Time series machine learning with shapelets

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A Hands-on Introduction to Time Series Classification and Regression

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Notation one-stop shop

Before diving in, let us define some notations :

- a **time series** of length m and of dimension d as a vector $X = [\mathbf{x}_1, \dots, \mathbf{x}_m]$ with $\mathbf{x}_i \in \mathbb{R}^d$.
- a time series dataset $\mathcal{X} = [X_1, \dots, X_n]$, this is equivalent to an array 3D X of shape : (n_samples, n_channels, n_timepoints)
- an **annotation vector** $Y = [y_1, \dots, y_n]$. For example, in a regression context, $y_i \in \mathbb{R}$ would represent the label of X_i .

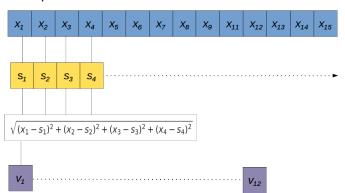
For simplicity, we'll consider time series with a fixed frequency. Additionally, we use:

- a blue color when talking about time series
- a yellow color when talking about shapelets
- a pink color when talking about annotation vectors
- a purple color when talking about distance vectors

What are shapelets?

A shapelet [4] is a subsequence $S = \{s_1, \dots, s_l\} \in \mathbb{R}^{l,d}$ representing a pattern of interest. The similarity between S and X is quantified using a distance vector V. Example with d = 1 (univariate):

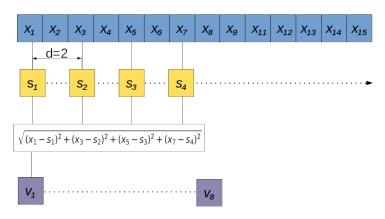
- $dist(X, S) = [v_1, \dots, v_{m-(l-1)}]$
- with $v_i = \sqrt{\sum_{j=1}^{I} (s_j x_{i+(j-1)})^2}$. (or any distance function)



What are (dilated) shapelets?

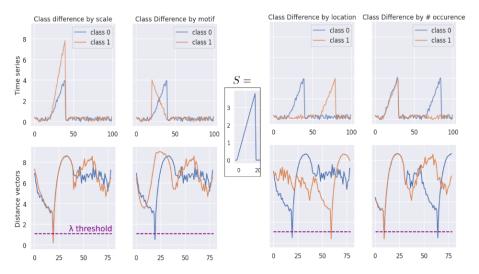
In [1], a parameter d is added to S. The intention is to :

- Allow matching non-contiguous, cyclic patterns (e.g. heartbeats)
- Increase receptive field.



What do we do with the distance vector?

We want to identify discriminative characteristics between S and X:



What do we do with the distance vector?

The following features are extracted to target these discriminative characteristics :

- min d(X, S): the feature most shapelet-based methods use. It allows discriminating by presence or absence of a pattern.
- **argmin** d(X, S): first used in [2] for shapelets, this allows to discriminate by **pattern location**.
- Shapelet occurrence [1]: $\sum_i I(v_i < \lambda)$, with I the identity function and a threshold λ , which is a shapelet parameter. It allows discriminating by number of λ -close occurrence of the pattern.

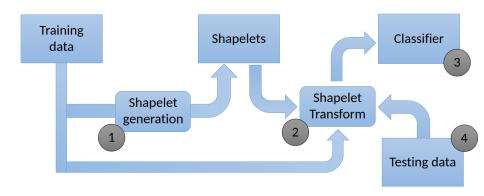
Standardizing S and every subsequence $[x_i, \ldots, x_{i+(l-1)}]$ independently makes v_i , and thus d(X, S) (and the extracted features), **scale invariant**. Not standardizing makes them **scale sensitive**.

Shapelet-based classification and regression

All Shapelet classifiers or regressors use the same methodology, given

$$\mathcal{X} = \{X_1, \dots, X_n\} \text{ and } Y = \{y_1, \dots, y_n\}$$
:

- **4** Generate $S = \{S_1, \dots, S_k\}$ the set of Shapelets
- ② Extract a matrix of features M, from \mathcal{X} using \mathcal{S}
- Learn a classifier based on M and Y



How do we generate Shapelets?

We can distinguish three approaches:

- Exhaustive search: extract all possible shapelet candidates and evaluate their quality to extract the best ones (information gain, F-statistic, . . .)
- Heuristic-based search: random or semi-random extraction from the training data, statistical or rule-based methods for reducing the number of shapelet candidates.
- **Shapelet generation**: Use of optimization methods (gradient descent, evolutionary algorithm) to generate shapelet values instead of extracting them from the input.

What is the Shapelet Transform?

Consider $\mathcal{X} = \{X_1, \dots, X_n\}$ and $\mathcal{S} = \{S_1, \dots, S_k\}$. The Shapelet Transform [3] algorithm is the same for all shapelet methods:

- Create a matrix M of size (n, k)
- ② $\forall i \in [1, n], \forall j \in [1, k], \text{ and extract features from } d(X_i, S_j)$

For example, if extracting the minimum value :

$$M = \begin{bmatrix} \min d(X_1, S_1) & \dots & \min d(X_n, S_1) \\ \vdots & \ddots & \vdots \\ \min d(X_1, S_k) & \dots & \min d(X_n, S_k) \end{bmatrix}$$

A column M_i of M describes X_i by the features extracted using S. The argmin can also be extracted, giving a matrix M of size (n, 2k).

References I



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