Structural Break Estimation for Non-Stationary Time Series Signals Richard A. Davis, Thomas C.M. Lee & Gabriel A. Rodriguez-Yam, Colorado State University

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1. Introduction

We consider the problem of modeling a non-stationary time series by segmenting the series into blocks of different autoregressive (AR) processes. The number of break points, denoted by *m*, as well as their location, and the order of the respective AR models are assumed to be unknown. We propose an automatic procedure for obtaining such an optimal partition called Auto-PARM for Automatic *P*iecewise Auto*R*egressive *M*odeling.

1.1 Piecewise Autoregressive processes

Setup: there exist *m* and $\tau_0 = 1 < \tau_1 < \ldots < \tau_m < \tau_{m+1} = n+1$ (n = sample size) such that $Y_t = \gamma_i + \phi_{i1}Y_{t-1} + \dots + \phi_{ip_i}Y_{t-p_i} + \sigma_i\varepsilon_t, \quad \text{if } \tau_{i-1} \le t < \tau_i,$

where $\{\varepsilon_t\}$ is IID(0,1).

Goal: Estimate

m = number of segments

 $\tau_i = \text{location of } j^{\text{th}} \text{ break point}$

 $\gamma_i = \text{level in } j^{\text{th}} \text{ epoch}$

 p_i = order of AR process in j^{th} epoch

 $(\phi_{i1},...,\phi_{ip_i}) = AR$ coefficients in *j*th epoch

 σ_i = scale in *j*th epoch

1.2 Motivation for using piecewise AR models

Piecewise AR is a special case of a *piecewise stationary process* (see

Adak 1998), $\widetilde{Y}_{t,n} = \sum_{i=1}^{m} Y_t^j I_{[\tau_{i-1},\tau_j)}(t/n),$

where $\{Y_t^j\}, j = 1, ..., m$ is a sequence of stationary processes. It is argued in Ombao et al. (2001) that if $\{Y_{t_n}\}$ is a locally stationary process (in the sense of Dahlhaus), then there exist a piecewise stationary process $\{\widetilde{Y}_{t,n}\}$ and a sequence m_n

 $m_n \rightarrow \infty$ and $m_n/n \rightarrow 0$, as $n \rightarrow \infty$,

that approximates $\{Y_{tn}\}$ (in average mean square).

Roughly speaking: $\{Y_{t_n}\}$ is a locally stationary process if it has a time-varying spectrum that is approximately $|A(t/n,w)|^2$, where A(u,w) is a continuous function in u.

2.1 Model selection using Minimum Description Length (MDL) The idea behind MDL is to choose the model which maximizes the compression of the data or, equivalently, select the model that minimizes the code length of the data (i.e., amount of memory required to encode the data).

2.2 The MDL applied to piecewise AR models

M = class of piecewise AR models for $y = (y_1, \ldots, y_n)$

$L_{\mathbf{F}}(\mathbf{y}) = \text{code length of } \mathbf{y} \text{ relative to } \mathbf{F} \in \mathbf{M}$

Best fitting MDL model is minimizer of

$$DL(m,(\tau_1, p_1), \dots, (\tau_m, p_m)) = \log_2 m + m \log_2 n + \sum_{j=1}^{j} \log_2 p_j$$

$$+\sum_{j=1}^{m} \frac{p_{j}+2}{2} \log_{2} n_{j} + \sum_{j=1}^{m} \frac{n_{j}}{2} \log_{2}(2\pi\hat{\sigma}_{j}^{2}) + \frac{n_{j}}{2}$$

where n_i is the length of the *j*-th segment and $\hat{\sigma}_i^2$ is the Yule-Walker estimate of σ^2 in the *j*-th segment.

2.3 Consistency

Μ

Assume there exist true values *m* and $0 < \lambda_1 < \ldots < \lambda_m < 1$ with $\tau_i = [\lambda_i n], i = 1, 2, ..., m.$

Theorem. For the piecewise AR model, if the number of breakpoints *m* is known, then

 $\hat{\lambda}_i \rightarrow \lambda_i$ a.s. j = 1, 2, ..., m.

3.1 Basics of the Genetic Algorithm (GA)

The GA is an optimization algorithms that mimics natural evolution.

- Start with an initial set of chromosomes, or population, of possible solutions to the optimization problem.
- · Parent chromosomes are randomly selected (proportional to the rank of their objective function values), and produce offspring using crossover or mutation operations.
- After a sufficient number of offspring are produced to form a second generation, the process then restarts to produce a third generation.
- Based on Darwin's *theory of natural selection*, the process should produce future generations that give a *smaller (or larger)* objective function.

3.2 Implementation of GA

A chromosome consists of n genes, each taking the value of -1 (no break) or *p* (order of AR process). Use natural selection to find a *near* optimal solution. An element $F \in M$ is mapped with a chromosome c by

$$(m,(\tau_1,p_1)...,(\tau_m,p_m)) \longleftrightarrow c = (\delta_1,...,\delta_n),$$

For example,

corresponds to

AR(2), t=1:5; AR(0), t=6:10; AR(0), t=11:14; AR(3), t=15:20

4.1 Piecewise stationary with dyadic structure $(9Y_{t-1} + \varepsilon_t, if 1 \le t < 513)$ $.9Y_{t-1} + \varepsilon_t$ $Y_t = \{ 1.69Y_{t-1} - .81Y_{t-2} + \varepsilon_t, \text{ if } 513 \le t < 769, \}$

 $1.32Y_{t-1} - .81Y_{t-2} + \varepsilon_t$, if $769 \le t \le 1024$, where $\{\varepsilon_t\} \sim \text{IID N}(0,1)$.

Sample realization.



Simulation Results based on 200 reps

# of	Auto-SLEX		Auto-PARM		
segments	%	Change Points	%	mean	std
2	0	1/2	0		
3	60.0	1/4 , 3/4	82.0	.500 .749	.006 .006
4	34.0	1/4, 2/4, 3/4	17.5	.476 .616 .761	.080 .110 .037
5	5.0	2/8, 4/8, 5/8, 6/8, 7/8	0		
≥6	1.0		0.5		

4.2 Slowly varying AR(2) model

Fitted

$Y_t = a_t Y_{t-1} - .81 Y_{t-2} + \varepsilon_t$ if $1 \le t \le 1024$

where $a_t = .8[1 - 0.5\cos(\pi t / 1024)]$, and $\{\varepsilon_t\} \sim \text{IID N}(0,1)$.



GA results: 3 pieces, breaks at τ_1 =293, τ_2 =615. Total run time 27.45 secs

model:		ϕ_1	ϕ_2	σ^2
	1- 292:	.365	-0.753	1.149
	293- 614:	.821	-0.790	1.176
	615-1024:	1.084	-0.760	0.960



Left: True model log-spectrogram; Center: Auto-PARM logspectrogram; Right: Average of Auto-PARM log-spectrogram based on 200 reps.

5. Speech signal segmentation Speech signal: GREASY, n = 5762 observations Auto-PARM results: m = 15 break points, run time = 18.02s



Bottom: Spectogram based on Auto-PARM model

6. Conclusions

- Introduced Auto-PARM (an automatic procedure for segmenting a time series signal into piecewise AR models).
- Model selection based on MDL (minimum description length) principle.
- A *genetic algorithm* was used to find a near optimal solution to the model selection problem based on MDL.
- Auto-PARM works well for both detecting segments and for estimating time-varying spectra.

7. References

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