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Data... over time.

# **Time series**

A univariate series  $s \in \mathcal{X}^n$  is a sequence  $s = [s_1, \ldots, s_n]$  of n values in a domain  $\mathcal{X}$ . A series is defined by:

- *Type*: discrete, e.g., nucleotide bases, or continuous, e.g., stock values in a financial market
- *Sampling rate*: How often values are sampled, e.g., daily
- *Amplitude*: Values sampled, e.g., value of the stock on a particular day



A time series.



### **Time series**



A univariate series  $s \in \mathcal{X}^K$  is a sequence  $s = [s_1, \ldots, s_K]$  of K values in a domain  $\mathcal{X}$  . A series is defined by:

- *Seasonality*: Series repeat (or almost repeat) over time, e.g., temperature
- *Period*: How much does it take for the series to repeat itself, e.g., length of a calendar year



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#### **Time series**

A multivariate time series *s* generalizes time series to multiple variables. Each instance is comprised of multiple time series, each representing a different feature.



#### **Time series statistics**

- Mean. Expected value  $\mathbb{E}[s]$  of s
- Variance. Variance of s
- Trend: slope  $\Delta$  of a linear model modeling s



A time series, and its trend (in beige).



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#### **Time series statistics**

- Interquantile ranges.  $q_a(s) q_b(s)$
- Skewness: is the distribution symmetric?  $\mathbb{E}[(rac{s-\mu}{\sigma})^3]$
- Kurtosis: what is the probability mass on the tails?  $\mathbb{E}[(rac{s-\mu}{\sigma})^4]$



Empirical distribution of a time series components, values color-coded by density.

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# **Time series: local behaviors**



Time series can be affected by:

- Seasonality: the series has some repeated periodic behavior, e.g., temperatures fall every winter
- *Trends*: the series tends to have a monotonic behavior, with values increasing/decreasing

We need to analyze time series on a local scale.

# **Time series statistics: rolling statistics**



Rolling indicates the act of extracting a series of consecutive subsequences of given dimensionalities  $w^1, \ldots, w^k$ . Each subseries gives a different view on s, and is thus named *window*. Given a series of windows, we can now *locally* describe a time series!

- Rolling mean
- Rolling variance
- ...

# Time series statistics: sliding statistics



Unlike rolling statistics, which compute a statistic over a subseries of a given series, *sliding* statistics slide a series t over a series s, computing a statistic between the two. The act of sliding a window through a function is called *convolution*.

#### Auto-covariance

How much does a component of a time series correlate with previous and future components? Slides a series over itself, computing covariance between the two components:

$$K_s = cov(s_{t:}, s_{t+\Delta:}) \sum_{i=1}^n s_i \cdot s_{i+\Delta}.$$

High autocovariances may indicate seasonality in the series.

# **Time series: sliding statistics**

#### **Cross-correlation**

Shifted pointwise correlation of the two series *a*, *b*, measured as a *sliding* inner product:

For univariate time series, the inner product is simply a multiplication.



$$CC_\Delta(a,b) = \sum_{i=1}^n a_i \cdot b_{i+\Delta}.$$

#### From analyzing to transforming

# **Representing by segmenting**



Given a set of time series  $S = \{s^1, \ldots, s^n\}$ , let  $S^{\subset}$  be a set of k subseries  $s_{\subset}^1, \ldots, s_{\subset}^k$ . We define an alphabet  $\Sigma$  of symbols, each symbol assigned to a subseries. Then, we can segment each series into a sequence of subseries, and represent each time series as a series of symbols in  $\Sigma^*$ .

$$s^i o \underbrace{[s^{i,1} \mid \cdots \mid s^{i,k_i}]}_{\in \Sigma^*}$$

We can tackle two problems independently:

- Segmentation: how to split a series into subseries?
- Transformation: what symbols do we use to define each segment?

### **Fixed window-based segmentation**

Each series  $s^i$  has an often intractable number of possible subseries, which makes exhaustive search unfeasible. Instead, we choose an arbitrary window w, and segment each series into a set of  $\frac{n}{w}$ subseries.

A fixed window of size w = 2, segmenting a time series in three non-overallping adjacent windows (color-coded).





### Learned segmentation

We *learn* segmentations which minimizing an approximation error.

Piecewise Linear Approximation (PLA) defines a segmentation minimizing segment-local linear models of the data.

- Given number of segments: segmentation which distributes approximation error as evenly as possible among segments
- Given error bound: segmentation which generates the *minimum* number of segments within given error bound  $\varepsilon$







# Learned equidistributional windows



We *learn* segmentations by their likelihood: given a desired number of segments, segments are learned to maximize their probability  $p(s_{\subset}^{i})$ , which results in segments as equiprobable as possible, ideally following a uniform distribution. By maximizing probability, we maximize entropy, and thus the diversity of the subseries.



# **Continuous transformations**

	Description	Туре	$f^{\Sigma}$	$\Sigma$
Piecewise Average (PA)	Average value of the segment	Continuous	$\mu_{s^i_\sub}$	$\mathbb{R}$
Piecewise Linear (PL)	Slope $ abla$ of the segment	Continuous	$ abla_{s^i_\sub}$	$\mathbb{R},\mathbb{R}^2$

 $\mathbb{R}^2$  given by storing both slope and intercept of the linear model.

# **Two-tier segmentations**



Window segmentations create series of symbols themselves! We can learn representations through a two-tier algorithm:

- continuous transformation, yielding segments  $S^{\sub}$  with symbols in  $\Sigma^{\sub}$
- transformation partitioning, mapping symbols in  $\Sigma^{\sub}$  to discrete symbols in  $\Sigma$

Symbols can be either categorical, or ordinal.

# Symbolic aggregate approximation (SAX)



Symbolic aggregate approximation (SAX) implements a two-tier transformation:

- fixed-window segmentation followed by piecewise average transformation in subsymbols  $(\Sigma^{\subset})^*$
- aggregation of subsymbols into equiprobable symbols in  $\Sigma$ : we partition the subsymbol distribution into equiprobable buckets, each defined by a symbol in  $\Sigma$

SAX symbols are discrete *and* ordinal!

# Symbolic aggregate approximation (SAX)



The three steps of SAX: subsymbol induction with fixed-window average segmentation, partitioning of the subsymbol distribution, and transformation.

# **Signal representations**



Segmentations can be tricky to handle, and simply offer a representation in a domain quite different from the original. Signal representations, on the other hand, aim to define a series in terms of other series. To stick with the signal processing literature, where they are most prevalent, we'll refer to series as *signals*.

- What other signals can we leverage?
- How do we combine them to represent the original series?



### **Fourier analysis**

Fourier analysis tackles **periodic** (also known as stationary) series, i.e., series which periodically repeat themselves.

- What other signals can we leverage? Sine and cosine signals at different frequencies
- How do we combine them to represent the original signal? Linear combination



# **Fourier analysis**



As a linear transformation, we need to learn a set of coefficients  $\alpha$  which map the basis to the signal s. In Fourier analysis, we constrain  $\alpha \in \mathbb{R}^+$ :

- basis signals which do not contribute to the signal have a 0 coefficient
- basis signals which do contribute do so with a positive coefficient  $lpha_j>0$

To compute  $\alpha$  we can thus use *inner products*, e.g., dot product, which are guaranteed to satisfy both.

#### The basis: sinusoids



First, we define the basis of the transformation. We use sinusoids, i.e.,  $\sin/\cos$  signals defined by a phasor  $\psi$ .



A phasor  $\psi$ , and the  $\sin$  series generated.

# The basis: sinusoids



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A phasor  $\psi$ , and the  $\cos$  series generated.

# Sinusoids define amplitude



For a phasor  $\psi$ , we can define the period  $t_0$  as the time required for the phasor to complete one rotation over the unit circle, and the frequency  $f_0 = t_0^{-1}$  as the rotations per unit of time. Then, at time t, the phasor defines a series component with amplitude





# Sinusoids and the complex unit circle

By Euler, we can map complex numbers to the complex unit circle.

$$e^{-i\theta} = \cos(\theta) + i\sin(\theta).$$



The complex unit circle and an imaginary number  $z = e^{-i\theta}$ . As in linear algebra, we can define a vector  $(z_{\mathbb{R}}, z_{\mathbb{C}})$  through the standard basis, in this case (1, 0), (0, i).



# Fourier, and the frequency domain

Phasors of different frequencies are the building block of the Fourier representation of the signal: we will map signals to a *frequency domain* populated by sinusoids of different frequencies.



Sinusoid signals of different frequencies define a basis in the frequency domain. An interactive visualization can be found **here**.

# Fourier, and the coefficients



Having defined a basis, we now need to learn the coefficients  $\alpha$ . Coefficients ought to measure the presence of a frequency, and its scaling. Thus,  $\alpha \in \mathbb{R}^{+n}$ , and  $\alpha_j = 0$  for frequencies not present in s. Inner products, e.g., dot products, satisfy both conditions: the coefficients will be given by the inner product of basis signals and the signal s:

$$lpha_j = s \cdot \psi_j.$$

# A small caveat: orthogonality... again?



Signals may be out orthogonally out of phase, thus inducing null products, leading to misses on the basis. To tackle this, we use the orthogonal components of the basis: the sin component of the phasor!

$$s_t = \sum_{
ho=1}^n lpha_
ho(\cos( heta_{
ho,t}) + i\sin( heta_{
ho,t}))$$

# **Discrete-Time Fourier Transform (DTFT)**



Signals may be out orthogonally out of phase, thus inducing null products, leading to misses on the basis. To tackle this, we use the orthogonal components of the basis: the sin component of the phasor!

$$s_t = \sum_{
ho=1}^n lpha_
ho(\cos( heta_{
ho,t}) + i\sin( heta_{
ho,t}))$$

Finally, we leverage Euler, and have

$$s_j = \sum_{
ho=1}^n lpha_
ho e^{ heta_{
ho,t}} = \sum_{
ho=1}^n lpha_
ho e^{-2\pi f_
ho t}$$

Remember: by definition, inner product satisfy  $a \cdot b = 0$  for a, b orthogonal!

# **Discrete-Time Fourier Transform**



- Quick:  $\mathcal{O}(n \log n)$
- Decomposition in separate and different signals

- Decomposition defined exclusively for sinusoidal series
- Decomposition of *periodic* series
- Decomposition exclusively in terms of frequency, not time

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#### Wavelets

Wavelet tackle several weaknesses of Fourier Transforms.

#### Fourier Transform: $\psi_\kappa$

- Decomposition defined exclusively for sinusoidal series
- Decomposition of *periodic* series
- Decomposition exclusively in terms of frequency, not time

Wavelets:  $\psi_{\kappa, au}$ 

VS

- Flexible decomposition defined by *mother wavelets*
- Decomposition of arbitrary series
- Wavelets parameterized in frequency **and** time



#### Wavelets



Wavelets (little waves) aim to replace sinusoidal phasors, and are both more general, and flexible enough for domain-specific application. A wavelet  $\psi_{\kappa,\tau} : \mathbb{R} \to \mathbb{R}$  is a function s.t.

- $\int_{-\infty}^{+\infty}\psi_{\kappa, au}(t)\,dt=0$  (zero mean)
- $\int_{-\infty}^{+\infty}\psi_{\kappa, au}(t)^2\,dt
  eq\infty$  (finite energy) or (compact support)

Finite energy makes it so a wavelet, unlike a sinusoidal function, is bounded: thus, by construction, wavelets **can be localized in time**!

$$-\sqrt{-}$$

Three wavelets: Mexican Hat, Morlet, Meyer.

# Surfing the series with wavelets



Localization is innate in wavelets by their own definition. A *mother wavelet*  $\psi_{K,T}$ defines a family of *daughter* wavelets defined by

- A frequency κ: shrink or stretch the daughter wavelet
- A shift  $\tau$ : pushes or pulls the daughter wavelet across time.

We define a daughter wavelet  $\psi_{\kappa, au}$  as

$$\psi_{\kappa, au}(t) = rac{1}{\sqrt{\kappa}}\psi(rac{t- au}{\kappa}).$$



Stretched wavelets: Mexican Hat wavelet (left column), and Meyer wavelet (right column).

#### **Wavelets Transforms**



Moving from sinuoidal phasors to wavelets is seamless in the formulation



Wavelets are convoluted across the series, producing a list of coefficients.  $\kappa$  and  $\tau$  are dyadic, i.e., they are taken as powers of 2:  $\tau = 2^{-i}$ ,  $\tau = k2^{-i}$ .

#### From representations to motifs

# **Time series: motifs**



Motifs tie strongly with subseries extraction for discretization: if representation algorithms extracted to *represent*, and thus describe, motif extraction algorithms instead extract subseries to *represent*, and thus *discriminate*. In other words, a *motif* is a subseries characteristic of a set of series. We can search motifs following the two views:

- descriptive: a motif is a reoccurring subseries in the set  ${\cal S}$
- discriminative: a motif is a reoccurring subseries in the set S... and not reoccurring in another set  $S^{\neq}$ . Also called a *shapelet*

# **Searching for Motifs**



The first step is already solved! We know how to extract a set of subseries  $S_{\subset}$  from series representation. We need to quantify the descriptive and discriminative power of candidate motifs.

**Descriptive power**. Distance of  $s_{\subset}^{i}$  with respect to all possible subseries of  $s^{j}$  yields distances  $D_{i,S} = \{d_{i,j}^{1}, \ldots, d_{i,j}^{k}\}$ . Descriptive power given by lower distances, e.g., min D**Discriminative power**. Comparison of descriptive power with respect to S and  $S^{\neq}$ , e.g., min  $D_{i,S}$ 

 $\min D_{i,S^{
eq}}$ 

# From motifs to shapelets



Since motifs are supposed to discriminate, why not directly measure their discriminative power? Partition subseries in  $S \cup S^{\neq}$  according to their distance from a given candidate  $s_{\subset}^{i}$  and a threshold  $\beta$ , obtaining two sets  $S_{\beta}^{\leftarrow}$ ,  $S_{\beta}^{\rightarrow}$ . Then, compute a discrimination measure, e.g., entropy, information gain, etc., on the two sets  $S_{\beta}^{\leftarrow}$ ,  $S_{\beta}^{\rightarrow}$ . The larger the measure, the higher the discriminative power!



A shapelet (in black), a threshold  $\beta$ , and the two sets  $S_{\beta}^{\leftarrow}$ ,  $S_{\beta}^{\rightarrow}$  (left and right, in blue and red, respectively).

# **Time series: alignment**

An alignment  $A = [(i, j)]^{\max n, m}$  of two time series l, u with components  $[l_1, \ldots, l_n], [u_1, \ldots, u_m]$  is an assignment of each component  $l_i$  of l to a component  $u_j$  of u. An alignment A induces an alignment  $cost C_A$  quantifying how unaligned the two series are.





Two possible alignments (in beige) of an *upper* series  $s^1$  (in red), and a *lower* series  $s^2$  (in blue).

We are going to assume equal time sampling for both series.

# Alignment: local cost



Alignment costs  $C_{\cdot}$  are based on two separate costs:

- local (component-wise, or point-wise) alignment cost  $c_{i,j}^A$ : defines the cost of aligning  $l_i$  with  $l_j$ . How much do I pay for this alignment?
- match alignment cost  $c_{i,j}^{\Sigma}$ : defines the cost of foregoing matching i with j, in favor of a lower-cost  $i' \neq i$ . How much would I pay for another alignment?

Alignment algorithms look to minimize a combined cost of the two.

# Alignment: straight match





A straight match alignment.

# Alignment: straight match



Trivial alignment: each point  $l_i$  is assigned to  $l_i$ . The alignment induces a pairwise alignment cost of  $c_{i,i}^a = || l_i - u_i ||^p$ . The alignment cost  $C_A$  is given by the sum of the pairwise alignments:

$$C_A = \sum_{i,i} c^a_{i,i} = \sum_i ~||~ l_i - u_i ~||^p.$$

Since i is always aligned with j, it follows that  $c_{i,j}^{\Sigma} = 0$ . This produces an alignment  $A = [(l_i, u_i)]^K$ , only applicable for m = n.

Applicable to series with the same sampling rate!



# Alignment: straight match

Two possible alignments of an *upper* series  $s^1$ , and a *lower* series  $s^2$ .

- Simple definition
- Minimizes norm cost
- Quick:  $\mathcal{O}(n)$

- Only applicable to equal-length series
- Assumes alignment

# Alignment: Shifted match



Not all time series are already aligned! Given a component  $l_i$ , we need to look for an aligning  $u_j$ , under a shift assumption.

#### Shifted alignment.

Assuming some subseries of either l, u are shifted, under a shifted alignment:

- Warp, Replication. Each component  $l_i$  can be assigned to any component  $u_j$
- **Planarity.** Assignments are monotonic: a successive point cannot be assigned backwards:  $orall i, j. \, (l_i, u_j) \in A \implies (l_{i+1}, u_{j-1}) 
  ot\in A$

# **Alignment: Shifted match**



If properly performed, a shifted match minimizes norm cost, and emulates a straight alignment... with additional aligned components.



A shifted match alignment as a straight match alignment.



# **Shifting as Warping**

#### **Warp, Replication.** Each component $l_i$ can be assigned to any component $u_j$ .

By *warping*, we replicate a component, warping it also further in the series. This allows us to replicate components, and emulate a straight match.

- *no warp*: the assignment  $(l_i, u_j)$  is attempted as is
- *lower warp*: the assignment  $(l_i, u_j)$  is attempted on  $(u_{i+1}, u_j)$ , replicating the component on the lower series u
- upper warp: the assignment  $(l_i, u_j)$  is attempted on  $(l_i, u_{j+1})$ , replicating the component on the upper series v

 $\Sigma$  stands for sum, unlike the alphabet used in SAX representation.

# $C^{\Sigma} \begin{bmatrix} 0.456 & 0.010 & 0.120 & 0.010 \\ 0.524 & 0.350 & 0.536 & 0.885 \\ 0.458 & 0.214 & 0.511 & 0.194 \end{bmatrix}$

 $0.618^{-1}$ 

0.429

A DTW alignment (top), and a cumulative alignment cost matrix  $C^{\Sigma}$  (bottom).

0.613

0.495

The (i, j) component holds the cumulative cost of aligning the  $l_i, u_j$ .

DTW divides the alignment problem in two steps:

- 1. compute a cumulative minimal alignment cost matrix  $C^{\Sigma}$ , defining the minimal alignment cost of every pair of components  $l_i, u_j$
- 2. search an alignment A on  $C^{\Sigma}$ , minimizing the alignment cost

# **Dynamic Time Warping (DTW)**





# Warping in Dynamic Time Warping



- $C^{\Sigma}$  computes a *minimal* and *cumulative*:  $c_{i,j}^{\Sigma}$  may be eq 0!
  - base case: the alignment  $(l_1, u_1)$  has minimal cost. This is trivially true, since there are no accumulated costs, i.e.,  $C_{1,1}^{\Sigma} = \mid\mid l_1 u_1 \mid\mid^p$
  - inductive case: i, j fall within the three warping categories, thus it must be one of three cases
    - $\circ~$  (no warp) i,j: the accumulated cost is  $c_{i,j}^{\Sigma}=C_{i-1,j-1}^{\Sigma}$
    - $\circ~$  (lower warp) i has warped: the accumulated cost is  $c_{i,j}^{\Sigma} = C_{i-1,j}^{\Sigma}$
    - $\circ~$  (upper warp) j has warped: the accumulated cost is  $c_{i,j}^{\Sigma}=C_{i,j-1}^{\Sigma}$

# **Computing warp cost**

Entries in 
$$C^{\Sigma}$$
 are computed as  $C_{i,j}^{\Sigma} = c_{i,j}^a + c_{i,j}^{\Sigma}$ 

alignment cost

$$= \underbrace{||l_i - u_j||^p}_{alignment\ cost} + \min\{\underbrace{C_{i-1,j-1}^{\Sigma}}_{no\ warp}, \underbrace{C_{i,j-1}^{\Sigma}}_{lower\ warp}, \underbrace{C_{i-1,j}^{\Sigma}}_{upper\ warp}\}$$

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	0.495	0.62	13	0.429	0	.618]
$C^{\Sigma}$	0.524	0.35	50	0.536	0	.885
	0.458	0.2	14	0.511	0	.194
		Γ-	$\leftarrow$	$\leftarrow$	$\leftarrow$	
	$C^{\leftarrow} =$	$\uparrow$	K	$\uparrow$	$\leftarrow$	
		L ↑	$\uparrow$	$\leftarrow$	$\leftarrow$ _	

Directions of minimal accumulated cost  $C^{\leftarrow}$  over  $C^{\Sigma}$ : entries indicate which alignment choice has produced the minimal cost.

# Searching for the lowest cost alignment



As a matrix of minimal accumulated alignment cost, we know that for each component i, j in  $C^{\Sigma}$ , we have, by construction, the minimal cost to align up to i, j. Thus, we can simply start from the last alignment, and follow  $C^{\leftarrow}$  backwards for the warps of minimal cost!

$$C^{\leftarrow} = egin{bmatrix} - & \leftarrow & \leftarrow \ \uparrow & \swarrow & \uparrow & \leftarrow \ \uparrow & \uparrow & \leftarrow & \leftarrow \end{bmatrix}, A = [(1,1),(2,2),(3,2),(3,3),(3,4)]$$

Directions of minimal accumulated cost  $C^{\leftarrow}$  over  $C^{\Sigma}$ : the red path from the last entry indicates induces the alignment of minimal cost. The alignment A is given by the indices of the path.

# **Dynamic Time Warping**





A Dynamic Time Warping alignment.

• Minimizes norm cost on shifted series

#### • High cost of $\mathcal{O}(mn)$

# **Constraining alignments**



Within  $C^{\Sigma}$ , we have several possible movements:

- alongside a row (*i* is constant): indicate a sequence of upper warps, as we are aligning one component of *l* with consecutive components of *u*
- alongside a column (j is constant): indicate a sequence of lower warps, as we are aligning one component of u with consecutive components of l
- diagonally (neither *i* nor *j* is constant): indicate a no-warp alignment

#### Sakoe-Chiba search



The optimal search we have highlighted can move almost arbitrarily over  $C^{\Sigma}$ , inducing arbitrarily large row- and column-segments, and thus arbitrarily large warps. Sakoe-Chiba search instead constraints the alignment A to be such that  $\forall i, j \in A. \mid i - j \mid < \gamma$  by introducing a search radius  $\gamma$ , and binding the maximum warp size of either series.



# Itakura parallelogram



The Itakura constraint instead binds the slope of segments, effectively binding consecutive warps.

For any two  $(i,j), (k,l) \in A.$   $k \geq i, l \geq j.$   $rac{l-j}{k-i} < lpha.$ 



#### From alignment to similarity

### **Time series: similarity**





#### References



Торіс	Reference
Fourier Transform	Digital Signals Theory e-book, Chapters 1, 2, 4, 5, 9
Wavelet Transform	Data-driven science and engineering. By S. L. Brunton, J. N. Kutz, 2nd edition. 2.5
Shapelets	Time series shapelets: a new primitive for data mining
Dynamic Time Warping	An introduction to Dynamic Time Warping

Credits for some images and animations to Romain Tavenard.