1 Repeated Interaction Systems (RIS)

Repeated Interaction Systems form a specific class of open quantum systems in which the environment \mathcal{E} has the following structure $\mathcal{E} = \mathcal{E}^1 + \mathcal{E}^2 + \dots + \mathcal{E}^n + \dots$ where $(\mathcal{E}^n)_n$ is a sequence of quantum subsystems with associated Hilbert spaces $\mathcal{H}_{\mathcal{E}^n}$ and free hamiltonians $H_{\mathcal{E}^n}$. \mathcal{S} will be called the *small system* and the \mathcal{E}^n 's will be called the *probes*. The Repeated Interaction dynamics consists in the joint evolution of \mathcal{S} and \mathcal{E}^1 for a duration τ_1 , immediately followed by the joint evolution of \mathcal{S} and \mathcal{E}^2 for a duration τ_2 , etc. For any $n, p \in \mathbb{N}^*$, \mathcal{E}^n and \mathcal{E}^p are disjoint and never interact directly. The interaction between \mathcal{S} and \mathcal{E}^n is described by the interaction operator V_n . Hence during the n -th interaction the coupled hamiltonian is

$$H_n := H_{\mathcal{S}} \otimes \mathbb{1} + \mathbb{1} \otimes H_{\mathcal{E}^n} + V_n,$$

and the unitary propagator for the coupled dynamics describing the N first interactions is thus given by $U_N U_{N-1} \dots U_1$ where $U_n := e^{-i\tau_n H_n}$, $n = 1, \dots, N$. Note that by construction the various $H_{\mathcal{E}^n}$ commute and that $[H_{\mathcal{E}^n}, H_p] = 0$ whenever $n \neq p$.

If \mathcal{S} is initially in the state ρ and the n -th probe is in the state $\rho_{\mathcal{E}^n}$ which we assume to be invariant for the free dynamics of \mathcal{E}^n , then the state of \mathcal{S} after N interactions is given by

$$\rho_N = \text{Tr}_{\mathcal{H}_{\mathcal{E}^1} \otimes \dots \otimes \mathcal{H}_{\mathcal{E}^N}} (U_N \dots U_1 \times \rho \otimes \rho_{\mathcal{E}^1} \otimes \dots \otimes \rho_{\mathcal{E}^N} \times U_1^* \dots U_N^*). \quad (3.1)$$

It is easy to see that the Repeated Interaction structure induces the following markovian behaviour, see e.g. [BJM3] for more details,

$$\rho_N = \text{Tr}_{\mathcal{H}_{\mathcal{E}^N}} (U_N \times \rho_{N-1} \otimes \rho_{\mathcal{E}^N} \times U_N^*) =: \mathcal{L}_N(\rho_{N-1}),$$

so that if, for any n , the map

$$\mathcal{L}_n(\rho) := \text{Tr}_{\mathcal{H}_{\mathcal{E}^n}} (U_n \times \rho \otimes \rho_{\mathcal{E}^n} \times U_n^*),$$

denotes the reduced dynamics map associated to the interaction of \mathcal{S} with the probe \mathcal{E}^n , we have

$$\rho_N = \mathcal{L}_N \circ \dots \circ \mathcal{L}_1(\rho). \quad (3.2)$$

In the simplest situation the probes are copies of an identical system \mathcal{E} , i.e. $\mathcal{H}_{\mathcal{E}^n} \equiv \mathcal{H}_{\mathcal{E}}$, $H_{\mathcal{E}^n} \equiv H_{\mathcal{E}}$, $V_n \equiv V$, $\tau_n \equiv \tau$, $\rho_{\mathcal{E}^n} \equiv \rho_{\mathcal{E}}$. Then the reduced dynamics maps coincide, $\mathcal{L}_n \equiv \mathcal{L}$, and the evolution of the system \mathcal{S} is given by $\rho_N = \mathcal{L}^N(\rho)$. We shall refer to this situation as an equilibrium situation, the various probes being considered as elements of a single reservoir. One may think of the initial states of the probes as Gibbs states at some common inverse temperature β .

3.2 Non-equilibrium RIS

In this paper we are interested in understanding RIS in a non-equilibrium situation. The picture one should have in mind is that the system \mathcal{S} is coupled to several reservoirs $\mathcal{R}_1, \dots, \mathcal{R}_M$ ($M \geq 2$) which are initially in thermal equilibrium but at possibly different temperatures. The various probes are then associated to one of these M reservoirs.

More precisely, we fix a finite set of quantum systems \mathcal{E}_j , $j = 1, \dots, M$, with the associated Hilbert spaces $\mathcal{H}_{\mathcal{E}_j}$, free hamiltonians $H_{\mathcal{E}_j}$, interaction operators V_j , interaction times τ_j and initial states $\rho_{\mathcal{E}_j}$. Then H_j , U_j and \mathcal{L}_j denote respectively the interacting hamiltonian, acting on $\mathcal{H}_{\mathcal{S}} \otimes \mathcal{H}_{\mathcal{E}_j}$, unitary propagator and reduced dynamics map as given by (2.2). Each probe will be a copy of one of these M systems.

Remark 3.1. To have a simple picture in mind the reader should think of all data identical, i.e. $\mathcal{H}_{\mathcal{E}_j} \equiv \mathcal{H}_{\mathcal{E}}$, $H_{\mathcal{E}_j} \equiv H_{\mathcal{E}}$, $V_j \equiv V$, $\tau_j \equiv \tau$, except for the initial states $\rho_{\mathcal{E}_j}$ which are Gibbs states at possibly different temperatures β_j^{-1} .

The sequence of probes is then described by a sequence $\mathbf{j} := (j_n)_{n \in \mathbb{N}^*} \in \{1, \dots, M\}^{\mathbb{N}^*}$ where j_n will specify which of the M systems \mathcal{E}_j 's the probe \mathcal{E}^n is a copy of, i.e. $\forall n \in \mathbb{N}^*$, $\mathcal{E}^n \equiv \mathcal{E}_{j_n}$. Thus \mathcal{S} interacts with M reservoirs $\mathcal{R}_1, \dots, \mathcal{R}_M$ where \mathcal{R}_j denotes the ‘‘union’’ of all the probes corresponding to the index j , i.e.

$$\mathcal{R}_j := \bigoplus_{n, j_n=j} \mathcal{E}^n.$$

If ρ is the initial state of \mathcal{S} then after n interactions \mathcal{S} is in the state

$$\rho_n(\mathbf{j}) := \mathcal{L}_{j_n} \circ \mathcal{L}_{j_{n-1}} \circ \dots \circ \mathcal{L}_{j_1}(\rho).$$

Accordingly, for an observable $A \in \mathcal{B}(\mathcal{H}_{\mathcal{S}})$, the corresponding Heisenberg evolution is

$$A_n(\mathbf{j}) := \mathcal{L}_{j_1}^* \circ \mathcal{L}_{j_2}^* \circ \dots \circ \mathcal{L}_{j_n}^*(A).$$

In this paper, we shall consider two specific and rather natural situation of non-equilibrium RIS whose reduced dynamics can be linked to a discrete quantum dynamical semigroup.

3.2.1 Cyclic case

A non-equilibrium RIS is called cyclic when \mathcal{S} interacts first with \mathcal{R}_1 , then $\mathcal{R}_2, \dots, \mathcal{R}_M$, then \mathcal{R}_1 again, etc. In other words the sequence \mathbf{j} describing the interactions is the M -periodic sequence $\mathbf{j}^{cy} = (j_n^{cy})_n$ where for any $k \in \{1, \dots, M\}$ and $n \in \mathbb{N}$ one has $j_{k+nM}^{cy} = k$. Consequently, the reduced dynamics of \mathcal{S} over the cycles can be described by the discrete semigroup $(\mathcal{L}_{cy}^n)_n$ where

$$\mathcal{L}_{cy} := \mathcal{L}_M \circ \mathcal{L}_{M-1} \circ \dots \circ \mathcal{L}_1. \quad (3.3)$$

Note that $\mathcal{L}_{cy}(\rho) = \text{Tr}_{\mathcal{H}_{\mathcal{E}_{cy}}}(U_{cy} \times \rho \otimes \rho_{cy} \times U_{cy}^*)$, where $\mathcal{E}_{cy} := \mathcal{E}_1 + \dots + \mathcal{E}_M$, $\mathcal{H}_{\mathcal{E}_{cy}} := \mathcal{H}_{\mathcal{E}_1} \otimes \dots \otimes \mathcal{H}_{\mathcal{E}_M}$, $U_{cy} := U_M \dots U_1$ and $\rho_{cy} := \rho_{\mathcal{E}_1} \otimes \dots \otimes \rho_{\mathcal{E}_M}$.

3.2.2 Random case

A non-equilibrium RIS is called random when the order in which the small system \mathcal{S} interacts with the $(\mathcal{R}_j)_{1 \leq j \leq M}$ is described by a random process, i.e. for each interaction the subsystem \mathcal{E}_j of which \mathcal{E}^n is a copy will be chosen randomly from $\mathcal{E}_1, \dots, \mathcal{E}_M$. The motivation to consider random RIS is related to the question of time-reversal invariance, see Remark 3.8 in Section 3.4.3. Let p denote the uniform probability measure on $\{1, \dots, M\}$ and \mathcal{P} the standard convoluted probability measure on $\{1, \dots, M\}^{\mathbb{N}^*}$ associated to p . For any $\mathcal{T} \subset \{1, \dots, M\}^{\mathbb{N}^*}$, $\mathcal{P}(\mathcal{T})$ is the probability for the random RIS to interact successively with

the reservoirs $\mathcal{R}_{j_1}, \mathcal{R}_{j_2}, \dots, \mathcal{R}_{j_n}, \dots$ with $\mathbf{j} = (j_n)_n \in \mathcal{T}$. We denote by $\mathbb{E}_{\mathcal{P}}$ the expectation value w.r.t. a measure \mathbb{P} . The choice of the measure \mathcal{P} reflects the fact that the order in which the reservoirs $\mathcal{R}_1, \dots, \mathcal{R}_M$ will interact with \mathcal{S} is chosen in an i.i.d. manner but on average \mathcal{S} will interact equally much with each of them. This is the simplest and only case we shall consider here. A generalization of our results to a Markovian situation will be considered in the forthcoming paper [BJP].

Then one easily gets that $\mathbb{E}_{\mathcal{P}}(\mathcal{L}_{j_n} \circ \mathcal{L}_{j_{n-1}} \circ \dots \circ \mathcal{L}_{j_1}) = \mathcal{L}_{ra}^n$ where

$$\mathcal{L}_{ra} := \mathbb{E}_{\mathcal{P}}(\mathcal{L}_j) = \frac{1}{M} (\mathcal{L}_1 + \dots + \mathcal{L}_M).$$

Thus, at least in expectation, the reduced dynamics of \mathcal{S} can be described by the semigroup $(\mathcal{L}_{ra}^n)_{n \in \mathbb{N}}$. We say at least because of the following theorem which was proven in [BJM2] (it is a particular case of Theorem 1.3 in that paper)

Theorem 3.1. *Suppose that there exists $1 \leq j \leq M$ such that \mathcal{L}_j is primitive. Then \mathcal{L}_{ra} is primitive. Moreover the random RIS dynamics converges almost surely and in the ergodic mean to the unique invariant state of \mathcal{L}_{ra} . More precisely, there is a subset $\mathcal{T} \subset \{1, \dots, M\}^{\mathbb{N}^*}$ such that $\mathcal{P}(\mathcal{T}) = 1$ and for any $\mathbf{j} \in \mathcal{T}$, any density matrix ρ , and any family of observables $\{1, \dots, M\} \ni j \mapsto A(j) \in \mathcal{B}(\mathcal{H}_{\mathcal{S}})$ one has*

$$\lim_{N \rightarrow +\infty} \frac{1}{N} \sum_{n=1}^N \text{Tr} (\mathcal{L}_{j_{n-1}} \circ \dots \circ \mathcal{L}_{j_1}(\rho) \times A(j_n)) = \text{Tr} (\rho_+^{ra} \times \mathbb{E}_{\mathcal{P}}(A)), \quad (3.4)$$

where ρ_+^{ra} denotes the unique invariant state of \mathcal{L}_{ra} and $\mathbb{E}_{\mathcal{P}}(A) = \frac{1}{M}(A(1) + \dots + A(M))$.

Remark 3.2. The observables $A(j)$ should be understood in the following sense. They represent the same physical quantity but their expression might vary depending on which probe \mathcal{S} is interacting with. They are called instantaneous observables in [BJM2]. A typical example is that of energy flux observables, see Section 4.1.

If in particular $A(j) \equiv A$, i.e. one considers a fixed observable, then (3.4) indeed traduces the ergodicity of the random RIS.

Notation. Throughout the paper when assumptions, results or identities are formulated for both the cyclic and random situation we will often use the symbol \sharp which will stand for either *cy* in the cyclic case or *ra* in the random case.

3.3 Assumptions

In this section we formulate the various assumptions that will be used in the paper.

3.3.1 Temperatures and thermal forces

The first assumption concerns the initial states of the probes. We will assume that they are initially in thermal equilibrium. Namely,

Assumption (KMS). *For any $1 \leq j \leq M$, $\rho_{\mathcal{E}_j} = \frac{e^{-\beta_j H_{\mathcal{E}_j}}}{\text{Tr} (e^{-\beta_j H_{\mathcal{E}_j}})}$ for some β_j .*

From now on, we will always suppose Assumption **(KMS)** holds.

Fix some $\beta_{\text{ref}} > 0$. If $\zeta_j := \beta_{\text{ref}} - \beta_j$ then $\zeta = (\zeta_1, \zeta_2, \dots, \zeta_M)$ denotes the vector of thermal forces. In the sequel, for any quantity A depending on ζ , we will write A_ζ when we need to stress its dependence on ζ , and only A otherwise. When all the temperatures are equal (but not necessarily to β_{ref}^{-1}), i.e. $\zeta = (\zeta, \dots, \zeta)$, we will write A_ζ instead of A_ζ .

Note that \mathcal{E}_j depends only on ζ_j . Hence we may write $\mathcal{L}_{j,\zeta} = \mathcal{L}_{j,\zeta_j}$, $\rho_{\mathcal{E}_j,\zeta} = \rho_{\mathcal{E}_j,\zeta_j,\dots}$

3.3.2 Ergodicity

The next assumption concerns spectral properties of the reduced dynamics maps $\mathcal{L}_{\sharp,\zeta}$, $\sharp = cy$ or ra , and the large time behaviour of \mathcal{S} (recall Proposition 2.2).

Assumption (\mathbf{ER}_\sharp) . *There exists $\zeta \in \mathbb{R}^M$ such that $\mathcal{L}_{\sharp,\zeta}$ is primitive.*

Proposition 3.2. *Suppose assumption (\mathbf{ER}_\sharp) holds. Then*

1. *For any $\zeta \in \mathbb{R}^M$, $\mathcal{L}_{\sharp,\zeta}$ is primitive. As a consequence, for any $\zeta \in \mathbb{R}^M$ and any ρ (not necessarily a state), $\lim_{n \rightarrow \infty} \mathcal{L}_{\sharp,\zeta}^n(\rho) = \text{Tr}(\rho)\rho_{+,\zeta}^\sharp$ where $\rho_{+,\zeta}^\sharp$ denotes the unique invariant faithful state of $\mathcal{L}_{\sharp,\zeta}$.*
2. *$\rho_{+,\zeta}^\sharp$ is infinitely differentiable w.r.t. $\zeta \in \mathbb{R}^M$.*

Proof. 1. Let λ_{jk} , φ_{jk} denote the eigenvalues and eigenvectors of $H_{\mathcal{E}_j}$ so that $H_{\mathcal{E}_j} = \sum_k \lambda_{jk} |\varphi_{jk}\rangle\langle\varphi_{jk}|$.

Then it follows from (2.2) that

$$\mathcal{L}_{j,\zeta_j}(\rho) = \sum_{k,k'} \frac{e^{-\lambda_{jk}(\beta_{\text{ref}} - \zeta_j)}}{Z_{\zeta_j}} V_{jkk'} \rho V_{jkk'}^*,$$

where $V_{jkk'} = \langle\varphi_{jk'}|e^{-i\tau_j H_j}|\varphi_{jk}\rangle$ does not depend on ζ . It follows immediately that the maps $\mathcal{L}_{\sharp,\zeta}$ have Kraus decomposition of the form $\mathcal{L}_{\sharp,\zeta} = \sum_i f_i(\zeta) V_i \rho V_i^*$ where the V_i do not depend on ζ and with $f_i(\zeta) > 0$. For any ζ, ζ' the maps $\mathcal{L}_{\sharp,\zeta}$ and $\mathcal{L}_{\sharp,\zeta'}$ therefore have the same Kraus decomposition up to positive scalar factors which proves 1. (recall Definition 2.1).

2. It follows from 1. and Proposition 2.1 that 1 is an isolated simple eigenvalue for any ζ so the result follows by regular perturbation theory [Ka]. \square

Finally, we have the following result which shows that it suffices to have information on one of the probes to get information on the entire non-equilibrium RIS.

Proposition 3.3. *Let $\zeta \in \mathbb{R}^M$. If \mathcal{L}_{j,ζ_j} is primitive for some $1 \leq j \leq M$ then (\mathbf{ER}_\sharp) holds for $\sharp = ra$, and if moreover \mathcal{L}_{j,ζ_j} is positivity improving then (\mathbf{ER}_\sharp) holds for $\sharp = cy$.*

The proof follows directly from Lemma 3.4 below in the cyclic case and from Theorem 3.1 in the random case.

Lemma 3.4. *If ρ is positive definite then so is $\mathcal{L}_j(\rho)$. As a consequence, if there exists $1 \leq j \leq M$ such that \mathcal{L}_{j,ζ_j} is positivity improving, then so is $\mathcal{L}_{cy,\zeta}$.*

Proof. If ρ is positive definite so is $\rho_{\mathcal{E}_j}$ and hence $\rho \otimes \rho_{\mathcal{E}_j}$ as well. Thus for any non-negative and non-zero A one has

$$\mathrm{Tr}(\mathcal{L}_j(\rho) A) = \mathrm{Tr}(\rho \otimes \rho_{\mathcal{E}_j} \times U_j^* A \otimes \mathbb{1} U_j) > 0,$$

so that $\mathcal{L}_j(\rho)$ is indeed positive definite. \square

3.3.3 Time-reversal

The next assumption concerns time-reversal invariance. A time reversal of a quantum system (\mathcal{H}, H) is an antiunitary involution θ on \mathcal{H} such that $\theta H = H\theta$. Given such a θ we denote by Θ the antilinear *-automorphism acting on $\mathcal{B}(\mathcal{H})$ as $\Theta(X) = \theta X \theta$. Note in particular that one has $\Theta(e^{-itH}) = e^{itH}$ for all $t \in \mathbb{R}$, or in an equivalent way if $\gamma_t(X) := e^{itH} X e^{-itH}$ is the Heisenberg evolution of an observable X then

$$\Theta \circ \gamma_t \circ \Theta = \gamma_{-t}, \quad \forall t \in \mathbb{R}. \quad (3.5)$$

In open systems one usually further specifies the structure of θ . Namely we assume that we are given time reversals θ_S and $\theta_{\mathcal{E}}$ associated to (\mathcal{H}_S, H_S) and $(\mathcal{H}_{\mathcal{E}}, H_{\mathcal{E}})$ respectively and such that $\theta_S \otimes \theta_{\mathcal{E}} V = V \theta_S \otimes \theta_{\mathcal{E}}$ where V is the interacting operator. Then $\theta = \theta_S \otimes \theta_{\mathcal{E}}$ is a time reversal for the coupled system $(\mathcal{H}_S \otimes \mathcal{H}_{\mathcal{E}}, H_S + H_{\mathcal{E}} + V)$.

A state ρ is then called time-reversal invariant for θ if $\Theta(\rho) = \rho$ which is equivalent to $\rho(\Theta(X)) = \rho(X^*)$ for all $X \in \mathcal{B}(\mathcal{H})$, and in particular $\rho(\Theta(X)) = \rho(X)$ when X is an observable hence self-adjoint. Finally, a quantum system (\mathcal{H}, H, ρ) is called time-reversal invariant iff there exists a time reversal θ on (\mathcal{H}, H) such that ρ is time-reversal invariant for θ . We consider the following assumption.

Assumption (TRI). *There exist antiunitary involutions θ and $(\theta_{\mathcal{E}_j})_{1 \leq j \leq M}$ acting on \mathcal{H}_S and $(\mathcal{H}_{\mathcal{E}_j})_{1 \leq j \leq M}$ such that $\theta H_S = H_S \theta$, $\theta_{\mathcal{E}_j} H_{\mathcal{E}_j} = H_{\mathcal{E}_j} \theta_{\mathcal{E}_j}$ and $\theta \otimes \theta_{\mathcal{E}_j} V_j = V_j \theta \otimes \theta_{\mathcal{E}_j}$.*

Since the $\rho_{\mathcal{E}_j}$ are KMS states **(TRI)** guarantees that all the probes $(\mathcal{H}_{\mathcal{E}_j}, H_{\mathcal{E}_j}, \rho_{\mathcal{E}_j})$ are time-reversal invariant systems. We shall come back to time-reversal invariance in Section 3.4.3.

3.4 The Non-Entanglement condition

There is a last assumption, see Assumption **(NE)** on p.14, which will play an important role in our paper and which is very specific to RIS. The purpose of this section is first to explain its origin and then to derive some of its consequences.

3.4.1 The Non-Entanglement condition: a signature of equilibrium

One of our goal is to understand linear response theory for RIS, which means how does the system respond to a small perturbation *from equilibrium*. It is therefore important to specify what do we mean by equilibrium. In this perspective the most ideal situation is certainly that of a single type of probes, i.e. $M = 1$, all initially at the same inverse temperature β . One of the usual features of equilibrium is that it is characterized by the vanishing of entropy production. During its interaction with a single probe \mathcal{E} the entropy production of the system is given by (2.4), and if the system is initially in an invariant state ρ we simply have

$$\sigma = \mathrm{Ent}(U \times \rho \otimes \rho_{\mathcal{E}} \times U^* | \rho \otimes \rho_{\mathcal{E}})$$

which vanishes iff $U \times \rho \otimes \rho_{\mathcal{E}} \times U^* = \rho \otimes \rho_{\mathcal{E}}$. It is thus natural to require that for each species of probes, and whatever is their initial temperature, there exists an invariant state ρ_j such that the triple $(U_j, \rho_j, \rho_{\mathcal{E}_j})$ satisfies the Non-Entanglement condition (2.6).

Consider now the general framework of Section 3.2 and more precisely the random situation of Section 3.2.2. When Assumption (**ER** \sharp) holds the entropy production in the system is given by, see Proposition 4.3 and also [BJM2, BJM3],

$$\sigma_{ra}^+ = - \sum_{j=1}^M \beta_j \rho_+^{ra}(\Phi_j), \quad (3.6)$$

where $\Phi_j = -\frac{1}{T} \text{Tr}_{\rho_{\mathcal{E}_j}}(U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j})$, $T = \tau_1 + \dots + \tau_M$, denotes the energy flux observable associated to the j -th type of probe (see Section 4.1 for more details about these Φ_j). A natural notion of equilibrium is that it leads to a vanishing of entropy production. The following proposition gives a simple characterization in terms of the individual probes.

Proposition 3.5. *If Assumption (**ER** \sharp) holds then $\sigma_{ra}^+ = 0$ if and only if*

$$U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* = \rho_+^{ra} \otimes \rho_{\mathcal{E}_j}, \quad \forall j \in \{1, \dots, M\}, \quad (3.7)$$

i.e. the states $\rho_+^{ra} \otimes \rho_{\mathcal{E}_j}$ are invariant states of the joint systems $\mathcal{S} + \mathcal{E}_j$ so that in particular $\mathcal{L}_j(\rho_+^{ra}) = \rho_+^{ra}$ for all j . In other words ρ_+^{ra} is a common invariant state for all the probes and the triples $(U_j, \rho_+^{ra}, \rho_{\mathcal{E}_j})$ all satisfy the Non-Entanglement condition (2.6).

Proof. One direction is obvious. Namely, if (3.7) holds for any j one easily computes

$$\rho_+^{ra}(\Phi_j) = -\frac{1}{T} \text{Tr} [\rho_+^{ra} \otimes \rho_{\mathcal{E}_j} (U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j})] = 0,$$

i.e. all the steady fluxes vanish, hence $\sigma_{ra}^+ = - \sum_{j=1}^M \beta_j \rho_+^{ra}(\Phi_j) = 0$.

Suppose now that $\sigma_{ra}^+ = 0$. Using (2.3) we have for all $j \in \{1, \dots, M\}$

$$\begin{aligned} & \text{Ent}(\mathcal{L}_j(\rho_+^{ra})) - \text{Ent}(\rho_+^{ra}) \\ &= \text{Ent}\left(U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* \mid \mathcal{L}_j(\rho_+^{ra}) \otimes \rho_{\mathcal{E}_j}\right) - \beta_j \text{Tr}\left((\rho_+^{ra} \otimes \rho_{\mathcal{E}_j})(U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j})\right) \\ &= \text{Ent}\left(U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* \mid \mathcal{L}_j(\rho_+^{ra}) \otimes \rho_{\mathcal{E}_j}\right) + \beta_j T \rho_+^{ra}(\Phi_j). \end{aligned}$$

Summing these identities over j and using (3.6) we get

$$\frac{1}{M} \sum_{j=1}^M \text{Ent}(\mathcal{L}_j(\rho_+^{ra})) - \text{Ent}(\rho_+^{ra}) = \frac{1}{M} \sum_{j=1}^M \text{Ent}\left(U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* \mid \mathcal{L}_j(\rho_+^{ra}) \otimes \rho_{\mathcal{E}_j}\right). \quad (3.8)$$

Relative entropies are non-negative quantities so that the right-hand side is non-negative. But the left-hand side is non-positive. Indeed von Neumann entropy is strictly concave so that, using $\mathcal{L}_{ra}(\rho_+^{ra}) = \rho_+^{ra}$, we have

$$\begin{aligned} & \frac{1}{M} \sum_{j=1}^M \text{Ent}(\mathcal{L}_j(\rho_+^{ra})) - \text{Ent}(\rho_+^{ra}) \\ & \leq \text{Ent}\left(\frac{1}{M} \sum_{j=1}^M \mathcal{L}_j(\rho_+^{ra})\right) - \text{Ent}(\rho_+^{ra}) = \text{Ent}(\mathcal{L}_{ra}(\rho_+^{ra})) - \text{Ent}(\rho_+^{ra}) = 0, \end{aligned}$$

with equality iff all the $\mathcal{L}_j(\rho_+^{ra})$ are equal. This proves that both sides of (3.8) vanish.

As a consequence all the $\mathcal{L}_j(\rho_+^{ra})$ are equal and, because $\mathcal{L}_{ra}(\rho_+^{ra}) = \rho_+^{ra}$, they have to be equal to ρ_+^{ra} , i.e. ρ_+^{ra} is a common invariant state for all the probes. Moreover since all the terms on the right-hand side of (3.8) are non-negative they all vanish, i.e.

$$\text{Ent}(U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* | \mathcal{L}_j(\rho_+^{ra}) \otimes \rho_{\mathcal{E}_j}) = 0 \quad \Rightarrow \quad U_j \times \rho_+^{ra} \otimes \rho_{\mathcal{E}_j} \times U_j^* = \mathcal{L}_j(\rho_+^{ra}) \otimes \rho_{\mathcal{E}_j} = \rho_+^{ra} \otimes \rho_{\mathcal{E}_j},$$

that is the triples $(U_j, \rho_+^{ra}, \rho_{\mathcal{E}_j})$ satisfy the Non-Entanglement condition (2.6). \square

The above proposition refers only to the vanishing of entropy production. If we want to consider the case where *all the temperatures are equal* as an *equilibrium situation*, whatever is this temperature, this immediately leads to the following

Assumption (NE). *There exists a function $\mathbb{R} \ni \zeta \mapsto \rho_{+,\zeta}$ such that, for any $j \in \{1, \dots, M\}$ and $\zeta \in \mathbb{R}$, the triple $(U_j, \rho_{+,\zeta}, \rho_{\mathcal{E}_j,\zeta})$ satisfies the Non-Entanglement condition*

$$U_j \times \rho_{+,\zeta} \otimes \rho_{\mathcal{E}_j,\zeta} \times U_j^* = \rho_{+,\zeta} \otimes \rho_{\mathcal{E}_j,\zeta}.$$

Remark 3.3. If moreover Assumption (ER \sharp) holds this obviously implies that, *when all the temperatures are equal*, the unique invariant states in the cyclic and random cases coincide with $\rho_{+,\zeta}$, i.e. $\rho_{+,\zeta}^{cy} = \rho_{+,\zeta}^{ra} = \rho_{+,\zeta}$.

Assumption (NE) may look quite restrictive at first sight. Indeed it requires that for any probe and at any temperature there is a non-entangled invariant state, and that this invariant state depends only on the temperature and not on the probe itself. Proposition 3.5 however shows that this is the natural condition if one wants to consider equal temperatures as an equilibrium situation, in the sense of vanishing of entropy production. Since linear response theory deals with a situation *close to equilibrium*, if Assumption (NE) does not hold one should therefore also take into account a departure from this assumption. It is however not clear how to quantify this or, said differently, what is the natural quantity one can associate to a generic RIS and the vanishing of which would correspond to the fulfilment of Assumption (NE). In particular, if Assumption (NE) fails, the Green-Kubo formulas (4.14) and (4.15) do not hold, at least in the form given in the present paper, and Onsager reciprocity relations are not valid.

On the other hand the large deviation principle for entropy production and the fluctuation relation are not directly related to any equilibrium situation. One can therefore expect that they hold true even without Assumption (NE) and this is indeed the case, see (5.10) and Theorem 5.8. However, when all temperatures are equal the vanishing of entropy production is also equivalent to energy conservation, see (3.6). Therefore, if Assumption (NE) does not hold this proves that, at least at equilibrium, energy conservation fails (which explains why we need it in Proposition 4.2). As we already mentioned in the introduction one should not be surprised. This comes from the fact that the RIS hamiltonian (1.1) is time-dependent. As a consequence the translation symmetry (5.11), which is directly related to energy conservation, also fails if Assumption (NE) does not hold.

In the rest of this section we derive two consequences of this Non-Entanglement assumption which will play an important role in our analysis.

3.4.2 Effective hamiltonian

The first consequence is the existence of a conserved quantity for the \mathcal{L}_j 's and which plays the role of an effective hamiltonian. This is directly related to the conservation of the entropy observable (2.7), see the end of Section 2.3.

Proposition 3.6. *Suppose **(ER \sharp)** and **(NE)** hold and let $H'_S = -\log\left(\rho_{+,0}^{1/\beta_{\text{ref}}}\right)$ so that $\rho_{+,0} = e^{-\beta_{\text{ref}}H'_S}$. Then for any $\zeta \in \mathbb{R}$ the state $\rho_{+,\zeta}$ is a Gibbs state at inverse temperature $\beta = \beta_{\text{ref}} - \zeta$ for the effective hamiltonian H'_S , i.e. $\rho_{+,\zeta} = \frac{e^{-\beta H'_S}}{\text{Tr}\left(e^{-\beta H'_S}\right)}$, $\beta = \beta_{\text{ref}} - \zeta$.*

Moreover, for any j , the observable $H'_S + H_{\mathcal{E}_j}$ is a conserved quantity of the interacting dynamics, namely

$$U_j^* (H'_S \otimes \mathbb{1} + \mathbb{1} \otimes H_{\mathcal{E}_j}) U_j = H'_S \otimes \mathbb{1} + \mathbb{1} \otimes H_{\mathcal{E}_j}. \quad (3.9)$$

Proof. Let $\zeta \in \mathbb{R}$ and $\beta = \beta_{\text{ref}} - \zeta$. Assumption **(ER \sharp)** together with Proposition 3.2 imply that $\rho_{+,0}$ is positive definite. Hence H'_S is well defined and Assumption **(NE)** guarantees that, for any $1 \leq k \leq M$, $[\rho_{+,0} \otimes e^{-\beta_{\text{ref}}H_{\mathcal{E}_k}}, U_k] = 0$. So

$$\left[\rho_{+,0}^{\frac{\beta}{\beta_{\text{ref}}}} \otimes e^{-\beta H_{\mathcal{E}_k}}, U_k \right] = \left[e^{-\beta H'_S} \otimes e^{-\beta H_{\mathcal{E}_k}}, U_k \right] = 0, \quad \forall \beta. \quad (3.10)$$

Hence $\frac{e^{-\beta H'_S}}{\text{Tr}\left(e^{-\beta H'_S}\right)}$ is an invariant state of $\mathcal{L}_{k,\zeta}$ for any k and hence of $\mathcal{L}_{\sharp,\zeta}$. Since the latter admits only one invariant state it coincides with $\rho_{+,\zeta}$.

Finally, (3.9) is a direct consequence of (3.10). \square

Remark 3.4. Note that the effective hamiltonian H'_S is intrinsic to the system and that changing the reference temperature β_{ref} only amounts to an irrelevant shift by a constant.

We also note the following lemma which can be seen as a sort of gauge invariance and whose proof is a straightforward computation left to the reader.

Lemma 3.7. *If Assumptions **(ER \sharp)** and **(NE)** hold and H'_S is as in Proposition 3.6, then for any state ρ , $j \in \{1, \dots, M\}$ and $t \in \mathbb{R}$ one has*

$$\mathcal{L}_j \left(e^{-itH'_S} \rho e^{itH'_S} \right) = e^{-itH'_S} \mathcal{L}_j(\rho) e^{itH'_S}.$$

Remark 3.5. Of course one has the same property in the Heisenberg picture,

$$\mathcal{L}_j^* \left(e^{itH'_S} X e^{-itH'_S} \right) = e^{itH'_S} \mathcal{L}_j^*(X) e^{-itH'_S}. \quad (3.11)$$

In particular, taking $X = H'_S$, we get that $[\mathcal{L}_j^*(H'_S), H'_S] = 0$ for all j .

3.4.3 Non-entanglement, time-reversal invariance and detailed balance

The second consequence of Assumption **(NE)** is related to time-reversal invariance. Assumption **(TRI)** is written in terms of the full interacting dynamics U_j . However the central objects in RIS are the reduced dynamics maps \mathcal{L}_j . It is therefore natural, and important, to understand what are the consequences of Assumption **(TRI)** on them.

Definition 3.1. *Let \mathcal{L} be a CPTP map and ρ an invariant state of \mathcal{L} . The pair (\mathcal{L}, ρ) is said to satisfied the standard quantum detailed balance (SQDB) condition with respect to the time-reversal Θ if $\Theta(\rho) = \rho$ and $\Theta \circ \mathcal{L}^* \circ \Theta = \mathcal{L}^{*\rho}$ where $\mathcal{L}^{*\rho}$ denotes the ρ adjoint of \mathcal{L}^* , i.e. its adjoint with respect to the inner product $\langle A, B \rangle_\rho := \text{Tr}(\rho A^* B)$.*

Remark 3.6. The above definition can be traced back to [Ag], see also [FU, Ma]. We mention that several other notions of quantum detailed balance which do not make reference to any time-reversal operator exist in the literature. The probably most common one is due to Kossakowski, Frigerio, Gorini and Verri [KFGV]. It is called GNS quantum detailed balance and it holds iff \mathcal{L}^* is ρ self-adjoint, i.e. $\mathcal{L}^{*\rho} = \mathcal{L}^*$. To enhance its dependence on the time-reversal operation the notion of quantum detailed balance as given in Definition 3.1 is sometimes called SQDB – Θ condition in the literature, see e.g. [FR].

Remark 3.7. In [JPW] Definition 3.1 is called time-reversal invariance. Indeed, if (\mathcal{H}, H, ρ) is a quantum system with ρ an invariant state and $\gamma_t(X) = e^{itH} X e^{-itH}$ denotes the Heisenberg evolution, it is easy to see that for all $t \in \mathbb{R}$ the ρ adjoint γ_t^ρ of γ_t is γ_{-t} so that (3.5) amounts to $\Theta \circ \gamma_t \circ \Theta = \gamma_t^\rho$. In the markovian description of open systems one then simply replaces the unitary evolution γ_t by the markovian one \mathcal{L}^* and the relation $\Theta \circ \mathcal{L}^* \circ \Theta = \mathcal{L}^{*\rho}$ can therefore also be understood as a signature of time-reversal invariance.

The following proposition shows that the Non-Entanglement condition allows one to make the connection between our Assumption **(TRI)** and Definition 3.1.

Proposition 3.8. *Let θ and $\theta_\mathcal{E}$ be time reversals for the quantum systems $(\mathcal{H}_\mathcal{S}, H_\mathcal{S})$ and $(\mathcal{H}_\mathcal{E}, H_\mathcal{E})$ respectively and $V \in \mathcal{B}(\mathcal{H}_\mathcal{S} \otimes \mathcal{H}_\mathcal{E})$ such that $\theta_\mathcal{S} \otimes \theta_\mathcal{E} V = V \theta_\mathcal{S} \otimes \theta_\mathcal{E}$. Let $\tau > 0$ and $\rho_\mathcal{E}$ be such that $(\mathcal{H}_\mathcal{E}, H_\mathcal{E}, \rho_\mathcal{E})$ is time-reversal invariant, and let \mathcal{L} be given by (2.2). If \mathcal{L} is primitive and its (unique) invariant state ρ is such that the triple $(U, \rho, \rho_\mathcal{E})$, $U = e^{-i\tau H}$ with $H = H_\mathcal{S} + H_\mathcal{E} + V$, satisfies the Non-Entanglement condition (2.6), then the pair (\mathcal{L}, ρ) satisfies the SQDB condition.*

In our framework of RIS we thus immediately get

Corollary 3.9. *If Assumptions **(ER \sharp)**, **(TRI)** and **(NE)** hold then for any $\zeta \in \mathbb{R}$ and any j the pair $(\mathcal{L}_{j,\zeta}, \rho_{+,\zeta})$ satisfies SQDB in the sense that $\Theta(\rho_{+,\zeta}) = \rho_{+,\zeta}$ and*

$$\Theta \circ \mathcal{L}_{j,\zeta}^* \circ \Theta = \mathcal{L}_{j,\zeta}^{*\rho_{+,\zeta}}. \quad (3.12)$$

Proof. Since the indices j and ζ do not play any role we omit them to alleviate the notation. For $X \in \mathcal{B}(\mathcal{H}_\mathcal{E})$ let $\Theta_\mathcal{E}(X) := \theta_\mathcal{E} X \theta_\mathcal{E}$. One immediately gets $\Theta \otimes \Theta_\mathcal{E}(U) = U^*$ so that

$$U \Theta(\rho) \otimes \rho_\mathcal{E} U^* = U (\Theta \otimes \Theta_\mathcal{E})(\rho \otimes \rho_\mathcal{E}) U^* = (\Theta \otimes \Theta_\mathcal{E})(U^* \rho \otimes \rho_\mathcal{E} U) = (\Theta \otimes \Theta_\mathcal{E})(\rho \otimes \rho_\mathcal{E}) = \Theta(\rho) \otimes \rho_\mathcal{E},$$

where we have used successively $\Theta_\mathcal{E}(\rho_\mathcal{E}) = \rho_\mathcal{E}$, $\Theta \otimes \Theta_\mathcal{E}(U) = U^*$, (2.6) and $\Theta_\mathcal{E}(\rho_\mathcal{E}) = \rho_\mathcal{E}$ again. Hence $\Theta(\rho)$ is an invariant state of \mathcal{L} .

We have proven that $\Theta(\rho)$ is an invariant state of \mathcal{L}_j for all j so $\Theta(\rho)$ is an invariant state of \mathcal{L}_{cy} and \mathcal{L}_{ra} . Since the latter are primitive by Assumption **(ER#)** this proves that $\Theta(\rho) = \rho$.

On the other hand, for any $A, B \in \mathcal{B}(\mathcal{H}_S)$ we have

$$\begin{aligned}
\text{Tr}(\rho \mathcal{L}^{*\rho}(A)^* B) &= \text{Tr}(\rho A^* \mathcal{L}^*(B)) \\
&= \text{Tr}(\rho A^* \otimes \rho_{\mathcal{E}} \times U^* \times B \otimes \mathbb{1} \times U) \\
&= \text{Tr}(B \rho \otimes \rho_{\mathcal{E}} \times U \times A^* \otimes \mathbb{1} \times U^*) \\
&= \overline{\text{Tr}(\Theta(B\rho) \otimes \Theta_{\mathcal{E}}(\rho_{\mathcal{E}}) \times U^* \times \Theta(A^*) \otimes \mathbb{1} \times U)} \\
&= \overline{\text{Tr}(\Theta(B\rho) \mathcal{L}^*(\Theta(A)^*))} \\
&= \text{Tr}(B \rho (\Theta \circ \mathcal{L}^* \circ \Theta(A))^*),
\end{aligned}$$

which proves that $\Theta \circ \mathcal{L}^* \circ \Theta = \mathcal{L}^{*\rho}$. Here we have used the definition of $\mathcal{L}^{*\rho}$ in line 1, of \mathcal{L}^* in line 2, cyclicity of the trace and (2.6) in line 3, antilinearity of $\Theta/\Theta_{\mathcal{E}}$ in line 4, $\Theta_{\mathcal{E}}(\rho_{\mathcal{E}}) = \rho_{\mathcal{E}}$ and definition of \mathcal{L}^* in line 5, and finally in line 6 that \mathcal{L}^* is completely positive so that $\mathcal{L}^*(X^*) = (\mathcal{L}^*(X))^*$. \square

Remark 3.8. Note that for cyclic RIS Assumption **(TRI)** makes the RIS time-reversal invariant only for the duration of the joint evolution with each individual probe, that is only locally in time. Indeed, the time-reversal operator Θ does not change the order of the interactions. This lack of global time-reversal invariance will however be resolved in the random model.

Remark 3.9. If **(ER#)**, **(TRI)** and **(NE)** hold it also follows from (3.9) that the effective hamiltonian H'_S is invariant under time reversal, i.e. $\Theta(H'_S) = H'_S$.

3.5 A toy example

Suppose that \mathcal{H}_S and $\mathcal{H}_{\mathcal{E}}$ are copies of \mathbb{C}^2 , $H_S = E a^* a$, $H_{\mathcal{E}} = E_0 b^* b$, and the interaction V has the form $V = \frac{\lambda}{2}(a^* \otimes b + a \otimes b^*)$. Here a/a^* , b/b^* are the usual annihilation/creation operators on \mathcal{H}_S and $\mathcal{H}_{\mathcal{E}}$ respectively, i.e. $a = b = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix}$. E and E_0 denote the excited energy levels of \mathcal{S} and \mathcal{E} respectively, and the interaction consists in an exchange of excitation between \mathcal{S} and \mathcal{E} . This model can be seen as a toy version of the Jaynes-Cummings hamiltonian describing the interaction between one mode of a quantized electro-magnetic field in a cavity and a two-level atom, see e.g. [CDG], and the corresponding RIS as a toy version of the one-atom maser model studied in [BP, Bru]. The spectral analysis and ergodic properties of this toy model can be found in [BJM3].

One easily sees that the total number operator $N_S \otimes \mathbb{1} + \mathbb{1} \otimes N_{\mathcal{E}} \equiv a^* a \otimes \mathbb{1} + \mathbb{1} \otimes b^* b$ is a conserved quantity. As a consequence, if $\rho_{\mathcal{E}} = \frac{e^{-\beta H_{\mathcal{E}}}}{\text{Tr}(e^{-\beta H_{\mathcal{E}}})}$, $H_{\mathcal{E}} = E_0 b^* b$, then $\rho = \frac{e^{-\beta H'_S}}{\text{Tr}(e^{-\beta H'_S})}$, $H'_S = E_0 N_S$, is an invariant state and the triple $(e^{-i\tau H}, \rho, \rho_{\mathcal{E}})$ satisfies condition (2.6).

Consider now the RIS where the probes \mathcal{E}_j are copies of \mathcal{E} with possibly different temperatures. We have just seen that Assumption **(NE)** indeed holds and leads to the effective hamiltonian $H'_S = E_0 N_S$. Assumption **(TRI)** is then obviously satisfied with θ and $\theta_{\mathcal{E}_j}$ the complex conjugation operations in the canonical bases of \mathbb{C}^2 .

Finally let $\nu_j := \sqrt{(E - E_0)^2 + \lambda_j^2}$. A simple calculation shows that if $\nu_j \tau_j$ is not a multiple of 2π then \mathcal{L}_j is primitive and that its unique invariant state is $\rho_+(\zeta_j) := \frac{e^{-\beta_j H'_S}}{\text{Tr}(e^{-\beta_j H'_S})}$ (see e.g. [BJM3] for more details). One thus infers that, if at least one of the $\nu_j \tau_j$'s is not a multiple of 2π , at equilibrium the maps \mathcal{L}_{cy} and \mathcal{L}_{ra} are primitive as well so that **(ER#)** holds.

Remark 3.10. It is easy to see that the true one-atom maser model based on the Jaynes-Cummings hamiltonian also satisfies Assumptions **(TRI)** and **(NE)**. Its ergodic properties are however much more delicate to study (here \mathcal{S} is equivalent to a harmonic oscillator, in particular $\mathcal{H}_{\mathcal{S}}$ has infinite dimension) and even the bare question of return to equilibrium, that is when all the probes are initially at the same temperature, turned out to be a difficult problem [BP, Bru].

4 Linear response of energy fluxes and entropy production

4.1 Energy flux observables

The energy flux observables describe the energy fluxes that get out of the reservoirs \mathcal{R}_j , as they are seen by the small system \mathcal{S} . Moreover we have to take into account the discrete-time nature of the RIS dynamics to define these fluxes. In other words we choose to study averaged fluxes, averaged over the duration of one interaction, instead of instantaneous ones.

Clearly the reservoir \mathcal{R}_j can exchange energy only when it interacts with \mathcal{S} . Moreover the typical time scale of the non-equilibrium RIS is $T = \tau_1 + \dots + \tau_M$ where we recall that τ_j is the duration of the interaction with probe \mathcal{E}_j . This leads to the following (see Section 2.2).

Definition 4.1. *The energy flux observable associated to \mathcal{R}_j is*

$$\Phi_j := -\frac{1}{T} \text{Tr}_{\rho_{\mathcal{E}_j}}(U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j}).$$

In the cyclic framework it will be convenient to also use a slightly different, but closely related, flux observable. If \mathcal{S} is in the state ρ , e.g. the steady state $\rho_{+\zeta}$ (see Remark 4.5), at the beginning of a cycle, i.e. before it interacts with \mathcal{R}_1 , then at the beginning of its interaction with \mathcal{R}_j it is in the state $\mathcal{L}_{j-1} \circ \dots \circ \mathcal{L}_1(\rho)$. Hence the corresponding expectation value of the flux observable is

$$\langle \Phi_j \rangle = \text{Tr}(\mathcal{L}_{j-1} \circ \dots \circ \mathcal{L}_1(\rho) \Phi_j).$$

It is thus natural to introduce

$$\Phi_j^{cy} := \mathcal{L}_1^* \circ \mathcal{L}_2^* \circ \dots \circ \mathcal{L}_{j-1}^*(\Phi_j), \quad (4.1)$$

so that $\langle \Phi_j \rangle = \text{Tr}(\rho \Phi_j^{cy})$. We shall call Φ_j^{cy} the cyclic flux observable associated to \mathcal{R}_j . It is easy to see that one also has

$$\Phi_j^{cy} = -\frac{1}{T} \text{Tr}_{\rho_{cy}}(U_{cy}^* H_{\mathcal{E}_j} U_{cy} - H_{\mathcal{E}_j}),$$

which is the mean energy variation observable in the reservoir \mathcal{R}_j during the entire cycle.

Remark 4.1. If we compare the definitions of Φ_j and Φ_j^{cy} we can see that the latter contains a supplementary information which is the order of the interactions in a cyclic RIS. Taking two distinct definitions allows us to obtain Green-Kubo formulas which are similar in both the cyclic and random situations, see (4.14), (4.15) and (4.17).

Remark 4.2. Note that the flux observables depend on the thermodynamic parameters but that Φ_j^{cy} depend on $\zeta_1, \zeta_2, \dots, \zeta_j$ whereas Φ_j depends only on ζ_j .

Notation. In the sequel the various results will often have a similar form when written in terms of Φ_j^{cy} in the cyclic case and of Φ_j in the random case. When we will write Φ_j^\sharp , $\sharp = cy$ or *ra*, the notation Φ_j^{ra} will therefore stand for Φ_j .

Time reversal plays an important role in linear response theory and the derivation of Green-Kubo formula and Onsager relations. Since we consider here averaged flux observables we need to consider also what we call the reversed-time flux observables.

Definition 4.2. *The reversed-time flux observable associated to \mathcal{R}_j is*

$$\Phi_{j,\text{rev}} := \frac{1}{T} \text{Tr}_{\rho_{\mathcal{E}_j}} (U_j H_{\mathcal{E}_j} U_j^* - H_{\mathcal{E}_j}), \quad U_j = e^{-i\tau_j H_j}.$$

Accordingly the cyclic reversed-time flux observable associated to \mathcal{R}_j is

$$\Phi_{j,\text{rev}}^{cy} := \mathcal{L}_{M,\text{rev}}^* \circ \dots \circ \mathcal{L}_{j+1,\text{rev}}^* (\Phi_{j,\text{rev}}) = \frac{1}{T} \text{Tr}_{\rho_{cy}} (U_{cy} H_{\mathcal{E}_k} U_{cy}^* - H_{\mathcal{E}_k}), \quad (4.2)$$

where $\mathcal{L}_{j,\text{rev}}$ is the reduced dynamics map associated to a time-reversed interaction, i.e.

$$\mathcal{L}_{j,\text{rev}}(\rho) := \text{Tr}_{\mathcal{H}_{\mathcal{E}_j}} (e^{i\tau_j H_j} \rho \otimes \rho_{\mathcal{E}_j} e^{-i\tau_j H_j}).$$

Remark 4.3. If **(ER \sharp)** and **(NE)** hold then for any $j = 1, \dots, M$ we actually have

$$\mathcal{L}_{j,\text{rev}}^* = \mathcal{L}_j^{*\rho_+, \zeta_j}, \quad (4.3)$$

the ρ_+, ζ_j -adjoint of \mathcal{L}_j^* , where we recall that ρ_+, ζ is the global invariant state as given in Assumption **(NE)**. The proof is very similar to the one of (3.12) and is left to the reader. This is of course related to the fact that in the Heisenberg picture the ρ adjoint of the dynamics $\gamma_t(\cdot) = e^{itH} \cdot e^{-itH}$ is γ_{-t} , see Remark 3.7. Note also that if Assumption **(TRI)** holds it follows directly from the definition of $\mathcal{L}_{j,\text{rev}}$ that

$$\mathcal{L}_{j,\text{rev}} = \Theta \circ \mathcal{L}_j \circ \Theta, \quad (4.4)$$

i.e. $\mathcal{L}_{j,\text{rev}}$ is the time-reversal of \mathcal{L}_j which is in agreement with (3.12).

Remark 4.4. 1) The sign discrepancy in the definitions of Φ_j and $\Phi_{j,\text{rev}}$ takes into account the fact that we have reversed the time. Assume for simplicity that all the interaction times τ_j are equal to τ . In the limit $\tau \rightarrow 0$ of instantaneous interactions it is easy to see that the flux observables Φ_j and $\Phi_{j,\text{rev}}$ both coincide with

$$\Phi_j^{\text{inst}} = \frac{1}{M} \text{Tr}_{\rho_{\mathcal{E}_j}} (-i[H_j, H_{\mathcal{E}_j}]).$$

2) When **(TRI)** holds the fluxes Φ_j and $\Phi_{j,\text{rev}}$ satisfy the relation

$$\Phi_{j,\text{rev}} = -\Theta(\Phi_j). \quad (4.5)$$

In the limit of instantaneous interactions one retrieves the usual relation $\Phi_j^{\text{inst}} = -\Theta(\Phi_j^{\text{inst}})$, i.e. flux observables are odd with respect to time reversal.

3) In the cyclic case, one can notice that $\Phi_{j,\text{rev}}^{\text{cy}}$ corresponds to a total time-reversal, where the order of the interactions is reversed as well (compare (4.1) and (4.2)).

When **(ER \sharp)** and **(NE)** hold then $H'_S + H_{\mathcal{E}_j}$ is a conserved quantity of the interacting system $\mathcal{S} + \mathcal{E}_j$ where we recall that H'_S is the effective hamiltonian, see Section 3.4.2. Hence Φ_j is naturally the flux observable associated to H'_S . Namely we have

Proposition 4.1. *Suppose **(ER \sharp)** and **(NE)** hold. Then*

$$\Phi_j = \frac{1}{T}(\mathcal{L}_j^*(H'_S) - H'_S), \quad (4.6)$$

i.e. Φ_j is the flux corresponding to the effective energy H'_S . Similarly one has

$$\Phi_{j,\text{rev}} = -\frac{1}{T}(\mathcal{L}_{j,\text{rev}}^*(H'_S) - H'_S), \quad (4.7)$$

Proof. Using (3.9) we have $U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j} = H'_S - U_j^* H'_S U_j$ hence

$$\Phi_j = \frac{1}{T} \text{Tr}_{\rho_{\mathcal{E}_j}}(U_j^* H'_S U_j - H'_S) = \frac{1}{T}(\mathcal{L}_j^*(H'_S) - H'_S),$$

where we have used that $H'_S \in \mathcal{B}(\mathcal{H}_S)$ and the definition of \mathcal{L}_j . This proves (4.6).

Then (4.7) follows from (4.6) using (4.3), (4.5) and Remark 3.9. \square

4.2 Steady fluxes, energy conservation and entropy production

The first central concepts in non-equilibrium systems are energy conservation (1st law) and the relation between entropy production and steady fluxes (entropy balance equation), together with non-negativity of entropy production.

Proposition 4.2. *[1st law] If **(ER \sharp)** and **(NE)** hold we have*

$$\sum_{j=1}^M \rho_+^\sharp(\Phi_j^\sharp) = 0, \quad \sharp = \text{cy or ra}. \quad (4.8)$$

Remark 4.5. In the cyclic framework, ρ_+^{cy} is the asymptotic state at the beginning of a cycle and not at the beginning of the interaction between \mathcal{S} and the j -th reservoir (which is thus $\mathcal{L}_{j-1} \circ \dots \circ \mathcal{L}_1(\rho_+^{\text{cy}})$). So the steady expectation value of the energy flux out of \mathcal{R}_j is indeed $\rho_+^{\text{cy}}(\mathcal{L}_1^* \circ \mathcal{L}_2^* \circ \dots \circ \mathcal{L}_{j-1}^*(\Phi_j)) = \rho_+^{\text{cy}}(\Phi_j^{\text{cy}})$.

Proof. Using (4.1), (4.6) and $\mathcal{L}_{\text{cy}}(\rho_+^{\text{cy}}) = \rho_+^{\text{cy}}$ we have

$$\sum_{j=1}^M \rho_+^{\text{cy}}(\Phi_j^{\text{cy}}) = T \sum_{j=1}^M \rho_+^{\text{cy}}(\mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^*(\mathcal{L}_j^*(H'_S) - H'_S)) = T \rho_+^{\text{cy}}(\mathcal{L}_{\text{cy}}^*(H'_S) - H'_S) = 0.$$

Similarly, recall $\mathcal{L}_{ra} = \frac{1}{M}(\mathcal{L}_1 + \dots + \mathcal{L}_M)$, so that using (4.6) and $\mathcal{L}_{ra}(\rho_+^{ra}) = \rho_+^{ra}$ we get

$$\sum_{j=1}^M \rho_+^{ra}(\Phi_j) = T \sum_{j=1}^M \rho_+^{ra} (\mathcal{L}_j^*(H'_S) - H'_S) = TM \rho_+^{ra} (\mathcal{L}_{ra}^*(H'_S) - H'_S) = 0.$$

□

Remark 4.6. Of course, as expected and mentioned at the beginning of the proof of Proposition 3.5, when all the temperatures are equal each steady flux vanishes, i.e. $\rho_+^\sharp(\Phi_j^\sharp) = 0$ for all j .

We now turn to entropy production. Recall that the entropy production during an interaction between \mathcal{S} and a system \mathcal{E} is given by (2.4). For a non-equilibrium RIS, and if $\mathbf{j} := (j_n)_{n \in \mathbb{N}^*}$ describes the sequence of interactions (see Section 3.2), the entropy production during the n^{th} interaction is thus

$$\sigma_n(\mathbf{j}) := \text{Ent}(U_{j_n} \times \rho_{n-1}(\mathbf{j}) \otimes \rho_{\mathcal{E}_{j_n}} \times U_{j_n}^* | \rho_n(\mathbf{j}) \otimes \rho_{\mathcal{E}_{j_n}}),$$

where $\rho_n(\mathbf{j}) = \mathcal{L}_{j_n} \circ \dots \circ \mathcal{L}_{j_1}(\rho)$ denotes the state of \mathcal{S} after n interactions. Using (2.3) we have the following entropy balance equation for the N first interactions

$$\sum_{n=1}^N \sigma_n(\mathbf{j}) = \text{Ent}(\rho_N(\mathbf{j})) - \text{Ent}(\rho) - T \sum_{n=1}^N \beta_{j_n} \text{Tr}(\mathcal{L}_{j_{n-1}} \circ \dots \circ \mathcal{L}_{j_1}(\rho) \times \Phi_{j_n}). \quad (4.9)$$

As a consequence in the random, resp. cyclic, cases we can define the entropy production associated to the N first interactions, resp. cycles, by

$$\sigma_{ra}(\mathbf{j}, N) := \sum_{n=1}^N \sigma_n(\mathbf{j}), \quad \text{resp. } \sigma_{cy}(N) := \sum_{n=1}^{NM} \sigma_n(\mathbf{j}). \quad (4.10)$$

For cyclic interactions (4.9) becomes

$$\sigma_{cy}(N) = \text{Ent}(\mathcal{L}_{cy}^N(\rho)) - \text{Ent}(\rho) - T \sum_{j=1}^M \sum_{n=1}^N \beta_j \text{Tr}(\mathcal{L}_{cy}^{n-1}(\rho) \times \Phi_j^{cy}). \quad (4.11)$$

Definition 4.3. We define the asymptotic entropy production rate of a cyclic or random Repeated Interaction System by

$$\sigma_{cy}^+ := \lim_{N \rightarrow +\infty} \frac{\sigma_{cy}(N)}{NT}, \quad \sigma_{ra}^+(\mathbf{j}) := \lim_{N \rightarrow +\infty} \frac{\sigma_{ra}(\mathbf{j}, N)}{N \frac{T}{M}},$$

provided the limits exist.

We then have the following entropy balance result.

Proposition 4.3. Assume (\mathbf{ER}^\sharp) holds. Then the asymptotic entropy productions exist and we moreover have

$$\sigma_{cy}^+ = - \sum_{j=1}^M \beta_j \rho_+^{cy}(\Phi_j^{cy}) \quad \text{and} \quad \sigma_{ra}^+(\mathbf{j}) = - \sum_{j=1}^M \beta_j \rho_+^{ra}(\Phi_j) \quad \mathcal{P} - a.s.$$

Proof. Since \mathcal{H}_S has finite dimension the von Neumann entropies $\text{Ent}(\mathcal{L}_{cy}^N(\rho))$ and $\text{Ent}(\rho_N(\mathbf{j}))$ are uniformly bounded hence give no contribution to the asymptotic entropy production. In the cyclic case the result thus follows directly from (4.11) and Proposition 3.2.

In the random case, using (4.9), Proposition 3.2 and Theorem 3.1 with $A(j) := \beta_j \Phi_j$ we get that $\sigma_{ra}^+(\mathbf{j})$ indeed exists \mathcal{P} -almost surely and is given by

$$\sigma_{ra}^+(\mathbf{j}) = -M\rho_+^{ra}(\mathbb{E}_p(\beta\Phi)) = -\sum_{j=1}^M \beta_j \rho_+^{ra}(\Phi_j).$$

Remark 4.7. In the cyclic case we have divided by T which is the total duration of a cycle while in the random case we have divided by $\frac{T}{M}$ which is the mean duration of an interaction, in agreement with (4.10). These different scalings will appear regularly in the sequel.

In the random case we could have equivalently divided by the total duration $\sum_{n=1}^N \tau_{j_n}$ instead of $N\frac{T}{M}$ because \mathcal{P} -almost surely $\frac{1}{N} \sum_{n=1}^N \tau_{j_n} \rightarrow \frac{T}{M}$ by the strong law of large numbers.

4.3 Green-Kubo formula and Onsager relations

The next step is linear response theory which is concerned with the response of the system to a small perturbation from equilibrium. In this section we derive the Green-Kubo fluctuation-dissipation formula which relates the transport coefficients of the system out of equilibrium to flux-flux correlations at equilibrium. Linear response theory will be completed in Section 5.4 by the Central Limit, *aka* Fluctuation-Dissipation, Theorem 5.9. As we have mentioned at the end of Section 3.4.1, and since they make reference to equilibrium, our results require that the Non-Entanglement Assumption (**NE**) holds.

In order to state the Green-Kubo formula we first recall the notion of dissipation function. Recall that the characteristic time of the system is T .

Definition 4.4. *The dissipation function associated to the reservoir \mathcal{R}_j is*

$$\mathcal{D}_j(X, Y) := \frac{1}{T} (\mathcal{L}_j^*(X^*Y) - \mathcal{L}_j^*(X^*)Y - X^*\mathcal{L}_j^*(Y) + X^*Y). \quad (4.12)$$

If the Kraus decomposition of \mathcal{L}_j is given by (2.1) it is easy to see that $\mathcal{D}_j(X, Y) = \frac{1}{T} \sum_{i \in I} [V_i, X]^* [V_i, Y]$.

In particular $\mathcal{D}_j(X, X) = \frac{1}{T} \sum_{i \in I} [V_i, X]^* [V_i, X]$ is non-negative and is zero only if X commutes with all the V_i 's.

Remark 4.8. The dissipation function has been introduced in [Li] for continuous-time quantum dynamical semigroups $(e^{tL})_{t \geq 0}$ as the sesquilinear map D acting on $\mathcal{B}(\mathcal{H}_S)$ and defined by

$$D(X, Y) = L(X^*Y) - L(X^*)Y - X^*L(Y).$$

Here L is the Lindblad generator of a semigroup of unital completely positive maps, i.e. corresponding to the Heisenberg picture. If we fix some characteristic time T , according to (4.12) the dissipation function associated to the unital completely positive map e^{TL} is

$$\mathcal{D}_T(X, Y) := \frac{1}{T} (e^{TL}(X^*Y) - e^{TL}(X^*)Y - X^*e^{TL}(Y) + X^*Y)$$

and it is easy to see that $\lim_{T \rightarrow 0} \mathcal{D}_T(X, Y) = D(X, Y)$.

Theorem 4.4 (GK formula and Onsager relations). *Suppose $(\mathbf{ER}_\#)$ and (\mathbf{NE}) hold.*

1) *The maps $\zeta \mapsto \rho_{+,\zeta}^\#$, $\zeta \mapsto \Phi_{j,\zeta}$ and $\zeta \mapsto \Phi_{j,\zeta}^{cy}$ are infinitely differentiable. The quantities*

$$L_{jk}^\# := \partial_{\zeta_k} \rho_{+,\zeta}^\#(\Phi_{j,\zeta}^\#) \Big|_{\zeta=0}, \quad \# = cy \text{ or } ra, \quad (4.13)$$

are called the kinetic coefficients.

2) *We have the following Green-Kubo formulas*

$$L_{jk}^{cy} = T \sum_{n=0}^{+\infty} \rho_+ \left(\mathcal{L}_{k+1}^* \circ \cdots \circ \mathcal{L}_M^* \circ \mathcal{L}_{cy}^{*n} \circ \mathcal{L}_1^* \circ \cdots \circ \mathcal{L}_{j-1}^* (\Phi_j) \times \Phi_{k,\text{rev}} \right) \quad (4.14)$$

$$+ \delta_{j>k} T \rho_+ \left(\mathcal{L}_{k+1}^* \circ \cdots \circ \mathcal{L}_{j-1}^* (\Phi_j) \times \Phi_{k,\text{rev}} \right) + \frac{1}{2} \delta_{jk} \rho_+ (\mathcal{D}_j(H'_S, H'_S)),$$

$$L_{jk}^{ra} = \frac{T}{M} \sum_{n=0}^{+\infty} \rho_+ (\mathcal{L}_{ra}^{*n} (\Phi_j) \Phi_{k,\text{rev}}) + \frac{1}{2} \delta_{jk} \rho_+ (\mathcal{D}_j(H'_S, H'_S)), \quad (4.15)$$

where all the quantities on the right-hand side are calculated at equilibrium $\zeta = 0$, e.g. ρ_+ stands for $\rho_+ = \rho_{+,0}^{cy} = \rho_{+,0}^{ra}$ (see Remark 3.3).

3) *If moreover Assumption (\mathbf{TRI}) holds then the following analogs of the Onsager reciprocity relations are satisfied:*

$$L_{jk}^{cy} = L_{kj}^{rcy} \quad \text{and} \quad L_{jk}^{ra} = L_{kj}^{ra}, \quad (4.16)$$

where L_{kj}^{rcy} denotes the kinetic coefficient associated to the cyclic RIS in which we reverse the order of the interactions, i.e. $\mathcal{L}^{rcy} = \mathcal{L}_1 \circ \cdots \circ \mathcal{L}_M$. In the specific case $M = 2$, we retrieve the usual Onsager reciprocity relations $L_{jk}^{cy} = L_{kj}^{cy}$ for the cyclic case too.

Let us comment on identities (4.14) and (4.15).

i) The sum in (4.14) describes the flux-flux correlation at equilibrium between reservoirs \mathcal{R}_j and \mathcal{R}_k . Since \mathcal{R}_j interacts only once during each cycle, using (4.1)-(4.2), it is easy to see that this sum can actually be written in terms of \mathcal{L}_{cy} leading to the more condensed form

$$\sum_{n=0}^{+\infty} \rho_+ \left(\mathcal{L}_{k+1}^* \circ \cdots \circ \mathcal{L}_M^* \circ \mathcal{L}_{cy}^{*n} \circ \mathcal{L}_1^* \circ \cdots \circ \mathcal{L}_{j-1}^* (\Phi_j) \times \Phi_{k,\text{rev}} \right) = \sum_{n=0}^{+\infty} \rho_+ (\mathcal{L}_{cy}^{*n} (\Phi_j^{cy}) \Phi_{k,\text{rev}}^{cy}). \quad (4.17)$$

ii) The sum (4.17) takes into account only the correlations when at least one cycle has been achieved. The second term in the right-hand side of (4.14) then takes into account the contribution of the flux-flux correlation between reservoirs \mathcal{R}_k and \mathcal{R}_j if less than a cycle has already occurred, which can happen only if $j > k$.

iii) In both (4.14) and (4.15) the δ_{jk} term takes into account the self-correlation of reservoir \mathcal{R}_j with itself during its first interaction with \mathcal{S} . It is non-negative and vanishes only if \mathcal{S} is not effectively coupled to \mathcal{E}_j in the following sense. If $\mathcal{L}_j = \sum_i V_i \cdot V_i^*$ then $\rho_+ (\mathcal{D}_j(H'_S, H'_S)) = 0$ iff $\mathcal{D}_j(H'_S, H'_S) = 0$ (the latter is non-negative and ρ_+ is positive definite) which in turn holds iff $[V_i, H'_S] = 0$ for all i . But this implies that $\mathcal{L}_j^*(H'_S) = \sum_i V_i^* H'_S V_i = \sum_i V_i^* V_i H'_S = H'_S$ because \mathcal{L}_j^* is unital, and hence $\Phi_j = 0$ by (4.6): whatever are the various temperatures and whatever is the state of \mathcal{S} there is no flux between \mathcal{S} and \mathcal{R}_j . We mention that a similar term appears in [JPW] in the framework of quantum dynamical semigroups.

iv) Finally the prefactors T and $\frac{T}{M}$ represent the “time-step”. In the cyclic case, \mathcal{L}_{cy} describes the evolution during a cycle while \mathcal{L}_{ra} is associated only to a time-step $\frac{T}{M}$.

Remark 4.9. The Onsager reciprocity relations are satisfied only for the random and the $M = 2$ cyclic cases. In the $M > 2$ cyclic case Assumption **(TRI)** reverses the time only locally, leaving the order of the interactions unchanged (see Remark 3.8). This explains why L_{jk}^{cy} has to be compared to the kinetic coefficient L_{kj}^{rcy} of the reversed-order cyclic model. This is similar to what happens when time reversal invariance is broken by the presence of a magnetic field and one has to replace Onsager relations by Onsager-Casimir relations.

Notation In the rest of the paper, any quantity with a subscript/superscript rcy should be understood as the quantity associated to the reversed-order cyclic RIS in which \mathcal{S} interacts first with \mathcal{R}_M , then \mathcal{R}_{M-1}, \dots . For example $\mathcal{L}_{rcy} = \mathcal{L}_1 \circ \dots \circ \mathcal{L}_M$, $\Phi_j^{rcy} = \mathcal{L}_M^* \circ \dots \circ \mathcal{L}_{j+1}^*(\Phi_j)$. Of course, any result which holds true for the cyclic RIS also holds for the reversed cyclic one.

5 Entropic fluctuation

The purpose of this section is to go beyond Proposition 4.3 and to study the statistical fluctuations of the entropy fluxes $\beta_j \Phi_j$ going out the reservoirs $(\mathcal{R}_j)_{1 \leq j \leq M}$.

5.1 Full counting statistics

To study the entropy fluctuation we consider the statistics of the increments of the entropy observable as given by a two time measurement protocol, also called Full Counting Statistics (FCS), and which we briefly recall for the convenience of the reader. This approach goes back to [Ku] and, independently, [Ta]. It has more recently been used in e.g. [JOPP, BDBP, BJPPP, BPP, BPR], see also the review [EHM].

Consider a quantum system \mathcal{S} with underlying finite dimensional Hilbert space \mathcal{H} and let A be the observable of interest. Suppose the system is in the state ρ when we perform a first measurement of A . The possible outcomes of the measurement are eigenvalues of A and $a \in \text{sp}(A)$ is observed with probability

$$\mathbb{P}(A = a) = \text{Tr}(\Pi_a(A)\rho\Pi_a(A))$$

where $\Pi_a(A)$ is the spectral projector of A associated to the eigenvalue a . After this first measurement, given that the outcome is a , the state of the system becomes $\frac{\Pi_a(A)\rho\Pi_a(A)}{\text{Tr}(\Pi_a(A)\rho\Pi_a(A))}$. Subsequently, if the evolution of the system during some time interval of length τ is described by some unitary operator U , a second measurement of A at time τ gives the value $a' \in \text{sp}(A)$ with probability

$$\frac{\text{Tr}(\Pi_{a'}(A)U\Pi_a(A)\rho\Pi_a(A)U^*\Pi_{a'}(A))}{\text{Tr}(\Pi_a(A)\rho)}.$$

The joint probability distribution of the two measurements is thus given by

$$\mathbb{P}(a, a') = \text{Tr}(\Pi_{a'}(A)U\Pi_a(A)\rho\Pi_a(A)U^*\Pi_{a'}(A)),$$

and the statistics of the increment ΔA of A as given by this protocol is therefore given by

$$\mathbb{P}(\Delta A = \delta) = \sum_{\substack{a, a' \in \text{sp}(A) \times \text{sp}(A) \\ a' - a = \delta}} \text{Tr}(\Pi_{a'}(A)U\Pi_a(A)\rho\Pi_a(A)U^*\Pi_{a'}(A)).$$

Note that if A commutes with the initial state ρ we simply have

$$\mathbb{P}(\Delta A = \delta) = \sum_{\substack{a, a' \in \text{sp}(A) \times \text{sp}(A) \\ a' - a = \delta}} \text{Tr}(\Pi_{a'}(A)U\Pi_a(A)\rho U^*) \quad (5.1)$$

so that the expectation of ΔA as given by this two-measurement protocol is

$$\mathbb{E}_{\mathbb{P}}(\Delta A) = \sum_{a, a' \in \text{sp}(A) \times \text{sp}(A)} (a' - a) \text{Tr}(\Pi_{a'}(A)U\Pi_a(A)\rho U^*) = \text{Tr}(\rho(U^*AU - A)), \quad (5.2)$$

and coincides with the expectation value of the flux observable $U^*AU - A$ associated to A .

Note also that, although it does not appear in the notation, the probability law \mathbb{P} depends on the initial state ρ of the system when the first measurement is performed.

In this section, we are interested in the full statistics of the increment of entropy observables of the probes, where the entropy observables are defined in Section 2.3. Since probes are initially in thermal equilibrium, up to an irrelevant constant the entropy observable of \mathcal{E}_j is $S_{\mathcal{E}_j} := \beta_j H_{\mathcal{E}_j}$. The corresponding entropy increment observable is therefore $U_j^* S_{\mathcal{E}_j} U_j - S_{\mathcal{E}_j} = \beta_j (U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j})$. We shall also consider the energy increment observables $U_j^* H_{\mathcal{E}_j} U_j - H_{\mathcal{E}_j}$ which up to a prefactor $\frac{1}{T}$ correspond to the flux observables considered in Section 4.1.

Let now $\mathbf{j} := (j_n)_{n \in \mathbb{N}^*} \in \{1, \dots, M\}^{\mathbb{N}^*}$ be the sequence of indices describing the sequence of probes with which \mathcal{S} interacts, and denote by $\mathcal{E}_{j_n}^n$ the n^{th} probe. Recall that $\mathcal{E}_{j_n}^n$ is a copy of \mathcal{E}_{j_n} . Then, according to (5.1), the probability distribution of the increment of entropy for the n -th interaction is given by

$$\mathbb{P}(\Delta \mathcal{S}_{\mathcal{E}_{j_n}^n} = \varsigma) = \sum_{s' - s = \varsigma} \text{Tr}(\Pi_{s'}(S_{\mathcal{E}_{j_n}^n})U_n \Pi_s(S_{\mathcal{E}_{j_n}^n}) \times \rho \otimes \rho_{\mathcal{E}_{j_n}^n} \times U_n^*),$$

where the sum runs over $s, s' \in \text{sp}(S_{\mathcal{E}_{j_n}^n})$ such that $s' - s = \varsigma$, ρ is the state of \mathcal{S} at the beginning of the interaction and where we have used that the observable $S_{\mathcal{E}_{j_n}^n}$ commutes with the ‘‘initial state’’ $\rho \otimes \rho_{\mathcal{E}_{j_n}^n}$.

To describe the statistics of entropy increments during the N first interactions we introduce some more notation. For any $\mathbf{s} := ((s_n, s'_n))_n \in (\mathbb{R}^2)^{\mathbb{N}^*}$ and $N \geq 1$ we define

$$\mathcal{U}_N^{\mathbf{s}}(\mathbf{j}) := U_{(s_N, s'_N)}^N(S_{\mathcal{E}_{j_N}^N}) \times \dots \times U_{(s_1, s'_1)}^1(S_{\mathcal{E}_{j_1}^1}), \quad \text{where } U_{(s, s')}^n(S) := \Pi_{s'}(S)U_n \Pi_s(S).$$

Note that clearly $U_{(s, s')}^n(S)$ is non-zero only when $s, s' \in \text{sp}(S)$. Then the statistics of entropy increments during the N first interactions is given by

$$\mathbb{P}(\Delta \mathcal{S}_{\mathcal{E}_{j_1}^1} = \varsigma_1, \dots, \Delta \mathcal{S}_{\mathcal{E}_{j_N}^N} = \varsigma_N) = \sum \text{Tr}(\mathcal{U}_N^{\mathbf{s}}(\mathbf{j}) \times \rho_{\text{tot}}^N(\mathbf{j}) \times \mathcal{U}_N^{\mathbf{s}}(\mathbf{j})^*),$$

where the sum runs over $\mathbf{s} \in (\mathbb{R}^2)^N$ such that $\forall 1 \leq n \leq N, s_n, s'_n \in \sigma(S_{\mathcal{E}_{j_n}^n}) \times \sigma(S_{\mathcal{E}_{j_n}^n})$ and $s'_n - s_n = \varsigma_n$, and where $\rho_{\text{tot}}^N(\mathbf{j})$ denotes the total initial state $\rho \otimes \rho_{\mathcal{E}_{j_1}^1} \otimes \dots \otimes \rho_{\mathcal{E}_{j_N}^N}$.

Definition 5.1. *The (random) vectors of entropy increments and energy increments after N interactions are $\mathcal{S}_{\mathcal{R}}^N(\mathbf{j}) := (\mathcal{S}_{\mathcal{R}_1}^N(\mathbf{j}), \dots, \mathcal{S}_{\mathcal{R}_M}^N(\mathbf{j}))$, $\mathcal{Q}_{\mathcal{R}}^N(\mathbf{j}) := (\mathcal{Q}_{\mathcal{R}_1}^N(\mathbf{j}), \dots, \mathcal{Q}_{\mathcal{R}_M}^N(\mathbf{j}))$ where*

$$\mathcal{S}_{\mathcal{R}_j}^N(\mathbf{j}) := \sum_{\substack{1 \leq n \leq N \\ j_n = j}} \Delta \mathcal{S}_{\mathcal{E}_{j_n}^n}, \quad \mathcal{Q}_{\mathcal{R}_j}^N(\mathbf{j}) := -\beta_j^{-1} \mathcal{S}_{\mathcal{R}_j}^N(\mathbf{j}).$$

Remark 5.1. The sign convention in the definition of $\mathcal{Q}_{\mathcal{R}_j}^N(\mathbf{j})$ is in agreement with the one of the flux observables Φ_j of Section 4.1.

We obviously have the following

Proposition 5.1. *The joint probability distribution of $\mathcal{S}_{\mathcal{R}}^N(\mathbf{j})$ is given by*

$$\mathbb{P}(\mathcal{S}_{\mathcal{R}}^N(\mathbf{j}) = \boldsymbol{\varsigma}) = \sum \text{Tr}(\mathcal{U}_N^{\mathbf{s}}(\mathbf{j}) \times \rho_{\text{tot}}^N(\mathbf{j}) \times \mathcal{U}_N^{\mathbf{s}}(\mathbf{j})^*), \quad (5.3)$$

where the sum runs over all \mathbf{s} such that for all $1 \leq n \leq N$, $s_n, s'_n \in \sigma(S_{\mathcal{E}_{j_n}^n}) \times \sigma(S_{\mathcal{E}_{j_n}^n})$ and for any $j \in \{1, \dots, M\}$ one has $\sum_{\substack{1 \leq n \leq N \\ j_n = j}} s'_n - s_n = \varsigma_j$ with $\boldsymbol{\varsigma} = (\varsigma_1, \dots, \varsigma_M)$.

In the case of cyclic RIS, in agreement with (4.11), we shall actually consider

$$\mathcal{S}_{\mathcal{R}}^N(\text{cy}) := \mathcal{S}_{\mathcal{R}}^{N \times M}(\mathbf{j}^{\text{cy}}), \text{ resp. } \mathcal{Q}_{\mathcal{R}}^N(\text{cy}) := \mathcal{Q}_{\mathcal{R}}^{N \times M}(\mathbf{j}^{\text{cy}}),$$

the vector of entropy, resp. energy, increments associated to the cyclic RIS after N cycles.

In the random case note that randomness is now twofold: randomness due to the random order of interactions and a “quantum” randomness due to the two time measurement protocol, and that the latter depends on the former. Indeed as already mentioned the probability law \mathbb{P} of a two time measurement protocol depends on the initial state of the (entire) system, hence on ρ but more importantly on the sequence \mathbf{j} of probes (via the $\rho_{\mathcal{E}_{j_n}^n}$ ’s). To be precise, given a sequence \mathbf{j} we should denote $\mathbb{P}_{\mathbf{j}}$ instead of \mathbb{P} and then the joint probability distribution with respect to the two alea is

$$\hat{\mathbb{P}}(\mathbf{j}, \mathbf{s}) = \mathcal{P}(\mathbf{j}) \times \mathbb{P}_{\mathbf{j}}(\mathbf{s}). \quad (5.4)$$

5.2 Moment generating function

We will analyze the large time limit, i.e. $N \rightarrow \infty$, statistics of the entropy and energy increments through their moment generating functions (MGF).

Definition 5.2 (Moment generating function). *Let $\boldsymbol{\alpha} = (\alpha_1, \dots, \alpha_M) \in \mathbb{C}^M$. We denote by $r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha})$ and $r_{N,\rho}^{\text{cy}}(\boldsymbol{\alpha})$ the respective MGF of the vectors $-\mathcal{S}_{\mathcal{R}}^N(\mathbf{j})$ and $-\mathcal{S}_{\mathcal{R}}^N(\text{cy})$ at $\boldsymbol{\alpha}$, i.e.*

$$r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha}) := \mathbb{E}_{\mathbb{P}_{\mathbf{j}}} \left(e^{-\sum_{j=1}^M \alpha_j \mathcal{S}_{\mathcal{R}_j}^N(\mathbf{j})} \right) \text{ and } r_{N,\rho}^{\text{cy}}(\boldsymbol{\alpha}) := \mathbb{E}_{\mathbb{P}} \left(e^{-\sum_{j=1}^M \alpha_j \mathcal{S}_{\mathcal{R}_j}^N(\text{cy})} \right).$$

Note that the above MGF are defined with respect to the “quantum” alea. We have also stressed their dependence on the initial state ρ of the small system. In the random RIS case we shall also consider the MGF with respect to the two alea.

Definition 5.3. *For $\boldsymbol{\alpha} \in \mathbb{C}^M$, let $r_{N,\rho}^{r\mathbf{a}}(\boldsymbol{\alpha}) := \mathbb{E}_{\mathcal{P}} \left(r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha}) \right) = \mathbb{E}_{\hat{\mathbb{P}}} \left(e^{-\sum_{j=1}^M \alpha_j \mathcal{S}_{\mathcal{R}_j}^N(\mathbf{j})} \right)$.*

To understand the MGF $r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha})$, $r_{N,\rho}^{\text{cy}}(\boldsymbol{\alpha})$ and $r_{N,\rho}^{r\mathbf{a}}(\boldsymbol{\alpha})$, following e.g. [JPW, BDBP, vHG, vHG2, BPP], we consider the following deformations of the reduced dynamics maps $(\mathcal{L}_j)_{1 \leq j \leq M}$.

Definition 5.4. For $\alpha = (\alpha_1, \dots, \alpha_M) \in \mathbb{C}^M$ we define

$$\mathcal{L}_j^{[\alpha]^*}(X) := \text{Tr}_{\mathcal{H}_{\mathcal{E}_j}} \left(\mathbb{1} \otimes \rho_{\mathcal{E}_j}^{1-\alpha_j} \times U_j^* \times X \otimes \rho_{\mathcal{E}_j}^{\alpha_j} \times U_j \right),$$

and

$$\mathcal{L}_{cy}^{[\alpha]^*} := \mathcal{L}_1^{[\alpha]^*} \circ \dots \circ \mathcal{L}_M^{[\alpha]^*}, \quad \mathcal{L}_{ra}^{[\alpha]^*} := \frac{1}{M} (\mathcal{L}_1^{[\alpha]^*} + \dots + \mathcal{L}_M^{[\alpha]^*}). \quad (5.5)$$

Remark 5.2. The notation $\mathcal{L}_j^{[\alpha]^*}$ is chosen in order to be consistent with the reduced dynamics maps \mathcal{L}_j^* . Indeed, $\mathcal{L}_j^{[0]^*} = \mathcal{L}_j^*$ and $\mathcal{L}_\#^{[0]^*} = \mathcal{L}_\#^*$, $\# = cy$ or ra . Of course $\mathcal{L}_j^{[\alpha]}$ will denote its dual map so that $\mathcal{L}_j^{[0]} = \mathcal{L}_j$. Note also that for any j the map $\mathcal{L}_j^{[\alpha]^*}$ only depends on α_j .

The connection between the MGF and the $\mathcal{L}_j^{[\alpha]^*}$'s is provided by the following proposition. For the convenience of the reader we briefly sketch its proof in Section 6.2.

Proposition 5.2. For any $\alpha \in \mathbb{C}^M$ and $N \geq 1$ one has

$$r_{N,\rho}^j(\alpha) = \rho \left(\mathcal{L}_{j_1}^{[\alpha]^*} \circ \dots \circ \mathcal{L}_{j_N}^{[\alpha]^*} (\mathbb{1}) \right). \quad (5.6)$$

As a consequence

$$r_{N,\rho}^{cy}(\alpha) = \rho \left(\left(\mathcal{L}_{cy}^{[\alpha]^*} \right)^N (\mathbb{1}) \right) \quad \text{and} \quad r_{N,\rho}^{ra}(\alpha) = \rho \left(\left(\mathcal{L}_{ra}^{[\alpha]^*} \right)^N (\mathbb{1}) \right). \quad (5.7)$$

Of course the same approach can be used to study the energy increments.

Definition 5.5. Let $\alpha \in \mathbb{C}^M$. We denote by $\tilde{r}_{N,\rho}^j(\alpha)$ and $\tilde{r}_{N,\rho}^{cy}(\alpha)$ the respective MGF of the vectors $\mathcal{Q}_{\mathcal{R}}^N(j)$ and $\mathcal{Q}_{\mathcal{R}}^N(cy)$ at α , i.e.

$$\tilde{r}_{N,\rho}^j(\alpha) := \mathbb{E}_{\mathbb{P}} \left(e^{\sum_{j=1}^M \alpha_j \mathcal{Q}_{\mathcal{R}_j}^N(j)} \right) \quad \text{and} \quad \tilde{r}_{N,\rho}^{cy}(\alpha) := \mathbb{E}_{\mathbb{P}} \left(e^{\sum_{j=1}^M \alpha_j \mathcal{Q}_{\mathcal{R}_j}^N(cy)} \right),$$

as well as, in the random case, the MGF with respect to the two alea, i.e.

$$\tilde{r}_{N,\rho}^{ra}(\alpha) := \mathbb{E}_{\mathcal{P}} \left(\tilde{r}_{N,\rho}^j(\alpha) \right) = \mathbb{E}_{\mathbb{P},\mathcal{P}} \left(e^{\sum_{j=1}^M \alpha_j \mathcal{Q}_{\mathcal{R}_j}^N(j)} \right).$$

It follows directly from Definition 5.1 that the entropy and energy MGF satisfy the relations

$$\tilde{r}_{N,\rho}^j(\alpha) = r_{N,\rho}^j \left(\frac{\alpha}{\beta} \right), \quad \tilde{r}_{N,\rho}^\#(\alpha) = r_{N,\rho}^\# \left(\frac{\alpha}{\beta} \right), \quad \# = cy, ra, \quad (5.8)$$

and where $\frac{\alpha}{\beta}$ denotes the vector $\left(\frac{\alpha_1}{\beta_1}, \dots, \frac{\alpha_M}{\beta_M} \right)$.

Although not mentioned explicitly all the above quantities depend on the vector $\zeta = (\zeta_1, \dots, \zeta_M)$ of thermal forces and, as for \mathcal{L}_j , the map $\mathcal{L}_j^{[\alpha]}$ actually only depends on ζ_j .

Proposition 5.3. If $(\mathbf{ER}_\#)$ holds then, for any $\zeta, \alpha \in \mathbb{R}^M$, the map $\mathcal{L}_{\#, \zeta}^{[\alpha]^*}$ is a primitive CP map. We shall denote by $r_\zeta^\#(\alpha) > 0$ its spectral radius.

Proof. The proof is the same as the one of Proposition 3.2. \square

One can then relate the large N behaviour of the moment generating function to the spectral radius of $\mathcal{L}_{\#}^{[\alpha]*}$ using Proposition 5.2.

Theorem 5.4. *Suppose $(\mathbf{ER}_{\#})$ holds.*

1) *For any initial state ρ and $\alpha, \zeta \in \mathbb{R}^M$ one has*

$$r_{\zeta}^{\#}(\alpha) = \lim_{n \rightarrow +\infty} r_{n,\rho}^{\#}(\alpha)^{\frac{1}{n}}. \quad (5.9)$$

2) *If (\mathbf{TRI}) holds we have the following version of the Evans-Searles symmetry:*

$$r_{\zeta}^{cy}(\mathbf{1} - \alpha) = r_{\zeta}^{rcy}(\alpha), \quad r_{\zeta}^{ra}(\mathbf{1} - \alpha) = r_{\zeta}^{ra}(\alpha) \quad (5.10)$$

where $r_{\zeta}^{rcy}(\alpha)$ is the spectral radius of $\mathcal{L}_{rcy,\zeta}^{[\alpha]*} := \mathcal{L}_{M,\zeta}^{[\alpha]*} \circ \dots \circ \mathcal{L}_{1,\zeta}^{[\alpha]*}$, the deformed reduced dynamics map corresponding to the cyclic RIS with a reversed order, and $\mathbf{1} = (1, \dots, 1)$.

3) *If (\mathbf{NE}) holds we have the following translation symmetry*

$$r_{\zeta}^{\#}(\alpha) = r_{\zeta}^{\#}(\alpha + \lambda \beta^{-1}), \quad \# = cy, ra, \quad (5.11)$$

for any $\alpha, \zeta \in \mathbb{R}^M$, $\lambda \in \mathbb{R}$ and where $\beta^{-1} = (\beta_1^{-1}, \dots, \beta_M^{-1})$.

In view of (5.9) we will call $\alpha \mapsto r_{\zeta}^{\#}(\alpha)$ the *large time moment generating function* of the entropy fluxes. Of course (5.8) immediately leads to a similar result for energy fluxes, i.e.

$$\tilde{r}_{\zeta}^{\#}(\alpha) := \lim_{n \rightarrow +\infty} \tilde{r}_{n,\rho}^{\#}(\alpha)^{\frac{1}{n}} = r_{\zeta}^{\#} \left(\frac{\alpha}{\beta} \right).$$

In particular the translation symmetry (5.11) becomes

$$\tilde{r}^{\#}(\alpha) = \tilde{r}^{\#}(\alpha + \lambda \mathbf{1}) \quad (5.12)$$

for any $\alpha \in \mathbb{R}^M$ and $\lambda \in \mathbb{R}$. This translation symmetry (5.12) is related to the energy conservation (4.8), see Remark 5.8 below. As we have already mentioned at the end of Section 3.4.1 it is thus not surprising that we need to assume (\mathbf{NE}) at this point.

Remark 5.3. We mention that analogous results are obtained in [AGMT, BPP] for systems continuously interacting with infinitely extended reservoirs. There the authors consider a two time measurement protocol with finitely extended reservoirs, then perform a thermodynamic limit which leads to the infinitely extended reservoirs and finally look at a large time limit. Their analysis requires some ultra-violet condition on the coupling between the small system \mathcal{S} and the reservoirs to control the large time limit (this ultraviolet assumption was initially missing in [AGMT] leaving a gap in the proof which was further corrected in [BPP]). The situation is somehow simpler here because at finite time only finitely many probes have already been coupled to \mathcal{S} . Said differently the thermodynamic and large time limit are in a sense taken simultaneously. On the other hand the RIS structure leads to a repeated two time measurement protocol (after n interactions $2n$ measurements have been done) instead of a single one.

5.3 First moments and link with linear response theory

Our next result concerns the first and second moments of the entropy/energy increments and makes the connection with the results presented in Section 4.3. We thus suppose throughout this section that **(ER_‡)** and **(NE)** hold. We also introduce the following maps which act on $\mathcal{B}(\mathcal{H}_S)$.

Definition 5.6. For any $j = 1, \dots, M$ let

$$\varphi_j(X) := \frac{1}{T} (\mathcal{L}_j^*(X H'_S) - \mathcal{L}_j^*(X) H'_S), \quad \varphi_j^{cy} := \mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^* \circ \varphi_j \circ \mathcal{L}_{j+1}^* \circ \dots \circ \mathcal{L}_M^*. \quad (5.13)$$

Remark 5.4. It follows from (4.1) and (4.6) that $\varphi_j(\mathbb{1}) = \Phi_j$ and $\varphi_j^{cy}(\mathbb{1}) = \Phi_j^{cy}$.

The maps φ_j appear naturally as the derivatives with respect to α of the deformed maps $\mathcal{L}_j^{[\alpha]^*}$, see Lemma 6.4. Of course, similarly to the flux observable Φ_j , the map φ_j depends on the thermodynamical force ζ_j . It then follows directly from (4.3) that for any $\zeta \in \mathbb{R}$ one has

$$\varphi_{j,\zeta}^{\rho_+^{\zeta}}(X) = \frac{1}{T} (\mathcal{L}_{j,\text{rev}}^*(X) H'_S - \mathcal{L}_{j,\text{rev}}^*(X H'_S))$$

so that in particular, using (4.7), we get

$$\varphi_{j,\zeta}^{\rho_+^{\zeta}}(\mathbb{1}) = \Phi_{j,\text{rev}}. \quad (5.14)$$

Theorem 5.5. Suppose **(ER_‡)** and **(NE)** hold. Then we have

1) For all $j = 1, \dots, M$,

$$\frac{\partial r^{cy}(\alpha)}{\partial \alpha_j} \Big|_{\alpha=0} = \beta_j T \rho_+^{cy}(\Phi_j^{cy}), \quad \frac{\partial r^{ra}(\alpha)}{\partial \alpha_j} \Big|_{\alpha=0} = \beta_j \frac{T}{M} \rho_+^{ra}(\Phi_j), \quad (5.15)$$

2) For all $j, k = 1, \dots, M$,

$$\begin{aligned} \frac{\partial^2 r^{cy}(\alpha)}{\partial \alpha_k \partial \alpha_j} \Big|_{\alpha=0} &= \beta_j \beta_k T^2 \sum_{n=0}^{\infty} \rho_+^{cy} \left(\varphi_j^{cy} \circ \mathcal{L}_{cy}^{*n} (\Phi_k^{cy} - \rho_+^{cy}(\Phi_k^{cy})) + \varphi_k^{cy} \circ \mathcal{L}_{cy}^{*n} (\Phi_j^{cy} - \rho_+^{cy}(\Phi_j^{cy})) \right) \\ &+ \delta_{k>j} \beta_j \beta_k T^2 \rho_+^{cy} (\mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^* \circ \varphi_j \circ \mathcal{L}_{j+1}^* \circ \dots \circ \mathcal{L}_{k-1}^* (\Phi_k)) \\ &+ \delta_{j>k} \beta_j \beta_k T^2 \rho_+^{cy} (\mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{k-1}^* \circ \varphi_k \circ \mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_{j-1}^* (\Phi_j)) \\ &+ \delta_{jk} \beta_j^2 T \rho_+^{cy} (\mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^* (\mathcal{D}_j(H'_S, H'_S))). \end{aligned} \quad (5.16)$$

and

$$\begin{aligned} \frac{\partial^2 r^{ra}(\alpha)}{\partial \alpha_k \partial \alpha_j} \Big|_{\alpha=0} &= \beta_j \beta_k \left(\frac{T}{M} \right)^2 \sum_{n=0}^{\infty} \rho_+^{cy} \left(\varphi_j \circ \mathcal{L}_{ra}^{*n} (\Phi_k - \rho_+^{ra}(\Phi_k)) + \varphi_k \circ \mathcal{L}_{ra}^{*n} (\Phi_j - \rho_+^{ra}(\Phi_j)) \right) \\ &+ \delta_{jk} \beta_j^2 \frac{T}{M} \rho_+^{ra} (\mathcal{D}_j(H'_S, H'_S)). \end{aligned} \quad (5.17)$$

Remark 5.5. In the above theorem all the quantities depend on the vector ζ of thermal forces and all the results hold for any value of ζ . We have not mentioned the dependence on ζ to make the formula not too heavy.

Remark 5.6. The prefactors T and T^2 , resp. $\frac{T}{M}$ and $\frac{T^2}{M^2}$, in (5.15)-(5.16), resp. (5.15) and (5.17), are due to the fact that the definitions of $r_n^\sharp(\boldsymbol{\alpha})$ correspond to the variation of entropy fluxes per interaction or cycle and not per unit time.

Remark 5.7. Of course (5.15) is in agreement with (5.2).

Remark 5.8. As mentioned at the end of the previous section (4.8) is a direct consequence of the translation symmetry (5.12) combined with (5.15).

Finally, we recall that the Green-Kubo formula (4.14)-(4.15) can also be obtained via the moment generating function $\tilde{r}(\boldsymbol{\alpha})$, see e.g. [JPR, JPW]. Indeed, we infer from (5.15) and (5.8) that

$$\frac{\partial^2 \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = \frac{T}{M} L_{jk}^{ra}, \quad \frac{\partial^2 \tilde{r}_\zeta^{cy}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = T L_{jk}^{cy}, \quad \frac{\partial^2 \tilde{r}_\zeta^{rcy}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = T L_{jk}^{rcy}.$$

If moreover **(TRI)** holds, then Evans-Searles symmetry gives for any $\boldsymbol{\alpha}, \boldsymbol{\zeta}$

$$\tilde{r}_\zeta^{rcy}(\boldsymbol{\alpha}) = \tilde{r}_\zeta^{cy}(\beta_{ref} \mathbf{1} - \boldsymbol{\zeta} - \boldsymbol{\alpha}), \quad \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha}) = \tilde{r}_\zeta^{ra}(\beta_{ref} \mathbf{1} - \boldsymbol{\zeta} - \boldsymbol{\alpha}),$$

and using translation symmetry with $\lambda = \beta_{ref}$ we get

$$\tilde{r}_\zeta^{rcy}(\boldsymbol{\alpha}) = \tilde{r}_\zeta^{cy}(-\boldsymbol{\zeta} - \boldsymbol{\alpha}), \quad \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha}) = \tilde{r}_\zeta^{ra}(-\boldsymbol{\zeta} - \boldsymbol{\alpha}).$$

Using the chain rule we therefore have

$$\frac{\partial^2 \tilde{r}_\zeta^{cy}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} + \frac{\partial^2 \tilde{r}_\zeta^{rcy}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = \frac{\partial^2 \tilde{r}_\zeta^{cy}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0}, \quad \frac{\partial^2 \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha})}{\partial \zeta_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = \frac{1}{2} \frac{\partial^2 \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0}.$$

The above computation is summarized in the following

Proposition 5.6. *If Assumptions **(ER \sharp)**, **(NE)** and **(TRI)** are satisfied then*

$$\frac{\partial^2 \tilde{r}_\zeta^{cy}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = T(L_{jk}^{cy} + L_{jk}^{rcy}) = T(L_{jk}^{cy} + L_{kj}^{cy}), \quad \frac{\partial^2 \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = \frac{2T}{M} L_{jk}^{ra},$$

where in the cyclic case the second equality follows from (4.16).

Note that similarly to what happens in continuous-time quantum dynamical semigroups [JPW], using (5.17) the above proposition gives another path to derive the Green-Kubo formula provided Assumption **(TRI)** holds. However due to the lack of global time-reversal invariance this does not allow us to retrieve (4.14) even if **(TRI)** holds. Nevertheless we have the following relations which hold without **(TRI)**.

Corollary 5.7. *If Assumptions **(ER \sharp)** and **(NE)** are satisfied, then*

$$\frac{\partial^2 \tilde{r}_\zeta^{cy}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = T(L_{jk}^{cy} + L_{kj}^{cy}), \quad \frac{\partial^2 \tilde{r}_\zeta^{ra}(\boldsymbol{\alpha})}{\partial \alpha_k \partial \alpha_j} \Big|_{\boldsymbol{\alpha}=\boldsymbol{\zeta}=0} = \frac{T}{M} (L_{jk}^{ra} + L_{kj}^{ra}).$$

Proof. Combining Remark 4.6, Eq. (4.3) and (5.14), and the fact that ρ_+ is an invariant state of all the \mathcal{L}_j 's, one easily gets that at equilibrium the various terms on the right-hand sides of (5.16)-(5.17) coincide with those in (4.14)-(4.15). \square

5.4 Fluctuation Theorem and fluctuation relation

Let $e_{n,\rho}^\sharp(\boldsymbol{\alpha})$ and $e^\sharp(\boldsymbol{\alpha})$ denote the cumulant generating functions of transient and large time entropy fluxes, i.e.

$$e_{n,\rho}^\sharp(\boldsymbol{\alpha}) := \frac{1}{\tau_\sharp} \log r_{n,\rho}^\sharp(\boldsymbol{\alpha}) = \frac{1}{\tau_\sharp} \log \mathbb{E}_{\mathbb{P}^\sharp} \left(e^{-\sum_{j=1}^M \alpha_j \mathcal{S}_{\mathcal{R}_j}^N(\sharp)} \right), \quad e^\sharp(\boldsymbol{\alpha}) := \frac{1}{\tau_\sharp} \log r^\sharp(\boldsymbol{\alpha}),$$

where $\tau_{cy} = T$ and $\tau_{ra} = \frac{T}{M}$, $\mathbb{P}^{cy} = \mathbb{P}$, $\mathbb{P}^{ra} = \hat{\mathbb{P}}$ and, with a slight abuse of notation, $\mathcal{S}_{\mathcal{R}}^N(\sharp)$ stands for $\mathcal{S}_{\mathcal{R}}^N(\mathbf{j})$ when $\sharp = ra$. Let also $I_\sharp(\boldsymbol{\varsigma}) := \sup_{\boldsymbol{\alpha} \in \mathbb{R}^M} (\boldsymbol{\alpha} \cdot \boldsymbol{\varsigma} - e^\sharp(-\boldsymbol{\alpha}))$ be the Fenchel-Legendre transform of $e^\sharp(-\boldsymbol{\alpha})$.

Theorem 5.8. *Suppose Assumption (ER $_\sharp$) holds. Then*

1) *The entropy fluxes satisfy a large deviation principle with rate function I_\sharp . Namely, for any Borel set $G \subset \mathbb{R}^M$,*

$$-\inf_{\boldsymbol{\varsigma} \in \overset{\circ}{G}} I_\sharp(\boldsymbol{\varsigma}) \leq \liminf_{N \rightarrow +\infty} \frac{1}{N\tau_\sharp} \log \mathbb{P}^\sharp \left(\frac{\mathcal{S}_{\mathcal{R}}^N(\sharp)}{N\tau_\sharp} \in \overset{\circ}{G} \right) \leq \limsup_{N \rightarrow +\infty} \frac{1}{N\tau_\sharp} \log \mathbb{P}^\sharp \left(\frac{\mathcal{S}_{\mathcal{R}}^N(\sharp)}{N\tau_\sharp} \in \overline{G} \right) \leq -\inf_{\boldsymbol{\varsigma} \in \overline{G}} I_\sharp(\boldsymbol{\varsigma}). \quad (5.18)$$

2) $\boldsymbol{\varsigma} \mapsto I_\sharp(\boldsymbol{\varsigma}) \in [0, +\infty]$ *is closed convex, with compact level sets and* $\inf_{\boldsymbol{\varsigma} \in \mathbb{R}^M} I_\sharp(\boldsymbol{\varsigma}) = 0$.

3) *The sequence of random vectors* $\left(\frac{\mathcal{S}_{\mathcal{R}_j}^N(\sharp)}{N\tau_\sharp} \right)_{N \in \mathbb{N}^*}$ *converges in probability and exponentially fast to* $\mathcal{S}_+(\sharp) := (-\beta_1 \rho_+^\sharp(\Phi_1^\sharp), \dots, -\beta_M \rho_+^\sharp(\Phi_M^\sharp))$. *Namely,*

$$\forall \epsilon > 0, \exists A(\epsilon) > 0, \forall N \in \mathbb{N}^*, \mathbb{P}^\sharp \left(\left\| \frac{\mathcal{S}_{\mathcal{R}}^N(\sharp)}{N\tau_\sharp} - \mathcal{S}_+(\sharp) \right\| > \epsilon \right) \leq e^{-NA(\epsilon)}. \quad (5.19)$$

Moreover $\frac{1}{\sqrt{N\tau_\sharp}} (\mathcal{S}_{\mathcal{R}}^N(\sharp) - \mathbb{E}(\mathcal{S}_{\mathcal{R}}^N(\sharp)))$ *converges in distribution towards a centered Gaussian* μ *whose covariance matrix is* $(\partial_{\alpha_j} \partial_{\alpha_k} e^\sharp(0))_{j,k}$.

4) *If Assumption (TRI) holds, we have the fluctuation relations*

$$\forall \boldsymbol{\varsigma} \in \mathbb{R}^M, \quad I_{cy}(-\boldsymbol{\varsigma}) - I_{cy}(\boldsymbol{\varsigma}) = I_{ra}(-\boldsymbol{\varsigma}) - I_{ra}(\boldsymbol{\varsigma}) = \sum_{j=1}^M \zeta_j, \quad (5.20)$$

where I_{rcy} is the analog of I_{cy} for the reversed-order cyclic RIS.

5) *If Assumption (NE) holds then*

$$\forall \boldsymbol{\varsigma} \in \mathbb{R}^M, \quad \sum_{j=1}^M \beta_j^{-1} \zeta_j \neq 0 \Rightarrow I_\sharp(\boldsymbol{\varsigma}) = +\infty. \quad (5.21)$$

In particular the gaussian measure μ in 3) is supported on the hyperplane $\sum_{j=1}^M \beta_j^{-1} \zeta_j = 0$.

Remark 5.9. Losely speaking the large deviation principle (5.18) can be written

$$\mathbb{P}^\sharp (\mathcal{S}_{\mathcal{R}}^N(\sharp) = \boldsymbol{\varsigma}) \approx e^{-N\tau_\sharp I_\sharp(\boldsymbol{\varsigma})},$$

as N goes to $+\infty$. So equation (5.20) can also be translated as

$$\frac{\mathbb{P}(\mathcal{S}_{\mathcal{R}}^N(cy) = -\varsigma)}{\mathbb{P}(\mathcal{S}_{\mathcal{R}}^N(rcy) = \varsigma)} \approx e^{-NT(\varsigma_1 + \dots + \varsigma_M)} \quad \text{and} \quad \frac{\hat{\mathbb{P}}(\mathcal{S}_{\mathcal{R}}^N(\mathbf{j}) = -\varsigma)}{\hat{\mathbb{P}}(\mathcal{S}_{\mathcal{R}}^N(\mathbf{j}) = \varsigma)} \approx e^{-N\frac{T}{M}(\varsigma_1 + \dots + \varsigma_M)}.$$

The latter is the original form of the fluctuation relation, see [ES, GC].

Remark 5.10. Recall from Definition 5.1 that

$$\mathcal{S}_{\mathcal{R}}^N(\#) = (\mathcal{S}_{\mathcal{R}_1}^N(\#), \dots, \mathcal{S}_{\mathcal{R}_M}^N(\#)) = (-\beta_1 \mathcal{Q}_{\mathcal{R}_1}^N(\#), \dots, -\beta_M \mathcal{Q}_{\mathcal{R}_M}^N(\#))$$

so that (5.19) is of course related to Proposition 4.3.

Remark 5.11. We would like to stress that except in point 5) the above theorem does not make use of Assumption **(NE)**. Those results concern what happens *far from equilibrium* hence there is indeed no reason that **(NE)**, which is related to the notion of equilibrium, plays any role.

Concerning 5), it is related to energy conservation (4.8). As we have argued at the end of Section 3.4.1 this explains why Assumption **(NE)** is needed here.

Obviously Theorem 5.8 has its analog for the energy flux variables $\mathcal{Q}_{\mathcal{R}}^N$. At equilibrium, 3. in the above theorem allows us to complete the linear response theory results of Section 4.3 with a Central Limit theorem on the large time behaviour of energy fluxes at equilibrium. Namely we have the following

Theorem 5.9 (Fluctuation-Dissipation). *If Assumptions **(ER $_{\#}$)**, **(NE)** and **(TRI)** hold, then the sequence of random vectors $\frac{1}{\sqrt{N\tau_{\#}}} (\mathcal{Q}_{\mathcal{R}}^N(\#) - \mathbb{E}(\mathcal{Q}_{\mathcal{R}}^N(\#)))_{N \in \mathbb{N}^*}$ converges in distribution to a centered Gaussian whose covariance matrix $(D_{jk}^{\#})_{1 \leq j, k \leq M}$ is given by $D_{jk}^{cy} = L_{jk}^{cy} + L_{jk}^{rcy}$ and $D_{jk}^{ra} = 2L_{jk}^{ra}$.*

Proof. Theorem 5.8 gives the convergence with $D_{jk}^{\#} = \partial_{\alpha_j \alpha_k}^2 \tilde{e}^{\#}(0)$ where $\tilde{e}^{\#}(\boldsymbol{\alpha}) = \frac{1}{\tau_{\#}} \log \tilde{r}^{\#}(\boldsymbol{\alpha})$. The result follows using Corollary 5.7 and the fact that at equilibrium $\partial_{\alpha_j} \tilde{r}(0) = \tau_{\#} \rho_{+}^{\#}(\Phi_j^{\#}) = 0$. \square

6 Proofs of the main results

6.1 Proof of Theorem 4.4

This proof is inspired by the one given in [LS] (see also [JPW]) for open quantum systems interacting with thermal reservoirs in the Van Hove weak coupling limit. In order to simplify the notation, all the quantities without any ζ or ζ parameter should be understood at equilibrium $\zeta = 0$.

1)-2) The differentiability of $\rho_{+, \zeta}^{\#}$ follows from Proposition 3.2 and the one of $\Phi_{j, \zeta}^{\#}$ is clear from its definition. Thus we have

$$L_{jk}^{\#} = \partial_{\zeta_k} \rho_{+, \zeta}^{\#} \Big|_{\zeta=0}(\Phi_j^{\#}) + \rho_{+} \left(\partial_{\zeta_k} \Phi_{j, \zeta}^{\#} \Big|_{\zeta=0} \right). \quad (6.1)$$

Since $\mathcal{L}_{\sharp, \zeta}(\rho_{+, \zeta}^{\sharp}) = \rho_{+, \zeta}^{\sharp}$ for any ζ we get $(\mathbb{1} - \mathcal{L}_{\sharp, \zeta})(\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp}) = (\partial_{\zeta_k} \mathcal{L}_{\sharp, \zeta})(\rho_{+, \zeta}^{\sharp})$, hence

$$(\mathbb{1} - \mathcal{L}_{\sharp, \zeta}^N)(\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp}) = \sum_{n=0}^{N-1} \mathcal{L}_{\sharp, \zeta}^n \circ \partial_{\zeta_k} \mathcal{L}_{\sharp, \zeta}(\rho_{+, \zeta}^{\sharp}), \quad \forall N \geq 1. \quad (6.2)$$

Using Proposition 3.2 we have $\lim_{N \rightarrow \infty} \mathcal{L}_{\sharp, \zeta}^N(\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp}) = \text{Tr}(\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp}) \rho_{+, \zeta}^{\sharp} = 0$ where we have used that $\text{Tr}(\rho_{+, \zeta}^{\sharp}) = 1$ for all ζ so that $\text{Tr}(\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp}) = 0$. Letting $N \rightarrow \infty$ in (6.2) we get

$$\partial_{\zeta_k} \rho_{+, \zeta}^{\sharp} \Big|_{\zeta=0} = \sum_{n=0}^{+\infty} \mathcal{L}_{\sharp}^n \circ \partial_{\zeta_k} \mathcal{L}_{\sharp, \zeta} \Big|_{\zeta=0}(\rho_+), \text{ and (6.1) becomes}$$

$$L_{jk}^{\sharp} = \sum_{n=0}^{+\infty} \rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{\sharp, \zeta}^* \Big|_{\zeta=0} \circ \mathcal{L}_{\sharp}^{*n}(\Phi_j^{\sharp}) \right) + \rho_+ \left(\partial_{\zeta_k} \Phi_{j, \zeta}^{\sharp} \Big|_{\zeta=0} \right). \quad (6.3)$$

Lemma 6.1. *For all k and $X \in \mathcal{B}(\mathcal{H}_S)$ one has $\rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{k, \zeta_k}^* \Big|_{\zeta_k=0}(X) \right) = T \rho_+ (X \Phi_{k, \text{rev}})$.*

Proof. Using $\partial_{\zeta_k} \rho_{\mathcal{E}_k, \zeta_k} = H_{\mathcal{E}_k} \rho_{\mathcal{E}_k, \zeta_k} - \text{Tr}(H_{\mathcal{E}_k} \rho_{\mathcal{E}_k, \zeta_k}) \rho_{\mathcal{E}_k, \zeta_k}$ and the definition of \mathcal{L}_{k, ζ_k} we have

$$\begin{aligned} & \rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{k, \zeta_k}^* \Big|_{\zeta_k=0}(X) \right) \\ &= \text{Tr}(\mathbb{1} \otimes H_{\mathcal{E}_k} \times \rho_+ \otimes \rho_{\mathcal{E}_k} U_k^* X \otimes \mathbb{1} U_k) - \text{Tr}(\rho_+ \otimes \rho_{\mathcal{E}_k} U_k^* X \otimes \mathbb{1} U_k) \times \text{Tr}(H_{\mathcal{E}_k} \rho_{\mathcal{E}_k}) \\ &= \text{Tr}(U_k \mathbb{1} \otimes H_{\mathcal{E}_k} U_k^* \times \rho_+ X \otimes \rho_{\mathcal{E}_k}) - \text{Tr}(\rho_+ X) \times \text{Tr}(H_{\mathcal{E}_k} \rho_{\mathcal{E}_k}) \\ &= \text{Tr}((U_k \mathbb{1} \otimes H_{\mathcal{E}_k} U_k^* - \mathbb{1} \otimes H_{\mathcal{E}_k}) \times \rho_+ X \otimes \rho_{\mathcal{E}_k}) \\ &= T \text{Tr}(\Phi_{k, \text{rev}} \rho_+ X), \end{aligned}$$

where we have used the cyclicity of the trace and Assumption **(NE)** in line 2. \square

Corollary 6.2. *For all $k = 1, \dots, M$ and $X \in \mathcal{B}(\mathcal{H}_S)$ one has*

$$\rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{cy, \zeta}^* \Big|_{\zeta=0}(X) \right) = T \rho_+ \left(\mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_M^*(X) \Phi_{k, \text{rev}} \right) = T \rho_+ \left(X \Phi_{k, \text{rev}}^{cy} \right), \quad (6.4)$$

$$\rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{ra, \zeta}^* \Big|_{\zeta=0}(X) \right) = \frac{T}{M} \rho_+ \left(X \Phi_{k, \text{rev}} \right). \quad (6.5)$$

Proof. Eq. (3.3) gives $\partial_{\zeta_k} \mathcal{L}_{cy, \zeta}^* \Big|_{\zeta=0} = \mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{k-1}^* \circ \partial_{\zeta_k} \mathcal{L}_{k, \zeta_k}^* \Big|_{\zeta_k=0} \circ \mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_M^*$. The first equality in (6.4) then follows from Lemma 6.1 and the fact that ρ_+ is a joint invariant state of \mathcal{L}_j , $j = 1, \dots, M$, while the second equality follows from (4.3) and Definition 4.2.

The second identity is immediate by definition of $\mathcal{L}_{ra, \zeta}$. \square

Inserting (6.4)-(6.5) into (6.3) leads to the infinite sums in (4.14)-(4.15)-(4.17). It thus remains to compute $\rho_+ \left(\partial_{\zeta_k} \Phi_{j, \zeta}^{\sharp} \Big|_{\zeta=0} \right)$. For $\sharp = ra$ it follows from (4.6) and Lemma 6.1 that

$$\rho_+ \left(\partial_{\zeta_k} \Phi_{j, \zeta} \Big|_{\zeta=0} \right) = \delta_{jk} \frac{1}{T} \rho_+ \left(\partial_{\zeta_j} \mathcal{L}_{j, \zeta_j}^* \Big|_{\zeta_j=0}(H'_S) \right) = \delta_{jk} \rho_+ \left(H'_S \Phi_{j, \text{rev}} \right),$$

while for $\sharp = cy$, using moreover (4.1) and the fact that ρ_+ is a joint invariant state of the \mathcal{L}_j 's, we have for $k < j$

$$\begin{aligned} \rho_+ \left(\partial_{\zeta_k} \Phi_{j, \zeta}^{cy} \Big|_{\zeta=0} \right) &= \rho_+ \left(\partial_{\zeta_k} \mathcal{L}_{k, \zeta_k}^* \Big|_{\zeta_k=0} \circ \mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_{j-1}^*(\Phi_j) \right) \\ &= T \rho_+ \left(\mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_{j-1}^*(\Phi_j) \times \Phi_{k, \text{rev}} \right), \end{aligned}$$

while $\rho_+ \left(\partial_{\zeta_k} \Phi_{j,\zeta}^{cy} \Big|_{\zeta=0} \right) = 0$ for $k > j$ and

$$\rho_+ \left(\partial_{\zeta_j} \Phi_{j,\zeta}^{cy} \Big|_{\zeta=0} \right) = \frac{1}{T} \rho_+ \left(\partial_{\zeta_j} \mathcal{L}_{j,\zeta_j}^* \Big|_{\zeta_j=0} (H'_S) \right) = \rho_+ (H'_S \Phi_{j,\text{rev}}).$$

Equations (4.14)-(4.15) finally follow because $\rho_+ (H'_S \Phi_{j,\text{rev}}) = \frac{1}{2} \rho_+ (\mathcal{D}_j(H'_S, H'_S))$. Indeed,

$$\begin{aligned} \rho_+ (H'_S \Phi_{j,\text{rev}}) &= \frac{1}{T} \rho_+ \left(H_S'^2 - H'_S \mathcal{L}_{j,\text{rev}}^* (H'_S) \right) \\ &= \frac{1}{T} \rho_+ \left(H_S'^2 - \mathcal{L}_j^* (H'_S) H'_S \right) \\ &= \frac{1}{2T} \rho_+ \left(H_S'^2 + \mathcal{L}_j^* (H_S'^2) - \mathcal{L}_j^* (H'_S) H'_S - H'_S \mathcal{L}_j^* (H'_S) \right) \\ &= \frac{1}{2} \rho_+ (\mathcal{D}_j(H'_S, H'_S)), \end{aligned}$$

where we have used (4.7) in line 1, (4.3) in line 2, and in line 3 that at equilibrium ρ_+ is \mathcal{L}_j invariant and $[\mathcal{L}_j^* (H'_S), H'_S] = 0$, see Remark 3.5.

3) We now prove the Onsager relations (4.16). We thus now assume that Assumption **(TRI)** also holds. It follows from (3.12) that $\Theta \circ \mathcal{L}_{ra}^* \circ \Theta = \mathcal{L}_{ra}^{*\rho_+}$, and $\rho_+ (\Theta(X)) = \rho_+ (X^*)$ for any $X \in \mathcal{B}(\mathcal{H}_S)$ because ρ_+ is Θ invariant. Hence using (4.5) we get for all n

$$\begin{aligned} \rho_+ (\mathcal{L}_{ra}^{*n} (\Phi_j) \Phi_{k,\text{rev}}) &= \rho_+ \left(\Theta (\Phi_{k,\text{rev}}) \times (\mathcal{L}_{ra}^{*\rho_+})^n (\Theta (\Phi_j)) \right) \\ &= \rho_+ \left(\Phi_k \times (\mathcal{L}_{ra}^{*\rho_+})^n (\Phi_{j,\text{rev}}) \right) \\ &= \rho_+ (\mathcal{L}_{ra}^{*n} (\Phi_k) \Phi_{j,\text{rev}}), \end{aligned} \tag{6.6}$$

from which $L_{jk}^{ra} = L_{kj}^{ra}$ follows.

Similarly, let

$$\mathcal{L}_{rcy} := \mathcal{L}_1 \circ \mathcal{L}_2 \circ \dots \circ \mathcal{L}_M, \quad \Phi_j^{rcy} := \frac{1}{T} \mathcal{L}_M^* \circ \dots \circ \mathcal{L}_{j+1}^* (\Phi_j) \quad \text{and} \quad \Phi_{j,\text{rev}}^{rcy} := \frac{1}{T} \mathcal{L}_{1,\text{rev}}^* \circ \dots \circ \mathcal{L}_{j-1,\text{rev}}^* (\Phi_{j,\text{rev}}),$$

denote the reversed-order analogs of \mathcal{L}_{cy} , Φ_j^{cy} and $\Phi_{j,\text{rev}}^{cy}$. The associated kinetic coefficients L_{jk}^{rcy} are thus given by

$$\begin{aligned} L_{jk}^{rcy} &= T \sum_{n=0}^{+\infty} \rho_+ (\mathcal{L}_{rcy}^{*n} (\Phi_j^{rcy}) \Phi_{k,\text{rev}}^{rcy}) \\ &\quad + \delta_{k>j} T \rho_+ \left(\mathcal{L}_{k-1}^* \circ \dots \circ \mathcal{L}_{j+1}^* (\Phi_j) \times \Phi_{k,\text{rev}} \right) + \frac{1}{2} \delta_{jk} \rho_+ (\mathcal{D}_j(H'_S, H'_S)), \end{aligned}$$

It follows from Lemma 3.12 and (4.4)-(4.5) that

$$\Theta \circ \mathcal{L}_{cy}^* \circ \Theta = \mathcal{L}_{rcy}^{*\rho_+}, \quad \Theta \left(\Phi_j^{cy} \right) = -\Phi_{j,\text{rev}}^{rcy} \quad \text{and} \quad \Theta \left(\Phi_{k,\text{rev}}^{cy} \right) = -\Phi_k^{rcy},$$

and the same reasoning than in (6.6) shows that $L_{jk}^{cy} = L_{kj}^{rcy}$.

6.2 Proof of Proposition 5.2

The idea of the proof is close to the one of an analogous statement in [vHG]. The main difference comes from the fact that we consider a double measurement with each probe while they only have a single post-interaction measurement.

Since $[H_{\mathcal{E}_j}, H_{\mathcal{E}_k}] = [H_{\mathcal{E}_j}, H_k] = 0$ for $j \neq k$, we have

$$\mathcal{U}_N^s(\mathbf{j}) = \Pi_{s'_N} \left(S_{\mathcal{E}_{j_N}}^N \right) \cdots \Pi_{s'_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \times U_N \cdots U_1 \times \Pi_{s_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \cdots \Pi_{s_N} \left(S_{\mathcal{E}_{j_N}}^N \right),$$

so that (5.3) becomes

$$\begin{aligned} & \mathbb{P}(S_{\mathcal{R}}^N(\mathbf{j}) = \varsigma) \\ &= \sum \text{Tr} \left(\Pi_{s'_N} \left(S_{\mathcal{E}_{j_N}}^N \right) \cdots \Pi_{s'_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \times U_N \cdots U_1 \times \Pi_{s_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \cdots \Pi_{s_N} \left(S_{\mathcal{E}_{j_N}}^N \right)^N \times \rho_{tot}^N(\mathbf{j}) \times U_1^* \cdots U_N^* \right), \end{aligned}$$

where the sum is as in (5.3). Hence

$$\begin{aligned} & r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha}) \\ &= \sum_{s_1, \dots, s_M} e^{-(\alpha_1 s_1 + \dots + \alpha_M s_M)} \mathbb{P}(S_{\mathcal{R}}^N(\mathbf{j}) = \varsigma) \\ &= \sum_{\mathbf{s}} \prod_{n=1}^N e^{-\alpha_{j_n} (s'_n - s_n)} \text{Tr} \left(\Pi_{s'_N} \left(S_{\mathcal{E}_{j_N}}^N \right) \cdots \Pi_{s'_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \times U_N \cdots U_1 \times \Pi_{s_1} \left(S_{\mathcal{E}_{j_1}}^1 \right) \cdots \Pi_{s_N} \left(S_{\mathcal{E}_{j_N}}^N \right) \right. \\ & \quad \left. \times \rho_{tot}^N(\mathbf{j}) \times U_1^* \cdots U_N^* \right) \\ &= \text{Tr} \left(e^{-\alpha_{j_N} S_{\mathcal{E}_{j_N}}^N} \cdots e^{-\alpha_{j_1} S_{\mathcal{E}_{j_1}}^1} \times U_N \cdots U_1 \times e^{\alpha_{j_1} S_{\mathcal{E}_{j_1}}^1} \cdots e^{\alpha_{j_N} S_{\mathcal{E}_{j_N}}^N} \times \rho_{tot}^N(\mathbf{j}) \times U_1^* \cdots U_N^* \right). \end{aligned}$$

Now, recall that $\rho_{tot}^N(\mathbf{j}) = \rho \otimes \rho_{\mathcal{E}_{j_1}} \otimes \cdots \otimes \rho_{\mathcal{E}_{j_N}}$ where $\rho_{\mathcal{E}_{j_n}} = \frac{e^{-S_{\mathcal{E}_{j_n}}}}{\text{Tr} \left(e^{-S_{\mathcal{E}_{j_n}}} \right)}$ for any n , and that

the $e^{-S_{\mathcal{E}_{j_n}}}$'s act on different probes. Hence we get

$$r_{N,\rho}^{\mathbf{j}}(\boldsymbol{\alpha}) = \text{Tr} \left(\mathbb{1}_S \otimes \rho_{\mathcal{E}_{j_1}}^{\alpha_{j_1}} \otimes \cdots \otimes \rho_{\mathcal{E}_{j_N}}^{\alpha_{j_N}} \times U_N \cdots U_1 \times \rho \otimes \rho_{\mathcal{E}_{j_1}}^{1-\alpha_{j_1}} \otimes \cdots \otimes \rho_{\mathcal{E}_{j_N}}^{1-\alpha_{j_N}} \times U_1^* \cdots U_N^* \right),$$

and (5.6) follows exactly in the same way than (3.2) follows from (3.1).

6.3 Proof of Theorem 5.4

1) The proof is standard, see e.g. [HMO, JPW, vHG], so we only briefly sketch it. Since $\mathcal{L}_{\sharp, \zeta}^{[\boldsymbol{\alpha}]^*}$ is completely positive, and primitive by Proposition 5.3, it follows from Perron-Frobenius theory for completely positive maps [EHK] that $r_{\zeta}^{\sharp}(\boldsymbol{\alpha})$ is a simple dominant eigenvalue with positive definite left and right eigenvectors and (5.9) follows then from (5.7). Indeed there exists $\gamma > 0$ (spectral gap) such that for all n one has $\left(\mathcal{L}_{\sharp}^{[\boldsymbol{\alpha}]^*} \right)^n = (r_{\zeta}^{\sharp}(\boldsymbol{\alpha}))^n |A\rangle\langle \nu| + O\left((r_{\zeta}^{\sharp}(\boldsymbol{\alpha}) - \gamma)^n \right)$ where ν and A denote the positive definite left and right eigenvectors of $\mathcal{L}_{\sharp, \zeta}^{[\boldsymbol{\alpha}]^*}$ normalized such that $\nu(A) = 1$. Thus

$$r_{n,\rho}^{\sharp}(\boldsymbol{\alpha}) = \rho \left(\left(\mathcal{L}_{\sharp}^{[\boldsymbol{\alpha}]^*} \right)^n (\mathbb{1}) \right) = (r_{\zeta}^{\sharp}(\boldsymbol{\alpha}))^n \left[\rho(A) \times \nu(\mathbb{1}) + O \left(\left(1 - \frac{\gamma}{r_{\zeta}^{\sharp}(\boldsymbol{\alpha})} \right)^n \right) \right],$$

and the result follows since both $\rho(A) > 0$ and $\nu(\mathbb{1}) > 0$.

2) The symmetry relies on the following Lemma which is a direct consequence of **(TRI)**.

Lemma 6.3. *For any $j = 1, \dots, M$ and $\zeta, \alpha \in \mathbb{R}^M$ one has $\Theta \circ \mathcal{L}_{j,\zeta}^{[\alpha]*} \circ \Theta = \mathcal{L}_{j,\zeta}^{[1-\alpha]}$. As a consequence*

$$\Theta \circ \mathcal{L}_{cy,\zeta}^{[\alpha]*} \circ \Theta = \mathcal{L}_{rcy,\zeta}^{[1-\alpha]} \quad \text{and} \quad \Theta \circ \mathcal{L}_{ra,\zeta}^{[\alpha]*} \circ \Theta = \mathcal{L}_{ra,\zeta}^{[1-\alpha]}.$$

The lemma indeed implies that the maps $\mathcal{L}_{ra,\zeta}^{[\alpha]*}$ and $\mathcal{L}_{ra,\zeta}^{[1-\alpha]}$, resp. $\mathcal{L}_{cy,\zeta}^{[\alpha]*}$ and $\mathcal{L}_{rcy,\zeta}^{[1-\alpha]}$, have the same spectral radius which proves (5.10).

Proof. Denote by Θ_j the time-reversal of \mathcal{E}_j as in Section 3.3.3. Then for any X, Y we have

$$\begin{aligned} & \text{Tr} \left(Y \times \Theta \circ \mathcal{L}_j^{[\alpha]*} \circ \Theta(X) \right) \\ &= \overline{\text{Tr} \left(\Theta(Y) \times \mathcal{L}_j^{[\alpha]*}(\Theta(X)) \right)} \\ &= \overline{\text{Tr} \left(\Theta(Y) \otimes \rho_{\mathcal{E}_j}^{1-\alpha_j} \times U_j^* \times \Theta(X) \otimes \rho_{\mathcal{E}_j}^{\alpha_j} \times U_j \right)} \\ &= \text{Tr} \left(Y \otimes \Theta_j(\rho_{\mathcal{E}_j}^{1-\alpha_j}) \times \Theta \otimes \Theta_j(U_j^*) \times X \otimes \Theta_j(\rho_{\mathcal{E}_j}^{\alpha_j}) \times \Theta \otimes \Theta_j(U_j) \right) \\ &= \text{Tr} \left(Y \otimes \rho_{\mathcal{E}_j}^{1-\alpha_j} \times U_j \times X \otimes \rho_{\mathcal{E}_j}^{\alpha_j} \times U_j^* \right) \\ &= \text{Tr} \left(Y \times \mathcal{L}_j^{[1-\alpha]}(X) \right), \end{aligned}$$

where we have used Assumption **(TRI)** in the 4-th line. \square

3) The argument is again of isospectral type and relies on the following lemma.

Lemma 6.4. *If **(ER \sharp)** and **(NE)** hold then for any $j = 1, \dots, M$, $\alpha = (\alpha_1, \dots, \alpha_M) \in \mathbb{R}^M$ and $\zeta = (\zeta_1, \dots, \zeta_M) \in \mathbb{R}^M$ and $X \in \mathcal{B}(\mathcal{H}_S)$ one has*

$$\mathcal{L}_j^{[\alpha]*}(X) = \mathcal{L}_j^* \left(X e^{\beta_j \alpha_j H'_S} \right) e^{-\beta_j \alpha_j H'_S} = e^{-\beta_j \alpha_j H'_S} \mathcal{L}_j^* \left(e^{\beta_j \alpha_j H'_S} X \right), \quad (6.7)$$

where β_j is the inverse temperature of \mathcal{E}_j , i.e. such that $\zeta_j = \beta_{ref} - \beta_j$.

Remark 6.1. Note that the mere existence of H'_S requires **(ER \sharp)** and **(NE)**.

Proof. For any $X \in \mathcal{B}(\mathcal{H}_S)$ we have

$$\begin{aligned} \mathcal{L}_j^{[\alpha]*}(X) &= \text{Tr}_{\mathcal{E}_j} \left(\mathbb{1} \otimes e^{\beta_j \alpha_j H_{\mathcal{E}_j}} \rho_{\mathcal{E}_j} \times U_j^* \times X \otimes e^{-\beta_j \alpha_j H_{\mathcal{E}_j}} \times U_j \right) \\ &= \text{Tr}_{\mathcal{E}_j} \left(\mathbb{1} \otimes \rho_{\mathcal{E}_j} \times U_j^* \times X e^{\beta_j \alpha_j H'_S} \otimes \mathbb{1} \times e^{-\beta_j \alpha_j (H'_S + H_{\mathcal{E}_j})} \times U_j \times \mathbb{1} \otimes e^{\beta_j \alpha_j H_{\mathcal{E}_j}} \right) \\ &= \text{Tr}_{\mathcal{E}_j} \left(\mathbb{1} \otimes \rho_{\mathcal{E}_j} \times U_j^* \times X e^{\beta_j \alpha_j H'_S} \otimes \mathbb{1} \times U_j \times e^{-\beta_j \alpha_j H'_S} \otimes \mathbb{1} \right) \\ &= \mathcal{L}_j^* \left(X e^{\beta_j \alpha_j H'_S} \right) e^{-\beta_j \alpha_j H'_S}, \end{aligned}$$

where we have used (3.9) in line 3. The second equality is a direct consequence of (3.11). \square

We finally use the above lemma to prove (5.11). For $\lambda \in \mathbb{R}$ let \mathcal{M}_λ denote the right-multiplication by $e^{\lambda H'_S}$ on $\mathcal{B}(\mathcal{H}_S)$. It follows from Lemma 6.4 that, for any j, α, ζ , we have $\mathcal{L}_{j,\zeta}^{[\alpha+\lambda\beta^{-1}]^*} = \mathcal{M}_\lambda^{-1} \circ \mathcal{L}_{j,\zeta}^{[\alpha]^*} \circ \mathcal{M}_\lambda$, so that

$$\mathcal{L}_{\sharp,\zeta}^{[\alpha+\lambda\beta^{-1}]^*} = \mathcal{M}_\lambda^{-1} \circ \mathcal{L}_{\sharp,\zeta}^{[\alpha]^*} \circ \mathcal{M}_\lambda, \quad \sharp = cy, ra.$$

The maps $\mathcal{L}_{\sharp,\zeta}^{[\alpha+\lambda\beta^{-1}]^*}$ and $\mathcal{L}_{\sharp,\zeta}^{[\alpha]^*}$ thus have the same spectral radius, i.e. $r_\zeta^\sharp(\alpha) = r_\zeta^\sharp(\alpha + \lambda\beta^{-1})$.

6.4 Proof of Theorem 5.5

The proof is an adaptation of the one in [JPW] for continuous time quantum dynamical semigroups. In all this section the thermodynamical parameter ζ is arbitrary but fixed, see Remark 5.5, and we shall omit it.

For $\alpha = 0$ the operator $\mathcal{L}_\sharp^{[0]^*} = \mathcal{L}_\sharp^*$ has a simple dominant eigenvalue 1. By perturbation theory there exists a small circle Γ centered at 1 such that for $\alpha \in \mathbb{R}^M$ sufficiently close to 0 the only point in $\text{sp}(\mathcal{L}_\sharp^{[\alpha]^*})$ inside or on Γ is its dominant eigenvalue $r_\sharp(\alpha)$. We shall further denote by $P_\sharp^{[\alpha]}$ the associated eigenprojection. Note that $P_\sharp^{[0]}(X) = \rho_+^\sharp(X)\mathbb{1}$.

For $n = 0, 1$ denote by $E_n^\sharp(\alpha)$ the quantity

$$E_n^\sharp(\alpha) := \oint_{\Gamma} (z-1)^n \rho_+^\sharp \left(\left(z - \mathcal{L}_\sharp^{[\alpha]^*} \right)^{-1} (\mathbb{1}) \right) \frac{dz}{2\pi i}. \quad (6.8)$$

Writing $E_n^\sharp(\alpha)$ as

$$E_n^\sharp(\alpha) := \oint_{\Gamma} (z-1)^n \rho_+^\sharp \left(\left(z - \mathcal{L}_\sharp^{[\alpha]^*} \right)^{-1} \circ (Id - P_\sharp^{[\alpha]})(\mathbb{1}) \right) \frac{dz}{2\pi i} + \oint_{\Gamma} \frac{(z-1)^n}{z - r_\sharp(\alpha)} \rho_+^\sharp \left(P_\sharp^{[\alpha]}(\mathbb{1}) \right) \frac{dz}{2\pi i},$$

the first term on the right-hand side is analytic inside Γ and it follows from Cauchy's integral formula that

$$E_0^\sharp(\alpha) = \rho_+^\sharp \left(P_\sharp^{[\alpha]}(\mathbb{1}) \right) \quad \text{and} \quad E_1^\sharp(\alpha) = (r_\sharp(\alpha) - 1) \rho_+^\sharp \left(P_\sharp^{[\alpha]}(\mathbb{1}) \right), \quad (6.9)$$

so that

$$r_\sharp(\alpha) = 1 + \frac{E_1^\sharp(\alpha)}{E_0^\sharp(\alpha)}. \quad (6.10)$$

Since $P_\sharp^{[0]}(\mathbb{1}) = \mathbb{1}$ and $r_\sharp(0) = 1$ it follows from (6.9) that $E_0^\sharp(0) = 1$, $E_1^\sharp(0) = 0$ hence $\partial_{\alpha_j} r_\sharp(0) = \partial_{\alpha_j} E_1^\sharp(0)$. Using (6.8) we get

$$\partial_{\alpha_j} E_n^\sharp(\alpha) = \oint_{\Gamma} (z-1)^n \rho_+^\sharp \left(\left(z - \mathcal{L}_\sharp^{[\alpha]^*} \right)^{-1} \circ \partial_{\alpha_j} \mathcal{L}_\sharp^{[\alpha]^*} \circ \left(z - \mathcal{L}_\sharp^{[\alpha]^*} \right)^{-1} (\mathbb{1}) \right) \frac{dz}{2\pi i}, \quad (6.11)$$

so that, because $\mathcal{L}_\sharp^*(\mathbb{1}) = \mathbb{1}$ and $\mathcal{L}_\sharp(\rho_+^\sharp) = \rho_+^\sharp$,

$$\partial_{\alpha_j} E_n^\sharp(0) = \oint_{\Gamma} (z-1)^{n-2} \rho_+^\sharp \left(\partial_{\alpha_j} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0](\mathbb{1}) \right) \frac{dz}{2\pi i},$$

hence

$$\partial_{\alpha_j} E_0^\sharp(0) = 0 \quad \text{and} \quad \partial_{\alpha_j} E_1^\sharp(0) = \rho_+^\sharp \left(\partial_{\alpha_j} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0(\mathbb{1})] \right). \quad (6.12)$$

Eq. (6.10) and (6.12) then give $\partial_{\alpha_k} \partial_{\alpha_j} r^\sharp(0) = \partial_{\alpha_k} \partial_{\alpha_j} E_1^\sharp(0)$, while from (6.11) we infer that

$$\begin{aligned} \partial_{\alpha_k} \partial_{\alpha_j} E_1^\sharp(0) &= \oint_{\Gamma} (z-1)^{-1} \rho_+^\sharp \left(\partial_{\alpha_k} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0] \circ (z - \mathcal{L}_\sharp^*)^{-1} \circ \partial_{\alpha_j} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0(\mathbb{1})] \right) \frac{dz}{2\pi i} \\ &\quad + \oint_{\Gamma} (z-1)^{-1} \rho_+^\sharp \left(\partial_{\alpha_j} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0] \circ (z - \mathcal{L}_\sharp^*)^{-1} \circ \partial_{\alpha_k} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0(\mathbb{1})] \right) \frac{dz}{2\pi i} \\ &\quad + \oint_{\Gamma} (z-1)^{-1} \rho_+^\sharp \left(\partial_{\alpha_j} \partial_{\alpha_k} \mathcal{L}_\sharp^{[\alpha]^*}[\alpha=0(\mathbb{1})] \right) \frac{dz}{2\pi i} \\ &=: I + II + III \end{aligned} \quad (6.13)$$

The next lemma is the main technical ingredient leading from (6.12)-(6.13) to (5.15)-(5.17).

Lemma 6.5. *For any $j, k = 1, \dots, M$ one has*

$$\partial_{\alpha_j} \mathcal{L}_{ra}^{[\alpha]^*}[\alpha=0(X)] = \beta_j \frac{T}{M} \varphi_j(X), \quad \partial_{\alpha_j} \mathcal{L}_{cy}^{[\alpha]^*}[\alpha=0(X)] = \beta_j T \varphi_j^{cy}(X), \quad (6.14)$$

$$\partial_{\alpha_k} \partial_{\alpha_j} \mathcal{L}_{ra}^{[\alpha]^*}[\alpha=0(\mathbb{1})] = \delta_{jk} \beta_j^2 \frac{T}{M} \mathcal{D}_j(H'_S, H'_S), \quad (6.15)$$

and

$$\partial_{\alpha_k} \partial_{\alpha_j} \mathcal{L}_{cy}^{[\alpha]^*}[\alpha=0(\mathbb{1})] = \begin{cases} \beta_j \beta_k T^2 \mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^* \circ \varphi_j \circ \mathcal{L}_{j+1}^* \circ \dots \circ \mathcal{L}_{k-1}^*(\Phi_k), & \text{if } k > j, \\ \beta_j \beta_k T^2 \mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{k-1}^* \circ \varphi_k \circ \mathcal{L}_{k+1}^* \circ \dots \circ \mathcal{L}_{j-1}^*(\Phi_j), & \text{if } k < j, \\ \beta_j^2 T \mathcal{L}_1^* \circ \dots \circ \mathcal{L}_{j-1}^*(\mathcal{D}_j(H'_S, H'_S)), & \text{if } j = k. \end{cases} \quad (6.16)$$

Proof. Using Lemma 6.4, by straightforward calculation we get

$$\partial_{\alpha_j} \mathcal{L}_j^{[\alpha]^*}(X) = \beta_j \left(\mathcal{L}_j^{[\alpha]^*}(X H'_S) - \mathcal{L}_j^{[\alpha]^*}(X) H'_S \right)$$

and

$$\partial_{\alpha_k} \partial_{\alpha_j} \mathcal{L}_j^{[\alpha]^*}(X) = \delta_{jk} \beta_j^2 \left(\mathcal{L}_j^{[\alpha]^*}(X H_S'^2) - 2 \mathcal{L}_j^{[\alpha]^*}(X H'_S) H'_S + \mathcal{L}_j^{[\alpha]^*}(X) H_S'^2 \right).$$

Using the definitions (5.13) of $\varphi_j(X)$ and (4.12) of $\mathcal{D}_j(X, Y)$ we thus obtain

$$\partial_{\alpha_j} \mathcal{L}_j^{[\alpha]^*}[\alpha=0(X)] = \beta_j T \varphi_j(X) \quad \text{and} \quad \partial_{\alpha_k} \partial_{\alpha_j} \mathcal{L}_j^{[\alpha]^*}[\alpha=0(\mathbb{1})] = \delta_{jk} \beta_j^2 T \mathcal{D}_j(H'_S, H'_S), \quad (6.17)$$

where we have used that $[\mathcal{L}_j^*(H'_S), H'_S] = 0$ in the second identity, see Remark 3.5.

Finally (6.14)-(6.16) follow directly from (6.17) and (5.5). \square

End of the proof of (5.15)-(5.17).

Since $\partial_{\alpha_j} r^\sharp(0) = \partial_{\alpha_j} E_1^\sharp(0)$ combining (6.12) and (6.14), recall also Remark 5.4, we get (5.15).

We now turn to (5.16)-(5.17). Using (6.14), the first term on the right-hand side of (6.13) becomes

$$I = \beta_j \beta_k \tau_{\sharp}^2 \oint_{\Gamma} (z-1)^{-1} \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ (z - \mathcal{L}_{\sharp}^*)^{-1} \left(\Phi_j^{\sharp} \right) \right) \frac{dz}{2\pi i},$$

where $\tau_{cy} = T$ and $\tau_{ra} = \frac{T}{M}$. Recall $P_{\sharp}^{[0]}$ denotes the eigenprojection of \mathcal{L}_{\sharp}^* associated to the eigenvalue 1. One has

$$\oint_{\Gamma} (z-1)^{-1} \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ (z - \mathcal{L}_{\sharp}^*)^{-1} \circ P_{\sharp}^{[0]} \left(\Phi_j^{\sharp} \right) \right) \frac{dz}{2\pi i} = \oint_{\Gamma} (z-1)^{-2} \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ P_{\sharp}^{[0]} \left(\Phi_j^{\sharp} \right) \right) \frac{dz}{2\pi i} = 0.$$

Hence

$$\begin{aligned} I &= \beta_j \beta_k \tau_{\sharp}^2 \oint_{\Gamma} (z-1)^{-1} \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ (z - \mathcal{L}_{\sharp}^*)^{-1} \circ (Id - P_{\sharp}^{[0]}) \left(\Phi_j^{\sharp} \right) \right) \frac{dz}{2\pi i} \\ &= \beta_j \beta_k \tau_{\sharp}^2 \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ (1 - \mathcal{L}_{\sharp}^*)^{-1} \circ (Id - P_{\sharp}^{[0]}) \left(\Phi_j^{\sharp} \right) \right), \end{aligned}$$

where we have used that $(z - \mathcal{L}_{\sharp}^*)^{-1} \circ (Id - P_{\sharp}^{[0]})$ is regular at $z = 1$. Moreover, since the spectral radius of \mathcal{L}_{\sharp}^* restricted to $\text{Ran}(Id - P_{\sharp}^{[0]})$ is strictly less than one due to Assumption **(ER)_{\sharp}** we can write

$$(1 - \mathcal{L}_{\sharp}^*)^{-1} \circ (Id - P_{\sharp}^{[0]}) \left(\Phi_j^{\sharp} \right) = \sum_{n=0}^{\infty} \mathcal{L}_{\sharp}^{*n} \left((Id - P_{\sharp}^{[0]}) \left(\Phi_j^{\sharp} \right) \right) = \sum_{n=0}^{\infty} \mathcal{L}_{\sharp}^{*n} \left(\Phi_j^{\sharp} - \rho_{+}^{\sharp} \left(\Phi_j^{\sharp} \right) \right),$$

so that

$$I = \beta_j \beta_k \tau_{\sharp}^2 \sum_{n=0}^{\infty} \rho_{+}^{\sharp} \left(\varphi_k^{\sharp} \circ \mathcal{L}_{\sharp}^{*n} \left(\Phi_j^{\sharp} - \rho_{+}^{\sharp} \left(\Phi_j^{\sharp} \right) \right) \right).$$

Proceeding in the same way with the second term II in (6.13) we obtain that $I + II$ indeed corresponds to the infinite sums in (5.16)-(5.17).

Finally, the third term on the right-hand side of (6.13) is $III = \rho_{+}^{\sharp} \left(\partial_{\alpha_j} \partial_{\alpha_k} \mathcal{L}_{\sharp}^{[\alpha]^*} [\alpha=0](\mathbb{1}) \right)$, and using (6.15)-(6.16) it is easy to see that it leads to the remaining terms in (5.16)-(5.17).

6.5 Proof of Theorem 5.8

The proof of 1)-3) is a direct application of the Gartner-Ellis and Bryc Theorems, see e.g. [DZ, El, Bry], and the following lemma, while (5.20) and (5.21) are direct consequences of (5.10) and (5.11) respectively.

Lemma 6.6. *Suppose Assumption **(ER)_{\sharp}** holds. Then $e^{\sharp}(\alpha)$ is a well defined real analytic function on \mathbb{R}^M and for any $\alpha \in \mathbb{R}^M$ and initial state ρ one has*

$$\lim_{n \rightarrow \infty} \frac{1}{n} e_{n,\rho}^{\sharp}(\alpha) = e^{\sharp}(\alpha). \quad (6.18)$$

Moreover there exists a neighborhood B of 0 in \mathbb{C}^M on which (6.18) holds.

The proof is standard, it is similar to the one given for similar results in e.g. [JPW, vHG]. We briefly sketch it for the convenience of the reader.

Proof. The maps $\alpha \mapsto \mathcal{L}_{\sharp}^{[\alpha]*}$ are clearly analytic on \mathbb{C}^M . Then Assumption **(ER $_{\sharp}$)** and Proposition 5.3 guarantee that, for any $\alpha \in \mathbb{R}^M$, $r^{\sharp}(\alpha)$ is a positive isolated simple eigenvalue of $\mathcal{L}_{\sharp}^{[\alpha]*}$. Hence regular perturbation theory, see e.g. [Ka], ensures that $r^{\sharp}(\alpha)$ defines a real analytic map on \mathbb{R}^M with positive values. This proves that $e^{\sharp}(\alpha) = \frac{1}{\tau_{\sharp}} \log r^{\sharp}(\alpha)$ is well defined and real analytic. Eq. (6.18) then follows from 1) in Theorem 5.4.

The extension to a complex neighborhood of 0 follows also by a standard perturbation theory argument, see also [JPW]. Indeed, $\mathcal{L}_{\sharp}^{[0]*}$ has a simple dominant eigenvalue $r^{\sharp}(0) = 1$ so there exists $\delta, \varepsilon > 0$ such that $\text{sp} \left(\mathcal{L}_{\sharp}^{[\alpha]*} \right) \setminus \{r^{\sharp}(\alpha)\} \subset \{z \in \mathbb{C} \mid |z| < |r^{\sharp}(\alpha)| - \delta\}$ for any $\alpha \in \mathbb{C}^M$, $|\alpha| < \varepsilon$. Hence for such α 's we have

$$r_{n,\rho}^{\sharp}(\alpha) = \rho \left(\left(\mathcal{L}_{\sharp}^{[\alpha]*} \right)^n (\mathbb{1}) \right) = r^{\sharp}(\alpha)^n \left[\rho \left(P_{\sharp}^{[\alpha]}(\mathbb{1}) \right) + O \left(\left(1 - \frac{\delta}{|r^{\sharp}(\alpha)|} \right)^n \right) \right],$$

where $P_{\sharp}^{[\alpha]}$ denotes the eigenprojection of $\mathcal{L}_{\sharp}^{[\alpha]*}$ associated to $r^{\sharp}(\alpha)$. In particular $P_{\sharp}^{[\alpha]}(\mathbb{1}) = P_{\sharp}^{[0]}(\mathbb{1}) + O(\varepsilon) = \mathbb{1} + O(\varepsilon)$, see Section 6.4, so that for all n

$$r_{n,\rho}^{\sharp}(\alpha) = r^{\sharp}(\alpha)^n \left[1 + O(\varepsilon) + O \left(\left(1 - \frac{\delta}{|r^{\sharp}(\alpha)|} \right)^n \right) \right]$$

which proves the result. □

References

- [Ag] Agarwal G.S.: Open Quantum Markovian Systems and the Microreversibility, *Z. Phys.* **258**, 409–422 (1973).
- [AGMT] Andrieux D., Gaspard P., Monnai T. and Tasaki S.: The fluctuation theorem for currents in open quantum systems, *New Journal of Physics* **11**, vol 4, 043014 (2009).
- [AJ] Attal S. and Joye A.: Weak coupling and continuous limits for repeated quantum interactions, *J. Stat. Phys.* **126**, 1241–1283 (2007).
- [AP] Attal S. and Pautrat, Y.: From repeated to continuous quantum interactions, *Ann. Henri Poincaré* **7**, 59–104 (2006).
- [BFFS] Ballesteros M., Fraas M., Fröhlich J. and Schubnel B.: Indirect acquisition of information in quantum mechanics, *J. Stat. Phys.* **162**, 924–958 (2016).
- [BBB] Bauer M., Bernard D. et Benoist T.: Repeated Quantum Non-Demolition Measurements: Convergence and Continuous Time Limit, *Ann. Henri Poincaré* **14**, 639–679 (2013).

- [BJPPP] Benoist T., Jakšić V., Panati A., Pautrat Y. and Pillet C.-A.: Full statistics of energy conservation in two-time measurement protocols, *Phys. Rev. E* **92**, vol 3, 032115 (2015).
- [BPP] Benoist T., Panati A. and Pautrat Y.: Heat conservation and fluctuations for open quantum systems in the Two-Time Measurement Picture, *J. Stat. Phys.* **178**, 893–925 (2020).
- [BPR] Benoist T., Panati A. and Raquépas R.: Control of fluctuations and heavy tails for heat variation in the two-time measurement framework, *Ann. Henri Poincaré* **20**, 631–674 (2019).
- [Bo] Bougron J.-F.: Etude des propriétés thermodynamiques hors-équilibre des systèmes quantiques en interactions répétées, PhD Thesis (2020).
- [BJP] Bougron J.-F., Joye A. and Pillet C.-A.: Markovian Random Repeated Interaction Systems, In preparation.
- [Bru] Bruneau L.: Mixing properties of the one-atom maser, *J. Stat. Phys.* **155**, 888–908 (2014).
- [BDBP] Bruneau L., De Bièvre S. and Pillet C.-A.: Scattering induced current in a tight-binding band, *J. Math. Phys.* **52**, 022109 (2011).
- [BJM1] Bruneau L., Joye A. and Merkli M.: Asymptotics of repeated interaction quantum systems, *J. Func. Anal.* **239**, 310–344 (2006).
- [BJM2] Bruneau L., Joye A. and Merkli M.: Random repeated interaction quantum systems, *Commun. Math. Phys.* **284**, 553–581 (2008).
- [BJM3] Bruneau L., Joye A. and Merkli M.: Repeated interactions in open quantum systems, *J. Math. Phys.* **55**, 075204 (2014).
- [BP] Bruneau L. and Pillet C.-A.: Thermal relaxation of a QED cavity, *J. Stat. Phys.* **134**, vol 5-6, 1071–1095 (2009).
- [Bry] Bryc W.: A remark on the connection between the large deviation principle and the central limit theorem, *Statistics and Probability Letters* **18**, 253–256 (1993).
- [CP] Carbone R. and Pautrat Y.: Irreducible decompositions and stationary states of quantum channels, *Rep. Math. Phys.* **77**, 293–313 (2016).
- [CDG] Cohen-Tannoudji C., Dupont-Roc J. and Grynberg G.: *Atom–Photon Interactions*, Wiley, New-York (1992).
- [DdrM] Dereziński J., de Roeck W. and Maes C.: Fluctuations of quantum currents and unravelings of master equations, *J. Stat. Phys.* **131**, 341–356 (2008).
- [dR] de Roeck W.: Large deviation generating function for currents in the Pauli-Fierz model, *Rev. Math. Phys.* **21**, 549–585 (2009).

- [dRM] de Roeck W. and Maes C.: Steady state fluctuations of the dissipated heat for a quantum stochastic model, *Rev. Math. Phys.* **18**, 619–653 (2006).
- [DS] Davies E.B. and Spohn H.: Open quantum systems with time-dependent Hamiltonians and their linear response, *J. Stat. Phys.* **19**, 511–523 (1978).
- [DZ] Dembo A. and Zeitouni O.: *Large Deviations Techniques and Applications*, Springer-Verlag, New-York (1998).
- [El] Ellis R. S.: *Entropy, Large Deviations and Statistical Mechanics*, Springer-Verlag (1985).
- [EHM] Esposito M., Harbola U. and Mukamel S.: Nonequilibrium fluctuations, fluctuation theorems, and counting statistics in quantum systems, *Rev. Mod. Phys.* **81**, vol 4, 1665–1702 (2009).
- [EHK] Evans D.E. and Høegh-Krohn R.: Spectral properties of positive maps on C^* -algebras, *J. London Math. Soc.* **17**, 345–355 (1978).
- [ES] Evans D.E. and Searles D.J.: Equilibrium microstates which generate second law violating steady states, *Phys. Rev. E* **50**, 1645–1648 (1994).
- [FR] Fagnola F. and Rebolledo R.: From classical to quantum entropy production, *QP-PQ: Quantum Probab. White Noise Anal.*, World Sci. Publ. **25**, 245–261 (2010).
- [FU] Fagnola F. and Umanità V.: Detailed balance, time reversal, and generators of quantum Markov semigroups, *Mathematical Notes* **84**, 108–115 (2008).
- [FJM] Filipowicz P., Javanainen J. and Meystre P.: Theory of a microscopic maser, *Phys. Rev. A* **34**, vol 4, 3077–3087 (1986).
- [Ga] Gallavotti G.: Extension of Onsager’s reciprocity to large fields and the chaotic hypothesis *Phys. Rev. Lett.* **77**, 4334–4337 (1996).
- [GC] Gallavotti G. and Cohen E.G.D.: Dynamical ensembles in stationary states, *J. Stat. Phys.* **80**, 931–970 (1995).
- [GK] Guță M. and Kiukas J.: Equivalence classes and local asymptotic normality in system identification for quantum Markov chains, *Commun. Math. Phys.* **335**, vol. 3, 1397–1428 (2015).
- [HJPR] Hanson E., Joye A., Pautrat Y. and Raquépas R.: Landauer’s Principle in Repeated Interaction Systems, *Commun. Math. Phys.* **349**, 285–327 (2017).
- [HBR] Haroche S., Brune M. and Raimond J.-M.: Experiments with single atoms in a cavity: entanglement, Schrödinger’s cats and decoherence, *Phil. Trans. Roy. Soc.* **355**, 2367–2380 (1997).
- [HMO] Hiai F., Mosonyi M., and Ogawa T.: Large deviations and Chernoff bound for certain correlated states on a spin chain, *J. Math. Phys.* **48**, 123301 (2007).

- [vHG] van Horssen M. and Guță M.: Sanov and Central limit theorems for output statistics of quantum Markov chains, *J. Math. Phys.* **56**, 022109 (2015).
- [vHG2] van Horssen M. and Guță M.: Large Deviations, Central Limit and dynamical phase transitions in the atom maser, arXiv preprint, arXiv:1206.4956 (2012).
- [JOP1] Jakšić V., Ogata Y. and Pillet C.-A.: The Green-Kubo formula and the Onsager reciprocity relations in quantum statistical mechanics, *Commun. Math. Phys.* **265**, vol 3, 721–738 (2006).
- [JOP2] Jakšić V., Ogata Y. and Pillet C.-A.: Linear response theory for thermally driven quantum open systems, *J. Stat. Phys.* **123**, vol 3, 547–569 (2006).
- [JOP3] Jakšić V., Ogata Y. and Pillet C.-A.: The Green-Kubo formula for the spin-fermion system, *Commun. Math. Phys.* **268**, vol 2, 369–401 (2006).
- [JOP4] Jakšić V., Ogata Y. and Pillet C.-A.: The Green-Kubo formula for locally interacting fermionic open systems, *Ann. Henri Poincaré* **8**, 1013–1036 (2007).
- [JOPP] Jakšić V., Ogata Y., Pautrat Y. and Pillet C.-A.: Entropic fluctuations in quantum statistical mechanics. An Introduction, *Quantum Theory from Small to Large Scales*, Lecture Notes of the Les Houches Summer School **95**, 213–410 (2010).
- [JPP] Jakšić V., Pautrat Y. and Pillet C.-A.: Central limit theorem for locally interacting Fermi gas, *Commun. Math. Phys.* **285**, 175–217 (2009).
- [JP] Jakšić V. and Pillet C.-A.: A note on the Landauer principle in quantum statistical mechanics, *J. Math. Phys.* **55**, 075210 (2014).
- [JPR] Jakšić V., Pillet C.-A. and Rey-Bellet L.: Entropic fluctuations in statistical mechanics I. Classical dynamical systems, *Nonlinearity* **24**, 699–763 (2011).
- [JPW] Jakšić V., Pillet C.-A. and Westrich M.: Entropic fluctuations of quantum dynamical semigroups, *J. Stat. Phys.* **54**, vol 1, 153–187 (2014).
- [Ka] Kato T.: *Perturbation theory for linear operators*, Springer-Verlag, Berlin (1980).
- [KFGV] Kossakowski A., Frigerio A., Gorini V. and Verri M.: Quantum detailed balance and KMS condition, *Commun. Math. Phys.* **57**, 91–110 (1977).
- [Kr] Kraus K.: *States, Effects and operations, fundamental notions of quantum theory*, Springer, Berlin (1983).
- [KM] Kümmerer B. and Maassen H.: A scattering theory for Markov chains, *Infin. Dimens. Anal. Quantum Probab. Relat. Top.* **3**, 161–176 (2000).
- [Ku] Kurchan J.: A quantum fluctuation theorem, arXiv preprint cond-mat/0007360 (2000).
- [LS] Lebowitz J. and Spohn H.: Irreversible thermodynamics for quantum systems weakly coupled to thermal reservoirs, *Adv. Chem. Phys.* **38**, 109–142 (1978).

- [Li] Lindblad G.: On the Generators of Quantum Dynamical Semigroups, *Commun. Math. Phys.* **48**, 119–130 (1976).
- [Ma] Majewski W.-A.: The detailed balance condition in quantum statistical mechanics, *J. Math. Phys.* **25**, 614–616 (1984).
- [MWM] Meschede D., Walther H. and Müller G.: One-Atom Maser, *Phys. Rev. Lett.* **54**, vol 6, 551–554 (1985).
- [NPe] Nechita I. and Pellegrini C.: Random repeated quantum interactions and random invariant states, *Prob. Th. and Rel. Fields.* **152**, 299–320 (2012).
- [Pe] Pellegrini C.: Markov Chain Approximations of Jump-Diffusion Stochastic Master Equations, *Ann. Inst. Henri Poincaré Prob. Stat.* **46**, 924–948 (2010).
- [PeP] Pellegrini C. and Petruccione F.: Non Markovian quantum repeated interactions: *J. Phys. A* **42**, 425304 (2009).
- [RBH] Raimond J.-M., Brune M. and Haroche S.: Manipulating quantum entanglement with atoms and photons in a cavity, *Rev. Mod. Phys.* **73**, 565–582 (2001).
- [Ta] Tasaki H.: Jarzynski Relations for Quantum Systems and Some Applications, arXiv preprint cond-mat/0009244 (2000).
- [WBKM] Wellens T., Buchleitner A., Kümmerer B. and Maassen H.: Quantum State Preparation via Asymptotic Completeness, *Phys. Rev. Lett.* **85**, 3361–3364 (2000).
- [Wo] Wolf M.M.: Quantum Channels and Operations: A Guided Tour, <https://www-m5.ma.tum.de/foswiki/pub/M5/Allgemeines/MichaelWolf/QChannelLecture.pdf>, Lecture notes based on a course given at the Niels-Bohr Institute (2012).