Temporal Infomax on Markov Chains with Input Leads to Finite State Automata

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Abstract

Information maximization between stationary input and output activity distributions of neural ensembles has been a guiding principle in the study of neural codes. We have recently extended the approach to the optimization of information measures that capture spatial *and* temporal signal properties. Unconstrained Markov chains that optimize these measures have been shown to be almost deterministic. In the present work we consider the optimization of stochastic interaction in constrained Markov chains where part of the units are clamped to prescribed probability distributions. Temporal Infomax in that case leads to finite state automata.

Key words: Markov model; Information maximization; Finite state automata;

1 Introduction

One of the most basic questions in computational neuroscience is that for the nature of the neural code. Experiments suggest a considerable interaction of neurons already on the level of spikes, e.g., expressed by spatio-temporal correlations [1,4,6,7]. A well-known measure that quantifies relations of interacting units is the so-called *mutual information*: The Kullback-Leibler divergence

$$I(p) := D(p || p_1 \otimes \cdots \otimes p_N) = \sum_{\nu=1}^N H(p_\nu) - H(p),$$
 (1)

where $H(\cdot)$ denotes the Shannon entropy and p_{ν} the ν 'th marginal of p, measures the "distance" of p from the factorized distribution $\Pi_{\nu=1...N}p_{\nu}$. It is a

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natural measure for "spatial" interdependence of N stochastic units and a starting point of many approaches to neural complexity [5,6]. In order to capture intrinsically temporal aspects of dynamic interaction, I in (1) has been extended by Ay [2] to the dynamical setting of Markov transitions, where it is referred to as (stochastic) interaction. In a previous paper we have shown that the optimization of stochastic interaction in Markov chains leads to almost deterministic dynamical systems [3]. This work neglegted external input into the considered systems. The present study, therefore, presents optimized Markov chains, where a part of the system is clamped to externally prescribed stochastic processes with varying degree of spatio-temporal correlations. Surprisingly, the optimized processes turn out to be finite state automata.

2 Dynamic Infomax on Constrained Markov Chains

Consider the set $V = \{1, ..., N\}$ of binary units with state sets $\Omega_{\nu} = \{0, 1\}$, $\nu \in V$. For a subsystem $A \subset V$, $\Omega_A := \{0, 1\}^A$ denotes the set of all configurations restricted to A, and $\bar{P}(\Omega_A)$ is the set of probability distributions on Ω_A . Given two subsets A and B, where B is non-empty, $\bar{K}(\Omega_B | \Omega_A)$ is the set of all Markov transition kernels from Ω_A to Ω_B . In the following we only consider the case A = B and use the abbreviation $\bar{K}(\Omega_A)$. For a probability distribution $p \in \bar{P}(\Omega_A)$ and a Markov kernel $K \in \bar{K}(\Omega_A)$, the conditional entropy of (p, K) is defined as

$$H(p,K) = -\sum_{\omega,\omega'\in\Omega_A} p(\omega) K(\omega' | \omega) \ln K(\omega' | \omega) . \qquad (2)$$

H(p, K) has been used in [2] to generalize (1) to Markov transitions. For that purpose, we define the marginal kernels $K_{\nu} \in K(\Omega_{\nu}), \nu \in V$, of K by

$$K_{\nu}(\omega_{\nu}' \mid \omega_{\nu}) := \frac{\sum_{\substack{\sigma, \sigma' \in \Omega_{V} \\ \sigma_{\nu} = \omega_{\nu}, \sigma_{\nu}' = \omega_{\nu}' \\ \sigma_{\nu} = \omega_{\nu}}} p(\sigma) K(\sigma' \mid \sigma)}{\sum_{\substack{\sigma \in \Omega_{V} \\ \sigma_{\nu} = \omega_{\nu}}} p(\sigma)}, \qquad \omega_{\nu}, \omega_{\nu}' \in \Omega_{\nu}.$$
 (3)

Then, the stochastic interaction measure of K with respect to p is given by

$$I(p,K) := \sum_{\nu \in V} H(p_{\nu}, K_{\nu}) - H(p,K). \tag{4}$$

Equation (4) provides the required generalization of (1) to Markov transitions. I(p, K) is large if the marginal transitions have high entropy, but that of the full transition is low. Thus, supposed the current state $\omega \in \Omega_V$ is known, the next state is predictable with high probability, but, conversely, not much information is gained from knowledge about single units, ω_{ν} .

In the sequel we consider Markov chains $X_n = (X_{\nu,n})_{\nu \in V}$, $n = 0, 1, 2, \ldots$, given by an initial distribution $p_0 \in \bar{P}(\Omega_V)$ and a kernel $K \in \bar{K}(\Omega_V)$. We further restrict attention to parallel Markov chains. A Markov kernel $K \in \bar{K}(\Omega_V)$ is called parallel if there exist kernels $K^{(\nu)} \in \bar{K}(\Omega_V | \Omega_V)$, $\nu \in V$, such that

$$K(\omega' | \omega) = \prod_{\nu \in V} K^{(\nu)}(\omega'_{\nu} | \omega), \quad \text{for all} \quad \omega, \omega' \in \Omega_V.$$
 (5)

Parallel Markov chains are a more natural assumption in neural modeling than general Markov chains, because the action potential generation of a neuron depends only on its own input and internal dynamics. Whence, each ω'_{ν} in (5) is given an individual kernel $K^{(\nu)}$. Spatial correlations in firing patterns may nonetheless result from common input, i.e., the global state ω in (5).

In [3] we studied parallel Markov chains which maximized I(p,K) under no further constraint regarding K. The optimized chains were shown to be almost deterministic and representable in phase space by more or less complex graphs consisting of transient trajectories and nested loops which correspond with repeteting firing patterns. These studies did not consider external input into the system. The observed activity patterns may therefore be envisaged as intrinsic modes of activity in strongly interacting systems. The present work, in contrast, shows simulations of parallel Markov chains which maximize I(p,K) under the additional constraint that the kernels $K^{(\nu)}$ of a subset of units in (5) are fixed during the optimization process. The simulations show that Markov chains with such clamped "input-units" converge to finite state automata.

3 Simulations

The simulations displayed in the sequel implement the usual Markov dynamics on a set of N binary units together with a random search scheme to optimize the stochastic interaction of the Markov chains. That is, the interaction measure, I, is computed with respect to an induced stationary probability distribution of a parallel Markov kernel, and starting from initial random values the kernel is iteratively perturbed such that I increases. Details of the optimization are described in [3]. In contrast to [3], however, the optimization here is not unconstrained, but B < N kernels $K^{(\nu)}$ in (5) are fixed during the optimization. In especially, these kernels are chosen to be independent of the state of the internal units. In the simplest case the "input units" are just B independent (discrete time) Poisson processes with fixed rates. Cluster- or burst-processes with $a_{\nu} := K^{(\nu)}(1|0) = 1 - K^{(\nu)}(0|0)$ and $b_{\nu} := K^{(\nu)}(0|1) = 1 - K^{(\nu)}(1|1)$ implement Poisson processes with rate $\lambda_{\nu} = a_{\nu}$ for $a_{\nu} = 1 - b_{\nu}$. For other $a_{\nu}, b_{\nu} \in]0,1[$ they generate burst-like firing patterns with a_{ν} setting the interburst intervals and $1-b_{\nu}$ the burst durations. Spatial correlations between input units can further be introduced by appropriate input kernels. Details of

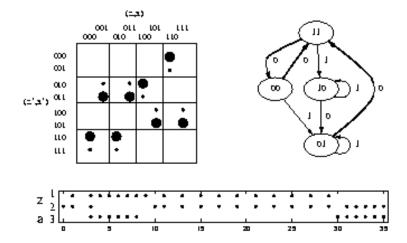


Fig. 1. Optimized example system with 2 internal units (1,2) and 1 input (3). Upper left: Markov kernel. The total kernel can be segregated into blocks defined by internal states (z, z') only. Right: corresponding automaton with internal states (z) as node- and input (a) as edge-symbols. Below: sample trajectory.

the simulations will be presented elsewhere. Because the optimization of I is algorithmically complex, simulations are restricted to small N.

Figure 3 shows an optimized systems with three units, where unit 3 has been clamped to a cluster process with $a_3 = 0.1$ and $b_3 = 0.1$. The unit has a firing probability of $\lambda = a_3/(a_3 + b_3) = 0.5$, and if it has fired the probability that it fires again in the next step is quite high: $K^{(\nu)}(1|1) = 1 - b_3 = 0.9$. The converged Markov transition kernel is displayed in the upper left of Fig.3 where circle area represents transition probability. Note, that the kernel can be divided into blocks according to the sub-states z, z' of the internal units, i.e., the first N - B bits of the total binary state representation: Given an internal state z and an input a (the B least significant bits of the total state), the internal target state z' in the next step is completely determined, only the next input a' is random. This means, that the Markov chain can be represented by a deterministic finite state automaton (DFA) with nodes labeled by the internal states and edges labeled by the (random) input patterns.

Figures 2 and 3 display examples with three units each and B=2. In Fig. 2 unit 2 is a Poisson process with $\lambda=0.1$ and unit 3 an independent but temporally correlated cluster process with $a_3=b_3=0.1$. In Fig. 3 the inputs (2,3) are spatially correlated near Poisson processes. The internal unit 1 combines activity from both input units. Both systems converge to comparable DFAs. Nonetheless, in Fig. 2 the externally induced temporal correlations are still expressed in the precise transition probabilites, e.g., three strong transitions to 000 because all units fire relatively seldom, or three strong transitions from states with unit 3 active to 101, because unit 3 fires long bursts but unit 2 is mostly silent. Similarly, spatial correlations are expressed in Fig. 3 by the dominating transitions $000 \rightarrow 000$ and $100 \rightarrow 100$, which reflect that input

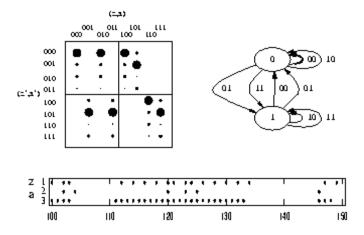


Fig. 2. Same as Fig. 3 but with two independent input units (2,3). Firing of the internal third unit (1) depends on both inputs but strongly reflects unit 3 activity.

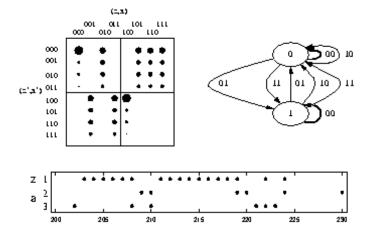


Fig. 3. Similar as Fig. 2 but with correlated inputs (2,3) favoring coactivation. The DFA is nonetheless similar to that in Fig. 2. Note also that the internal unit 1 acts as a latch (hysteresis, memory) activated from state 0 by input from unit 3.

units 2 and 3 are simultaneously silent in most steps. Activation of one input unit, however, often induces transitions to states where the same or other input unit fires, since both units excite themselves and each other.

Fig. 4 reveals an example with 4 units and 2 inputs. Here, as in the previous DFA-graphs, transitions for zero input, $a=0\cdots 0$, are drawn as thick solid edges. The respective subgraphs determine the system behavior in the absence of input and can be interpreted as intrinsic dynamic modes with which external input interacts. These modes appear in an almost identical way also under the unconstrained optimization described in [3]. Finally Fig. 5 shows that the optimization does not always converge to deterministic finite state machines: The arrow in Fig. 5B indicates 2 possible transitions from internal state z=01 to z'=10 and z'=11. Thus, the resulting finite state machine can make non-deterministic transitions given the same internal state and external input.

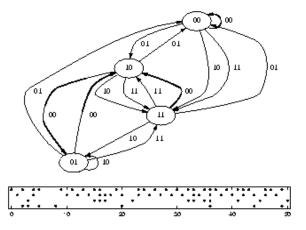


Fig. 4. A more complex example with N=4, B=2. Intrinsic modes (bold edges, input = 00) consist of a 1-cycle, a 2-cycle, and a transient transition from 11 to 10.

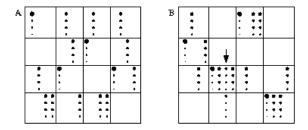


Fig. 5. A) Markov matrix corresponding with Fig. 4. B) Simulation as in (A) but the converged process contains non-deterministic internal transitions (arrow).

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