Order Determination for Functions of Finite Markov Chains

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1. Introduction

The simplest modeling problem for stochastic processes can be described as follows. We observe a scalar, finite valued, stationary stochastic process $\{Y_t\}$ (FVP) and try to model it as a probabilistic function of a finite Markov chain. The underlying assumption is that the probabilistic mechanism generating $\{Y_t\}$ can be reasonably well explained in the following way. There is an (unobserved) finite stationary Markov chain $\{X_t\}$ such that:

(1.1)
$$P\left[X_{t+1}, Y_{t+1} | X^t, Y^t\right] = P\left[X_{t+1}, Y_{t+1} | X_t\right]$$

Without loss of generality we can assume that the chain $\{X_t\}$ satisfying (1.1) is such that:

$$(1.2) P[X_{t+1}, Y_{t+1}|X_t] = P[Y_{t+1}|X_{t+1}] P[X_{t+1}|X_t]$$

Once $\{X_t\}$ has been found it is easy to describe $\{Y_t\}$ as a deterministic function of a Markov chain. The standard construction is to define the new process $S_t := (X_t, Y_t)$ with values in the cartesian product $\mathcal{S} := \mathcal{X} \times \mathcal{Y}$ of the state spaces of $\{X_t\}$ and $\{Y_t\}$ and the function $f:\mathcal{S}\to\mathcal{Y}$ as $f(i,\varepsilon)=\varepsilon$. It follows from (1.1) that $\{S_t\}$ is a Markov chain and $Y'_t := f(S_t)$ has the same law as $\{Y_t\}$. We can now give the following interpretation: the process we observe, $\{Y_t\}$, is just a component of a process, $\{S_t\}$, which is Markovian. As one would expect the "completion" of $\{Y_t\}$ to a Markov chain $\{S_t\}$ is inherently non unique. Define as size of the model the cardinality of the state space of $\{X_t\}$. Between all possible models for $\{Y_t\}$ we would like to select one of minimal size (order). The status of Realization Theory for FVP is rather unsatisfactory. The main result here is the characterization of the class of FVP that admit realization [1], [2] but characterization and construction of minimal ralizations are still poorly understood aspects of the theory [3]. The lack of a Realization Theory for FVP makes therefore impossible to follow Kalman's prescription [4]: Identification = Realization + Parametrization. To surmount the impasse and solve the identification problem we must follow the classical approach: select a parametric class of models P containing the observed process and estimate parameters on the basis of the data $\{y_1, y_2, \dots, y_T\}$. In section 2 we define Pas a class of probabilistic functions of Markov chains (the order considered as a parameter) satisfying conditions sufficient to guarantee identifiability. In section 3 we show how the order can be consistently estimated from the data. We present the results without proofs here. Further details and proofs can be found in [9]. Let $\{Y_t\}$ be a FVP. Denote by r its state space, where $\mathbf{r} := (1, 2, \dots, r)$ and by \mathbf{r}^* the set of finite sequences of elements of r. We assume that r is known a priori (at most this will cost the elimination of a null subset from the state space of the observed process) and denote by $\hat{\mathcal{P}}$ the set of pdf of stationary r-valued sequences.

2. The class of models

Our objective is to obtain a finite dimensional parametric representation for $\hat{\mathcal{P}}$. Finite dimensionality of the parameter space forces us to restrict attention to a subclass of \hat{P} .

Definition 2.1

$$\mathcal{P} := \{p_{\theta}; \ \theta \in \Theta\}$$

where -

$$\Theta := \left\{ \begin{array}{ll} \theta := & (k, A, B, \pi); k \in \mathbb{N}, \ A \in \mathbb{R}^{k \times k}, \\ & B \in \mathbb{R}^{k \times r}, \ \pi \in \mathbb{R}^k \\ & A, B, \pi \ \text{ stochastic and } \pi = \pi A \end{array} \right\}$$

and, defining $\forall arepsilon \in \mathbf{r}$

$$B_{\varepsilon} := \operatorname{diag} \{b_{1\varepsilon}, b_{2\varepsilon}, \dots, b_{k\varepsilon}\},\$$

the pdf

$$p_{\theta}: \mathbf{r}^* \to [0,1]$$
 is given by:

$$(2.1)\,\varepsilon_1\varepsilon_2\ldots\varepsilon_t\longmapsto p_\theta(\varepsilon_1\varepsilon_2\ldots\varepsilon_t):=\pi AB_{\varepsilon_1}AB_{\varepsilon_2}\ldots AB_{\varepsilon_t}e$$

where
$$e := (1, 1, \dots 1) \in I\!\!R^k$$
.

Remark 2.2

In terms of the underlying Markov chain $\{X_t\}$ we can make the following identifications:

 $k = \text{cardinality of the state space of } \{X_t\}.$

 $A = \text{transition matrix of } \{X_t\}$

 π = invariant measure of $\{X_t\}$

$$b_{i\epsilon} = P[Y_t = \epsilon | X_t = j]$$

For notational convenience we define $M(\varepsilon):=AB_{\varepsilon}.$ The interpretation of $M(\varepsilon)$ is:

(2.2)
$$M(\varepsilon)_{ij} = P[Y_{t+1} = \varepsilon, X_{t+1} = j | X_t = i].$$

As one would expect the class P just defined is not identifiable.

Assumption 2.3

A and B in definition (2.1) are such that:

- i) $a_{ij} > 0 \quad \forall i, j$
- ii) $b_{i\epsilon} > 0 \quad \forall j, \epsilon$
- A is invertible
- iv) B_{ε} has distinct diagonal elements for some $\varepsilon \in \mathbf{r}$ Definition 2.4

$$\Theta' := \{\theta \in \Theta; \text{ assumption 2.3 is satisfied}\}\$$

$$P' = \{p_{\theta}; \theta \in \Theta'\}$$

The main result of this section is theorem 2.9 which guarantees the identifiability of a subset of Θ' of full (Lebesgue) measure.

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Definition 2.5

Let $p_{\theta} \in \mathcal{P}$. Define the set of compound sequence matrices of p_{θ} relative to $\varepsilon \in \mathbf{r}$ as:

$$C(\theta, \varepsilon) := \begin{cases} M; \ m \in N, \ M \in \mathbb{R}^{m \times m}, \\ M_{ij} := p_{\theta}(s_{i}\varepsilon t_{j}) \ s_{i}, t_{j} \in \mathbf{r}^{*} \end{cases}.$$

Definition 2.6

Let $p_{\theta} \in \mathcal{P}$. Define the rank of p_{θ} relative to $\varepsilon \in \mathbf{r}$ as:

$$r_{ heta}(arepsilon) := \sup_{M \in C(heta, arepsilon)} \ (\mathit{rank}\ M)$$

Remark 2.7 If $p_{\theta} \in \mathcal{P}$ and $\theta = (k, A, B, \pi)$ then:

$$p_{\theta}(s_i \varepsilon t_j) = \pi M(s_i) M(\varepsilon) M(t_j) e$$

and, from Sylvester's inequality, $r_{\theta}(\varepsilon) \leq k$.

Definition 2.8

Let $p_{\theta} \in \mathcal{P}$ and $\theta = (k, A, B, \pi)$.

 p_{θ} is said to be regular if $r_{\theta}(\varepsilon) = k$, $\forall \varepsilon \in \mathbf{r}$.

 θ is regular if $\theta \in \Theta'$ and p_{θ} is regular.

Definitions 2.6 and 2.8 are analogous to notions defined in [5] for deterministic functions of Markov chains.

Theorem 2.9

Regular points of Θ' are identifiable modulo permutations of the states.

Remark 2.10

It is interesting to observe that under the same hypotheses theorem 2.9 has been proved in [6] in the case of fixed order of the model. For identification it is convenient to have compactness of the parameter space. Define, for $\delta > 0$:

$$\Theta'_{\delta} := \{ \theta \in \Theta'; \ a_{ij} \geq \delta, \ b_{j\epsilon} \geq \delta, \ \forall i, j, \epsilon \}$$

In what follows we will assume that the observed process $\{Y_t\}$ has a regular pdf $p_Y(\cdot) \in \mathcal{P}'$, i.e. there exists a regular point $\theta_0 = (n, A^0, B^0) \in \Theta'$ such that $p_Y(\cdot) = p_{\theta_0}(\cdot)$.

In practice the choice of δ is reduced to the choice of an upper bound K to the order n of the observed process. If $n \leq K$ and K is large enough taking $\delta := \frac{1}{K}$ we have $\theta_0 \in \Theta'_{\delta}$

3. Order Determination

We want to analyze the following question: given observations $(y_1, y_2, \dots y_T)$ with T arbitrarily large, can we determine the order n of $\{Y_t\}$? The answer is yes if $p_{\theta_0} \in \mathcal{P}'_{\delta}$ and is regular.

It is convenient to decompose Θ'_{δ} as follows: $\Theta'_{\delta} = U^K_{k=1} \Theta'_{\delta,k}$ where $\Theta'_{\delta,k}$ is the section of Θ'_{δ} along k. We identify $\theta = (k, A, B) \in \Theta'_{\delta,k}$ with its projection $\theta = (A, B)$.

Define the random variables on r*:

$$\begin{aligned} \forall \theta \in \Theta_{\delta}' & \quad f_1(\theta, Y(\cdot)) := P_{\theta}[Y_0] \\ \forall \theta \in \Theta_{\delta}' & \text{and} \quad T > 2 & \quad f_T(\theta, Y(\cdot)) := P_{\theta}[Y_0|Y_{-T-1}^{-1}] \end{aligned}$$

Lemma 3.1[7]

$$f(\theta, Y(\cdot)) := \lim_{T \to \infty} f_T(\theta, Y(\cdot))$$

exists $\forall Y(\cdot) \in \mathbf{r}^*$ and is continuous on $\Theta'_{\delta,k}$ $\forall k \in \mathbf{K}$.

Define:

$$\begin{split} h_k(\cdot) : \Theta_{\delta,k}' I\!\!R, &\rightarrow h_k(\theta) := E_{\theta_0}[\log f(\theta, Y(\cdot))] \\ h_{k,T}(\cdot) : \Theta_{\delta,k}' &\rightarrow I\!\!R, h_{k,T}(\theta) := \frac{1}{T} \log P[Y_1^T | \theta] \\ h_k : \mathbf{K} &\rightarrow I\!\!R, h_k := \sup_{\theta \in \Theta_{\delta,k}} h_k(\theta) \\ h_{k,T} : \mathbf{K} \times N &\rightarrow I\!\!R, h_{k,T} := \sup_{\theta \in \Theta_{\delta,k}} h_{k,T}(\theta) \end{split}$$

Lemma 3.2[7]

- i) $h_k(\cdot)$ is continuous on $\Theta'_{\delta,k}$ $\forall k \in \mathbf{K}$
- ii) $h_k(\cdot) = \lim_{T\to\infty} h_{k,T}(\cdot) \quad \forall \theta \in \Theta'_{\delta,k} \text{ and a.e. } P_{\theta_0}$

Theorem 3.3

- i) $h_k(\theta) \leq h_n(\theta_0) \quad \forall k \in \mathbf{K}, \ \theta \in \Theta'_{\theta,k}$
- ii) $h_k(\theta) = h_n(\theta_0)$ iff k = n

This is an important consequence of theorem 2.9. Compactness of Θ'_{δ} allows us to conclude with:

Corollary 3.4
$$h_n > h_k$$
 $\forall k \in \mathbf{K}, k \neq n$

From the abstract point of view this solves the problem of order determination. To compute the sequence h_k we must start from the data and therefore from the sequence of normalized maximum log-likelihoods $h_{k,T}$. In our situation the asymptotic behavior of $h_{k,T}$ is easy to analyze.

Lemma 3.5: $\delta^T \leq P\left[Y_1^T | \theta\right] \leq (1 - \delta)^T \quad \forall \theta \in \Theta_{\delta}'$ and therefore:

Lemma 3.6: $h_{k,\infty} := \lim_{T\to\infty} h_{k,T}$ exists $\forall k \in \mathbf{K}$

To be able to apply corollary 3.4 we need $h_{k,\infty} = h_k$ and this is true if:

(3.1)
$$\lim_{T \to \infty} \sup_{\theta \in \Theta_{k,k}^T} h_{k,T}(\theta) = \sup_{\theta \in \Theta_{k,k}^T} \lim_{T \to \infty} h_{k,T}(\theta)$$

The interchange of limit operations (3.1) is allowed under uniform convergence (for $T \to \infty$) of $h_{k,T}(\theta)$ on $\Theta'_{\delta,k}$. Using the previous results one can verify that actually the sequence $h_{k,T}(\theta)$ satisfies Dini's criterion for uniform convergence.

Observe that (3.1) is verified for every trajectory $Y_t(\omega)$ and from Lemma 3.2 we conclude that $h_{k,T} \longrightarrow h_k$ for almost (P_{θ_o}) every trajectory.

Theorem 3.7

- i) $arg \ max_k\{h_{k,T}\}$ is a strongly consistent estimator of n.
- ii) For (almost) every trajectory there exists T_o such that: $n = arg \max_{k} \{h_{k,T}\}$ $(\forall T \geq T_o)$

Remark The practical computation of $h_{k,T}$ (and of the maximizing $\theta = (A, B)$) can be done using the Baum-Eagon algorithm; see [8] for a description and the analysis of its numerical properties. p

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