Machine Learning for Time Series

Lecture 5: Change-point and Anomaly Detection

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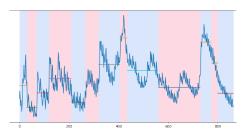
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Discovering events in time series

- ► Time series are widely used for monitoring: finance, industry, healthcare, meteorology...
- ▶ When recorded for hours, days, weeks... data is likely to be redundant
- ► Two fundamental questions:
 - ▶ Were there significant changes in my data across time?
 - ▶ Was there something new or unusual in my data?
- ▶ Two ill-posed problems: what is a significant change? What is new or unusual?

Problem 1: Change-Point Detection

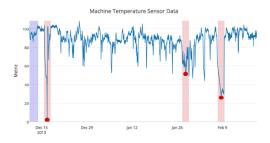


Change-Point Detection

Given a time series x, retrieve the times (t_1, \ldots, t_K) where a significant change occurs

- ► Necessitates to estimate both the change-points but also the number of changes *K*
- Highly depends on the meaning given to change

Problem 2: Anomaly Detection



Anomaly Detection

Given a time series x, retrieve the set of samples $\mathcal T$ that corresponds to unusual phenomenon

- May include isolated or contiguous samples (see Lecture 4 on outlier detection/removal)
- Highly depends on the meaning given to usual/unusual

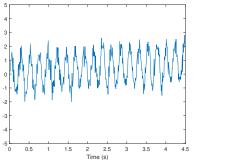
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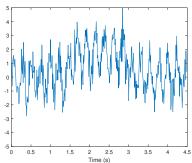
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Importance of stationarity

- As seen in Lecture 2, stationarity (e.g. at the wide sense) is a fundamental assumption when processing time series
- Necessity when using DFT, autocorrelation function or extracting features
- When observed during a long period of time, the system behaviour monitored by time series is likely to change over time, either smoothly or abruptly
- Several strategies can be used to deal with non-stationary time series, either simple or complex
- In some context, knowledge on these abrupt changes also carries relevant information on the system

Example





Smooth evolution vs. abrupt change

How to bypass the problem

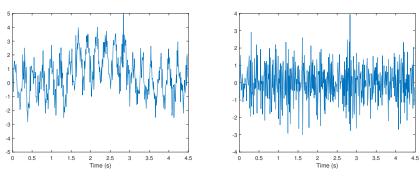
In a first order approximation, several strategies can be used to bypass the problem:

- If changes are smooth, the task can be seen as a detrending task: remove the slow phenomenon and only keep the seasonality that may be more stationary
- Divide the signal into small frames on which the signal is assumed to be stationary (see Lecture 2 on spectrogram)
- ► Instead of working on the original signal x[n], we can work on the signal derivative

$$x'[n] = x[n] - x[n-1]$$

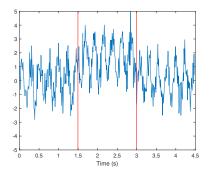
which in general has nicer stationarity properties Careful! This can imply to re-integrate the signal after processing, which can be a source of errors!

Example



Use of derivation to make the signal more stationary Left: original signal / Right: first order derivative

Problem statement

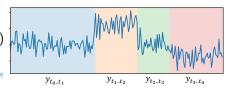


- When the changes are abrupt or when the estimation of the change-points is relevant in the context, we can use change-point detection methods
- Let assume that signal x[n]undergoes abrupt changes at times

$$\mathcal{T}^* = (t_1^*, \ldots, t_{K^*}^*)$$

- Goal: retrieve the number of change-points K^* and their times
- One assumption: offline segmentation (but can easily be adapted to online setting) [Truong et al., 2020]

$$(\hat{t}_1,\ldots,\hat{t}_K) = \underset{(t_1,\ldots,t_K)}{\operatorname{argmin}} \sum_{k=0}^K c(x[t_k:t_{k+1}])$$

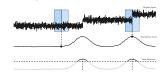


Cost function c(.)

- Measures the homogeneity of the segments
- Choosing c(.) conditions the type of change-points that we want to detect
- Often based on a probabilistic model for the data

Problem solving

- Optimal resolution with dynamic programming
- Approximate resolution (sliding windows...)



Cost function

$$(\hat{t}_1, \dots, \hat{t}_K) = \underset{(t_1, \dots, t_K)}{\operatorname{argmin}} \sum_{k=0}^K c(x[t_k : t_{k+1}])$$

Convention :
$$t_0 = 0$$
, $t_{K+1} = N$
 $a : b = [a, a + 1, ..., b - 1]$

- Function c(.) is characteristic of the notion of homogeneity
- The most common cost functions are linked to parametric probabilistic models: in this case change-points are defined as changes in the parameters of a probability density function [Basseville et al., 1993]
- Non-parametric cost functions can also be introduced when no model is available

Maximum likelihood estimation

Given a parametric family of distribution densities $f(\cdot|\theta)$ parametrized with $\theta \in \Theta$, a cost function can be derived:

$$c_{ML}(x[a:b]) = -\sup_{\theta} \sum_{n=a+1}^{b} \log f(x[n]|\theta)$$

- Corresponds to the assumption that on a regime, samples are i.i.d. according to a parametric distribution density
- On each regime, the parameters are estimated through maximum likelihood estimation
- This model can be adapted to several situations: change in mean, change in variance, change in both mean and variance...

Change in mean

The most popular is indubitably the L2 norm [Page, 1955]

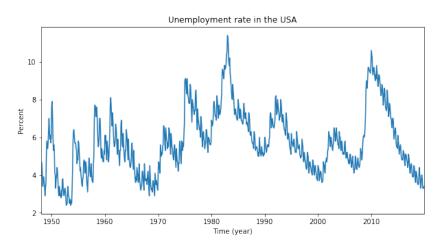
$$c_{L_2}(x[a:b]) = \sum_{n=a+1}^{b} ||x[n] - \mu_{a:b}||_2^2$$

where $\mu_{a:b}$ is the empirical mean of the segment x[a:b].

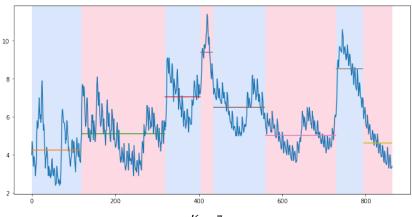
- Particular case of c_{ML} with Gaussian model with fixed variance
- Allows to detect changes in mean



Example

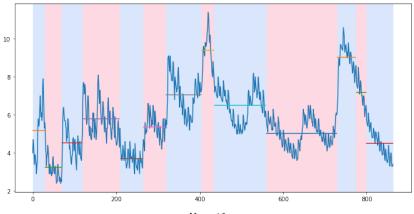


Example: Change-Point Detection with c_{L_2}



K = 7

Example: Change-Point Detection with c_{L_2}



K = 12

Change in mean and variance

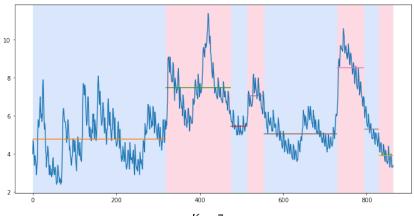
When the mean and variance change over time the cost function becomes

$$c_{\Sigma}(x[a:b]) = (b-a)\log \sigma_{a:b}^2 + \frac{1}{\sigma_{a:b}^2} \sum_{n=a+1}^b \|x[n] - \mu_{a:b}\|_2^2$$

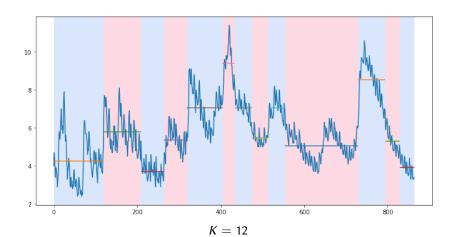
where $\mu_{a;b}$ and $\sigma_{a;b}^2$ are the empirical mean / variance of the segment x[a:b].

- Particular case of c_{MI} with Gaussian model with unknown mean and variance
- Can be adapted to multivariate time series by replacing the variance by the covariance matrix: in this case, changes of correlations between dimensions can also be detected [Lavielle, 1999]

Example: Change-Point Detection with c_{Σ}



K = 7



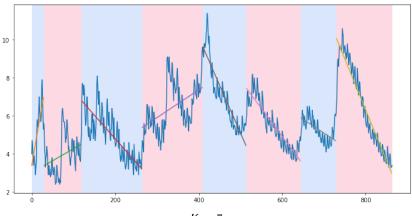
Change in slope and intercept

Change in slope and intercept can be handled in the general context of piecewise linear regression

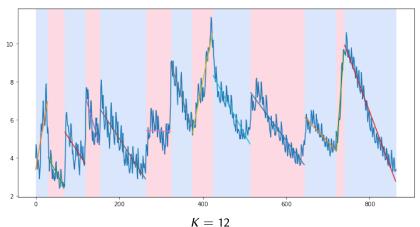
$$c_{linear}(x[a:b]) = \min_{\alpha} \sum_{n=a+1}^{b} \left\| x[n] - \sum_{i=1}^{M} \alpha_i \beta_i[n] \right\|_2^2$$

- **r** Functions $\beta_1[n], \ldots, \beta_M[n]$ are covariate functions and we seek for changes in the regression parameters
- Allows to detect changes in trend, seasonality, etc... [Bai et al., 1998]
- For slope and intercept, we choose $\beta_1[n] = 1$ and $\beta_2[n] = n$

Example: Change-Point Detection with c_{linear}



K = 7



Rank-based cost functions

In order to remove the need for a parametric probability model, one trick is to work on the notion of rank instead of the whole signal

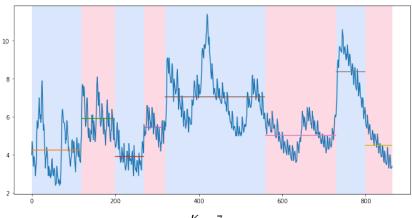
$$r[n] = \text{number of } i \text{ such that } x[i] < x[n]$$

- ▶ Robust and invariant with respect to amplitude changes: r[n] corresponds to the rank of sample x[n] in the time series **x**
- Cost functions can be derived by detecting changes in mean and/or variance in the rank signal [Lung-Yut-Fong et al., 2015]

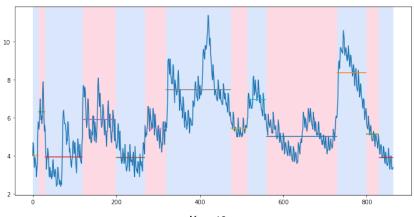
$$c_{rank}(x[a:b]) = \sum_{n=a+1}^{b} ||r[n] - \mu_{a:b}^{r}||_{2}^{2}$$

where $\mu_{a:b}^r$ is the empirical mean of the rank signal r[a:b]

Example: Change-Point Detection with c_{rank}



K = 7



K = 12

$$(\hat{t}_1,\ldots,\hat{t}_K) = \underset{(t_1,\ldots,t_K)}{\operatorname{argmin}} \sum_{k=0}^K c(x[t_k:t_{k+1}])$$

Convention :
$$t_0 = 0$$
, $t_{K+1} = N$

- Several methods can be used to solve this problem with a fixed K
- Optimal resolution with dynamic programming: find the true solution of the problem (but costly)
- Approximated resolution with windows: test for one unique change-point on a window

Optimal resolution

By denoting

$$\mathcal{V}(\mathcal{T},\mathbf{x}) = \sum_{k=0}^{K} c(x[t_k:t_{k+1}])$$

we can see that

$$\min_{|\mathcal{T}|=K} V(\mathcal{T}, \mathbf{x}) = \min_{0=t_0 < t_1 < \dots < t_K < t_{K+1} = N} \sum_{k=0}^{K} c(x[t_k : t_{k+1}])$$

$$= \min_{t \le N-K} \left[c(x[0 : t]) + \min_{t_0 = t < t_1 < \dots < t_{K-1} < t_K = N} \sum_{k=0}^{K-1} c(x[t_k : t_{k+1}]) \right]$$

$$= \min_{t \le N-K} \left[c(x[0 : t]) + \min_{|\mathcal{T}|=K-1} V(\mathcal{T}, x[t : N]) \right]$$

- Recursive problem (just like DTW in Lecture 1): resolution with dynamic programming [Bai et al., 2003]
- Two steps: computation of the cumulative costs + determination of the change-points

Algorithm 1 Algorithm Opt

```
Input: signal \{y_t\}_{t=1}^T, cost function c(\cdot), number of regimes K \geq 2.
for all (u, v), 1 \le u < v \le T do
    Initialize C_1(u, v) \leftarrow c(\{y_t\}_{t=u}^v).
end for
for k = 2, ..., K - 1 do
    for all u, v \in \{1, \dots, T\}, v - u \ge k do
        C_k(u, v) \leftarrow \min_{u+k-1 < t < v} C_{k-1}(u, t) + C_1(t+1, v)
    end for
end for
Initialize L, a list with K elements.
Initialize the last element: L[K] \leftarrow T.
Initialize k \leftarrow K.
while k > 1 do
    s \leftarrow L(k)
    t^* \leftarrow \operatorname{argmin}_{k-1 \leq t < s} C_{k-1}(1,t) + C_1(t+1,s)
    L(k-1) \leftarrow t^*
    k \leftarrow k - 1
end while
Remove T from L
```

Output: set L of estimated breakpoint indexes.

Complexity of $\mathcal{O}(KN^2)$

Approximated resolution

- Main limitation of optimal resolution: high complexity. Prohibitive for long time series...
- Approximated resolution methods exist, which are based on the single change-point detection, which is way easier to perform
- ► Idea: consider a sliding window of length 2w and for each position, determine if there is a change or not
- How to detect a single change by using the cost functions?

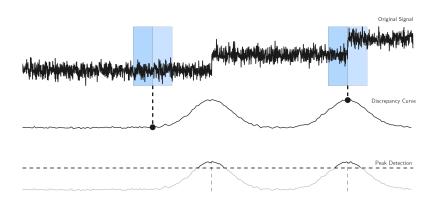
Discrepancy function

Given a window of length 2w centered on sample n, we compute the discrepancy function

$$d[n] = c(x[n-w:n+w]) - c(x[n:n+w]) - c(x[n-w:n])$$

- \triangleright The discrepancy function d[n] allows to compare
 - ► The homogeneity of the whole window c(x[n-w:n+w])
 - The homogenities of the right/left windows c(x[n:n+w]), c(x[n-w:n])
- Intuitively, if a change-point occurs at time n and if the window length w is well adapted, both subsegments x[n-w:n] and x[n:n+w] will be homogeneous (i.e. small values) and the whole segment x[n-w:n+w] will be heterogeneous (i.e. large values)
- Large values for d[n] suggests that a change-point is likely to appear at time n

Discrepancy function



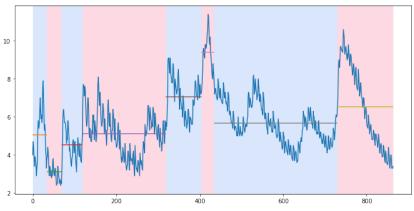
Sliding window approximated resolution

```
Algorithm 2 Algorithm Win
```

```
Input: signal \{y_t\}_{t=1}^T, cost function c(\cdot), half-window width w, peak search procedure PKSearch. Initialize Z \leftarrow [0,0,\dots] a T-long array filled with 0. \triangleright Score list. for t=w,\dots,T-w do p \leftarrow (t-w)..t. q \leftarrow t..(t+w). r \leftarrow (t-w)..(t+w). Z[t] \leftarrow c(y_r) - [c(y_p) + c(y_q)]. end for L \leftarrow \text{PKSearch}(Z) \triangleright Peak search procedure. Output: set L of estimated breakpoint indexes.
```

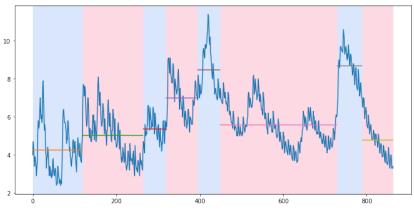
- ► Computation of the discrepancy function + peak search procedure to detect the *K* largest peaks
- \triangleright Complexity of $\mathcal{O}(N)$

Example: Sliding window CPD with c_{L_2}



$$K = 7, w = 20$$

Example: Sliding window CPD with c_L ,



K = 7, w = 40

How to set w

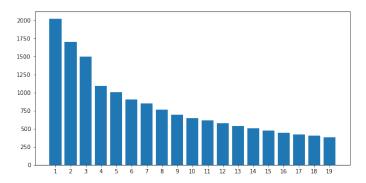
- Parameter w should correspond to the smallest length of stationarity: the discrepancy function makes sense if the two subsegments (right and left) are homogeneous
- Window length is also necessarily smaller than the smallest regime length
- But careful! In order to be relevant, the window length should be large enough so that there is enough samples to properly estimate the homogeneity
- Other vision: statistical tests between two sets of samples, each containing w samples. Good estimation requires a sufficient number of samples.

Finding the number of change points

- In all previously described algorithms, the number of change-point *K* was supposed to be known
- In practice, this parameter is difficult to set: as such, the total cost $\mathcal{V}(\mathcal{T}, \mathbf{x})$ will always decrease when K increases...
- Three solutions
 - Use heuristics by testing several values of K
 - Use a penalized formulation of the CPD problem to seek for a compromise between reconstruction error and complexity
 - Use supervised approaches from annotated signals

Heuristics for finding the number of change-points

- One easy solution is to test a set of change-points number K from 1 to K_{max} and to compute the sum of costs $\mathcal{V}(\mathcal{T}, \mathbf{x})$
- The *optimal* number of change-points can be estimated by searching for an elbow on the curve of $\mathcal{V}(\mathcal{T}, \mathbf{x})$ as a function of K



Penalized Change-Point Detection

- Intuitively, the optimal number of change-points is the one that allows the best compromise between the sum of costs $\mathcal{V}(\mathcal{T}, \mathbf{x})$ and the number of ruptures $|\mathcal{T}|$
- ▶ We actually had the same problem in various tasks: order estimation in AR models (Lecture 3), number of atoms in dictionary learning (Lecture 3 & 4), etc... In all cases the higher the order (and the number of parameters), the better the reconstruction
- **Model selection problem**: find the best model among a class of models
- Penalized change-point detection

$$(\hat{t}_1,\ldots,\hat{t}_{\hat{K}}) = \operatorname*{argmin}_{(t_1,\ldots,t_K),K} \sum_{k=0}^K c(x[t_k:t_{k+1}]) + \beta K$$

Penalized change-point detection

$$(\hat{t}_1,\ldots,\hat{t}_{\hat{K}}) = \underset{(t_1,\ldots,t_K),K}{\operatorname{argmin}} \sum_{k=0}^{K} c(x[t_k:t_{k+1}]) + \beta K$$

- Joint estimation of the change-point times and the number of change-points
- Parameter β penalizes the introduction of a new change-point in the model: an additional change-point should decrease the sum of costs $\mathcal{V}(\mathcal{T}, \mathbf{x})$ by at least β
- Luckily, this problem is even easier to solve than the original one with fixed *K*!

Pruning strategy

Given two times s and t such that t < s < N, remark that if

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:s]) \geq \min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right]$$

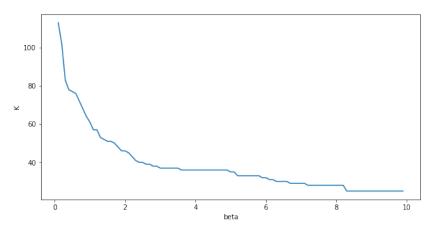
then t cannot be the last change point prior to N (demo in the last slides).

- Considerable speed-up since most times will not satisfy this criterion
- **Pruned Exact Linear Time (PELT)** algorithm: under the assumption that regime lengths are randomly drawn from a uniform distribution, the complexity of PELT is $\mathcal{O}(N)$ [Killick et al., 2012]
- Optimal algorithm: exact solution

PELT algorithm

Algorithm 3 Algorithm Pelt

```
Input: signal \{y_t\}_{t=1}^T, cost function c(\cdot), penalty value \beta. Initialize Z a (T+1)-long array; Z[0] \leftarrow -\beta. Initialize L[0] \leftarrow \emptyset. Initialize \chi \leftarrow \{0\}. \diamond Admissible indexes. for t=1,\ldots,T do \hat{t} \leftarrow \mathop{\rm argmin}_{s \in \chi} \left[Z[s] + c(y_{s..t}) + \beta\right]. Z[t] \leftarrow \left[Z[t] + c(y_{t..t}) + \beta\right] L[t] \leftarrow L[t] \cup \{\hat{t}\}. \chi \leftarrow \{s \in \chi : Z[s] + c(y_{s..t}) \le Z[t]\} \cup \{t\} end for Output: set L[T] of estimated breakpoint indexes.
```



 β can be difficult to set: no explicit formula between β and K

Model selection criterion

Two popular criteria can be used to estimate the relevance of a model

▶ Bayesian information criterion (BIC) [Schwarz, 1978]

$$BIC = k \log N - 2 \log \hat{L}$$

Akaike information criterion (AIC) [Akaike, 1974]

$$AIC = 2k - 2\log \hat{L}$$

where

- k is the number of parameters
- *N* is the number of samples
- \hat{L} is the maximum value of the likelihood function for the model

Standard criterion

In the context of change-point detection with L2 cost function, these criteria provide estimates for the β parameter:

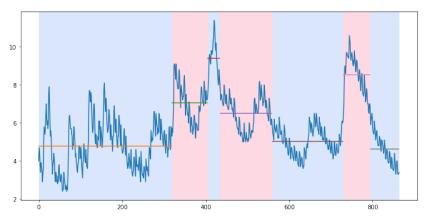
Bayesian information criterion (BIC) for L2 change-point detection

$$\beta = 4\sigma^2 \log N$$

Akaike information criterion (AIC) for L2 change-point detection

$$\beta = 4\sigma^2$$

Example



Results obtained with the BIC criterion

Supervised Change-Point Detection

- Parameter β can also be learned from a collection of annotated signals $\mathbf{x}^{(1)}, \dots, \mathbf{x}^{(M)}$ with annotations $\mathcal{T}_*^{(1)}, \dots, \mathcal{T}_*^{(M)}$ [Truong et al., 2017]
- By denoting

$$V_eta(\mathcal{T},\mathbf{x}) = \sum_{k=0}^K \ c(x[t_k:t_{k+1}]) + eta K$$

the relevance of a value of β can be assessed by computing the excess penalized risk

$$E(\mathbf{x}^{(\ell)}, \beta) = V_{\beta}(\mathcal{T}_{*}^{(\ell)}, \mathbf{x}^{(\ell)}) - \min_{\mathcal{T}} V_{\beta}(\mathcal{T}, \mathbf{x}^{(\ell)})$$

This quantity reflects how well we can approach the annotated segmentation with a given value of β

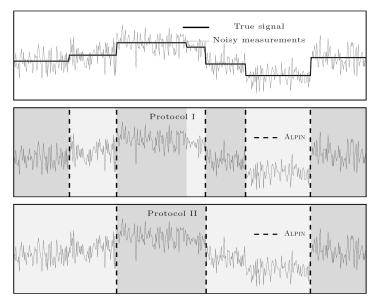
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Supervised Change-Point Detection

- ▶ The function $\beta \mapsto E(\mathbf{x}^{(\ell)}, \beta)$ is a convex function of β , which can be easily optimized with off-the-shelf solvers
- The final optimization problem writes

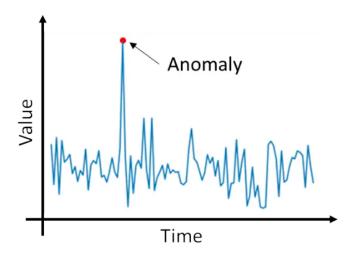
$$eta_{opt} = \operatorname{argmin}_{eta>0} rac{1}{M} \sum_{\ell=1}^M E(\mathbf{x}^{(\ell)}, eta)$$

Experimental results show that with only a few annotated examples, it is possible to find an adequate range for β

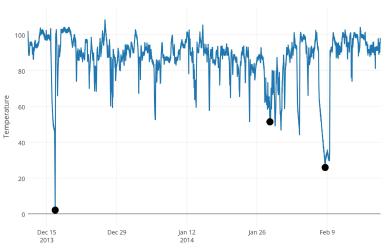


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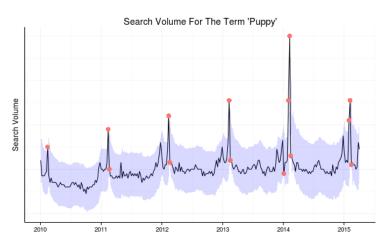
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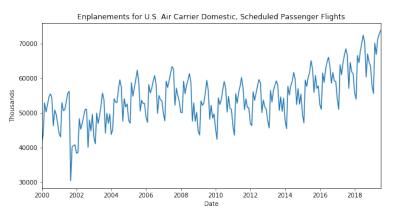
Easy: an anomaly is a too small or too large value (outlier)



More complex: some small/large values are anomalies



Anomalies depend in the previous values



Anomalies correspond to unusual events

Anomaly Detection

Anomalies can take various forms and have different meanings [Chandola et al., 2009]:

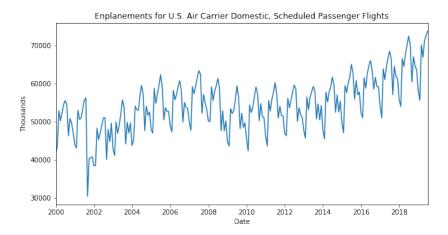
- Outliers, i.e. isolated samples with exceptionally large/low values
- Bursts of outliers, i.e. segments that do not coincide with what is observed usually in the time series (in terms of values)
- Unusual events that breaks the regularity within the time series

Outlier detection

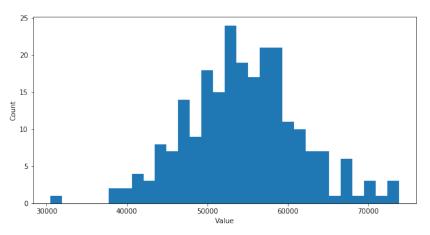
Simple anomalies (isolated samples or contiguous samples) can be detected with techniques already described in Lecture 4:

- Statistical methods:
 - ► Global: Histogram visualization to detect aberrant values (see Lecture 4)
 - Adaptive: Threshold-based methods on sliding windows (mean/standard deviation or median)
- Model-based methods:
 - Residual and prediction error (trend+seasonality, sinusoidal or AR model)

Example

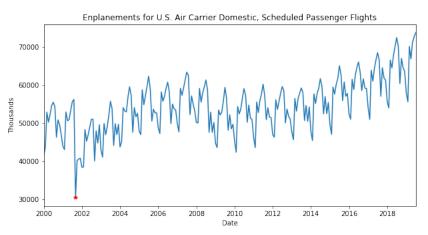


Example: Histogram



One outlier can be considered as an anomaly

Example: Histogram



Only one detected anomaly

Adaptive statistical methods

- ▶ The main idea is to use sliding window and to perform a statistical test for outlier detection
- Contrary to histogram, these methods allow to take into account the local context but careful, time information is lost! Only the distribution of values is used for detection.
- Multitude tests can be used but the most common are
 - ► Mu/sigma [Roberts, 2000]:

$$|x[n] - \mu_n| > \lambda \sigma_n$$

where μ_n and σ_n are respectively the local mean/standard deviation around sample n and λ a threshold.

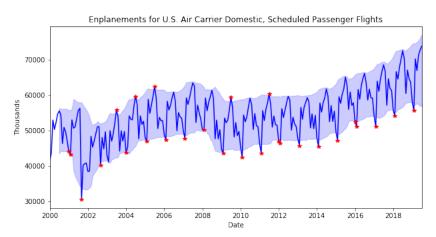
Under i.i.d. Gaussian assumption, $\lambda = 1 \rightarrow 68\%$, $\lambda = 2 \rightarrow 95\%$, $\lambda = 3 \rightarrow 99.7\%$

► Median/median absolute deviation [Leys et al., 2013]:

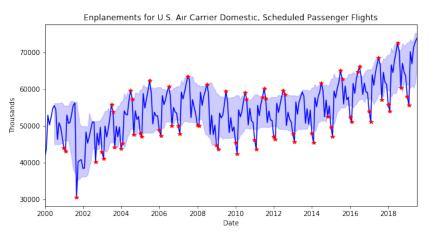
$$|x[n] - \mathsf{med}_n| > \lambda \, \mathsf{mad}_n$$

where med_n and mad_n are respectively the local median/median absolute deviation around sample n and λ a threshold.

Example: Mu-Sigma



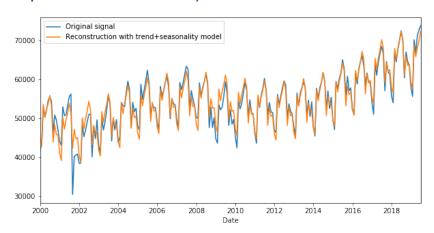
Mu-Sigma, $\lambda = 1.5$, window length of 12 samples



Med-Mad, $\lambda = 1.5$, window length of 12 samples

Model-based anomaly detection

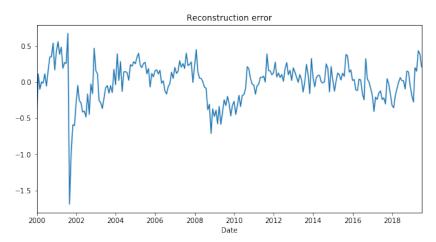
- ► Idea: use a time series model to detect anomaly [Yamanishi et al., 2002; Hill et al., 2010]
- Advantage: truly takes into account the temporal aspects
- ► Three steps:
 - 1. Choose an adequate model and learn the parameters
 - 2. Compute the prediction/signal reconstruction
 - 3. Anomalies are samples that diverge from the model



- ► Trend: polynomial of degree 4
- Seasonality: cosine/sine functions with fundamental frequencies multiples of

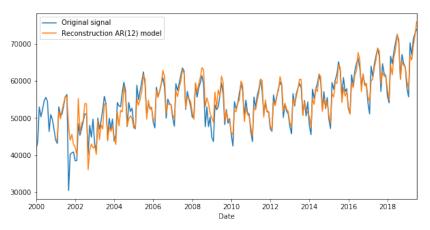
 $\frac{1}{12}$

Example: trend+seasonality



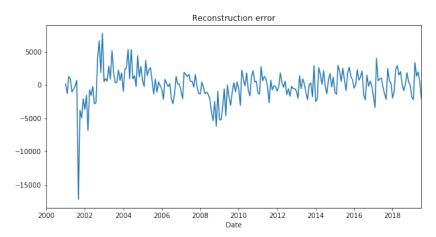
Not only large/small values but also temporal progression

Example: AR model



AR model with p = 12

Example: AR model



Anomaly also changes the prediction of the next *p* samples

Distance-based methods

- Some anomalies may be more complex to detect as they are not characterized by aberrant values but by a new behavior that was not previously seen in the time series
- In this case, anomalies can only be defined as a divergence from a normal behavior
- ▶ This task is the dual of the task already seen in Lecture 1 (Pattern Detection/Extraction), and the same techniques can therefore be used
- Instead of searching for repetitive patterns, we are searching for non-repetitive events!

Unsupervised anomaly detection

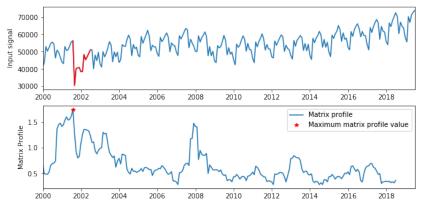
Reminder: Matrix profile [Yeh et al., 2016]: given a pattern length L, compute

$$m[n] = \min_{i > n+L \text{ or } i < n-L} d(x[n:n+L-1], x[i:i+L-1])$$

- Small matrix profiles values indicate that the subsequence has been found elsewhere in the time series, suggesting that it could be a pattern
- Efficient computation with normalized Euclidean distance (see Lecture 1)

What about large values in the matrix profile?

Example: matrix profile



Matrix profile with window of length L = 12 months

- By examining large values on the matrix profile, anomalies can be detected
- ► Subsequences that are *far* from all subsequences in the signal: likely to correspond to new behaviors
- Advantages: no need for a parametric model
- Necessitates to have a rough idea of the scale of the anomaly (parameter *L*)

Other distance-based approaches

Clustering approaches can be used for detecting anomalies:

- 1. Divide the signal into (possibly) overlapping subsequences
- 2. Perform clustering on the subsequences (k-Means, spectral clustering etc...)
- 3. Subsequences that are far from their centroids/medoids are likely to be outliers

More details in [Schmidl et al., 2022; Boniol et al., 2022]

A word on supervised approaches

- Semi-supervised approaches: use knowledge on normality
 - Learn a model on normality and detect anomalies in the residual or as a derivation from normality
 - Example: use annotated templates representing normal behavior, retrieve them in the signal up to a measure of fit (see Lecture 1), and detect all segments in the time series that do not correspond to a known pattern as anomalies
- ➤ Supervised approaches: supervised classification techniques can also be used: SVM, random forests, neural networks...

Contents

- 1. Problem statement
- Change-Point Detection
- 3. Anomaly Detection
- 4. Evaluation of event detection methods

Evaluation of event detection methods

- Several time series ML tasks are event detection tasks: pattern recognition (Lecture 1), change-point detection, anomaly detection...
- In order to benchmark the tested methods, one need to use
 - An annotated dataset where the events of interested have been highlighted
 - Some relevant metrics of evaluation
- ▶ What does a *good detection* mean?

Point-based vs. range-based

When the events consist of single points, the method can be assessed with the standard precision/recall metric:

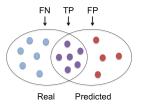
$$precision = \frac{TP}{TP + FP}$$

$$recall = \frac{TP}{TP + FN}$$

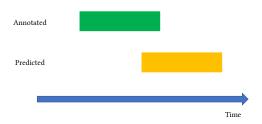
where *TP* is the number of true positive, *FP* the number of false positive and *FN* the number of false negative

These metrics are comprised between 0 and 1, and we can plot the precision/recall curve to benchmark the methods

What if the events are range-based?



Range-based detection: example



- Event is correctly detected
- ▶ BUT the detection is a bit delayed: would not be suitable for e.g. anomaly detection in industrial monitoring

Range-based detection: example



- Event is correctly detected
- BUT the duration of the event is poorly estimated

Range-based detection: example



- Event is correctly detected
- ▶ BUT the two annotated events are detected as one single event

Metrics for event detection

Several principles can be taken into account [Tatbul et al., 2018]:

- Existence: Catching the existence of the event (even by predicting only a single point), by itself, might be valuable for the application.
- ➤ *Size:* The larger the size of the correctly predicted portion of the event, the higher the recall score.
- ▶ *Position:* In some cases, not only size, but also the relative position of the correctly predicted portion of the event might matter to the application.
- Cardinality: Detecting the event with a single prediction range may be more valuable than doing so with multiple different ranges in a fragmented manner.

Formulation

▶ We consider a set of predicted intervals $P = \{P_1, \dots, P_{N_P}\}$ and a set of real intervals $R = \{R_1, \dots, R_{N_R}\}$, the recall can be computed as:

$$recall = \frac{1}{N_R} \sum_{i=1}^{N_R} recall(R_i, P)$$

- ▶ The term $recall(R_i, P)$ will be defined as a weighted sum of several terms that will assess how well event R_i has been detected
- ▶ The same definition can be computed for the precision, but this time as

$$precision = \frac{1}{N_P} \sum_{i=1}^{N_P} precision(R, P_i)$$

Formulation

Existence (only used for recall):

$$\mathsf{existence}(R_i, P) = \begin{cases} 1 & \text{if } \sum_{j=1}^{N_p} |R_i \cap P_j| \ge 1 \\ 0 & \text{elsewhere} \end{cases}$$

Size/position (used for precision and recall):

$$size_position(R_i, P) = \sum_{j=1}^{N_p} w(R_i, R_i \cap P_j)$$

where w(A, B) is an overlap score (between 0 and 1) between $R_i \cap P_j$ and the *desirable* portion of R_i (can voluntary introduce a bias if e.g. we wish to detect the event R_i in advance, or the middle part of R_i , etc...)

Cardinality (used for position and recall):

cardinality(
$$R_i, P$$
) =
$$\begin{cases} 1 & \text{if } R_i \text{ overlaps with at most one } P_j \\ \gamma(R_i, P) & \text{elsewhere} \end{cases}$$

where γ is a penalty function

Formulation

Recall:

$$\operatorname{recall}(R_i, P) = \alpha \operatorname{existence}(R_i, P) + (1 - \alpha) \operatorname{cardinality}(R_i, P) \times \operatorname{size_position}(R_i, P)$$

Precision:

$$precision(R, P_i) = cardinality(R, P_i) \times size_position(R, P_i)$$

- Note that different functions w and γ can be used for precision and recall, and several choices can be used (see [Tatbul et al., 2018] and associated mini-project for details)
- If all events in R and P are single-point, $\alpha=0,\gamma(.,.)=1$ and w(.,.) is the percentage of common points, then this definition is compliant with the regular precision/recall definition

How to choose the parameters

- Several possible parametrizations are provided in the original article: depends on the usecases
- ➤ One simpler solution is to compute the Intersection Over Union (IoU) metric between the detected segment and the true segment

$$loU = \frac{|P_i \cap R_j|}{|P_i \cup R_j|}$$

and use a threshold value (e.g. 25%, 50%, 75%...) as a detection criteria

▶ Performances can be provided with different threshold values so as to give a better idea of the accuracy of the detection method

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List of possible topics/projects

- How to contribute to the ruptures package (see with C. Truong) https://centre-borelli.github.io/ruptures-docs/
- Truong, C., Oudre, L., & Vayatis, N. (2017). Penalty learning for changepoint detection. In 2017 25th European Signal Processing Conference (EUSIPCO) (pp. 1569-1573). IEEE.
 - How to learn the penalty for change point detection
- Fearnhead, P., & Rigaill, G. (2019). Changepoint detection in the presence of outliers. Journal of the American Statistical Association, 114(525), 169-183.

 How to detect change-points in presence of outliers
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 How to use change point detection to study soccer games
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 How to introduce a sparsity penalty into change-point detection problems
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 How to introduce graph constraints into change-point detection problems
- Jewell, S. W., Hocking, T. D., Fearnhead, P., & Witten, D. M. (2020). Fast nonconvex deconvolution of calcium imaging data. Biostatistics, 21(4), 709–726.
 How to apply change-point detection to biology

List of possible topics/projects

- Chin, S. C., Ray, A., & Rajagopalan, V. (2005). Symbolic time series analysis for anomaly detection: A comparative evaluation. Signal Processing, 85(9), 1859-1868.
 - How to use symbolic representation for detecting anomalies
- Chandola, V., Banerjee, A., & Kumar, V. (2010). Anomaly detection for discrete sequences: A survey. IEEE transactions on knowledge and data engineering, 24(5), 823-839.
- How to detect anomalies in discrete time series
- Boniol, P., Linardi, M., Roncallo, F., & Palpanas, T. (2020, April). Automated Anomaly Detection in Large Sequences. In 2020 IEEE 36th International Conference on Data Engineering (ICDE) (pp. 1834–1837). IEEE.
 How to use detect anomalies in large time series
- Nakamura, T., Imamura, M., Mercer, R., & Keogh, E. MERLIN: Parameter-Free Discovery of Arbitrary Length Anomalies in Massive Time Series Archives. In Proc. 20th IEEE Intl. Conf. Data Mining.

 How to detect anomalies with different lengths
- ► Tatbul, N., Lee, T. J., Zdonik, S., Alam, M., & Gottschlich, J. (2018). Precision and recall for time series. arXiv preprint arXiv:1803.03639. How to assess event detection techniques
- Schmidl, S., Wenig, P., & Papenbrock, T. (2022). Anomaly detection in time series: a comprehensive evaluation. Proceedings of the VLDB Endowment, 15(9), 1779–1797.
 Wonderful article with tons of references, implementations, etc...
- Boniol, P., Paparrizos, J., Kang, Y., Palpanas, T., Tsay, R. S., Elmore, A. J., & Franklin, M. J. (2022). Theseus: navigating the labyrinth of time-series anomaly detection. Proceedings of the VLDB Endowment, 15(12), 3702-3705.
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PELT pruning strategy

$$(\hat{t}_1,\ldots,\hat{t}_{\hat{K}}) = \operatorname*{argmin}_{(t_1,\ldots,t_K),K} \sum_{k=0}^K c(x[t_k:t_{k+1}]) + \beta K$$

We want to show that given two times s and t such that t < s < N, if

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:s]) \ge \min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right]$$

then t cannot be the last change point prior to N

• Given two times s and t such that t < s < N, assume that

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:s]) \ge \min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right]$$

▶ By adding $c(x[s:N]) + \beta$ to this equation, and using the fact that $c(x[t:s]) + c(x[s:N]) \le c(x[t:N])$ (adding a breakpoint decreases the sum of costs), it goes that

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:N]) + \beta \ge \min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right] + c(x[s:N]) + \beta$$

PELT pruning strategy

Now, if t < s is the last change point prior to N, then

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right] = \min_{t} \left[\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:s]) + \beta \right]$$

But we also have

$$\min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:t]) + \beta |\mathcal{T}| \right] + c(x[t:N]) + \beta \ge \min_{\mathcal{T}} \left[V(\mathcal{T}, x[0:s]) + \beta |\mathcal{T}| \right] + c(x[s:N]) + \beta |\mathcal{T}|$$

► This shows that s is a better choice than t in the first equation, and thus t cannot be the last change point prior to N