Non-equilibrium steady states: maximization of the Shannon entropy associated with the distribution of dynamical trajectories in the presence of constraints

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Abstract. Filyokov and Karpov (1967 Inzh.-Fiz. Zh. 13 624) have proposed a theory of non-equilibrium steady states in direct analogy with the theory of equilibrium states: the principle is to maximize the Shannon entropy associated with the probability distribution of dynamical trajectories in the presence of constraints, including the macroscopic current of interest, via the method of Lagrange multipliers. This maximization leads directly to the generalized Gibbs distribution for the probability distribution of dynamical trajectories, and to some fluctuation relation of the integrated current. The simplest stochastic dynamics where these ideas can be applied are discrete-time Markov chains, defined by transition probabilities $W_{i \to j}$ between configurations $i$ and $j$: instead of choosing the dynamical rules $W_{i \to j}$ a priori, one determines the transition probabilities and the associate stationary state that maximize the entropy of dynamical trajectories with the other physical constraints that one wishes to impose. We give a self-contained and unified presentation of this type of approach, both for discrete-time Markov chains and for continuous-time master equations. The obtained results are in full agreement with the Bayesian approach introduced by Evans (2004 Phys. Rev. Lett. 92 150601) under the name ‘Non-equilibrium Counterpart to detailed balance’, and with the ‘invariant quantities’ derived by Baule and Evans (2008 Phys. Rev. Lett. 101 240601), but provide a slightly different perspective via the formulation in terms of an eigenvalue problem.
Non-equilibrium steady states

Keywords: stationary states, current fluctuations, statistical inference, optimization under uncertainty

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1. **Introduction**

In contrast to the statistical physics of equilibrium, which is based on a few general principles, everybody complains that no global theory has emerged to describe out-of-equilibrium systems. A natural explanation of this fact is that the problem is ill defined for many reasons. The first reason is that there are actually very different ways to be out-of-equilibrium: a system can be ‘out-of-equilibrium’ because it has a very slow dynamics that never converges towards some stationary state, or because it is in a non-equilibrium steady state, characterized by some currents that are imposed by the environment. It is clear that in the first case, one really needs a non-trivial dynamical description, whereas in the second case, one may hope that the ‘steady state’ property could help to develop some theory generalizing the equilibrium steady state. The second reason is that even within the field of ‘non-equilibrium steady states’, very different goals have been pursued. On one hand, many theorists of statistical physics consider that the best strategy consists in solving specific models with chosen dynamical rules, in order to see if some general principles emerge in the end: in this direction, the most studied models are one-dimensional exclusion processes, for which many remarkable results have been obtained over the years (see the recent short review [1] and references therein). On the other hand, it is clear that the success of the theory of equilibrium is based on the fact that one does not specify the details of the dynamical evolution but that on the contrary, one keeps only its conservation properties (for instance the energy conservation) and then maximizes the Shannon entropy associated with the equilibrium state. So there have been various attempts to formulate general principles based on some notion of ‘entropy’ for non-equilibrium steady states. Although the principles of ‘minimum entropy production’ and of ‘maximum entropy production’ seem popular, in particular in applications to related areas such as chemistry, biology or climatology, the proposed justifications have not been widely accepted, either because they are at the level of ‘generalized thermodynamics’ rather than statistical physics, or because they are limited to ‘near-equilibrium’ conditions.

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**1. Introduction**

In contrast to the statistical physics of equilibrium, which is based on a few general principles, everybody complains that no global theory has emerged to describe out-of-equilibrium systems. A natural explanation of this fact is that the problem is ill defined for many reasons. The first reason is that there are actually very different ways to be out-of-equilibrium: a system can be ‘out-of-equilibrium’ because it has a very slow dynamics that never converges towards some stationary state, or because it is in a non-equilibrium steady state, characterized by some currents that are imposed by the environment. It is clear that in the first case, one really needs a non-trivial dynamical description, whereas in the second case, one may hope that the ‘steady state’ property could help to develop some theory generalizing the equilibrium steady state. The second reason is that even within the field of ‘non-equilibrium steady states’, very different goals have been pursued. On one hand, many theorists of statistical physics consider that the best strategy consists in solving specific models with chosen dynamical rules, in order to see if some general principles emerge in the end: in this direction, the most studied models are one-dimensional exclusion processes, for which many remarkable results have been obtained over the years (see the recent short review [1] and references therein). On the other hand, it is clear that the success of the theory of equilibrium is based on the fact that one does not specify the details of the dynamical evolution but that on the contrary, one keeps only its conservation properties (for instance the energy conservation) and then maximizes the Shannon entropy associated with the equilibrium state. So there have been various attempts to formulate general principles based on some notion of ‘entropy’ for non-equilibrium steady states. Although the principles of ‘minimum entropy production’ and of ‘maximum entropy production’ seem popular, in particular in applications to related areas such as chemistry, biology or climatology, the proposed justifications have not been widely accepted, either because they are at the level of ‘generalized thermodynamics’ rather than statistical physics, or because they are limited to ‘near-equilibrium’ conditions.
or because they involve other additional questionable assumptions (see for instance the discussions in [2]–[5] and references therein). But we feel that instead of focusing on the notion of ‘entropy production’ to formulate general principles, it is much more natural to consider the ‘Shannon entropy associated with the probability distribution of dynamical trajectories’ which we will call equivalently ‘dynamical entropy’ to have a shorter name. This notion can be introduced for any stochastic dynamics, or for any deterministic chaotic dynamics that becomes effectively ‘stochastic’, as follows: to characterize the dynamics during some time interval $[0,t]$, one considers all possible dynamical trajectories $\Omega_{[0,t]}$ and their probabilities $P(\Omega_{[0,t]})$ normalized to

$$\sum_{\Omega_{[0,t]}} P(\Omega_{[0,t]}) = 1.$$  

(1)

The Shannon entropy associated with this distribution $P(\Omega_{[0,t]})$ over dynamical trajectories reads

$$S_{\text{dyn}}(t) \equiv - \sum_{\Omega_{[0,t]}} P(\Omega_{[0,t]}) \ln P(\Omega_{[0,t]}).$$  

(2)

For many stochastic process of interest, this entropy becomes extensive in the time $t$ in the large-time limit $t \to +\infty$: it is then convenient to introduce the entropy rate per unit time, called the Kolmogorov–Sinai entropy in the context of chaotic dynamical systems (see for instance the book [6] and references therein)

$$h_{\text{KS}} \equiv \lim_{t \to +\infty} \frac{S_{\text{dyn}}(t)}{t}.$$  

(3)

More than forty years ago, Filyokov and Karpov [7]–[9] proposed to base the theory of non-equilibrium steady states on the maximization of the dynamical entropy of equation (2) in the presence of some constraints via the method of Lagrange multipliers. We are not aware of direct continuations of these works, except the very recent work of Favretti [10]. However, closely related ideas have been proposed independently by Evans [11] under the name ‘Non-equilibrium Counterpart to detailed balance’ and have been applied to shear flow [12]–[14]. Also among recent attempts to ‘prove’ the principle of maximum entropy production by Dewar [15], one can see that the starting point is precisely the maximization of the dynamical entropy of equation (2). We thus feel that if one wishes to formulate a general principle based on the maximization of some notion of entropy for non-equilibrium steady states, the dynamical entropy of equation (2) is definitely the most natural and could be the basis of some consensus.

Let us now mention various recent works that are related to some of these notions:

- For random walks on arbitrary networks, the optimization of the dynamical entropy of equation (2) leads to the ‘maximal entropy random walk’ that was introduced by Burda, Duda, Luck and Waclaw [16]: it is different from the usual random walk that maximizes the entropy only locally, whenever there exist some fluctuations in the coordination numbers of nodes of the network (but of course these two types of random walks coincide on regular networks).
- For non-equilibrium steady states, even within the point of view where the dynamical rules are given a priori, the dynamical entropy of equation (2) plays a major role in the
thermodynamics of histories’ that has been developed by Lecomte, Appert-Rolland and van Wijland [17], and that has been then applied to various glassy systems [18].

- The method of Lagrange multipliers has been much used recently for non-equilibrium quantum systems [19]–[22], but we should stress that the physical meaning is completely different: in these quantum studies, the dynamics is governed by some given quantum Hamiltonian, and the Lagrange multiplier associated with a macroscopic current is added to the Hamiltonian, whereas in the Filyokov–Karpov approach [7]–[10] or in the Evans approach [11]–[14] the dynamical rules are not given at the beginning but on the contrary they are determined by the optimization of the dynamical entropy of equation (2) in the presence of appropriate constraints.

The aim of this paper is to give a self-contained and unified presentation of this type of approach, to discuss the various results that can be obtained, and to make the links with other recent developments in the field of non-equilibrium statistical physics.

The paper is organized as follows. In section 2, we describe how the maximization of the dynamical entropy of equation (2) leads to generalized Gibbs distribution for the probability distribution $P(\Omega_{[0,t]})$ of dynamical trajectories, and to some fluctuation relation of the integrated current. In the remainder of the paper, we discuss in detail how this general formal argument can be given a more precise meaning when one considers well-defined stationary stochastic dynamics. In section 3, we recall the important properties of discrete-time Markov chains, which constitute the simplest framework to describe stochastic dynamics with a time-extensive dynamical entropy (equations (2) and (3)). In section 4, we explain how the maximization of the dynamical entropy of equation (2) in the presence of a constraint concerning the energy of dynamical trajectories allows the recovery of the usual Boltzmann–Gibbs distribution of equilibrium. In section 5, we add a supplementary constraint concerning the flux of some observable and derive the form of the optimal solution for the stationary distribution and for the transition rates. We then focus on the problem of finding effective local stochastic rules, and explain how a consistent formulation can be obtained by the maximization of the relative dynamical entropy with respect to the equilibrium trajectories: we describe the case of discrete-time Markov chains in section 6 and the case of continuous-time master equations in section 7. In both cases, we make the link with the Evans approach called ‘Non-equilibrium Counterpart to detailed balance’ [11]–[14]. Our conclusions are summarized in section 8. Appendix A contains a reminder on the maximization of static entropy for the equilibrium, to make the comparison with the maximization of the dynamical entropy discussed in the text. In appendix B, we explain the technical simplifications that occur if one considers an alternate Markov chain as done originally [7]–[10].

2. General idea: generalized Gibbs distribution for dynamical trajectories

As recalled in appendix A, the usual Boltzmann–Gibbs distribution for equilibrium configurations can be obtained from the maximization of the static entropy in the presence of constraints introduced via Lagrange multipliers. This simple derivation of the equilibrium has been introduced by Jaynes [23], and can be found nowadays in most textbooks on statistical physics. Although Jaynes has often proposed to apply the same strategy to non-equilibrium phenomena, in particular in [24], we have not been able to
find in his articles the expression of equation (2) for the dynamical entropy; whereas equation (2) is the clear starting point of some ‘maximization’ in the works of Filyokov and Karpov [7]–[9] and in the Evans approach [11]–[14]. (Note that equation (2) is also the starting point in the works of Dewar [15], but with the different goal of justifying the ‘maximum production entropy principle’.)

2.1. Maximization of the dynamical entropy with constraints

For non-equilibrium steady states characterized by some macroscopic current $J_0$, the general idea is the following: one wishes to optimize the dynamical entropy of equation (2) with constraints concerning the normalization of equation (1)

$$\mathcal{N} \equiv \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) = 1$$

(4)

the averaged energy of the trajectory

$$E_{\text{dyn}}(t) \equiv \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) E(\Omega_{[0,t]}) = tE_0$$

(5)

and the averaged current of the trajectory

$$J_{\text{dyn}}(t) \equiv \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) J(\Omega_{[0,t]}) = tJ_0.$$  (6)

The optimization of the Lagrange functional

$$\psi = S_{\text{dyn}}(t) - \rho (\mathcal{N} - 1) - \beta \left( E_{\text{dyn}}(t) - tE_0 \right) + \nu \left( J_{\text{dyn}}(t) - tJ_0 \right)$$

$$= - \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) \ln \mathcal{P}(\Omega_{[0,t]}) - \rho \left( \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) - 1 \right) - \beta \left( \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) E(\Omega_{[0,t]}) - tE_0 \right) + \nu \left( \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) J(\Omega_{[0,t]}) - tJ_0 \right)$$

(7)

with respect to $\mathcal{P}(\Omega_{[0,t]})$

$$0 = \frac{\delta \psi}{\delta \mathcal{P}(\Omega_{[0,t]})} = - \ln \mathcal{P}(\Omega_{[0,t]}) - 1 - \rho - \beta E(\Omega_{[0,t]}) + \nu J(\Omega_{[0,t]})$$

(8)

directly leads to the following generalized Gibbs distribution for dynamical trajectories

$$\mathcal{P}(\Omega_{[0,t]}) = e^{-1-\rho-\beta E(\Omega_{[0,t]})+\nu J(\Omega_{[0,t]})}$$

(9)

where the Lagrange multipliers $\rho$, $\beta$ and $\nu$ are fixed respectively by the constraints of equations (4)–(6).
2.2. Maximization of the relative dynamical entropy with respect to some equilibrium reference process

Another formulation that turns out to be more appropriate in many cases consists in considering the relative dynamical entropy with respect to the corresponding known equilibrium distribution $P_{\text{eq}}(\Omega_{[0,t]})$ of trajectories in the absence of the current

$$S_{\text{rel}}^{\text{dyn}}(t) \equiv - \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) \ln \frac{\mathcal{P}(\Omega_{[0,t]})}{P_{\text{eq}}(\Omega_{[0,t]})}. \quad (10)$$

The idea is then to maximize this relative entropy in the presence of the constraints concerning the normalization of equation (4) and the imposed current (equation (6)) to obtain

$$\mathcal{P}(\Omega_{[0,t]}) = P_{\text{eq}}(\Omega_{[0,t]}) e^{-1 - \rho + \nu J(\Omega_{[0,t]})} \quad (11)$$

instead of equation (9).

The notion of relative entropy with respect to some reference, also called the Kullback–Leibler entropy, is well known in many areas. At a naive level, the idea is that the probability distribution used as reference constitutes some ‘prior’ for the problem at hand: in the absence of any other constraint, the relative entropy is maximal for the ‘prior’ distribution. At a more fundamental level, the importance of the notion of relative entropy comes from its essential role in the Sanov theorem concerning the theory of large deviations when formulated at the so-called ‘level-2’ (see for instance the review [27]).

The interest to consider relative entropies in various areas of statistical physics has been discussed recently in [26]. In the present context of non-equilibrium steady states, the idea to take the equilibrium dynamics as the reference process is the basis of the Evans approach [11]–[14].

2.3. Consequence: fluctuation relation concerning the integrated current

It is clearly impossible to summarize here all the recent developments concerning the various ‘fluctuation relations’ that have been established in the field of non-equilibrium dynamics, and we refer the interested reader to some recent reviews [28]–[34] and references therein. Here we simply mention the minimum for our present purpose.

The probability distribution of the integrated current is expected to satisfy some large deviation principle: the probability to have a given time-averaged value $J(\Omega_{[0,t]})/t = j$ behaves at large time $t$ as

$$\text{Prob} \left( \frac{J(\Omega_{[0,t]})}{t} = j \right) \propto e^{tG(j)} \quad (12)$$

where $G(j) \leq 0$ is called the large deviation function. The typical value $j_{\text{typ}}$ corresponding to the point where $G(j_{\text{typ}}) = 0$ is fixed here to $J_0$ by the constraint of equation (64). Equivalently, the generating function of the integrated current $J(\Omega_{[0,t]})$ behaves for large time $t$ as

$$\langle e^{\lambda J(\Omega_{[0,t]})} \rangle \propto \int dj \, e^{t(j + G(j))} \propto e^{t\mu(\lambda)} \quad (13)$$

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where $\mu(\lambda)$ is the Legendre transform of $G(j)$

$$
\mu(\lambda) = \max_j [\lambda j + G(j)].
$$

(14)

To characterize the irreversibility of the dynamics, it is interesting to compare the probabilities of a given trajectory $\Omega_{[0,t]}$ and of its associate trajectory $\tilde{\Omega}_{[0,t]}$ obtained by time-reversal. In our present case, these two trajectories have the same energy $E(\Omega_{[0,t]}) = E(\tilde{\Omega}_{[0,t]})$ but have opposite currents $J(\Omega_{[0,t]}) = -J(\tilde{\Omega}_{[0,t]})$. So equation (9) yields that the ratio of the probabilities between a trajectory and its time-reversed trajectory reads

$$
\frac{\mathcal{P}(\Omega_{[0,t]})}{\mathcal{P}(\tilde{\Omega}_{[0,t]})} = e^{2\nu J(\Omega_{[0,t]})}.
$$

(15)

More generally, any equilibrium dynamics has the property that two reversed-time trajectories occur with the same probability $\mathcal{P}^{eq}(\Omega_{[0,t]}) = \mathcal{P}^{eq}(\tilde{\Omega}_{[0,t]})$. As a consequence, equation (11) also leads to equation (15).

A direct consequence of equation (15) is that the generating function of the integrated current defined as

$$
\langle e^{\lambda J(\Omega_{[0,t]})} \rangle \equiv \sum_{\Omega_{[0,t]}} e^{\lambda J(\Omega_{[0,t]})} \mathcal{P}(\Omega_{[0,t]})
$$

(16)

will satisfy the following symmetry (using the antisymmetry $J(\tilde{\Omega}_{[0,t]}) = -J(\Omega_{[0,t]})$ and the one-to-one change of variables between $\Omega_{[0,t]}$ and $\tilde{\Omega}_{[0,t]}$)

$$
\langle e^{\lambda J(\Omega_{[0,t]})} \rangle = \sum_{\tilde{\Omega}_{[0,t]}} e^{-(\lambda+2\nu) J(\tilde{\Omega}_{[0,t]})} \mathcal{P}(\tilde{\Omega}_{[0,t]}) = \langle e^{-(\lambda+2\nu) J(\Omega_{[0,t]})} \rangle
$$

(17)

that can be rewritten for the function $\mu(\lambda)$ introduced in equation (13)

$$
\mu(\lambda) = \mu(-\lambda - 2\nu)
$$

(18)

or equivalently after the Legendre transform of equation (14)

$$
G(j) = G(-j) + 2\nu j.
$$

(19)

In summary, the generalized Gibbs distribution of dynamical trajectories of equation (9) or equation (11) directly leads to some fluctuation relation for the integrated current, as already discussed by Baule and Evans [12]. Independently of the context of the maximization of the dynamical entropy, the fact that fluctuation relations have actually for origin some generalized Gibbs distribution of space-time trajectories has been proposed by Maes [35].

2.4. Entropy production

The notion of entropy ‘production’ is of course subtle. Since it is expected to measure the ‘irreversibility’ of the dynamics, it is natural to relate it to the probabilities between a trajectory and its reversed-time trajectory already introduced in equation (15): a possibility [35, 36] is thus to define the entropy production $S^{prod}(\Omega_{[0,t]})$ of a dynamical

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trajectory $\Omega_{[0,t]}$ as the logarithm of the ratio of equation (15)

$$S_{\text{prod}}(\Omega_{[0,t]}) \equiv \ln \frac{\mathcal{P}(\Omega_{[0,t]})}{\mathcal{P}(\tilde{\Omega}_{[0,t]})}. \quad (20)$$

Such a definition, which is in agreement with the usual expressions of the averaged entropy production derived for Markov chains [37, 38], has for advantage that the fluctuation relation for the entropy production

$$\frac{\text{Prob}(S_{\text{prod}})}{\text{Prob}(-S_{\text{prod}})} = e^{S_{\text{prod}}} \quad (21)$$

is then a direct consequence of equation (20) because

$$\text{Prob}(S_{\text{prod}}) \equiv \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) \delta \left( S_{\text{prod}} - \ln \frac{\mathcal{P}(\Omega_{[0,t]})}{\mathcal{P}(\tilde{\Omega}_{[0,t]})} \right)$$

$$= \sum_{\tilde{\Omega}_{[0,t]}} \mathcal{P}(\tilde{\Omega}_{[0,t]}) e^{S_{\text{prod}}} \delta \left( S_{\text{prod}} + \ln \frac{\mathcal{P}(\tilde{\Omega}_{[0,t]})}{\mathcal{P}(\Omega_{[0,t]})} \right)$$

$$= e^{S_{\text{prod}}} \text{Prob}(-S_{\text{prod}}). \quad (22)$$

For our present case, the definition of equation (20) leads with equation (15) to

$$S_{\text{prod}}(\Omega_{[0,t]}) \equiv \ln \frac{\mathcal{P}(\Omega_{[0,t]})}{\mathcal{P}(\tilde{\Omega}_{[0,t]})} = 2\nu J(\Omega_{[0,t]}) \quad (23)$$

i.e. it is simply proportional to the time-extensive trajectory current $J(\Omega_{[0,t]})$ that breaks the time asymmetry.

2.5. Discussion

In this section, we have described the general idea of maximization of the dynamical entropy with constraints that leads to generalized Gibbs distributions for dynamical trajectories, and to some fluctuation relation for the integrated current. However, the derivation presented above remains a bit formal, since the space of dynamical trajectories has not been precisely defined. In particular, the important notion of ‘stationarity’ of the dynamics has been imposed only implicitly by the time-extensive constraints of equations (5) and (6), but no direct condition of stationarity has been explicitly imposed on $\mathcal{P}(\Omega_{[0,t]})$ (and we do not see how one could impose such a stationarity condition at this formal level). In the following sections, we thus consider well-defined stochastic stationary dynamics generated by discrete-time Markov chains or by continuous-time master equations to study the consequences of the maximization of the dynamical entropy.

3. Dynamical observables for a stationary Markov chain

Discrete-time Markov chains constitute the simplest formulation of stochastic dynamics presenting a dynamical entropy (equation (2)) that is extensive in time (equation (3)). In this section, we introduce the notations that will be useful in the remainder of the paper.
3.1. Discrete-time Markov chain

Let us consider a system where possible microscopic configurations are indexed by $i$, and where the probability $P_t(i)$ to be in configuration $i$ at time $t$ evolves according to the discrete-time Markov chain

$$P_{t+1}(j) = \sum_i P_t(i)W_{i\rightarrow j}$$

(24)

where $W_{i\rightarrow j}$ represents the transition probability from $i$ to $j$ during the unit time interval, with the following normalization for each $i$

$$\sum_j W_{i\rightarrow j} = 1.$$  

(25)

The stationary state $P^{\text{st}}(i)$ associated with this Markov chain satisfies

$$P^{\text{st}}(i) = \sum_j P^{\text{st}}(j)W_{j\rightarrow i}$$

(26)

and the normalization

$$\sum_i P^{\text{st}}(i) = 1.$$  

(27)

3.2. Probability of a dynamical trajectory in the stationary regime

Suppose one starts at time $t = 0$ in the stationary state $P^{\text{st}}$ of equation (26). The probability of a trajectory $\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\}$ to be in state $i_0$ at time $t = 0$, in state $i_1$ at time $t = 1$, in state $i_2$ at time $t = 2, \ldots$, and in state $i_t$ at time $t$ reads

$$P(\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\}) = P^{\text{st}}(i_0)W_{i_0\rightarrow i_1}W_{i_1\rightarrow i_2}\cdots W_{i_{t-1}\rightarrow i_t}.$$  

(28)

The normalization of equation (1)

$$\sum_{i_0} \sum_{i_1} \sum_{i_2} \cdots \sum_{i_t} P(\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\}) = \sum_{i_0} P^{\text{st}}(i_0)\sum_{i_1} W_{i_0\rightarrow i_1} \sum_{i_2} W_{i_1\rightarrow i_2} \cdots \sum_{i_t} W_{i_{t-1}\rightarrow i_t} = 1$$  

(29)

can be easily checked using equations (25) and (27).

3.3. Dynamical entropy associated with the distribution of trajectories

Using the normalization and stationarity properties of equations (25) and (26), the dynamical entropy of equation (2) reads for the distribution $P(\Omega_{[0,t]})$ of equation (28)

$$S^{\text{dyn}}(t) = -\sum_{\Omega_{[0,t]}} P(\Omega_{[0,t]}) \ln P(\Omega_{[0,t]})$$

$$= -\sum_{i_0} \sum_{i_1} \cdots \sum_{i_t} P^{\text{st}}(i_0)W_{i_0\rightarrow i_1} \cdots W_{i_{t-1}\rightarrow i_t}$$

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The physical meaning of this formula is clear: it is the average, over the possible configurations $i$, so that its averaged value over the trajectories $\Omega$ is extensive in time, and the ratio $E^{\text{dyn}}(t)/t$ corresponds to the static averaged energy computed from the stationary distribution as it should.

3.4. Energy associated with trajectories

In statistical physics, one usually wishes to fix some constraint concerning the energy. For instance at equilibrium, the Boltzmann–Gibbs distribution can be obtained from the optimization of the static entropy in the presence of a constraint fixing the average energy (see appendix A for a short reminder). In our present dynamical context, it is natural to associate with each trajectory $\Omega_{[0,t]}$ the integrated energy over the time interval

$$E(\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\}) \equiv E_{i_0} + E_{i_1} + \cdots + E_{i_{t-1}}$$

(33)

so that its averaged value over the trajectories $\Omega_{[0,t]}$ (using equations (25) and (26))

$$E^{\text{dyn}}(t) \equiv \sum_{\Omega_{[0,t]}} P(\Omega_{[0,t]}) E(\Omega_{[0,t]})$$

$$= \sum_{i_0} P^\text{st}(i_0) E_{i_0} + \sum_{i_1} P^\text{st}(i_1) E_{i_1} + \cdots + \sum_{i_{t-1}} P^\text{st}(i_{t-1}) E_{i_{t-1}} = t \sum_i P^\text{st}(i) E_i$$

(34)

is extensive in time, and the ratio $E^{\text{dyn}}(t)/t$ corresponds to the static averaged energy computed from the stationary distribution as it should.
3.5. Flux associated with trajectories

In the field of non-equilibrium steady states, one is interested in particular in situations where a flux or current of some observable is imposed on the system. It can be a flux of particles, a flux of energy, or a flux of something else depending on the system under study and on the environment.

In our present framework, we will consider that each transition \( i \rightarrow j \) is characterized by some contribution \( J_{i \rightarrow j} \) to this current, with the antisymmetry property
\[
J_{i \rightarrow j} = -J_{j \rightarrow i}.
\]

(35)

The integrated current \( J(\Omega_{[0,t]}) \) associated with a trajectory \( \Omega_{[0,t]} \) simply reads
\[
J(\Omega_{[0,t]} = \{i_0,i_1,\ldots,i_t\}) \equiv J_{i_0 \rightarrow i_1} + J_{i_1 \rightarrow i_2} + \cdots + J_{i_{t-1} \rightarrow i_t}
\]

so that its averaged value over the trajectories \( \Omega_{[0,t]} \) (using equations (25) and (26))
\[
J^{\text{dyn}}(t) \equiv \sum_{\Omega_{[0,t]}} \mathcal{P}(\Omega_{[0,t]}) J(\Omega_{[0,t]})
\]

\[
= \sum_{i_0} \sum_{i_1} \cdots \sum_{i_t} P^\text{st}(i_0) W_{i_0 \rightarrow i_1} \cdots W_{i_{t-1} \rightarrow i_t} [J_{i_0 \rightarrow i_1} + J_{i_1 \rightarrow i_2} + \cdots + J_{i_{t-1} \rightarrow i_t}]
\]

\[
= \sum_{i_0} P^\text{st}(i_0) \sum_{i_1} W_{i_0 \rightarrow i_1} J_{i_0 \rightarrow i_1} + \cdots + \sum_{i_{t-1}} P^\text{st}(i_{t-1}) \sum_{i_t} W_{i_{t-1} \rightarrow i_t} J_{i_{t-1} \rightarrow i_t}
\]

\[
= t \sum_i P^\text{st}(i) \sum_j W_{i \rightarrow j} J_{i \rightarrow j}
\]

(37)

is again extensive in time as expected.

4. Recovering equilibrium from the maximization of the dynamical entropy

The minimal requirement for any approach concerning non-equilibrium steady states is of course to be compatible with the theory of the equilibrium. In this section, we thus describe how the maximization of the dynamical entropy of equation (2) allows one to recover the equilibrium. As a comparison, we recall in appendix A how the equilibrium is usually obtained via the maximization of the static entropy.

4.1. Lagrange functional taking into account the constraints

We consider all possible stationary distributions \( P^\text{st}(i) \) satisfying the normalization
\[
\mathcal{N} \equiv \sum_i P^\text{st}(i) = 1
\]

(38)

and all Markov chains defined by some transition probabilities \( W_{i \rightarrow j} \), that satisfy the normalization condition of equation (25) for all \( i \)
\[
\mathcal{N}_i \equiv \sum_j W_{i \rightarrow j} = 1
\]

(39)
and the stationarity condition of equation (26) for all $i$

$$
\Sigma_i \equiv \sum_j P_{\text{st}}(j) W_{j\rightarrow i} - P_{\text{st}}(i) = 0.
$$

(40)

We also wish to fix the time-averaged energy $E_{\text{dyn}}(t)/t$ (see equation (A.5)) to some value $E_0$

$$
\frac{E_{\text{dyn}}(t)}{t} \equiv \sum_i P_{\text{st}}(i) E_i = E_0.
$$

(41)

To optimize the dynamical entropy $S_{\text{dyn}}(t)$ of equation (31) in the presence of all these constraints, we introduce Lagrange multipliers $\rho$, $\beta$, $\lambda_i$ and $\mu_i$ and we consider the functional

$$
\Psi \equiv \frac{S_{\text{dyn}}(t)}{t} - \rho (N - 1) - \beta \left( \frac{E_{\text{dyn}}(t)}{t} - E_0 \right) - \sum_i \lambda_i (N_i - 1)
- \sum_i \mu_i (\Sigma_i) = -\sum_i P_{\text{st}}(i) \sum_j W_{i\rightarrow j} \ln W_{i\rightarrow j} - \rho \left( \sum_i P_{\text{st}}(i) - 1 \right)
- \beta \left( \sum_i E_i P_{\text{st}}(i) - E_0 \right) - \sum_i \lambda_i \left( \sum_j W_{i\rightarrow j} - 1 \right)
- \sum_i \mu_i \left( \sum_j P_{\text{st}}(j) W_{j\rightarrow i} - P_{\text{st}}(i) \right).
$$

(42)

4.2. Solving the optimization equations

We wish to optimize the functional of equation (42) both over the stationary probabilities $P_{\text{st}}(i)$

$$
0 = \frac{\delta \Psi}{\delta P_{\text{st}}(i)} = -\sum_j W_{i\rightarrow j} \ln W_{i\rightarrow j} - \rho - \beta E_i + \mu_i - \sum_j \mu_j W_{i\rightarrow j}
$$

(43)

and over the transition probabilities $W_{i\rightarrow j}$

$$
0 = \frac{\delta \Psi}{\delta W_{i\rightarrow j}} = P_{\text{st}}(i) [- \ln W_{i\rightarrow j} - 1 - \mu_j] - \lambda_i.
$$

(44)

The Lagrange multipliers have then to be determined by imposing the various constraints. Equation (44) yields

$$
W_{i\rightarrow j} = e^{-1-\mu_j-\lambda_i/P_{\text{st}}(i)}.
$$

(45)

The normalization constraint of equation (39) for all $i$

$$
1 = \sum_j W_{i\rightarrow j} = e^{-1-\lambda_i/P_{\text{st}}(i)} \sum_j e^{-\mu_j}
$$

(46)
implies that the Lagrange multipliers $\lambda_i$ are given by

$$\lambda_i = P^{\text{st}}(i) \ln \left( \sum_j e^{-\mu_j} \right)$$  \hspace{1cm} (47)$$

so that the rates of equation (45) can be rewritten as

$$W_{i \rightarrow j} = \frac{e^{-\mu_j}}{\sum_k e^{-\mu_k}}.$$  \hspace{1cm} (48)$$

The stationarity constraint of equation (40) yields that for all $j$

$$P^{\text{st}}(j) = \sum_i P^{\text{st}}(i) W_{i \rightarrow j} = \frac{e^{-\mu_j}}{\sum_k e^{-\mu_k}}$$  \hspace{1cm} (49)$$

so that the optimum found in equation (48) simply corresponds to [8,10]

$$W_{i \rightarrow j} = P^{\text{st}}(j).$$  \hspace{1cm} (50)$$

Equation (43) becomes, using equation (49) and equation (38)

$$0 = -\rho - \beta E_i - \ln P^{\text{st}}(i)$$  \hspace{1cm} (51)$$

so that using the normalization constraint of equation (38) one recovers the Boltzmann–Gibbs distribution

$$P^{\text{st}}(j) = \frac{e^{-\beta E_j}}{\sum_i e^{-\beta E_i}} \equiv P^{\text{eq}}(j)$$  \hspace{1cm} (52)$$

where $\beta$ is fixed by the constraint of equation (41).

4.3. Consequence for the probabilities of dynamical trajectories

With the optimal solution obtained in equations (50) and (52), the probability of a trajectory $\Omega_{[0,t]}$ of equation (28) takes the simple form

$$P^{\text{eq}}(\Omega_{[0,t+1]}) = \left\{ i_0, i_1, \ldots, i_t \right\} = P^{\text{st}}(i_0) W_{i_0 \rightarrow i_1} W_{i_1 \rightarrow i_2} \cdots W_{i_{t-1} \rightarrow i_t}$$

$$= \frac{e^{-\beta E(i_0)} e^{-\beta E(i_1)} \cdots e^{-\beta E(i_t)}}{[Z(\beta)]^{t+1}}$$  \hspace{1cm} (53)$$

where

$$Z(\beta) = \sum_i e^{-\beta E_i}$$  \hspace{1cm} (54)$$

is the usual static partition function. So within the present approach, the probability distribution over dynamical trajectories is simply the Boltzmann–Gibbs distribution with respect to the trajectory energy $E(\Omega_{[0,t+1]})$ of equation (34)

$$P^{\text{eq}}(\Omega_{[0,t+1]}) = \frac{e^{-\beta E(\Omega_{[0,t+1]})}}{[Z(\beta)]^{t+1}}.$$  \hspace{1cm} (55)$$

The appearance of $E(\Omega_{[0,t+1]})$ (instead of $E(\Omega_{[0,t]})$ found in equation (9) via the formal argument) is a consequence of the choice of equation (34) for the energy for the discrete
Markov chain: the energy $E(i_t)$ of the last state $i_t$ will contribute to the trajectory energy only during the next time interval $[t, t + 1]$, but it nevertheless appear in the probability of the trajectory $\Omega_{[0,t]}$ ending in state $i_t$. This slight problem thus comes from the discrete nature of the dynamics, and is not expected to play an important role: it is only a finite boundary term that will become subleading in the large-time limit $t \to +\infty$ with respect to the time-extensive energy of the whole trajectory.

An essential property of this distribution of trajectories is the symmetry by time-reversal: the probabilities of a given trajectory $\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_{t-1}, i_t\}$ and of its associate trajectory $\tilde{\Omega}_{[0,t]} = \{i_t, i_{t-1}, \ldots, i_1, i_0\}$ obtained by time-reversal are equal

$$\frac{P(\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\})}{P(\tilde{\Omega}_{[0,t]} = \{i_t, i_{t-1}, \ldots, i_1, i_0\})} = 1.$$ (56)

4.4. Discussion

In this section, we have described how the maximization of the dynamical entropy allows one to recover that the stationary distribution follows the Boltzmann–Gibbs distribution (equation (52)). However the result of equation (50) concerning the transition probabilities $W_{i \to j}$ may be surprising at first sight: this solution means that the new configuration $j$ is chosen with probability $P_{\text{st}}(j)$ and is completely independent of the initial state $i$. Physically this corresponds to a very coarse-grained dynamics with no memory. It is thus clear that equation (50) does not represent an effective ‘microscopic’ dynamics, but represents an effective dynamics on some macroscopic time $\tau$. The corresponding entropy rate per unit time of equation (3)

$$h_{\text{KS}} = - \left( \sum_i P_{\text{st}}(i) \right) \sum_j P_{\text{st}}(j) \ln P_{\text{st}}(j) = - \sum_j P_{\text{st}}(j) \ln P_{\text{st}}(j)$$ (57)

then exactly coincides with the static entropy of equation (A.1), which seems natural if one wishes this approach to be equivalent to the maximization of the static entropy that one uses for the equilibrium (see appendix A). This discussion suggests the following interpretation of this type of computation.

4.5. Final formulation of the physical meaning of this approach

The statistical physics theory of equilibrium is usually based on some ‘ergodic’ hypothesis, stating that time-averages of observables $A(i)$ will converge in the infinite-time limit towards averages computed with respect to the Boltzmann–Gibbs distribution $P_{\text{eq}}(i)$ of equation (52)

$$\lim_{t \to +\infty} \frac{1}{t} \int_0^t dt A(i(t)) = \sum_i A(i) P_{\text{eq}}(i).$$ (58)

Since the right-hand side contains only the Boltzmann–Gibbs distribution $P_{\text{eq}}(i) = e^{-\beta E_i}/Z(\beta)$ and no other information about the dynamics except the conserved energy, this means that the precise form of the ‘true’ microscopic dynamics becomes completely irrelevant at large time. Loosely speaking, this means that there exists some finite macroscopic correlation time $\tau_{\text{correl}}$, beyond which the dynamical correlations have been
lost in practice, so that the large-time interval $t$ can be decomposed into $t/\tau_{\text{correl}}$ quasi-independent time intervals, and equation (58) becomes possible. On the contrary, if the correlation time $\tau_{\text{correl}}$ of the ‘true’ microscopic dynamics is infinite, equation (58) cannot really be satisfied, since other dynamical information besides the conservation of energy remains relevant forever.

This discussion suggests that the analysis presented above based on the maximization of the dynamical entropy, with the result of equation (50), actually describes what happens at this coarse-grained macroscopic scale $\tau_{\text{correl}}$, i.e. the unit time of the Markov chain of equation (24) should be interpreted as $\tau_{\text{correl}}$. Then we have shown above that one recovers the Boltzmann–Gibbs distribution, and thus all the ‘static’ properties that can be derived from it.

However, in the statistical physics of equilibrium, once one has understood the properties of the Boltzmann–Gibbs measure over configurations, one can become interested in the ‘equilibrium dynamics’. However, one does not wish to return to the ‘true’ deterministic microscopic dynamics, which is usually very complicated for systems with a very large number (like $10^{23}$) degrees of freedom, because all the details of the dynamics have proven to be irrelevant at large time scale. So one introduces an effective stochastic microscopic dynamics that is compatible with the known properties on large time scales. For definiteness, let us consider a microscopic Markov chain

$$ p_{t+\Delta t}(j) = \sum_i p_t(i) w_{i \rightarrow j} \quad (59) $$

where $w_{i \rightarrow j}$ represents the transition probability from $i$ to $j$ during the microscopic time interval $\Delta t$, so that $w_{i \rightarrow j}$ are now non-zero only if the configurations $i$ and $j$ are sufficiently close in configuration space. For instance in a system of $N$ spins with $2^N$ configurations, one may require that $w_{i \rightarrow j}$ is non-zero only if $j$ can be obtained from $i$ by the flip of a single spin. More generally, this notion of elementary moves has to be defined in an appropriate way for each type of models. Then the question is: beyond this locality requirement, how should these microscopic transition probabilities $w_{i \rightarrow j}$ be chosen to be compatible with the properties known on large time scales? The first obvious requirement is that the microscopic Markov chain of equation (59) should have for stationary state $p^\text{st}(i)$ the Boltzmann–Gibbs distribution which is known to be the stationary state on macroscopic times

$$ p^\text{st}(i) = P^\text{eq}(i). \quad (60) $$

However this is not the only constraint, since one also wishes to reproduce the essential time-reversibility property of dynamical trajectories of equation (56). For this, it is sufficient to impose that the probability of the elementary microscopic trajectory $\omega_{\Delta t} = \{i_0, i_1\}$

$$ \text{Prob}(\omega_{\Delta t} = \{i_0, i_1\}) = p(i_0) w_{i_0 \rightarrow i_1} \quad (61) $$

is equal to the probability of the reversed trajectory $\tilde{\omega}_{\Delta t} = \{i_1, i_0\}$

$$ 1 = \frac{\text{Prob}(\omega_{\Delta t} = \{i_0, i_1\})}{\text{Prob}(\tilde{\omega}_{\Delta t} = \{i_1, i_0\})} = \frac{p(i_0) w_{i_0 \rightarrow i_1}}{p(i_1) w_{i_1 \rightarrow i_0}}. \quad (62) $$

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Taking into account equation (60), this leads to the well-known detailed balance condition

\[ \frac{w^{eq}_{i_0 \rightarrow i_1}}{w^{eq}_{i_1 \rightarrow i_0}} = \frac{p(i_1)}{p(i_0)} = \frac{P^{eq}(i_1)}{P^{eq}(i_0)} = e^{-\beta(E(i_1) - E(i_0))}, \] (63)

Besides this detailed balance constraint, there is still some freedom to choose the effective microscopic transition probabilities \( w_{i \rightarrow j} \), but one usually considers that they are equally valid, in the sense that they will all reproduce the essential large-time properties of convergence towards the Boltzmann–Gibbs distribution while preserving the time-reversibility of dynamical trajectories.

4.6. Is it possible to add locality constraints within the maximization procedure?

As explained above, the requirements on effective local Markov chain dynamics of equation (59) are derived from the macroscopic equilibrium properties. A natural question here is whether these requirements can be instead directly derived from a maximization of the dynamical entropy in the presence of locality constraints on the \( w_{i \rightarrow j} \). We have now the following picture: configurations are the nodes of a network, and the links are present between configurations that are related by an elementary local move. As a first example, in a spin models with \( N \) spins and \( 2^N \) configurations, one may consider a single spin-flip dynamics, where each configuration has \( N \) neighbors (corresponding to the flip of one of its \( N \) spins). As a second example, in lattice gases models with hard-core interactions, the number of neighbors of a configuration will be given by the number of possible local moves from this configuration, and will thus depend on the configuration.

Whenever the connectivities (i.e. the numbers of neighbors) are configuration-dependent, it is clear that one cannot recover the equilibrium from the maximization of the dynamical entropy in the presence of locality constraints, as can be easily understood on the simple infinite temperature case \( \beta = 0 \):

(i) At infinite temperature, the problem of the maximization of the dynamical entropy is completely equivalent to the maximal entropy random walk (MERW) on arbitrary networks studied in detail in [16], with the following main conclusion: in networks where the connectivities of the various nodes are not all the same, the MERW tends to visit more the sites with higher connectivities than the sites with lower connectivities, because sites with higher connectivities are associated with a bigger number of choices for the next step of the random walk.

(ii) On the other hand, in the infinite temperature case \( \beta = 0 \), the Boltzmann–Gibbs distribution simply corresponds to the uniform distribution over all configurations, independently of their connectivities from the point of view of the local dynamics.

Our conclusion is thus that the choice of the connectivities that one wishes to impose on the local dynamics has too strong an effect on the dynamical entropy that one wishes to maximize, whereas the maximization of the dynamical entropy at a macroscopic time scale allows one to recover the equilibrium, as explained above in section 4.5. In the non-equilibrium case, this discussion on local dynamics will be the subject of sections 6 and 7, but in section 5, we first describe what happens in the absence of locality constraints.
5. Non-equilibrium steady state with an imposed current

5.1. Lagrange functional taking into account the constraints

With respect to the previous section 4.1, we now add another constraint concerning the flux of equation (37)

\[ \frac{J_{\text{dyn}}(t)}{t} \equiv \sum_i P_{\text{st}}(i) \sum_j W_{i \rightarrow j} J_{i \rightarrow j} = J_0. \]  

(64)

So we introduce this constraint with a new Lagrange multiplier \( \nu \) into the functional of equation (42) to obtain

\[
\Psi \equiv \frac{S_{\text{dyn}}(t)}{t} - \rho(N - 1) - \beta \left( \frac{E_{\text{dyn}}(t)}{t} - E_0 \right)
- \sum_i \lambda_i (N_i - 1) - \sum_i \mu_i (\Sigma_i) + \nu \left( \frac{J_{\text{dyn}}(t)}{t} - J_0 \right)
- \beta \left( \sum_i E_i P_{\text{st}}(i) - E_0 \right)
- \sum_i \lambda_i \sum_j W_{i \rightarrow j} - 1
- \sum_i \mu_i \left( \sum_j P_{\text{st}}(j) W_{j \rightarrow i} - P_{\text{st}}(i) \right) + \nu \left( \sum_i P_{\text{st}}(i) \sum_j W_{i \rightarrow j} J_{i \rightarrow j} - J_0 \right).
\]

(65)

5.2. Solving the optimization equations

Equations (43) and (44) are modified into

\[
0 = \frac{\delta \Psi}{\delta P_{\text{st}}(i)} = -\sum_j W_{i \rightarrow j} \ln W_{i \rightarrow j} - \rho - \beta E_i + \mu_i - \sum_k \mu_k W_{i \rightarrow k} + \nu \sum_j W_{i \rightarrow j} J_{i \rightarrow j}
\]

(66)

and

\[
0 = \frac{\delta \Psi}{\delta W_{i \rightarrow j}} = P_{\text{st}}(i) \left[ -\ln W_{i \rightarrow j} - 1 - \mu_j + \nu J_{i \rightarrow j} \right] - \lambda_i.
\]

(67)

Equation (67) yields

\[
W_{i \rightarrow j} = e^{-1 - \mu_j + \nu J_{i \rightarrow j} - \lambda_i / P_{\text{st}}(i)}.
\]

(68)

The normalization constraint of equation (39) for all \( i \)

\[
1 = \sum_j W_{i \rightarrow j} = e^{-1 - \lambda_i / P_{\text{st}}(i)} \sum_j e^{-\mu_j + \nu J_{i \rightarrow j}}
\]

(69)
leads to the Lagrange multipliers
\[ \lambda_i = P^\text{st}(i) \ln \left( \sum_j e^{-\mu_j + \nu J_{i \rightarrow j}} \right) \]  
so that the rates of equation (45) become
\[ W_{i \rightarrow j} = \frac{e^{-\mu_j + \nu J_{i \rightarrow j}}}{\sum_k e^{-\mu_k + \nu J_{i \rightarrow k}}} \]  

To analyze the stationarity constraint of equation (40) for all \( j \)
\[ P^\text{st}(j) = \sum_i P^\text{st}(i) W_{i \rightarrow j} = \sum_i P^\text{st}(i) \frac{e^{-\mu_j + \nu J_{i \rightarrow j}}}{\sum_k e^{-\mu_k + \nu J_{i \rightarrow k}}} \]  
it is convenient to introduce the notations
\[ z_i \equiv \sum_k e^{-\mu_k + \nu J_{i \rightarrow k}} \]  
and
\[ y_j \equiv \sum_i \frac{P^\text{st}(i)}{z_i} e^{\nu J_{i \rightarrow j}} \]  
to rewrite equation (72) as
\[ e^{-\mu_j} = \frac{P^\text{st}(j)}{y_j} \]  
so we may now eliminate all \( \mu_j \) in terms of the \( y_j \). In particular
\[ z_i = \sum_k \frac{P^\text{st}(k)}{y_k} e^{\nu J_{i \rightarrow k}} \]  
is somewhat the ‘dual’ of equation (74). The transition probability of equation (71) now reads
\[ W_{i \rightarrow j} = \frac{P^\text{st}(j)}{z_i y_j} e^{\nu J_{i \rightarrow j}} \]  
and equation (66) becomes
\[ 0 = -\rho - \beta E_i - \ln P^\text{st}(i) + \ln(y_i) + \sum_j W_{i \rightarrow j} [-\ln W_{i \rightarrow j} - \mu_j + \nu J_{i \rightarrow j}] \]
\[ = -\rho - \beta E_i - \ln P^\text{st}(i) + \ln(y_i) + \ln(z_i) \]  
leading to
\[ P^\text{st}(i) = y_i z_i e^{-\rho - \beta E_i}. \]  
Plugging equation (79) into equations (74) and (76) leads to
\[ y_j = \sum_i y_i e^{-\rho - \beta E_i} e^{\nu J_{i \rightarrow j}} \]  
and
\[ z_i = \sum_k z_k e^{-\rho - \beta E_k} e^{\nu J_{i \rightarrow k}}. \]
5.3. Optimal stationary distribution \(P^{\text{st}}(i)\) and transition probabilities \(W_{i\rightarrow j}\)

In summary, we have obtained that the optimal stationary distribution \(P^{\text{st}}(i)\) and transition probabilities \(W_{i\rightarrow j}\) follow the form

\[
P^{\text{st}}(i) = y_i z_i e^{-\rho - \beta E_i}
\]
\[
W_{i\rightarrow j} = \frac{z_j}{z_i} e^{-\rho - \beta E_i + \nu J_{i\rightarrow j}}
\]

where the \(y_i\) and \(z_i\) are positive variables satisfying respectively the equations

\[
y_j = \sum_i y_i e^{-\rho - \beta E_i} e^{\nu J_{i\rightarrow j}}
\]

that correspond to the stationarity constraints \(P^{\text{st}}(j) = \sum_i P^{\text{st}}(i) W_{i\rightarrow j}\) and the equations

\[
z_i = \sum_j e^{\nu J_{i\rightarrow j}} e^{-\rho - \beta E_j} z_j
\]

that correspond to the normalizations \(1 = \sum_j W_{i\rightarrow j}\), and

\[
1 = \sum_i y_i z_i e^{-\rho - \beta E_i}
\]

that correspond to the normalization \(1 = \sum_i P^{\text{st}}(i)\). To see more clearly what this structure means, it is convenient to introduce the bra and ket notations and to set

\[
y_i \equiv e^{(\beta/2)E_i} \langle L | i \rangle
\]
\[
z_i \equiv e^{(\beta/2)E_i} \langle i | R \rangle.
\]

Then in terms of the matrix \(M\) defined by the positive matrix elements

\[
\langle i | M | j \rangle \equiv e^{-(\beta/2)E_j} e^{\nu J_{i\rightarrow j}} e^{-(\beta/2)E_i}
\]

Equations (83) and (84) mean that \(\langle L | \) and \(| R \rangle\) are respectively positive left eigenvector and right eigenvector of the non-symmetric matrix \(M\)

\[
e^\rho \langle L | = \langle L | M |
e^\rho | R \rangle = M | R \rangle
\]

associated with the highest eigenvalue \(e^\rho\) of the matrix \(M\) (Perron–Frobenius), with the normalization given by equation (85)

\[
e^\rho = \langle L | R \rangle.
\]

The equilibrium case discussed in the previous section 4 corresponds to \(\nu = 0\), \(y_i = 1 = z_i\), \(\langle L | i \rangle = \langle i | R \rangle = e^{-(\beta/2)E_i}\). In the following to discuss what happens for \(\nu > 0\), we will use the notations \(y_i\) and \(z_i\) that appear in equations (82).
5.4. Consequence for the probability distribution of dynamical trajectories

From the optimal solution of equation (82), one obtains that the probability of a trajectory \(\Omega_{[0,t]}\) of equation (28) takes the form

\[
\mathcal{P}(\Omega_{[0,t]} = \{i_0, i_1, \ldots, i_t\}) = P^{st}(i_0)W_{i_0 \rightarrow i_1}W_{i_1 \rightarrow i_2} \cdots W_{i_{t-1} \rightarrow i_t} \\
= y_{i_0}z_{i_0}e^{-\beta E_{i_0}}\frac{z_{i_1}}{z_{i_0}}e^{-\beta E_{i_1} + \nu J_{i_0 \rightarrow i_1}}\frac{z_{i_2}}{z_{i_1}}e^{-\beta E_{i_2} + \nu J_{i_1 \rightarrow i_2}} \cdots \frac{z_{i_t}}{z_{i_{t-1}}}e^{-\beta E_{i_t} + \nu J_{i_{t-1} \rightarrow i_t}} \\
= y_{i_0}z_{i_t}e^{-\beta E_{i_0} + \nu J_{i_0 \rightarrow i_1} + \nu J_{i_1 \rightarrow i_2} + \cdots + \nu J_{i_{t-1} \rightarrow i_t}}e^{-\beta E_{i_t} + \nu J_{i_{t-1} \rightarrow i_t}}e^{-(t+1)\nu}.
\]

(90)

So it is not exactly as simple as equation (9): besides the expected Boltzmann–Gibbs factors involving the trajectory energy \(E(\Omega_{[0,t+1]})\) (already found for the equilibrium case in equation (55)) and the trajectory current \(J(\Omega_{[0,t]}),\) and besides the normalization ensured by the choice of \(\rho\) in the last factor, there remains the non-trivial boundary prefactor \(y_{i_0}z_{i_t}:\) we believe that this term comes from the stationary constraint on the dynamics that we could impose within the Markov chain framework, whereas we were not able to impose this stationary constraint at the formal level discussed in section 2. However, this supplementary prefactor is a ‘boundary term’ containing only the initial and the final configurations. As a consequence, it is not expected to grow in time, and in the large-time limit \(t \rightarrow +\infty,\) it will become subleading with respect to the time-extensive terms present in the exponential, since the trajectory energy and the trajectory current are extensive in time by the imposed constraints. (Note however that in some cases with unbounded phase space, the ‘boundary terms’ may diverge and remain important even in the large-time limit, see section 5.3 of the review [29] and references therein for more details.)

Our conclusion is thus that the formal solution of equation (9) neglects only the boundary terms of equation (90) and thus captures correctly the dominant terms in the large-time limit \(t \rightarrow +\infty.\) In particular, it contains the factor that is responsible for the fluctuation relation of the integrated current discussed in section 2.3 and for the dominant term of the entropy production discussed in section 2.4. We will obtain the same conclusion in the presence of locality constraints on the dynamics that we consider in section 6.

6. Effective stochastic microscopic dynamics in the non-equilibrium case

As explained in section 4.6, there exists some difficulty to recover the equilibrium if one tries to maximize the dynamical entropy associated with a Markov chain containing locality constraints, as a consequence of the possible configuration-dependent connectivities that one wishes to impose on the possible elementary moves. However, as recalled in section 4.5, the requirements on effective stochastic microscopic dynamics (equation (59)) to describe the equilibrium dynamics are well understood, and lead to the detailed balance condition of equation (63). In this section, we will thus follow the point of view of the Evans approach [11]–[14]: we consider that a local equilibrium dynamics generated by some Markov chain \(w^\text{eq}_{i \rightarrow j}\) satisfying detailed balance is known, and we wish to determine the non-equilibrium appropriate modified Markov chain \(w_{i \rightarrow j}\) in the presence of an imposed current. Instead of maximizing the ‘full’ dynamical entropy of equation (2),

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we will thus maximize the relative dynamical entropy with respect to the equilibrium dynamics, which we have introduced in equation (10).

6.1. Maximization of the relative dynamical entropy with respect to the equilibrium

For the microscopic Markov chain

\[ p_{t+\Delta t}(j) = \sum_i p_t(i) w_{i\rightarrow j} \]

the relative dynamical entropy of equation (10) with respect to the equilibrium dynamical trajectory corresponding to \( w_{eq}^{i\rightarrow j} \) has for time-extensive behavior

\[ S_{rel}^{dyn}(t) \equiv -\sum_{\Omega(0,t)} P(\Omega(0,t)) \ln \frac{P(\Omega(0,t))}{P_{eq}(\Omega(0,t))} \sim -\frac{t}{\Delta t} \sum_i p_{st}(i) \sum_j w_{i\rightarrow j} \ln \left( \frac{w_{i\rightarrow j}}{w_{eq}^{i\rightarrow j}} \right). \]  

The maximization of this relative entropy in the presence of the constraints concerning

the normalizations

\[ \sum_i p_{st}(i) = 1 \]

\[ \sum_j w_{i\rightarrow j} = 1 \]

the stationarity condition

\[ \sum_j p_{st}(j) w_{j\rightarrow i} - p_{st}(i) = 0 \]

and the current

\[ J_{dyn}(t) = \frac{1}{t} \sum_i p_{st}(i) \sum_j w_{i\rightarrow j} J_{i\rightarrow j} = J_0 \]

can be done with Lagrange multipliers, with steps similar to the ones detailed in section 6.

To write the final result, it is convenient to introduce the positive non-symmetric matrix

\[ \langle i | M | j \rangle = w_{i\rightarrow j}^{eq} e^\nu J_{i\rightarrow j} \]

its largest eigenvalue \( e^\rho \), and the corresponding right and left positive eigenvectors \( |R\rangle \) and \( \langle L | \)

\[ M | R \rangle = e^\rho | R \rangle \]

\[ \langle L | M = e^\rho \langle L | \]

normalized with

\[ \langle L | R \rangle = e^\rho. \]
With these notations, the optimal solution reads

\[ p^{st}(i) = e^{-\rho} \langle i| R \rangle \langle L|i \rangle \]

\[ w_{i \rightarrow j} = w_{i \rightarrow j}^{eq} e^{-\nu J_{i \rightarrow j} - \rho} = e^{-\rho} \langle i| M |j \rangle \langle j|R \rangle. \tag{99} \]

In the equilibrium case \( \nu = 0 \), the matrix \( M \) reduces to the generator of the equilibrium Markov chain \( M_{ij} = w_{i \rightarrow j}^{eq} \), so the maximal eigenvalue corresponds to \( e^{\rho_{\text{eq}} = 1} \), the right and the left eigenvectors are simply \( \langle i| R_{\text{eq}} \rangle = 1 \) and \( \langle L_{\text{eq}}|i \rangle = p^{eq}(i) \).

6.2. Consequences for the probability distribution of dynamical trajectories

From the solution of equation (99), one obtains that the probability of a trajectory \( \omega = \{i_0, i_1, \ldots, i_n\} \) takes the form

\[
\mathcal{P}(\omega = \{i_0, i_1, \ldots, i_n\}) = p^{st}(i_0)w_{i_0 \rightarrow i_1}w_{i_1 \rightarrow i_2} \cdots w_{i_{n-1} \rightarrow i_n} \\
= e^{-\rho \langle i_0| R \rangle \langle L|i_0 \rangle} w_{i_0 \rightarrow i_1}^{eq} e^{-\rho + \nu J_{i_0 \rightarrow i_1} - \rho} \cdots w_{i_{n-1} \rightarrow i_n}^{eq} e^{-\rho + \nu J_{i_{n-1} \rightarrow i_n} - \rho} \langle i_n| R \rangle \langle i_{n-1}| R \rangle \\
= \left[ \frac{\langle L|i_0 \rangle \langle i_n| R \rangle}{p^{eq}(i_0)} \right] \mathcal{P}^{eq}(\omega = \{i_0, i_1, \ldots, i_n\}) e^{\nu J(\omega) - (n+1)\rho}. \tag{100} \]

Again, it is not exactly as simple as equation (11): besides the equilibrium trajectory probability, the expected Boltzmann–Gibbs factors involving the trajectory current \( J(\Omega_{[0,t]} \rangle) \), and besides the normalization insured by the choice of \( \rho \) in the last factor, there remains a boundary prefactor that depends only on the initial and final configurations. But this boundary factor is expected to become subleading in the large-time limit \( t \to +\infty \), as already discussed in section 5.4.

6.3. Link with the Bayesian approach of Evans [11]

In [11], Evans has introduced an approach called ‘Non-equilibrium Counterpart to detailed balance’. The general idea is clearly the same as in the present paper, namely the maximization of the dynamical entropy in the presence of the appropriate constraints. However, the way of reasoning is slightly different and thus gives other useful points of view: Bayes theorem is used to analyze the properties of an elementary trajectory belonging to a macroscopic trajectory satisfying the flux constraint with respect to the equilibrium dynamics. The outcome is that the transitions \( w_{i \rightarrow j}^{\text{driven}} \) in the driven case should have the following form (see equation (24) of [11])

\[ w_{i \rightarrow j}^{\text{driven}} = w_{i \rightarrow j}^{eq} e^{\nu J_{i \rightarrow j} + q_j(\nu) - q_i(\nu) - Q(\nu) \Delta t}. \tag{101} \]

From the comparison with equation (99), one obtains the following correspondence: \( Q(\nu) \Delta t \) corresponds to the normalization factor \( \rho \), whereas \( q_j(\nu) \) corresponds to \( \ln \langle j| R \rangle \). The physical interpretation proposed by Evans is that the factor \( q_j(\nu) \) measures the ‘willingness’ of state \( j \) to accept a future flux. We refer the reader to the very interesting series of articles [11]–[14] to have more detailed explanations and to see various examples of application.
7. Case of continuous-time master equation

Up to now we have considered the case of discrete-time Markov chains, which constitute the simplest framework to define the probabilities of dynamical trajectories and their Shannon entropy of equation (2). However, many studies on non-equilibrium systems prefer to consider stochastic dynamics that are generated by some continuous-time master equation

$$\partial_t p_t(j) = \sum_{i \neq j} p_t(i) k_{i \rightarrow j} - p_t(j) \left[ \sum_{i \neq j} k_{j \rightarrow i} \right]. \quad (102)$$

In this section, we describe how the results of the previous sections should be adapted for this case.

7.1. Dynamical entropy for the continuous-time master equation

Let us first recall how the master equation (102) can be obtained as the continuous limit $\Delta t \to 0$ of the Markov chain of equation (91). In the limit where the elementary time step becomes small $\Delta t \to 0$, it is natural to assume that the transition probability $w_{i \rightarrow j}$ from $i$ to another state $j \neq i$ becomes proportional to $\Delta t$

$$w_{i \rightarrow j} \simeq \Delta t \rightarrow 0 k_{i \rightarrow j} \Delta t \quad (103)$$

where $k_{i \rightarrow j}$ represents the transition rate per unit time from $i$ to $j$. The normalization of equation (25) yields that the transition probability to remain on site $i$ takes the form

$$w_{i \rightarrow i} = 1 - \sum_{j \neq i} w_{i \rightarrow j} \simeq 1 - \Delta t k_{\text{out}}(i) \quad (104)$$

where

$$k_{\text{out}}(i) \equiv \sum_{j \neq i} k_{i \rightarrow j} \quad (105)$$

represents the total exit rate out of state $i$. In this limit $\Delta t \to 0$, the finite-time Markov chain of equation (91) thus becomes the master equation (102).

However as explained in [17], the continuous-time description leads to some difficulties when one considers the dynamical entropy of equation (2). Indeed, the naive continuous limit of equation (32) becomes

$$S_{\text{dyn}}(t) = -\frac{1}{\Delta t} \sum_i p^{st}(i) \sum_j w_{i \rightarrow j} \ln w_{i \rightarrow j} \quad \simeq \Delta t \to 0 -\frac{1}{\Delta t} \sum_i p^{st}(i) \left[ \sum_{j \neq i} k_{i \rightarrow j} \Delta t \ln(k_{i \rightarrow j} \Delta t) + (1 - \Delta t k_{\text{out}}(i)) \ln(1 - \Delta t k_{\text{out}}(i)) \right]$$

$$\simeq \Delta t \to 0 - \sum_i p^{st}(i) \left[ \sum_{j \neq i} k_{i \rightarrow j} \ln(k_{i \rightarrow j} \Delta t) - k_{\text{out}}(i) \right] \quad (106)$$

i.e. the elementary time $\Delta t$ remains in the argument of the logarithm to have a non-dimensional quantity. This difficulty to define the ‘absolute’ entropy in problems involving
continuous variables can be usually circumvented by considering the ‘relative’ entropy with respect to some reference. In our present framework, the reference will be the corresponding equilibrium master equation defined by transition rates $k_{eq}^{i\rightarrow j}$ satisfying the detailed balance condition (equation (63))

$$p^{eq}(i)k_{eq}^{i\rightarrow j} = p^{eq}(j)k_{eq}^{j\rightarrow i}. \quad (107)$$

The continuous-time limit $\Delta t \to 0$ of the relative dynamical entropy of equation (92) reads

$$S_{rel}^{dyn}(t) = -\frac{1}{\Delta t} \sum_i p^{st}(i) \sum_j w_{i\rightarrow j} \ln \left(\frac{w_{i\rightarrow j}}{w_{eq}^{i\rightarrow j}}\right)$$

$$= -\frac{1}{\Delta t} \sum_i p^{st}(i) \left[ \sum_{j \neq i} k_{i\rightarrow j} \Delta t \ln \left(\frac{k_{i\rightarrow j}}{k_{eq}^{i\rightarrow j}}\right) + (1 - \Delta t k_{out}^{eq}(i)) \ln \left(\frac{1 - \Delta t k_{out}^{eq}(i)}{1 - \Delta t k_{eq}^{i\rightarrow j}}\right) \right]$$

$$\sim \Delta t \to 0 \sum_i p^{st}(i) \left[ -\sum_{j \neq i} k_{i\rightarrow j} \ln \left(\frac{k_{i\rightarrow j}}{k_{eq}^{i\rightarrow j}}\right) + k_{out}^{eq}(i) - k_{eq}^{i\rightarrow j} \right]$$

$$\sim \Delta t \to 0 \sum_i p^{st}(i) \left[ -\sum_{j \neq i} k_{i\rightarrow j} \ln \left(\frac{k_{i\rightarrow j}}{k_{eq}^{i\rightarrow j}}\right) + \sum_{j \neq i} k_{i\rightarrow j} - \sum_{j \neq i} k_{eq}^{i\rightarrow j} \right]. \quad (108)$$

So this relative dynamical entropy is well defined for continuous-time master equations, and can be used in maximization procedures.

7.2. Maximization of the relative entropy with the appropriate constraints

We wish to maximize the relative dynamical entropy of equation (108) with respect to all stationary states $p^{st}(i)$ satisfying the normalization (as in equation (38))

$$N \equiv \sum_i p^{st}(i) = 1 \quad (109)$$

and with respect to all possible transition rates $k_{i\rightarrow j}$ with $j \neq i$ (note that here there is no normalization equation as equation (39), since it has already been taken into account in equations (104) and (105)) that have $p^{st}(i)$ as stationary distribution (equivalent of equation (40))

$$\Sigma_i \equiv \sum_{j \neq i} p^{st}(j)k_{j\rightarrow i} - p^{st}(i) \sum_{j \neq i} k_{i\rightarrow j} = 0 \quad (110)$$

in the presence of the following flux constraint (equivalent of equation (95))

$$\frac{J_{dyn}(t)}{t} \equiv \sum_i p^{st}(i) \sum_{j \neq i} k_{i\rightarrow j} J_{i\rightarrow j} = J_0. \quad (111)$$
So we introduce the following functional with appropriate Lagrange multipliers

\[ \Psi \equiv \frac{S_{\text{rel}}^\text{dyn}(t)}{t} - \rho(N - 1) - \sum_i \mu_i (\Sigma_i) + \nu \left( \frac{J^\text{dyn}(t)}{t} - J_0 \right) \]

\[ = \sum_i p^\text{st}(i) \left[ -\sum_{j\neq i} k_{i\rightarrow j} \ln \left( \frac{k_{i\rightarrow j}^\text{eq}}{k_{i\rightarrow j}} \right) + \sum_{j\neq i} k_{i\rightarrow j} - \sum_{j\neq i} k_{i\rightarrow j}^\text{eq} \right] - \rho \left( \sum_i p^\text{st}(i) - 1 \right) \]

\[ - \sum_i \mu_i \left( \sum_{j\neq i} p^\text{st}(j) k_{j\rightarrow i} - p^\text{st}(i) \sum_{j\neq i} k_{i\rightarrow j} \right) + \nu \left( \sum_i p^\text{st}(i) \sum_{j\neq i} k_{i\rightarrow j} J_{i\rightarrow j} - J_0 \right). \]

(112)

The optimization with respect to \( k_{i\rightarrow j} \) with \( i \neq j \)

\[ 0 = \frac{\delta \Psi}{\delta k_{i\rightarrow j}} = p^\text{st}(i) \left[ -\ln \left( \frac{k_{i\rightarrow j}^\text{eq}}{k_{i\rightarrow j}} \right) - \mu_j + \mu_i + \nu J_{i\rightarrow j} \right] \]

yields

\[ k_{i\rightarrow j} = k_{i\rightarrow j}^\text{eq} e^{\mu_i - \mu_j + \nu J_{i\rightarrow j}} \]

(114)

which is the appropriate continuous limit of the Markov chain result of equation (99) and of the Evans result cited in equation (101).

The optimization with respect to \( p^\text{st}(i) \)

\[ 0 = \frac{\delta \Psi}{\delta p^\text{st}(i)} = \sum_{j\neq i} k_{i\rightarrow j} \left[ -\ln \left( \frac{k_{i\rightarrow j}^\text{eq}}{k_{i\rightarrow j}} \right) + 1 - \mu_j + \mu_i + \nu J_{i\rightarrow j} \right] - \sum_{j\neq i} k_{i\rightarrow j}^\text{eq} - \rho \]

(115)

can be simplified using equation (113) in

\[ 0 = \sum_{j\neq i} k_{i\rightarrow j} - \sum_{j\neq i} k_{i\rightarrow j}^\text{eq} - \rho \]

(116)

or equivalently in terms of the total exit rates introduced in equation (105), one obtains

\[ k_{\text{out}}(i) = k_{\text{out}}^\text{eq}(i) + \rho. \]

(117)

Physically, this means that for all configurations \( i \), the exit rate is shifted by the same amount \( \rho \) with respect to the equilibrium case. Equation (117) has been previously obtained by Baule and Evans in [12] within their slightly different perspective, in order to obtain a set of invariant quantities for shear flows and to compute the non-equilibrium rates via some network rules (see [12] for more details).

In contrast equations (113) and (117) can appear as ‘new’ with respect to the discrete-time Markov chain result of equation (99). Another apparent difference is that the discrete-time Markov chain result of equation (99) contains both the stationary distribution \( p^\text{st}(i) \) and the transition probabilities \( w_{i\rightarrow j} \) that are determined together, whereas here the two types of equations (equation (113) and equation (117)) concern only the transition rates, and \( p^\text{st}(i) \) no longer appears. However, this absence of \( p^\text{st}(i) \) is only apparent, because \( p^\text{st}(i) \) is of course directly determined by the rates \( k_{i\rightarrow j} \) via the stationarity condition of equation (110). To make clearer the correspondence with the discrete-time Markov chain result of equation (99), it is thus useful to reformulate the continuous-time master equation solution as follows.

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7.3. Reformulation of the solution in terms of an eigenvalue problem

The aim of this section is to reformulate the solution found above in equations (113) and (117) in terms of an eigenvalue problem to see the similarity with the discrete-time Markov chain solution of equation (99).

To replace the matrix $M$ of equation (96), it is convenient to introduce the matrix $N$ defined by the matrix elements

$$
N_{ij} = \begin{cases} 
N_{ij}^{\text{eq}} & \text{if } j \neq i \\
- \sum_{j \neq i} N_{ij}^{\text{eq}} & \text{if } i = j
\end{cases} = -k_{\text{out}}^{\text{eq}}(i).
$$

Setting $e^{-\mu_j} = \langle j | R \rangle$, equation (113) becomes

$$
k_{i \rightarrow j} = N_{ij} \langle j | R \rangle \langle i | R \rangle
$$

whereas equation (117) becomes

$$
\rho = k_{\text{out}}(i) - k_{\text{eq}}^{\text{out}}(i) = \sum_{j \neq i} N_{ij} \frac{\langle j | R \rangle}{\langle i | R \rangle} + N_{ii} + \sum_{j \neq i} N_{ij} \frac{\langle j | R \rangle}{\langle i | R \rangle} = \frac{\langle i | N | R \rangle}{\langle i | R \rangle}
$$

meaning that $|R\rangle$ is a right eigenvector of the matrix $N$ corresponding to the eigenvalue $\rho$

$$
N |R\rangle = \rho |R\rangle.
$$

Let us now rewrite the stationary equation of equation (110) using equations (118), (119) and (117)

$$
0 = \sum_{i \neq j} p^\text{st}(i) k_{i \rightarrow j} - p^\text{st}(j) k_{\text{out}}(j) = \sum_{i \neq j} p^\text{st}(i) N_{ij} \frac{\langle j | R \rangle}{\langle i | R \rangle} - p^\text{st}(j) (k_{\text{out}}^{\text{eq}}(j) + \rho).
$$

It is thus convenient to set

$$
p^\text{st}(i) = \langle i | R \rangle \langle L | i \rangle
$$

and rewrite equation (122) as

$$
0 = \sum_{i \neq j} \langle L | i \rangle N_{ij} + \langle L | j \rangle (N_{jj} - \rho)
$$

$$
= \sum_i \langle L | i \rangle \langle i | N | j \rangle - \rho \langle L | j \rangle = \langle L | N | j \rangle - \rho \langle L | j \rangle
$$

meaning that $\langle L \rangle$ is a left eigenvector of the matrix $N$ corresponding to the eigenvalue $\rho$

$$
\langle L | N = \rho \langle L \rangle.
$$

In summary, the solution of equations (113) and (117) for the continuous-time master equation can be written in terms of an eigenvalue problem for the matrix $N$ introduced

$$
doi:10.1088/1742-5468/2011/03/P03008
$$
in equation (118) as follows:

\[ p^\text{st}(i) = \langle i|R\rangle \langle L|i \rangle \]
\[ k_{i\rightarrow j} = N_{ij} \frac{\langle j|R \rangle}{\langle i|R \rangle} \]  

(126)

where \( |R\rangle \) and \( \langle L| \) are the right and the left positive eigenvectors associated with the largest eigenvalue \( \rho \) of the matrix \( N \), with the following normalization

\[ 1 = \langle L|R \rangle. \]  

(127)

The correspondence with the discrete-time Markov chain result of equation (99) is now clear. As a final remark, let us mention that in the equilibrium case \( \nu = 0 \), the matrix \( N \) is the generator of the equilibrium dynamics, so the maximal eigenvalue corresponds to \( \rho_{\text{eq}} = 0 \), the corresponding right and the left eigenvectors being simply \( \langle i|R_{\text{eq}} \rangle = 1 \) and \( \langle L_{\text{eq}}|i \rangle = p_{\text{eq}}(i) \).

7.4. Relation with the ‘constrained dynamics’ introduced to study large deviations

The reformulation in terms of an eigenvalue problem described in section 7.3 is also useful to make the link with the ‘constrained dynamics’ that have been introduced in various studies concerning large deviations of stochastic processes described by master equations (see for instance the recent works [39]–[42] and references therein). In this context, the point of view is as follows: one considers some ‘true’ dynamics and one is interested in large deviation properties of some observable like the current. To compute the probability of rare events giving rise to an anomalous current for the ‘true’ dynamics, it is useful to introduce an ‘auxiliary’ dynamics that takes into account the conditioning to produce a given anomalous current (see [39]–[42] for more details). The transition rules of this auxiliary dynamics and its stationary state are given in terms of left/right eigenvectors of a modified operator via the same formula written above, see for instance equations (2.15) (2.17) (2.18) (2.19) in [39] or equations (2.16) (2.18) and (2.19) in [40]: the exchange of the roles of the right and left eigenvectors with respect to the present notations comes from the different choice in the writing of matrices, since [39,40] have chosen the ‘quantum mechanical’ convention where the initial state is on the right and the final state is on the left. The fact that exactly the same equations appear can be understood as follows. In the present paper, we have described how the maximization of the dynamical entropy yields equation (11) at a formal level, and give the results of the previous section when applied to an explicit master equation. But of course another possibility could be to take the formal solution of equation (11) as a starting point, and to derive the consequences of this reweighting of trajectories for the specific case of a master equation: this is exactly the route followed in the large deviation studies mentioned above [39]–[42] and so one should indeed obtain the same results by consistency. However, even if the equations are exactly the same, the physical interpretation of this modified dynamics is different: in the large deviation studies [39]–[42], this modified dynamics is usually presented only as a useful technical tool to better understand the large deviations of the ‘true’ dynamics, whereas in the approach summarized in the present paper, one considers that the modified dynamics is the ‘real’ dynamics in the presence of an imposed current.
An interesting output of this comparison with large deviation studies is that it provides specific studies of this modified dynamics in models different from the examples considered in the works of Evans [11]–[14]: the auxiliary dynamics corresponding to the large deviations of the current of the asymmetric simple exclusion process has been studied in various situations or regimes in [39, 41, 42], whereas constrained dynamics for the Glaube–Ising chain has been considered in [40]. Since the change of measure is a standard tool in the large deviation theory, many other examples can actually be found in the huge number of works concerning large deviations for stochastic processes.

8. Conclusions and perspectives

In this paper, we have argued that if one wishes to formulate a general principle based on the maximization of some notion of ‘entropy’ for non-equilibrium steady states, the Shannon entropy associated with the probability distribution of dynamical trajectories is definitely the most natural, as first proposed by Filyokov and Karpov in 1967. The general idea is then to maximize the dynamical entropy of equation (2) in the presence of appropriate constraints, including the macroscopic current of interest, via the method of Lagrange multipliers. We have tried to give a self-contained and unified presentation of this type of approach. We have first described at the formal level how this maximization leads to a generalized Gibbs distribution for the probability distribution of dynamical trajectories, and to some fluctuation relation of the integrated current. We have then discussed in detail the cases of well-defined stochastic dynamics generated either by discrete-time Markov chains or by continuous-time master equations. In the cases where the use of the ‘full’ dynamical entropy of equation (2) leads to some difficulties, we have shown how the use of the ‘relative’ dynamical entropy of equation (10) allows one to solve the problems. The obtained results are in full agreement with the Evans approach called ‘Non-equilibrium Counterpart to detailed balance’ [11]–[14], but give a slightly different perspective. In particular, we have obtained that the stationary distribution and the transition probabilities or transition rates could be obtained in various cases from an eigenvalue problem. Finally, we have explained the link with the constrained dynamics that are often introduced in large deviation studies of stochastic processes.

In this paper, we have remained at a very general level with a dynamics visiting a series of configurations, to see more clearly what general properties could emerge, since it is at this level of generality that the statistical physics theory of equilibrium is formulated. However it is clear that it would be interesting in the future to study more precisely how this type of approach can be applied in various models of interest, and to discuss whether the obtained dynamics can be considered as ‘real’. An important issue is of course to compare with other approaches that solve exactly some non-equilibrium dynamics. In the field of quantum models, the non-equilibrium steady states have been found to be generalized Gibbs states, but they involve an infinite number of conserved quantities coming from the integrability of the considered models [43, 44]. For non-integrable models, one may thus hope that the non-equilibrium steady states correspond to much simpler generalized Gibbs states.

As a final remark, we should stress that the idea of maximizing some dynamical entropy to describe non-equilibrium steady states has been argued here to hold for ‘physical systems’ for which, in the absence of any imposed current, the notion of
equilibrium exists and corresponds to the maximization of the usual ‘static’ entropy of statistical physics, and for which effective stochastic models satisfying detailed balance are well accepted to describe the equilibrium dynamics. But of course, in the field of non-equilibrium, many models of interest are inspired not by physics, but by biology (such as predator–prey models), sociology (such as road traffic models), politics (such as voter models) etc. In all these cases, it is clear that one can choose arbitrarily the microscopic rates to model the considered phenomenon as one wishes.

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Appendix A. Reminder on the maximization of the static entropy for the equilibrium state

At equilibrium, one is interested into the equilibrium probabilities $P^{\text{eq}}(i)$ of occupations of configurations $i$. The Shannon entropy associated with the equilibrium distribution is

\[
S^{\text{eq}} = -\sum_i P^{\text{eq}}(i) \ln P^{\text{eq}}(i). \tag{A.1}
\]

(i) If the only constraint is the normalization,

\[
N \equiv \sum_i P^{\text{eq}}(i) = 1 \tag{A.2}
\]

the maximization of the entropy of equation (A.1) with the constraint of equation (A.2) can be obtained by introducing a Lagrange multiplier $\lambda$ and the functional

\[
\Phi_\lambda = S^{\text{eq}} - \lambda(N - 1) = -\sum_i P^{\text{eq}}(i) \ln P^{\text{eq}}(i) - \lambda \left( \sum_i P^{\text{eq}}(i) - 1 \right). \tag{A.3}
\]

The optimization with respect to $P^{\text{eq}}(j)$ yields

\[
0 = \frac{\delta \Phi_\lambda}{\delta P^{\text{eq}}(j)} = -\ln P^{\text{eq}}(j) - 1 - \lambda \tag{A.4}
\]

leading to $P^{\text{eq}}(j) = e^{-1-\lambda}$, where $\lambda$ is determined by the normalization constraint of equation (A.2). So $P^{\text{eq}}(j)$ is simply uniform over all states.

(ii) If in addition to the normalization of equation (A.2), one imposes also a constraint for the averaged energy

\[
\langle E \rangle \equiv \sum_i E_i P^{\text{eq}}(i) = E_0 \tag{A.5}
\]
Non-equilibrium steady states

one has to introduce another Lagrange multiplier $\beta$ and the functional

$$\Phi_{\lambda, \beta} = S^{eq} - \lambda(N - 1) - \beta(\langle E \rangle - E_0)$$

$$= -\sum_i P^{eq}(i) \ln P^{eq}(i) - \lambda \left( \sum_i P^{eq}(i) - 1 \right) - \beta \left( \sum_i E_i P^{eq}(i) - E_0 \right).$$

(A.6)

The optimization with respect to $P^{eq}(j)$ yields

$$0 = \frac{\delta \Phi_{\lambda, \beta}}{\delta P^{eq}(j)} = - \ln P^{eq}(j) - 1 - \lambda - \beta E_j$$

(A.7)

leading to $P^{eq}(j) = e^{-1-\lambda-\beta E_j}$, where $\lambda$ is determined by the normalization constraint of equation (A.2), and $\beta$ by equation (A.5). Then $P^{eq}(j)$ follows the Boltzmann–Gibbs distribution

$$P^{eq}(j) = \frac{e^{-\beta E_j}}{\sum_i e^{-\beta E_i}}.$$  

(A.8)

This derivation of the Boltzmann–Gibbs distribution, introduced by Jaynes [23], has the advantage of being very simple and really showing what the principle of maximum entropy contains. For more details concerning the consideration of statistical physics from the point of view of the Shannon information entropy, we refer the reader to the ‘old’ book [46] and to recent presentations [47] and references therein.

Appendix B. Technical simplifications for an alternate Markov chain

In their original paper [7], Filyokov and Karpov did not consider the homogeneous Markov chain of equation (24), but focused instead on an alternate Markov chain, which has been also reconsidered recently by Favretti [10]. In this appendix, we show how this alternate Markov chain framework yields technical simplifications with respect to the solution described in section 5 concerning the homogeneous Markov chain.

B.1. Alternate Markov chain

In this section, we consider the alternate Markov chain, where the transition probabilities $W_{i \rightarrow j}$ take alternately two sets of values $A_{i \rightarrow j}$ and $B_{i \rightarrow j}$, so that the probability of a trajectory of equation (28) now becomes

$$P(\Omega_{0,2t} = \{i_0,i_1,\ldots,i_{2t}\}) = P^{even}(i_0)A_{i_0 \rightarrow i_1}B_{i_1 \rightarrow i_2} \cdots A_{i_{2t-2} \rightarrow i_{2t-1}}B_{i_{2t-1} \rightarrow i_{2t}}$$

(B.1)

with the normalizations (as in equation (25))

$$N^A_i \equiv \sum_j A_{i \rightarrow j} = 1$$

$$N^B_i \equiv \sum_j B_{i \rightarrow j} = 1.$$  

(B.2)
The stationary states \( P_{\text{even}}(i) \), \( P_{\text{odd}}(i) \) at even and odd times associated with this alternate Markov chain satisfy

\[
\Sigma_{i}^{\text{even}} \equiv \sum_{j} P_{\text{odd}}(j) B_{j\rightarrow i} - P_{\text{even}}(i) = 0 \tag{B.3}
\]

\[
\Sigma_{i}^{\text{odd}} \equiv \sum_{j} P_{\text{even}}(j) A_{j\rightarrow i} - P_{\text{odd}}(i) = 0
\]

with the normalizations

\[
\mathcal{N}^{\text{even}} \equiv \sum_{i} P_{\text{even}}(i) = 1 \tag{B.4}
\]

\[
\mathcal{N}^{\text{odd}} \equiv \sum_{i} P_{\text{odd}}(i) = 1.
\]

The dynamical entropy of equation (31) becomes

\[
S_{\text{dyn}}(2t) = -t \left[ \sum_{i} P_{\text{even}}(i) \sum_{j} A_{i\rightarrow j} \ln A_{i\rightarrow j} + \sum_{j} P_{\text{odd}}(i) \sum_{j} B_{i\rightarrow j} \ln B_{i\rightarrow j} \right]. \tag{B.5}
\]

The averaged energy is now fixed by the constraint

\[
\frac{E_{\text{dyn}}(2t)}{2t} = \frac{1}{2} \left[ \sum_{i} P_{\text{even}}(i) E_{i} + \sum_{i} P_{\text{odd}}(i) E_{i} \right] = E_{0} \tag{B.6}
\]

that replaces equation (41).

### B.2. Simplification concerning the current constraint

In [7, 10], the interest was in an energy current flowing through the system: the advantage of this formulation with an alternate Markov chain is that one may consider that during the evolution described by \( A \), there is a flow of energy entering the system, whereas during the evolution described by \( B \), there is a flow of energy coming out of the system. To remain in a stationary state, these two flows have to be equal, and can be simply expressed in terms of the difference of averaged energies between the even and odd stationary states

\[
J_{\text{in}}^{E} = \sum_{i} P_{\text{even}}(i) \sum_{j} A_{i\rightarrow j} (E(j) - E(i)) = \sum_{j} E(j) P_{\text{odd}}(j) - \sum_{i} E(i) P_{\text{even}}(i)
\]

\[
J_{\text{out}}^{E} = \sum_{i} P_{\text{odd}}(i) \sum_{j} B_{i\rightarrow j} (E(i) - E(j)) = \sum_{i} E(i) P_{\text{odd}}(i) - \sum_{j} E(j) P_{\text{even}}(j) = J_{\text{in}}^{E}. \tag{B.7}
\]

In the following, we consider more generally the case of a current \( J_{K} \) that can be written similarly as the difference between some observable \( K \) between the even and odd stationary states (\( K \) is the ‘charge’ associated with the current \( J_{K} \))

\[
J_{\text{in}}^{K} = \sum_{i} P_{\text{even}}(i) \sum_{j} A_{i\rightarrow j} (K(j) - K(i)) = \sum_{j} K(j) P_{\text{odd}}(j) - \sum_{i} K(i) P_{\text{even}}(i)
\]

\[
J_{\text{out}}^{K} = \sum_{i} P_{\text{odd}}(i) \sum_{j} B_{i\rightarrow j} (K(i) - K(j)) = \sum_{i} K(i) P_{\text{odd}}(i) - \sum_{j} K(j) P_{\text{even}}(j) = J_{\text{in}}^{K}. \tag{B.8}
\]
so that the constraint of equation (64) can be replaced by
\[
J_{\text{dyn}}(2t) \equiv \sum_i K(i) \left[ P^{\text{odd}}(i) - P^{\text{even}}(i) \right].
\] (B.9)

**B.3. Optimization of the Lagrange functional**

So the functional of equation (65) is replaced by (some factors of 2 have been added to simplify slightly the notations)
\[
\Psi \equiv \frac{S_{\text{dyn}}(2t)}{t} - \rho_{\text{even}}(N_{\text{even}}^{A} - 1) - \rho_{\text{odd}}(N_{\text{odd}}^{B} - 1) - \beta \left( \frac{E_{\text{dyn}}(t)}{t} - E_0 \right) - \sum_i \lambda_i^{A} (N_i^{A} - 1) - \sum_i \lambda_i^{B} (N_i^{B} - 1) - \sum_i \mu_{i}^{\text{even}} (\Sigma_i^{\text{even}}) - \sum_i \mu_{i}^{\text{odd}} (\Sigma_i^{\text{odd}}) + \nu \left( J_{\text{dyn}}(2t) - J_0 \right)
\]
\[
= - \left[ \sum_i P^{\text{even}}(i) \sum_j A_{i \rightarrow j} \ln A_{i \rightarrow j} + \sum_j P^{\text{odd}}(i) \sum_j B_{i \rightarrow j} \ln B_{i \rightarrow j} \right] - \rho_{\text{even}} \left( \sum_i P^{\text{even}}(i) - 1 \right) - \rho_{\text{odd}} \left( \sum_i P^{\text{odd}}(i) - 1 \right) - \beta \left( \sum_i P^{\text{even}}(i) E_i + \sum_i P^{\text{odd}}(i) E_i \right) - 2E_0 \right) - \sum_i \lambda_i^{A} \left( \sum_j A_{i \rightarrow j} - 1 \right) - \sum_i \lambda_i^{B} \left( \sum_j B_{i \rightarrow j} - 1 \right) - \sum_i \mu_{i}^{\text{even}} \left( \sum_j P^{\text{odd}}(j) B_{j \rightarrow i} - P^{\text{even}}(i) \right) - \sum_i \mu_{i}^{\text{odd}} \times \left( \sum_j P^{\text{even}}(j) A_{j \rightarrow i} - P^{\text{odd}}(i) \right) + \nu \left( \sum_i K(i) \left[ P^{\text{odd}}(i) - P^{\text{even}}(i) \right] - J_0 \right).
\] (B.10)

The optimization with respect to \( A_{i \rightarrow j} \)
\[
0 = \frac{\delta \Psi}{\delta A_{i \rightarrow j}} = P^{\text{even}}(i) \left[ - \ln A_{i \rightarrow j} - 1 - \lambda_i^{A} - \mu_{j}^{\text{odd}} \right]
\] (B.11)
yields
\[
A_{i \rightarrow j} = e^{-1-\lambda_i^{A}+\mu_{j}^{\text{odd}}}. \] (B.12)
The normalization of equation (B.2) yields that $\lambda_i^A$ is independent of $i$ and given by
\[ e^{1+\lambda_i^A} = \sum_j e^{-\mu_j^{\text{odd}}}. \] (B.13)

The stationary equation of equation (B.3) yields, using equation (B.4)
\[ P^{\text{odd}}(j) = \sum_i P^{\text{even}}(i) A_{i\rightarrow j} = \left( \sum_i P^{\text{even}}(i) \right) \frac{e^{-\mu_j^{\text{odd}}}}{\sum_k e^{-\mu_k^{\text{odd}}}} = \frac{e^{-\mu_j^{\text{odd}}}}{\sum_k e^{-\mu_k^{\text{odd}}}}. \] (B.14)

The optimization with respect to $B_{i\rightarrow j}$ yields similar equations and solutions, so that finally one has the simple forms
\[ A_{i\rightarrow j} = P^{\text{odd}}(j) = \frac{e^{-\mu_j^{\text{odd}}}}{\sum_k e^{-\mu_k^{\text{odd}}}} \] (B.15)
\[ B_{i\rightarrow j} = P^{\text{even}}(j) = \frac{e^{-\mu_j^{\text{even}}}}{\sum_k e^{-\mu_k^{\text{even}}}}. \] (B.15)

The optimization of equation (B.10) with respect to $P^{\text{even}}(i)$ yields
\[ 0 = \frac{\delta\Psi}{\delta P^{\text{even}}(i)} = -\sum_j A_{i\rightarrow j} \ln A_{i\rightarrow j} - \rho^{\text{even}} - \beta E_i + \mu_i^{\text{even}} - \sum_j \mu_j^{\text{odd}} A_{i\rightarrow j} - \nu K(i) \] (B.16)
yields using equation (B.15)
\[ -\mu_i^{\text{even}} = -\rho^{\text{even}} - \beta E_i - \nu K(i) + \ln \left( \sum_k e^{-\mu_k^{\text{odd}}} \right) \] (B.17)
equation (B.15) then yields
\[ P^{\text{even}}(i) = \frac{e^{-\mu_i^{\text{even}}}}{\sum_k e^{-\mu_k^{\text{even}}}} = \frac{e^{-\beta E_i - \nu K_i}}{\sum_k e^{-\beta E_k - \nu K_k}}. \] (B.18)

Similarly, the optimization with respect to $P^{\text{even}}(i)$ yields
\[ P^{\text{odd}}(i) = \frac{e^{-\beta E_i + \nu K_i}}{\sum_k e^{-\beta E_k + \nu K_k}}. \] (B.19)

So here, in contrast to the case of the homogeneous Markov chain discussed in section 5, one obtains two Boltzmann–Gibbs distributions, without any prefactor like the functions $(y_i, z_i)$ in equation (82). This simple result was found on the special case of the energy flow $K_i = E_i$ by Favretti [10] (who has used the correct expression for the Kolmogorov–Sinai entropy of the alternate Markov chain, whereas an erroneous expression was actually used in the initial work [7]): the even and odd stationary distributions of equation (B.18) and (B.19) are then two Boltzmann–Gibbs distributions at two different temperatures that are fixed by the constraints on the averaged energy and on the averaged energy flow.

Despite the technical simplifications of the alternate Markov chain, we find that it is rather artificial from a physical point of view. Since the result means that the system oscillates between two distinct Boltzmann–Gibbs distributions, we feel that it could only correspond to physical situations where the energy is added or removed instantaneously.
at even and odd times, and that the system then relaxes to its new equilibrium during the macroscopic time $\tau$ of the Markov chain. Otherwise, if the energy were added and removed continuously during the time intervals, this would mean that one should somehow have thermal equilibrium at all times with an adiabatic change of temperature, which seems extremely restrictive. This is why, in the present paper, we have chosen to put the main focus on the case of the homogeneous Markov chain.

References

[1] Derrida B, 2010 arXiv:1012.1136
[2] Martyushev L M and Seleznev V D, 2006 Phys. Rep. 426 1
[3] Bruers S, 2006 arXiv:cond-mat/0604482
[4] Martyushev L M, Nazarova A S and Seleznev V D, 2007 J. Phys. A: Math. Theor. 40 371
[5] Bruers S, Maes C and Netocny K, 2007 J. Stat. Phys. 129 725
[6] Gaspard P, 1998 Chaos, Scattering and Statistical Mechanics (Cambridge: Cambridge University Press)
[7] Filyokov A A and Karpov V Ya, 1967 Inzh.-Fiz. Zh. 13 624
[8] Filyokov A A, 1967 Inzh.-Fiz. Zh. 13 798
[9] Filyokov A A, 1968 Inzh.-Fiz. Zh. 14 814
[10] Favretti M, 2009 Entropy 11 675
[11] Evans R M L, 2004 Phys. Rev. Lett. 92 150601
[12] Baule A and Evans R M L, 2008 Phys. Rev. Lett. 101 240601
[13] Simha A, Evans R M L and Baule A, 2008 Phys. Rev. E 77 031117
[14] Evans R M L, Simha R A, Baule A and Olmsted P D, 2010 Phys. Rev. E 81 051109
[15] Dewar R, 2003 J. Phys. A: Math. Gen. 36 631
[16] Burda Z, Duda J, Luck J M and Waclaw B, 2009 Phys. Rev. Lett. 102 160602
[17] Lecomte V, Appert-Rolland C and van Wijland F, 2005 Phys. Rev. Lett. 95 010601
[18] Garrahan J P, Jack R L, Lecomte V, Pitard E, van Duijvendijk K and van Wijland F, 2007 Phys. Rev. Lett. 98 195702
[19] Antal T, Racz Z and Sasvari L, 1997 Phys. Rev. Lett. 78 167
[20] Antal T, Racz Z, Rakos A and Schutz G M, 1998 Phys. Rev. E 57 5184
[21] Antal T, Racz Z, Rakos A and Schutz G M, 1999 Phys. Rev. E 59 4912
[22] Eisler V, Racz Z and van Wijland F, 2003 Phys. Rev. E 67 056129
[23] Eisler V and Zimors Z, 2005 Phys. Rev. A 71 042318
[24] Jaynes E T, 1957 Phys. Rev. 106 620
[25] Koralov L B and Sinai Y G, 2007 Theory of Probability and Random Processes (Berlin: Springer)
[26] Banavar J R and Maritan A, 2007 arXiv:cond-mat/0703622
[27] Banavar J R, Maritan A and Volkov I, 2010 J. Phys.: Condens. Matter 22 063101
[28] Touchette H, 2009 Phys. Rep. 478 1
[29] Derrida B, 2007 J. Stat. Mech. P07023
[30] Harris R J and Schütz G M, 2007 J. Stat. Mech. P07020
[31] Kurchan J, 2007 J. Stat. Mech. P07005

doi:10.1088/1742-5468/2011/03/P03008
Non-equilibrium steady states

[31] Sevick E M, Prabhakar R, Williams S R and Searles D J, 2008 Ann. Rev. Phys. Chem. 59 603
[32] Zia R K P and Schmittmann B, 2007 J. Stat. Mech. P07012
[33] Maes C, Netocny K and Shergelashvili B, A selection of nonequilibrium issues, 2009 Methods of Contemporary Mathematical Statistical Physics (Lecture Notes in Mathematics vol 1970) ed R Kotecký p 247
[34] Chetrite R, 2008 PhD Thesis available at http://perso.ens-lyon.fr/raphael.chetrite/indexfra.html
[35] Maes C, 1999 J. Stat. Phys. 95 367
[36] Maes C and Redig F, 2000 J. Stat. Phys. 101 3
Maes C, Netocny K and Wynants B, 2008 Markov Proc. Rel. Fields 14 445
[37] Schnakenberg J, 1976 Rev. Mod. Phys. 48 571
[38] Lebowitz J L and Spohn H, 1999 J. Stat. Phys. 95 333
[39] Simon D, 2009 J. Stat. Mech. P07017
[40] Jack R L and Sollich P, 2010 Prog. Theor. Phys. Suppl. 184 304
[41] Popkov V, Schutz G M and Simon D, 2010 J. Stat. Mech. P10007
Popkov V and Schutz G M, 2010 arXiv:1011.3913
[42] Simon D, 2011 arXiv:1011.3590
[43] Ogata Y, 2002 Phys. Rev. E 66 016135
Ogata Y, 2002 Phys. Rev. E 66 066123
[44] Karevski D and Platini T, 2009 Phys. Rev. Lett. 102 207207
[45] Herbert C, Paillard D and Dubrulle B, 2010 Earth Syst. Dyn. Discuss. 1 325
Herbert C, Paillard D, Kageyama M and Dubrulle B, 2011 Earth Syst. Dyn. at press, arXiv:1101.3173
[46] Brillouin L, 1956 Science and Information Theory (New York: Academic)
[47] Balian R, 1999 Am. J. Phys. 67 1078
Balian R, 2005 Stud. Hist. Phil. Mod. Phys. 36 323

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