## Time Series

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## Importance of Time Series Analysis

Ample of time series data is being generated from a variety of fields. And hence the study of time series analysis holds a lot of applications

Time series analysis is important in different areas:

- Economics
- Finance
- Healthcare
- Environmental Science
- Sales Forecasting

## Why & where Time Series is used?

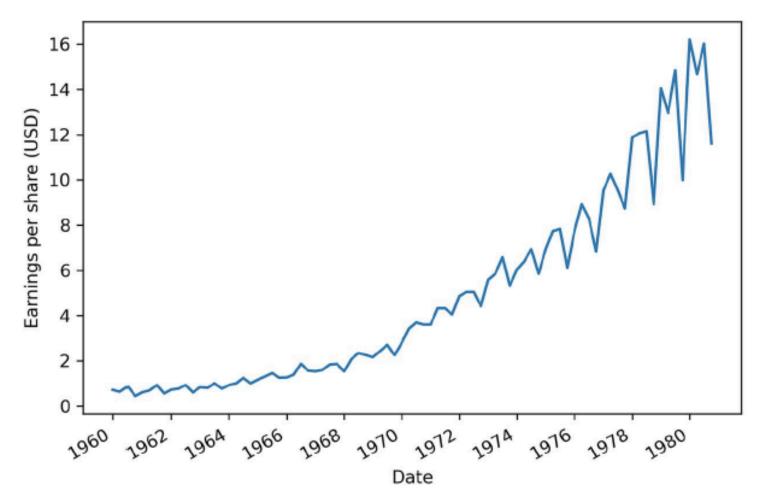
Time series data can be analysed in order to extract meaningful statistics and other characteristics

It's used in at least in the four scenarios:

- Business Forecasting
- Understanding past behaviour
- Plan the future
- Evaluate current accomplishment

### Time Series

• Time Series is simple a set of data points ordered in time intervals (usually equal time intervals)



## Components of a Time Series

Time series decomposition is a process by which we separate a time series into its components:

#### Trend

Represents the slow-moving changes in a time series. It is responsible for making the series gradually increase or decrease over time

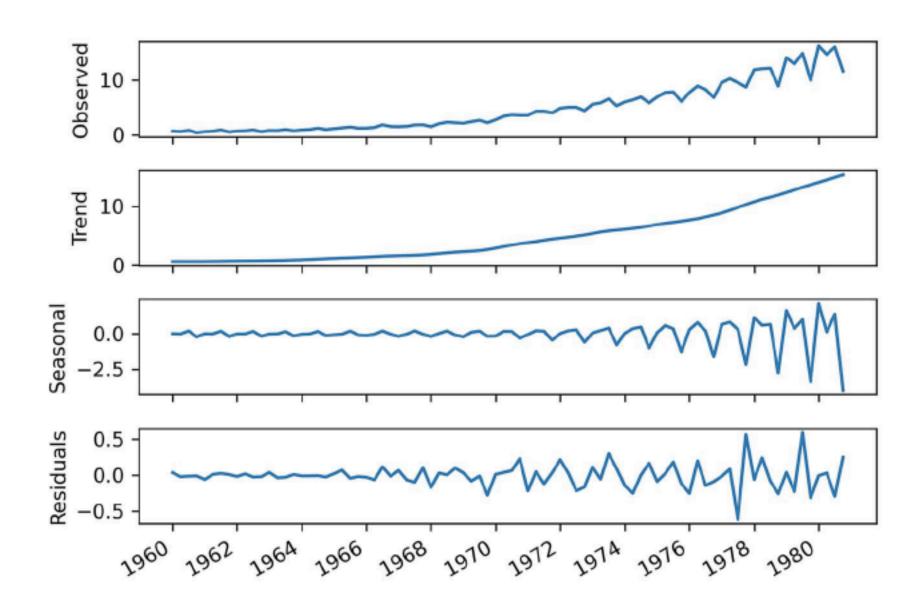
#### Seasonality

Represents the seasonal pattern in the series. The cycles occur repeatedly over a fixed period of time

#### Residuals

Represent the behaviour that cannot be explained by the trend and seasonality components. They correspond to random errors, also termed white noise

## Components of a Time Series



## Baseline models

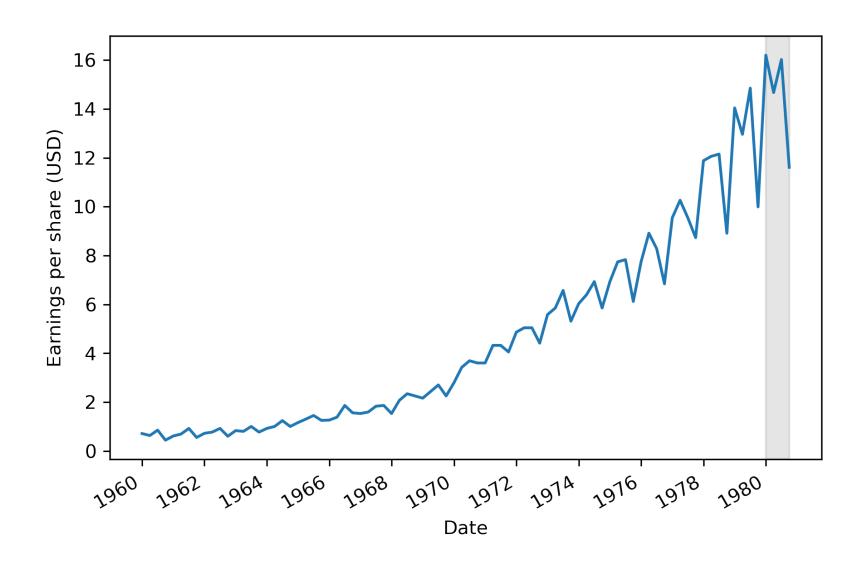
#### Baseline model

- A baseline model is a trivial solution for forecasting a problem
- It relies on heuristics or simple statistics and is usually the simplest solution
- It does not require model fitting, and it is easy to implement

#### Examples of baseline models:

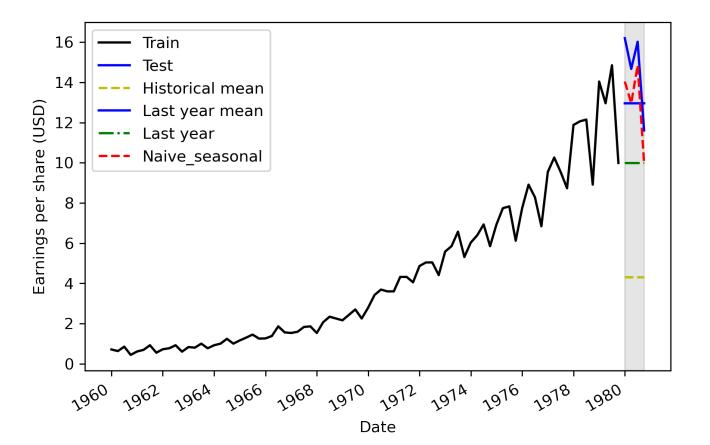
- historical mean
- last year's mean
- last known value
- naive seasonal forecast

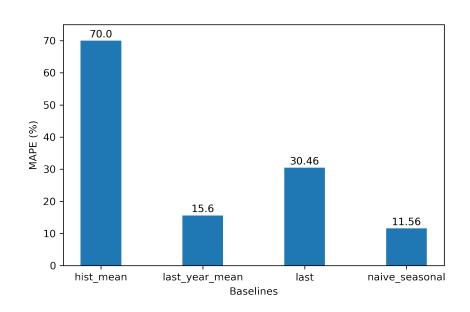
## Predict the last four quarters with baseline models



## Predict the last four quarters with baseline models

	date	data	hist_mean	last_year_mean	last	naive_seasonal
80	1980-01-01	16.20	4.3085	12.96	9.99	14.04
81	1980-04-01	14.67	4.3085	12.96	9.99	12.96
82	1980-07-02	16.02	4.3085	12.96	9.99	14.85
83	1980-10-01	11.61	4.3085	12.96	9.99	9.99





## Random Walk

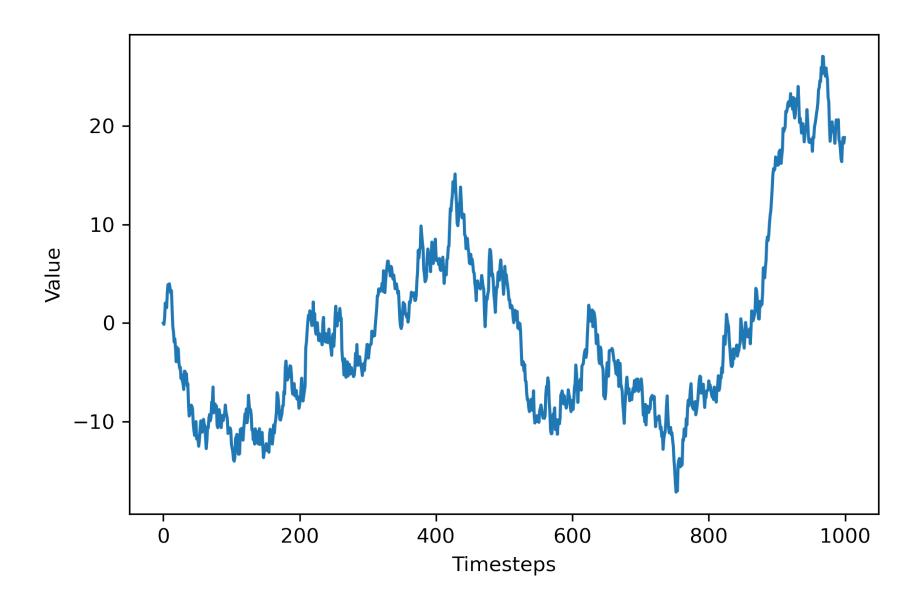
#### Random Walk

- A random walk is a process in which there is an equal chance of going up or down by a random number
- Random walks often expose long periods where a positive or negative trend can be observed. They are also often accompanied by sudden changes in direction

A random walk can be mathematically express with the following equation

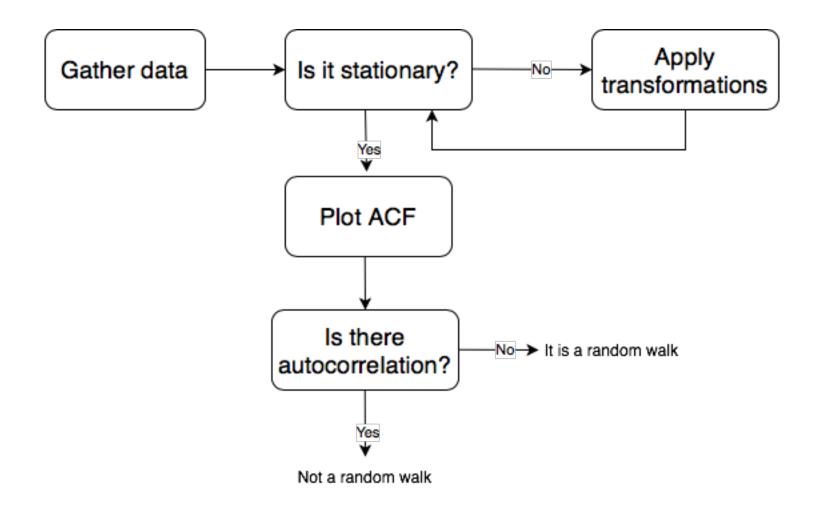
$$y_t = C + y_{t-1} + \epsilon_t$$

## Random walk example



## Identifying a Random Walk

A random walk is a series whose first difference is stationary and uncorrelated



## Stationarity

Before applying any statistical model on a time series, the series has to be stationary, which means that, over different time periods:

- 1. it should have constant mean
- 2. It should have constant variance or standard deviation
- 3. Autocorrelation do not change on time

## How to make the time series stationary

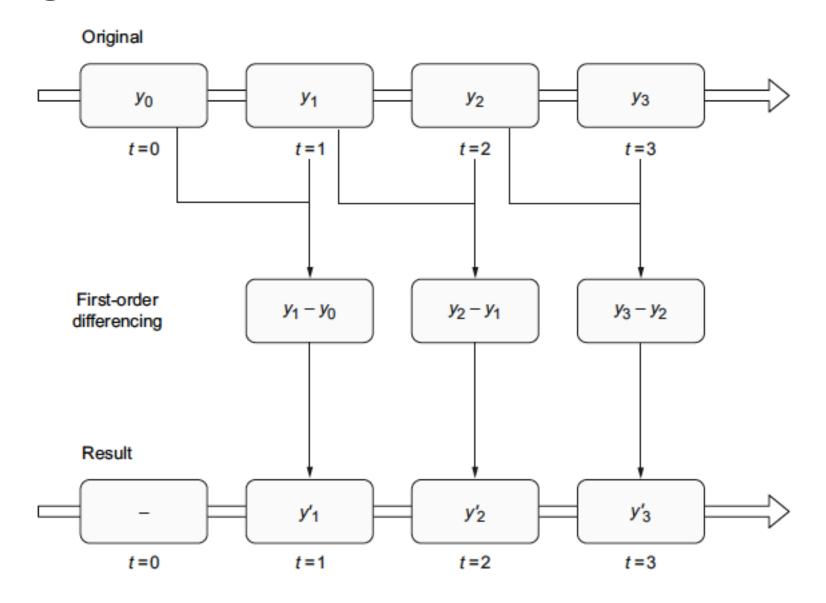
- Differencing is a transformation that calculates the change from one timestep to another
- This transformation helps stabilize the mean, which in turn removes or reduces the trend and seasonality effects
- Differencing involves calculating the series of changes from one timestep to another

$$y_t' = y_t - y_{t-1}$$

It is possible to difference a time series many times:

- taking the difference once is applying a first-order differencing
- taking it a second time would be a second-order differencing

## Visualizing a first-order difference



## Test to check if a series is stationary

#### **ADCF Test - Augmented Dickey-Fuller test**

Null hypothesis: says that the time series is non-stationary

The result of this test is the ADF statistic, which is a negative number. The more negative it is, the stronger the rejection of the null hypothesis

If the p-value is less than 0.05, we can also reject the null hypothesis and say the series is stationary

## ADCF Test to check if a series is stationary

```
from statsmodels.tsa.stattools import adfuller

ADF_result = adfuller(random_walk)

print('ADF Statistic:', round(ADF_result[0],3))
print('p-value:', round(ADF_result[1],3))

ADF Statistic: -0.966
p-value: 0.765
```

Since the series is not stationary a first-order differencing will be applied

```
diff_random_walk = np.diff(random_walk, n=1)

ADF_result = adfuller(diff_random_walk)

print('ADF Statistic:', round(ADF_result[0],3))
print('p-value:', round(ADF_result[1],3))

ADF Statistic: -31.789
p-value: 0.0
```

## Autocorrelation function (ACF)

The autocorrelation function (ACF) measures the linear relationship between lagged values of a time series

It measures the correlation of the time series with itself

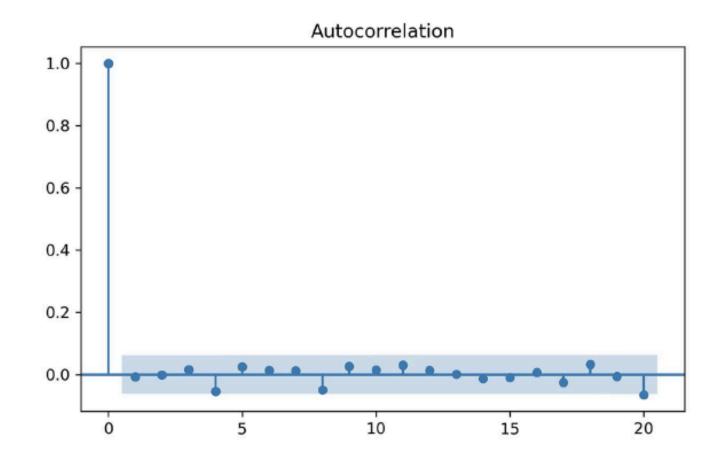
The ACF calculates the autocorrelation coefficient between:

$$y_t$$
 and  $y_{t-1} : r_1$   
 $y_t$  and  $y_{t-2} : r_2$ 

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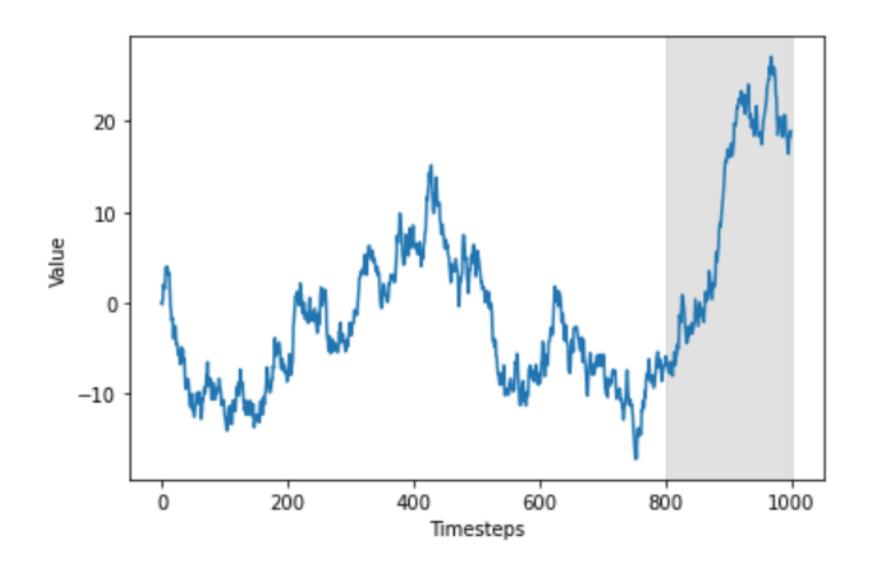
In the ACF plot the coefficient is the dependent variable, while the lag is the independent variable

## ACF plot of the random walk



There are no significant coefficients after lag 0, which is a clear indicator of a random walk - can be described as white noise

## Forecasting a random walk



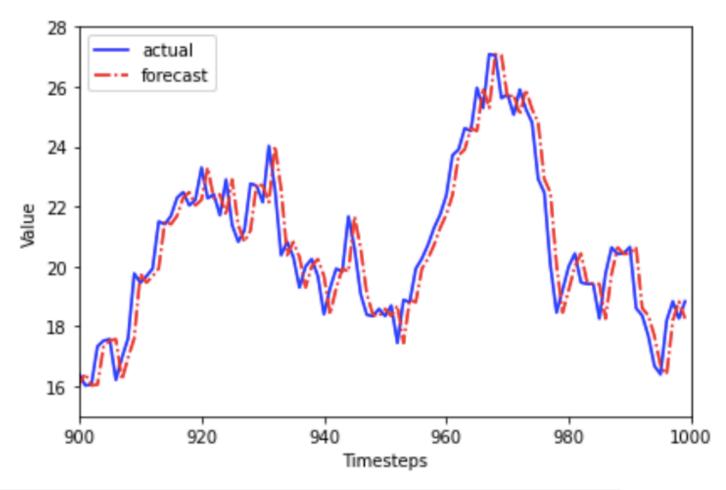
## Forecasting a random walk

- To predict a random walk, we can only use naive forecasting methods or baseline methods
- Forecasting a random walk on a **long horizon** does not make sense the randomness portion is magnified in a long horizon where many random numbers are added over the course of many timesteps

• In a random walk, it is only possible to forecast the **next timestep** 

• The present observed value is used as a forecast for the next timestep. Once a new value is recorded, it will be used as a forecast for the following timestep

## Forecasting a random walk



```
In [39]: mse_one_step = mean_squared_error(test['value'], df_shift[800:])
    mse_one_step
```

Out[39]: 0.9256876651440581

# Forecasting with statistical models

## Statistical models for time series forecasting

- MA(q) models
- AR(p) models
- ARMA(p,q) models
- ARIMA(p,d,q) models for non-stationary time series
- SARIMA(p,d,q)(P,D,Q)<sub>m</sub> for seasonal time series
- SARIMAX models to include external variables in the forecast
- VAR(p) model for predicting many time series at once

## Stationary time series

## Moving Average MA(q)

## Moving Average model

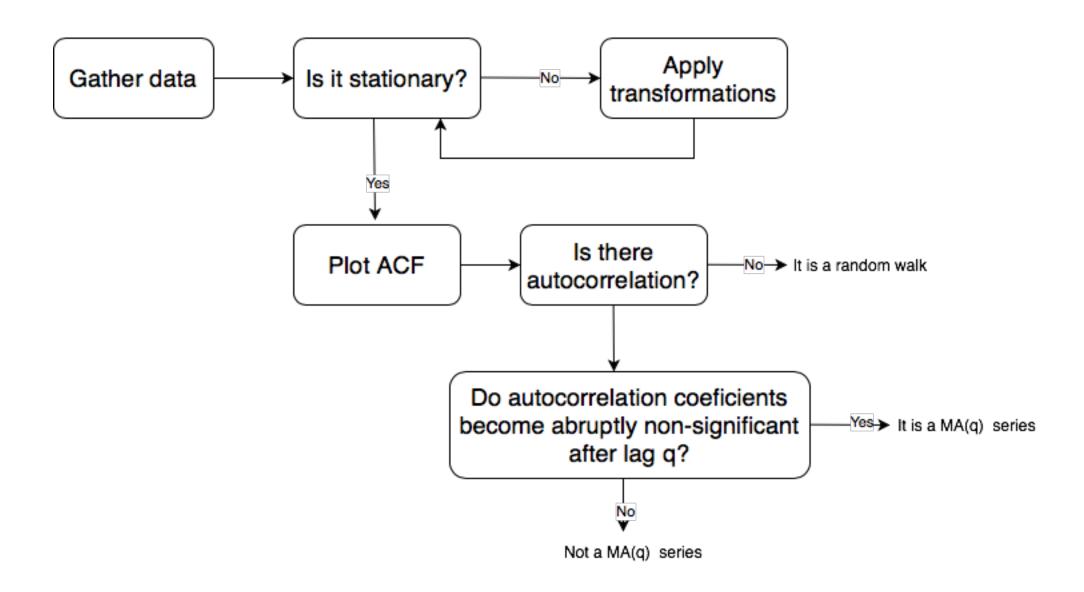
- In a moving average (MA) model, the current value depends linearly on the mean of the series, the current error term, and past error terms
- The moving average model is denoted as MA(q), where q is the order

The general expression of an MA(q) model is

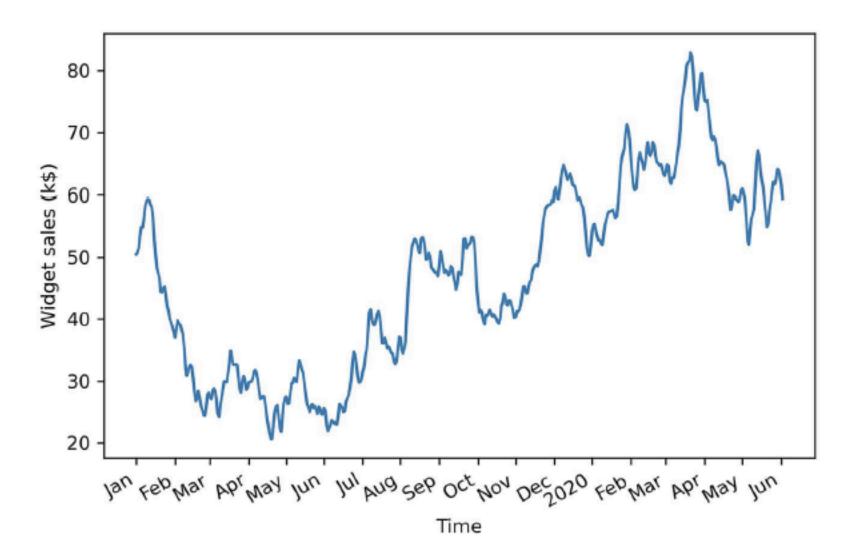
$$y_t = \mu + e_t + \theta_1 e_{t-1} + \theta_2 e_{t-2} + \dots + \theta_q e_{t-q}$$

The order q of the moving average model determines the number of past error terms that affect the present value

## Identifying a Moving Average series



## Moving Average series example



Volume of sales for a Company over 500 days, starting on January 1, 2019

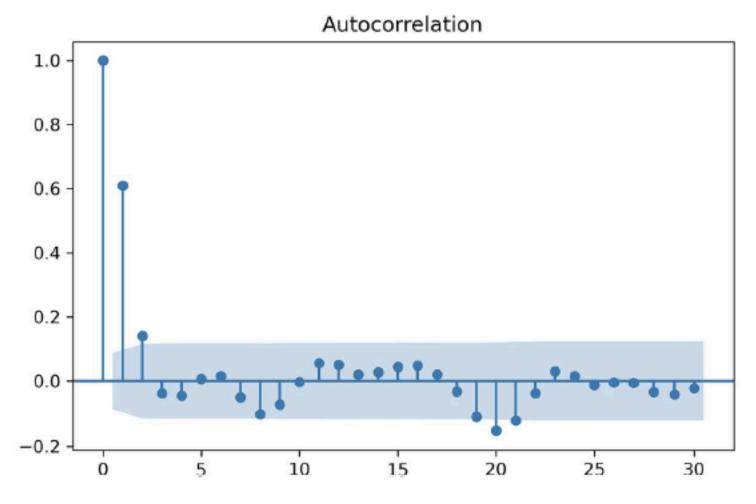
## Identifying the order of a Moving Average series

**Test for stationarity** 

```
# Test for stationarity
from statsmodels.tsa.stattools import adfuller
ADF result = adfuller(df['widget sales'])
print(f'ADF Statistic: {ADF result[0]}')
print(f'p-value: {ADF_result[1]}')
ADF Statistic: -1.5121662069359012
p-value: 0.5274845352272624
# first-order differencing to make it stationary
widget sales diff = np.diff(df['widget sales'], n=1)
ADF_result = adfuller(widget_sales_diff)
print(f'ADF Statistic: {ADF result[0]}')
print(f'p-value: {ADF_result[1]}')
ADF Statistic: -10.576657780341957
p-value: 7.076922818587346e-19
```

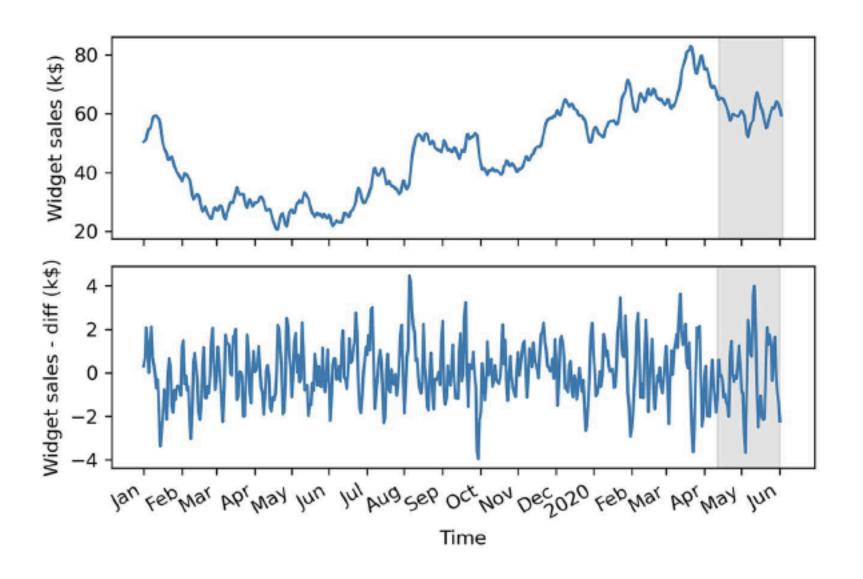
## Identifying the order of a Moving Average series

Plot the autocorrelation function - ACF



Stationary moving average series of order 2 - MA(2)

## Forecasting a Moving Average series



## Forecasting using the MA(q) model

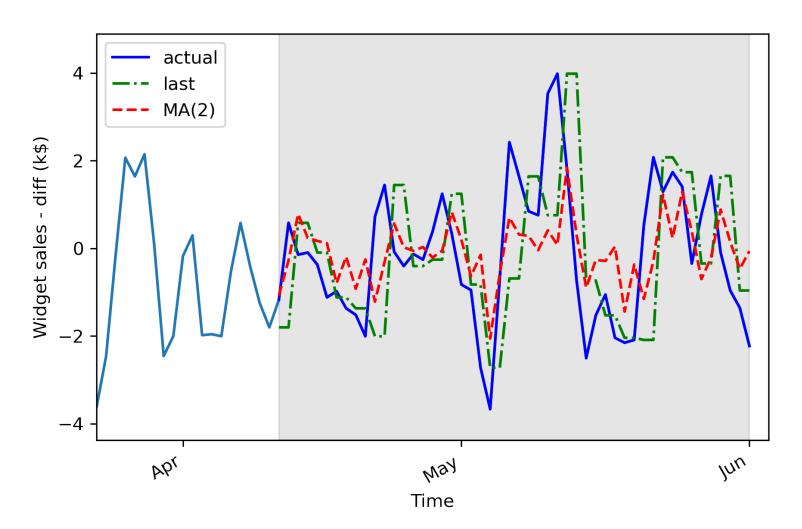
- When using an MA(q) model, forecasting **beyond q steps** into the future will simply return the mean there are no error terms to estimate beyond q steps
- It is possible to use *rolling forecasts* to predict up to q steps at a time
- In a dataset with 500 steps, to predict the last 50 steps:
  - First pass: train on the first 449 timesteps to predict timesteps 450 and 451
  - Second pass: train on the first 451 timesteps to predict timesteps 452 and 453
  - ...
  - This is repeated until the values at timesteps 498 and 499 are predicted

## A function for rolling forecasts on a horizon

return pred MA

```
from statsmodels.tsa.statespace.sarimax import SARIMAX
def rolling forecast(df: pd.DataFrame, train len: int, horizon: int, window: int) -> list:
   total len = train len + horizon
   pred MA = []
   for i in range(train len, total len, window):
       model = SARIMAX(df[:i], order=(0,0,2))
       res = model.fit(disp=False)
       predictions = res.get_prediction(0, i + window - 1)
       oos pred = predictions.predicted mean.iloc[-window:]
       pred MA.extend(oos pred)
```

### Forecasting the MA(2) series



MSE last value: 3.249

MSE MA(2): 1.948

### Differencing to obtain the series to the original scale

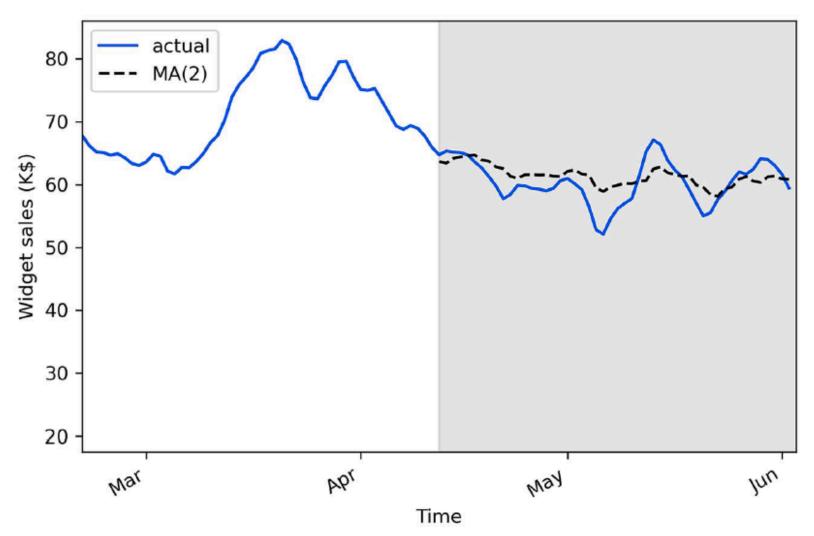
In order to reverse the first-order difference it is necessary to add an initial value  $y_0$  to the first differenced value  $y_1'$ 

$$y_1 = y_0 + y_1'$$

Then  $y_2$  can be obtained using a cumulative sum of the differenced values

$$y_2 = y_0 + y_1' + y_2'$$

### Inverse-transformed MA(2) series



MAE MA(2): 2.32

# AutoRegressive AR(p)

### Autoregressive model

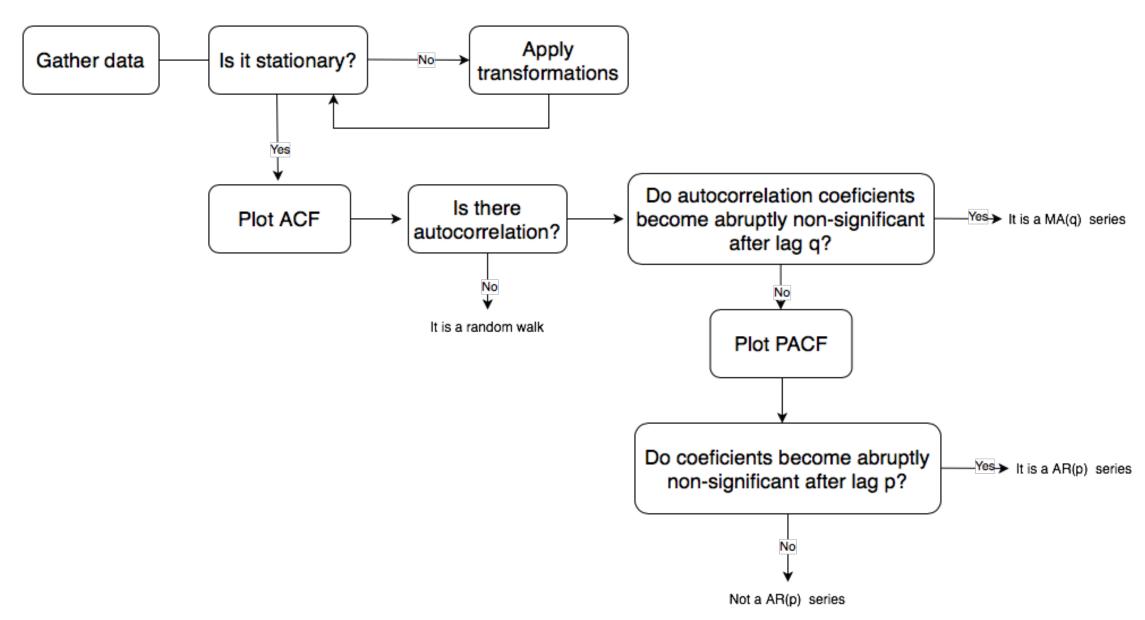
- An autoregressive model is a regression of a variable against itself. In a time series, this means that the present value is linearly dependent on its past values
- The autoregressive process is denoted as AR(p), where p is the order

The general expression of an AR(p) model is

$$y_t = C + \phi_1 y_{t-1} + \phi_2 y_{t-2} + ... + \phi_p y_{t-p} + \epsilon_t$$

Similar to the MA(q) series, the order p of an autoregressive process determines the number of past values that affect the present value.

### Identifying a Autoregressive series



### Partial Autocorrelation function (PACF)

In a second-order autoregressive series or AR(2)

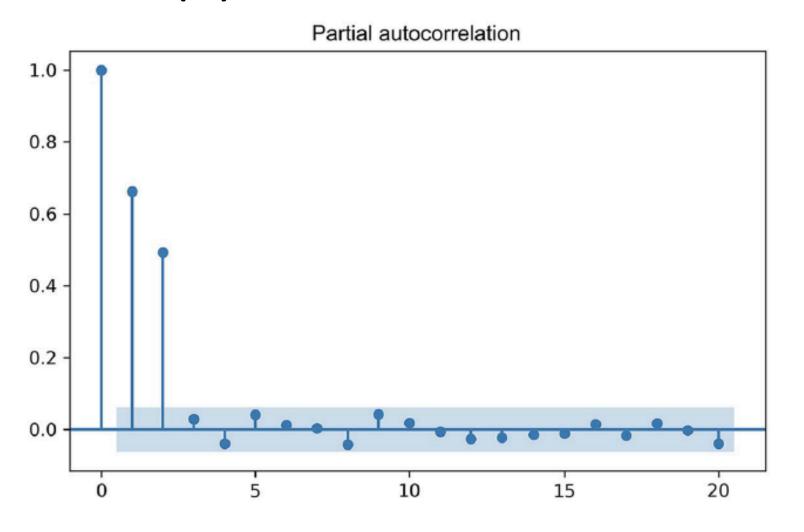
$$y_t = C + \phi_1 y_{t-1} + \phi_2 y_{t-2} + \epsilon_t$$

The autocorrelation between yt and yt-2 using the ACF does not take into account the fact that  $y_{t-1}$  has an influence on both  $y_t$  and  $y_{t-2}$ 

To do so, it is necessary to remove the effect of  $y_{t-1}$ . Thus, measuring the partial autocorrelation between  $y_t$  and  $y_{t-2}$ 

The partial autocorrelation function measures the correlation between lagged values in a time series when the influence of correlated lagged values in between are removed

### PACF plot of a AR(2) series



The partial autocorrelation function can be used to determine the order of a stationary AR(p) series - the coefficients will be non-significant after lag p

## AutoRegressive Moving Average ARMA(p,q)

### Autoregressive moving average series

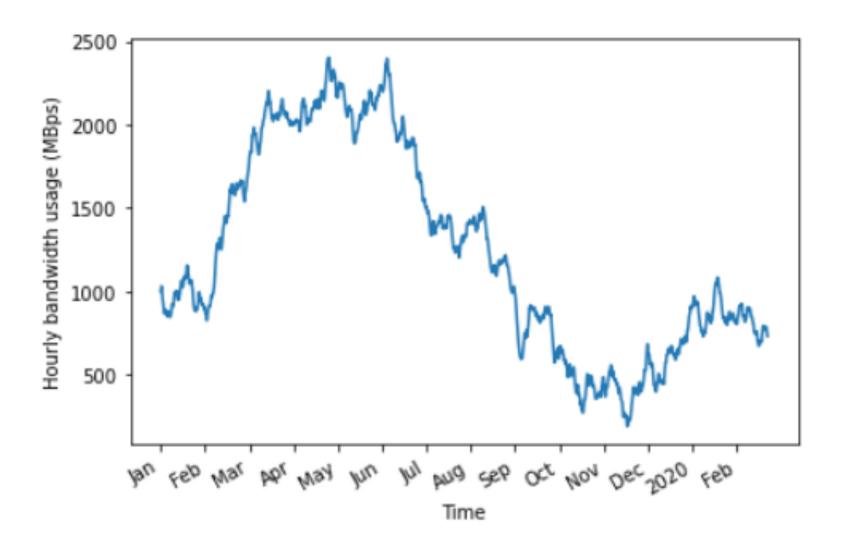
- An autoregressive moving average series is a combination of the autoregressive and the moving average series
- It is denoted as ARMA(p,q), where p is the order of the autoregressive process, and q is the order of the moving average process

The general equation of the ARMA(p,q)

$$y_t = C + \phi_1 y_{t-1} + \phi_2 y_{t-2} + \dots + \phi_p y_{t-p} + \mu + \epsilon_t + \theta_1 e_{t-1} + \theta_2 e_{t-2} + \dots + \theta_q e_{t-q}$$

ARMA(0,q)  $\approx$  MA(q), since the order p = 0 cancels the AR(p) portion ARMA(p,0)  $\approx$  AR(p), since the order q = 0 cancels the MA(q) portion

### Autoregressive moving average series example

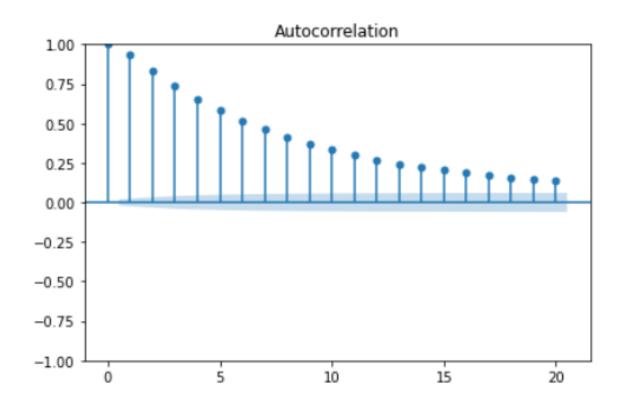


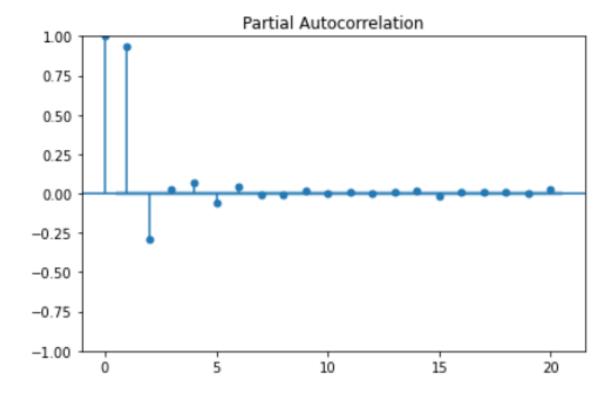
Bandwidth usage for a large data center in bits per second (bps)

### Identifying a stationary ARMA series

- If the series is a stationary ARMA(p,q) process both the ACF and PACF plots show a decaying or sinusoidal pattern
- The ACF and PACF plots cannot be used to determine the orders q and p of an ARMA(p,q) process
- The solution is to fit many ARMA(p,q) models with various combinations of values for p and q, then choosing a model using the **Akaike information** criteria

### ACF and PACF plots





### Akaike information criterion (AIC)

• Is a metric that aims to find a balance between a model's maximum likelihood and a model's simplicity

The AIC is a function of the number of parameters k in a model and the maximum value of the likelihood function  $\hat{L}$ :

$$AIC = 2k - 2\ln(\hat{L})$$

- The AIC quantifies the relative amount of information lost by the model
- The **better the model**, the **lower the AIC value** and the less information is lost

### Function to fit several ARMA(p,q) models

```
def optimize_ARMA(data, order_list) -> pd.DataFrame:
    results = []
    for order in order list:
        try:
            model = SARIMAX(data, order=(order[0], 0, order[1]), simple_differencing=False)
        except:
            continue
        aic = model.aic
        results.append([order, aic])
    result df = pd.DataFrame(results)
    result_df.columns = ['(p,q)', 'AIC']
    #Sort in ascending order, lower AIC is better
    result_df = result_df.sort_values(by='AIC', ascending=True).reset_index(drop=True)
    return result df
```

### Residual analysis

The residuals of a model are simply the difference between the predicted values and the actual values

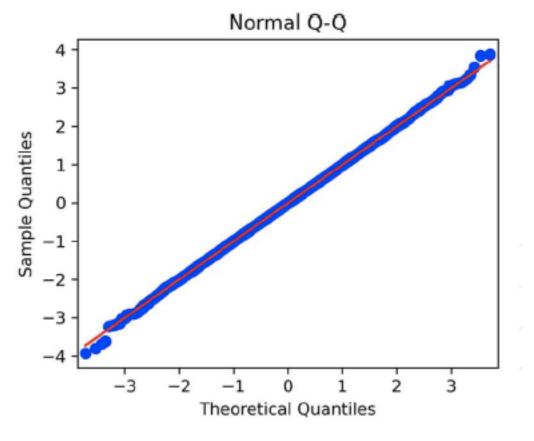
If the model has captured all predictive information from a dataset, the residuals of the model are **white noise**; there is only a random fluctuation left that cannot be modelled

To have a good model for making forecasts, the **residuals** must be **uncorrelated** and have a **normal distribution** 

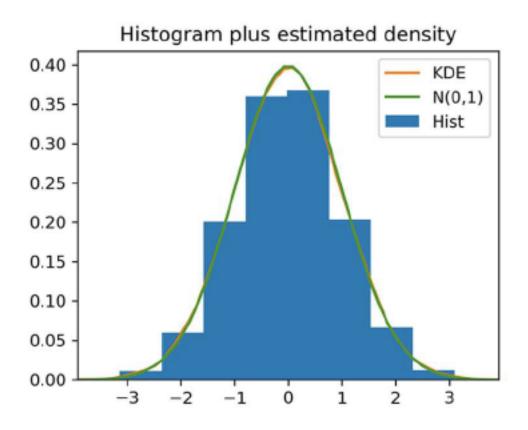
### Two residual analysis

- A qualitative analysis through the study of the quantile-quantile plot (Q-Q plot), for verifying if the model's residuals are normally distributed
- A quantitative analysis applying the *Ljung-Box test* to demonstrate that the residuals are uncorrelated

### Two residual analysis: quantile-quantile plot



Q-Q plot of the residuals



Histogram of the residuals

### Two residual analysis: Ljung-Box test

```
from statsmodels.stats.diagnostic import acorr_ljungbox
# run the Ljung-Box test on the residuals for the first 10 lags
residuals = model fit.resid
residuals test = acorr ljungbox(residuals, np.arange(1, 11, 1))
residuals_test['lb_pvalue'].describe()
count
         10.000000
       0.923847
mean
         0.057180
std
min
     0.811247
25%
         0.915579
50%
         0.942076
75%
         0.961415
          0.981019
max
Name: 1b pvalue, dtype: float64
```

All the returned p-values exceed 0.05, the residuals are uncorrelated

### Non-stationary time series

# AutoRegressive Integrated Moving Average ARIMA(p,d,q)

### Autoregressive integrated moving average series

- An autoregressive integrated moving average series is denoted as ARIMA(p,d,q), where p is the order of the AR(p) process, d is the order of integration, and q is the order of the MA(q) process
- Integration is the reverse of differencing, and the order of integration d is equal to the number of times the series has been differenced to be rendered stationary

The general equation of the ARIMA(p,d,q)

$$y'_{t} = C + \varphi_{1}y'_{t-1} + ... + \varphi_{p}y'_{t-p} + \theta_{1}\epsilon'_{t-1} + ... + \theta_{q}\epsilon'_{t-q} + \epsilon_{t}$$

### General modelling procedure for using the ARIMA(p,d,q)

- Gather data
- If data is not stationary
  - d = the minimum number of times the series must be differenced to become stationary
- Fit several combinations ARIMA(p, d, q)
- Select the model with lowest AIC
- Make the residual analysis of the model:
  - with Q-Q plot to evaluate whether the residuals are normally distributed
  - with Ljung-Box test to determine whether the residuals are correlated or not
- If the ARIMA(p,d,q) model has passed all the checks on the residual analysis it can be used to forecast the series

# Seasonal AutoRegressive Integrated Moving Average SARIMA(p,d,q)(P,D,Q)<sub>m</sub>

### Seasonal autoregressive integrated moving average series

- The SARIMA(p,d,q)(P,D,Q) $_{\rm m}$  model adds seasonal parameters to the ARIMA(p,d,q) model
- It has four new parameters: P, D, Q, and m
- The parameter m is the frequency
- P is the order of the seasonal AR(P) process, D is the seasonal order of integration, Q is the order of the seasonal MA(Q) process

A SARIMA(p,d,q)(0,0,0)m model is equivalent to an ARIMA(p,d,q) model

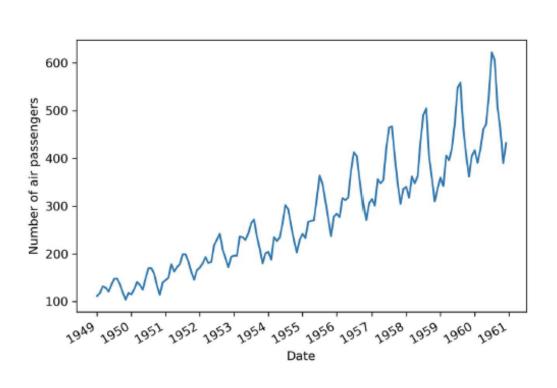
### Parameter m — Frequency

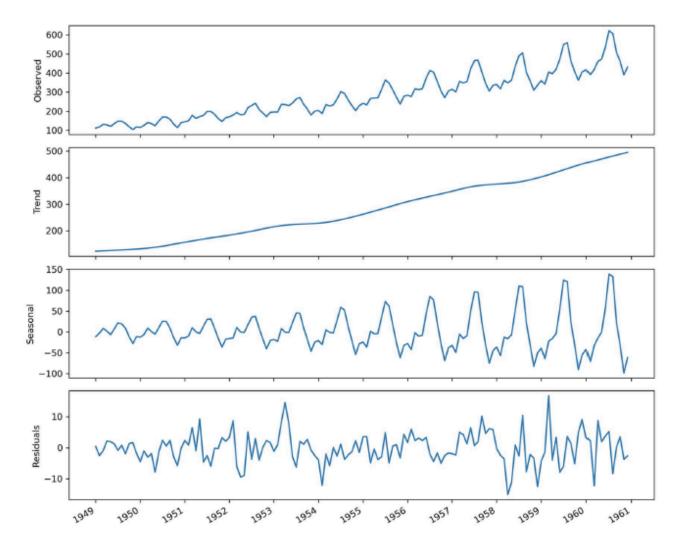
- The frequency is defined as the number of observations per cycle
- The length of the cycle will depend how the data was recorded

	Frequency	Frequency		
Data collection		day	week	year
Annual	1			
Quarterly	4			
Monthly	12			
Weekly	52			
Daily		1	7	365
Hourly		24	168	8766

### Identifying seasonal patterns in a time series

By plotting the time series or using time series decomposition





### Decomposition of a time series

The modelling of the decomposed components can be:

**Additive**: when the original time series can be reconstructed by adding all three components

$$y_t = T_t + S_t + R_t$$

**Multiplicative**: the time series can be reconstructed by multiplying all three components

$$y_t = T_t \times S_t \times R_t$$

### Four new parameters: P, D, Q, and m

#### Example where m = 12

- If P=2, this means that we will be include two past values of the series at a lag that is a multiple of m, the values at  $y_{t-12}$  and  $y_{t-24}$
- If D=1, this means that one seasonal difference makes the series stationary. The seasonal difference would be expressed as  $y'_t = y_t y_{t-12}$
- If Q=2, past error terms at lags that are a multiple of m will be included, the errors  $\epsilon_{t-12}$  and  $\epsilon_{t-24}$

### General modelling procedure for using the SARIMA

- Gather data
- Until the data is not stationary
  - make the difference of the data
  - make the seasonal difference of data
- When data is stationary d and D are defined
- Fit several combinations of p, q, P, D of SARIMA(p, d, q)(P,D,Q)<sub>m</sub>
- Select the model with lowest AIC
- Make the residual analysis of the model
- If the SARIMA(p,d,q)(P,D,Q) $_{\rm m}$  model has passed all the checks on the residual analysis it can be used to forecast the series

### SARIMAX(p,d,q)(P,D,Q)<sub>m</sub> X – exogenous variables

#### SARIMAX model

- Extends the SARIMA(p,d,q)(P,D,Q) $_{\rm m}$  model by adding the effect of exogenous variables
- The present value  $y_t$  is expressed as a SARIMA(p,d,q)(P,D,Q)<sub>m</sub> model to which there are added any number of exogenous variables  $X_t$

The general equation of the SARIMAX(p,d,q)(P,D,Q) $_{\rm m}$ 

$$y_t = SARIMA(p, d, q)(P, D, Q)_m + \sum_{i=1}^n \beta_i X_i^t$$

### SARIMAX model: exogenous variables

- The exogenous variables may or may not be good predictors
- However, it is not necessary to perform feature selection
- The SARIMAX model will attribute a coefficient close to 0 for exogenous variables that are not significant in predicting the target

### SARIMAX model: number of timesteps to forecast

- The **SARIMAX model** uses the SARIMA(p,d,q)(P,D,Q)<sub>m</sub> model and a linear combination of exogenous variables to predict **one timestep into the future**
- To predict more timesteps into the future the SARIMAX model requires to forecast the exogenous variables too
- There is no clear recommendation to predict only one timestep. It is dependent on the exogenous variables available:
  - If the exogenous variables can be accurately predicted, it is possible to forecast many timesteps into the future
  - if the forecast of exogenous variables have some error associated they can ampliated the prediction error of the target, as more timesteps are predicted into the future

### General modelling procedure for using the SARIMAX

- Gather data
- Until the data is not stationary
  - make the difference of the data
  - make the seasonal difference of data
- When data is stationary d and D are defined
- Fit several combinations of p, q, P, D of SARIMAX(p, d, q)(P,D,Q)<sub>m</sub>
- Select the model with lowest AIC
- Make the residual analysis of the model
- If the SARIMA(p,d,q)(P,D,Q) $_{\rm m}$  model has passed all the checks on the residual analysis it can be used to forecast the series

## Vector AutoRegression (VAR) Multivariate Forecasting

# Vector Autoregression model

- The vector autoregressive model VAR(p) is a generalization of the AR(p) model that allows the forecast of multiple time series
- In this model, each time series has an impact on the others, the past values of one time series affect the other time series, and vice versa
- The order p of the VAR(p) model determines how many lagged values impact the present value of the series

# Vector Autoregression model

For two time series, the general equation for the VAR(p) model is a linear combination of a vector of constants, past values of both time series, and a vector of error terms:

$$\begin{bmatrix} y_{1,t} \\ y_{2,t} \end{bmatrix} = \begin{bmatrix} C_1 \\ C_2 \end{bmatrix} + \begin{bmatrix} \phi_{1,1}^1 & \phi_{1,2}^1 \\ \phi_{2,1}^1 & \phi_{2,2}^1 \end{bmatrix} \begin{bmatrix} y_{1,t-1} \\ y_{2,t-1} \end{bmatrix} + \begin{bmatrix} \phi_{1,1}^2 & \phi_{1,2}^2 \\ \phi_{2,1}^2 & \phi_{2,2}^2 \end{bmatrix} \begin{bmatrix} y_{1,t-2} \\ y_{2,t-2} \end{bmatrix} + \dots + \begin{bmatrix} \phi_{1,1}^p & \phi_{1,2}^p \\ \phi_{2,1}^p & \phi_{2,2}^p \end{bmatrix} \begin{bmatrix} y_{1,t-p} \\ y_{2,t-p} \end{bmatrix} + \begin{bmatrix} \epsilon_{1,t} \\ \epsilon_{2,t} \end{bmatrix}$$

The time series must be stationary to apply the VAR model

## Granger causality test

- The VAR model can only be used if each of the time series are predictive of each other
- The **Granger causality test** determines whether one time series is predictive of another
- This statistical test helps to determine if past values of a time series  $y_{2,t}$  can help to forecast time series  $y_{1,t}$  => If true  $y_{2,t}$  Granger-causes  $y_{1,t}$
- The Granger causality test tests causality only in one direction
- In order for the VAR model to be valid, the test we must be repeated to verify that  $y_{1,t}$  also Granger causes  $y_{2,t}$

## Modelling procedure for the VAR(p) model

- Gather data
- If the series are not stationary
  - must be differentiated the number of times needed to become stationary
- Fit many VAR(p) models to select the one with the smallest AIC the p with the lowest AIC value is the order of VAR(p) model
- Make the Granger causality test for both series
  - If p-value less than 0.05, the null hypothesis can be rejected, meaning that the variables are Granger cause each other
- Make the residual analysis of the two series:
  - with **Q-Q plot** to evaluate whether the residuals are normally distributed
  - with Ljung-Box test to determine whether the residuals are correlated or not
- If the VAR(p) model passed the residual analysis it can be used to forecast the series

# Time series forecasting with Machine Learning

# When to use ML/DL for time series forecasting

- When statistical models take too much time to fit or when they result in correlated residuals that do not approximate white noise
- This can be due to the fact that data has multiple seasonal periods that cannot be considered in the model, or simply because there is a nonlinear relationship between the features and the target
- In those cases, machine learning/deep learning models can be used to capture this nonlinear relationship
- These models also have the added advantage of being faster to train

# Types of ML models for time series forecasting

#### Single-step model

Predicts one step into the future for one variable

#### Multi-step model

Predicts many steps into the future for one variable

#### Multi-output model

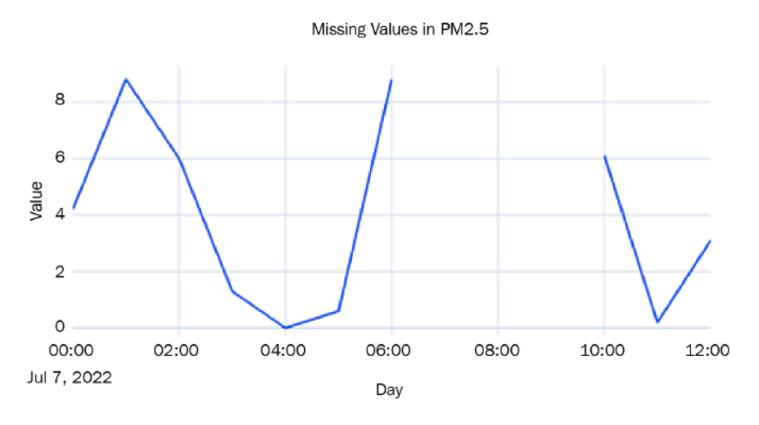
• Has more than one time dependent variable. Each variable not only depends on its past values but also has some dependency on other variables

### Transform time series into a two-dimensional format

- Explore the time series: trend, seasonality
- Define its frequency
- Treat missing values
- Split data into train/test
- Scaling data
- Put the time series in a tabular format (slide window)

## Handling missing values

Missing values are represented in the time series by a **sequence gap** in the time stamp variable or in other values



The first thing to do is to evaluate if the data in question is indeed missing or not

## Techniques to fill the missing values

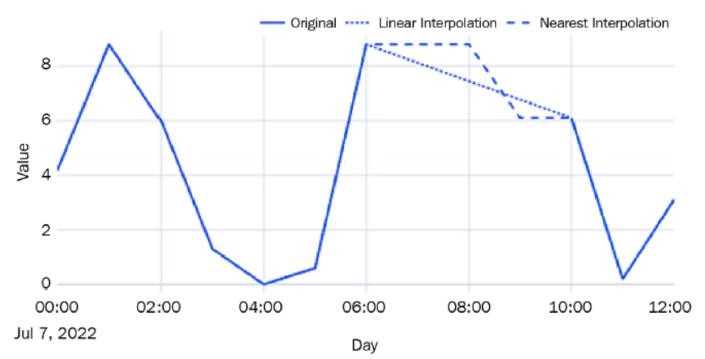
#### Simple techniques:

- Forward Fill: fill in the missing values with the last observed value until it finds the next observation
- Backward Fill: takes the next observation and backtracks to fill in all the missing values
- Mean Value Fill: fill in the missing values with the mean value of the entire series

## Techniques to fill the missing values

Imputation techniques covers interpolation:

- Linear Interpolation: draws a line between two observed points and fills the missing values so that they lie on this line
- Nearest Interpolation: is a combination of the forward and backward fill. For each missing value, the closest observed value is used to fill in

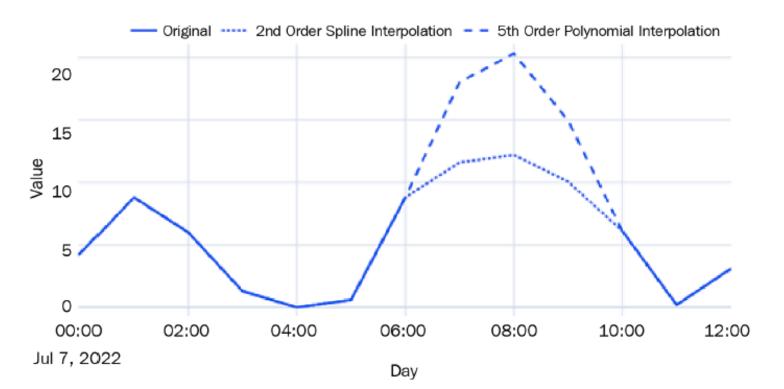


## Techniques to fill the missing values

Non-linear interpolation techniques:

• **Spline, Polynomial**: fit a spline/polynomial of a given order to the data and use that to fill in missing values

The higher the order, the more flexible the function is to fit the observed points



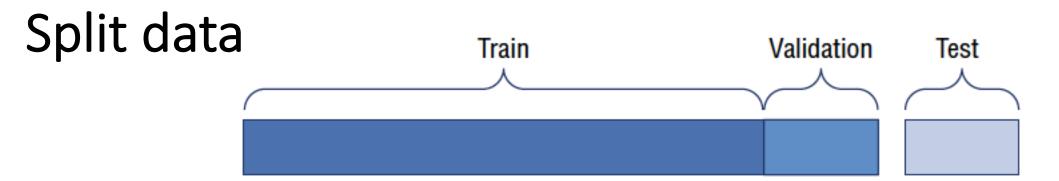
## Handling longer periods of missing data



- Imputing with the previous day
- Hourly average profile
- The hourly average for each weekday

#### **Seasonal interpolation**

- 1. Calculate the seasonal profile
- 2. Subtract the seasonal profile and apply any interpolation technique
- 3. Add the seasonal profile back to the interpolated series



#### Split by Time

Since time matters in time series data, data should be split based on time, rather than randomly

#### **Split Points**

- Training set: must be the earlier portion of the time series, 70-80% of the data
- Validation set: the next 10-15% of the data after the training set
- Test set: the remaining data, the last 15-20%

If data has any seasonality or specific patterns for example weekly patterns, the splits must respect the pattern

# Time Series Scaling/Normalizing

- Scaling
  - Min-Max
- Normalization
  - Sigmoid

## Time series data transformation

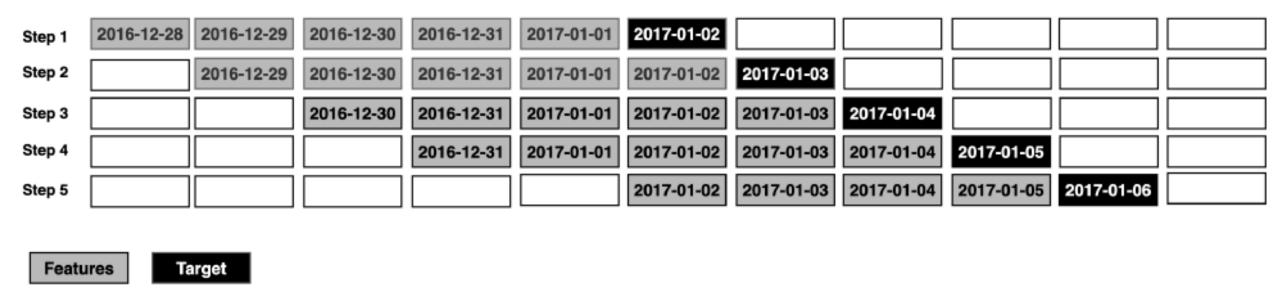
Time series to be processed by ML algorithms must be transformed to contain independent and dependent variables

For that, it is necessary to define a **sliding window** to convert the data into a window of inputs

For example, a **five-period sliding window** in a univariate time series, will produce five independent variables,  $x_1$ ,  $x_2$ ,  $x_3$ ,  $x_4$ ,  $x_5$ , which are lagged versions of the dependent variable, to be used to predict the current period  $x_6$ 

This representation of multiple inputs (a sequence) to one output is referred to as a single-step forecast

# Five-period sliding window for single-step forecast



This technique was described in Machine Learning Strategies for Time Series Forecasting, Lecture Notes in Business Information Processing. Springer Berlin Heidelberg (https://doi.org/10.1007/978-3-642-36318-4\_3)

## Multi-step forecast

- Predicts more than one step forecast into the future
- Forecast horizon is the number of time steps into the future to forecast at any point in time

#### Example:

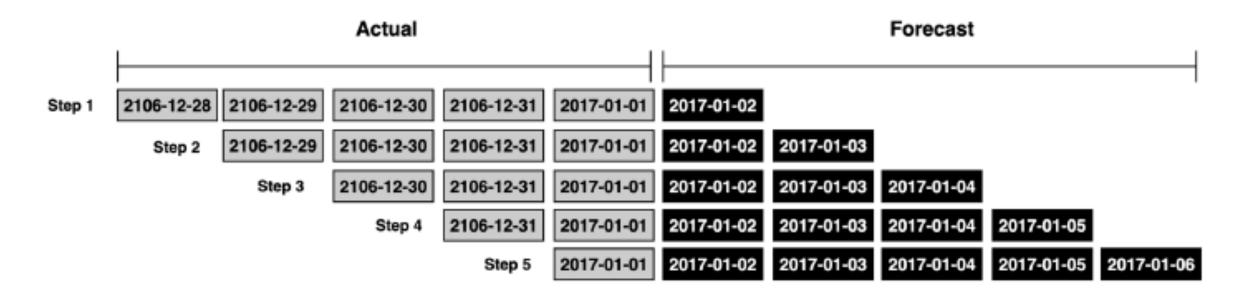
Prediction of three months into the future  $(x_{11}, x_{12}, x_{13})$  based on a sequence of 10 previous months  $(x_1, ..., x_{10})$ 

Four different ways to transform a series for multi-step forecast:

- Recursive strategy
- Direct strategy
- DirRec (Direct-Recursive) strategy

## Recursive multi-step forecast strategy

Is based on one-step forecasts that are reused (recursively) to make the next one-step prediction, and the process continues until all the future steps are produced



## Slide window

Window = 6 Horizon = 3

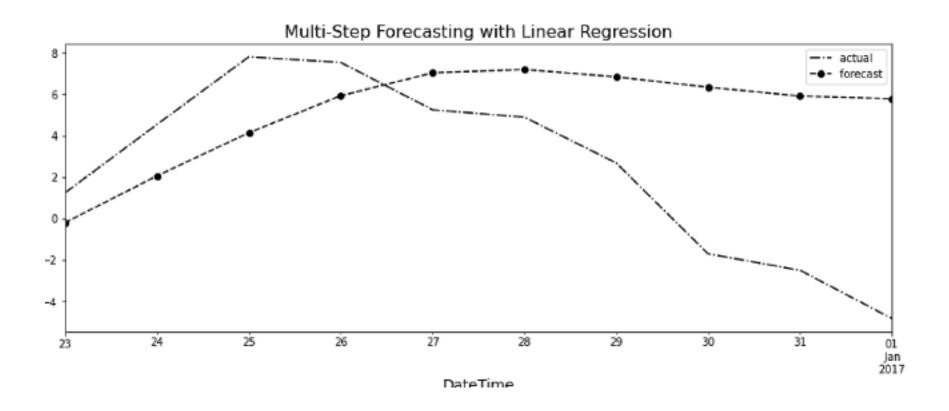
	date	temperature
0	2009-01-01	-6.810629
1	2009-01-02	-3.360486
2	2009-01-03	5.435694
3	2009-01-04	7.283889
4	2009-01-05	12.690069

	x1	x2	х3	x4	х5	х6	y1	y2	у3
date									
2009-01-01	-6.810629	-3.360486	5.435694	7.283889	12.690069	15.201597	20.121875	18.864792	21.289722
2009-01-02	-3.360486	5.435694	7.283889	12.690069	15.201597	20.121875	18.864792	21.289722	11.937847
2009-01-03	5.435694	7.283889	12.690069	15.201597	20.121875	18.864792	21.289722	11.937847	3.210903
2009-01-04	7.283889	12.690069	15.201597	20.121875	18.864792	21.289722	11.937847	3.210903	3.682431
2009-01-05	12.690069	15.201597	20.121875	18.864792	21.289722	11.937847	3.210903	3.682431	-1.678194

## Recursive multi-step forecast strategy

Some steps further, the recursive approach will only rely on the estimated values and any related estimation errors

The accumulation of these errors can turn the forecast highly biased



## Other multi-step forecast strategies

#### **Direct strategy**

- creates multiple independent models for each future step
- this independence turns impossible to identify relationships that would emerge between different predictions
- this strategy can suffer from high variance

#### **DirRec (Direct-Recursive) strategy**

• is a hybrid of direct and recursive strategies and a way to mitigate their individual shortcomings

## Multiple output model

- A multivariate time series has more than one time dependent variable
- Each variable not only depends on its past values but also has some dependency on other variables
- Example, forecasting the temperature and the wind speed

## Forecasting Time Series with Machine Learning Models

Regression Algorithms

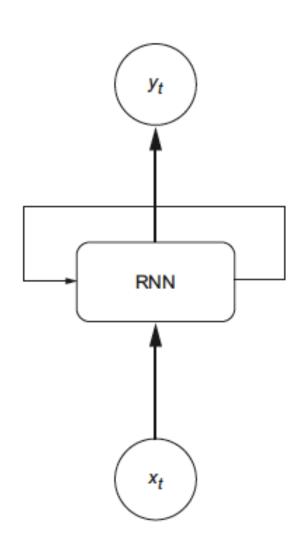
# Time series forecasting with Deep Learning

## Reasons to use Deep Learning in time series forecasting

Deep learning neural networks have three main intrinsic capabilities:

- are capable of automatically learning and extracting features from raw and imperfect data
- supports multiple inputs and outputs
- Recurrent neural networks are especially useful with sequential data because each neuron or unit can use its internal memory to maintain information about the previous input
  - —**long short-term memory** (LSTM) networks
  - —gated recurrent units (GRUs)
    - are good at extracting patterns in input data that span over relatively long sequences

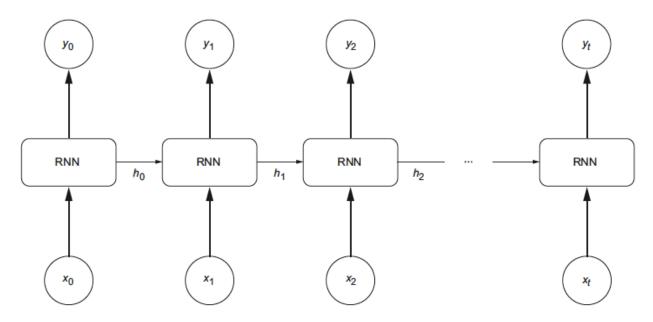
## Recurrent Neural Networks (RNNs)



- When an element of the sequence  $(x_t)$  it is fed to the RNN, it computes a hidden state  $(h_t)$
- This hidden state acts as memory. It is computed for each element of the sequence and fed back to the RNN as an input
- That way, the RNN uses past information computed for previous elements of the sequence to inform the output for the next element of the sequence

## Recurrent Neural Networks (RNNs)

• RNN units are described as recurrent units because the type of dependence of the current value on the previous event is recurrent and can be thought of as multiple copies of the same node, each transmitting a recurrent message to a successor



This is how it replicates the concept of memory

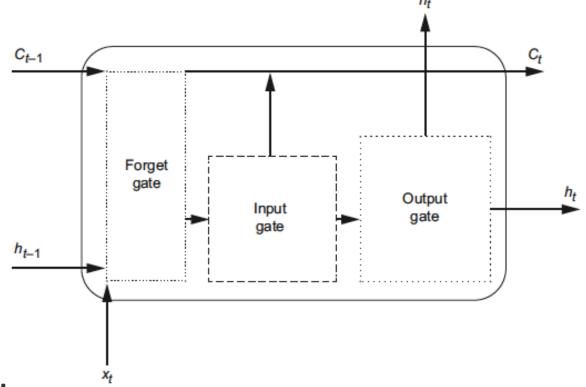
## RNN vanishing gradient problem

- RNNs suffer from **short-term memory** information from an early element in the sequence will stop having an impact further into the sequence
- During backpropagation RNNs suffer from short-term memory due to the vanishing gradient problem the change in gradient becomes very small, sometimes close to 0 and the weights of the network do not get updated, and the network stops learning

## Long Short-term Memory (LSTM)

- LSTMs are a particular type of RNN that addresses the short-term memory problem
- LSTMs introduce additional elements in the RNN architecture, called gates and cell state
- The output of the LSTM network is modulated by the state cells that act as long-term or short-term memory cells these state cells allows for past information to flow through the network for a longer period of time
- Instead of mapping inputs to outputs alone, the network can learn a mapping function of inputs over time to an output

## LSTM gates



The LSTM has three gates:

- forget gate determines what information from past steps is still relevant
- input gate determines what information from the current step is relevant to update the network's memory
- output gate determines what information is passed on to the next element of the sequence or as a result to the output layer

# Gated Recurrent Unit (GRU)

- GRU can be considered as a variation on the LSTM because both are designed similarly
- Unlike LSTM, GRU networks contain only two gates and do not maintain an internal cell state
- The information that is stored in the internal cell state in an LSTM is incorporated into one of the gates
- A GRU can be trained to keep information from long ago, without removing it through time, or delete information that is irrelevant to the prediction

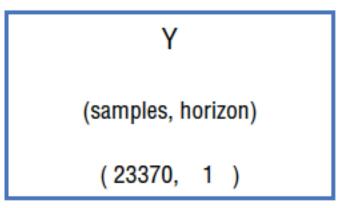
## **GRU** gates

- **Update gate** helps the network to determine how much of the past information (from previous time steps) needs to be passed along to the future
- Reset gate decides how much past information to forget
- cell state is incorporated into the reset gate
  - is used to initiate some nonlinearity into the input
  - is also used to decrease the impact that previous information has on the current information that is being passed into the future

### Transform time series into a three-dimensional structure

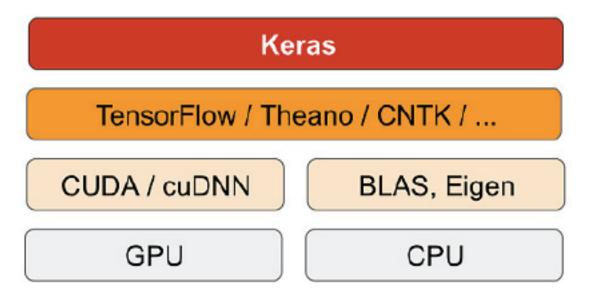
- 1. Index the data on time stamp for time-based filtering
- 2. Treat missing values
- 3. Split data into train/validation/test
- 4. Scaling data
- 5. Put the time series in a tabular format (slide window)
- 6. Transform time series data into two tensors

```
X
(samples, time steps, features)
( 23370, 6 , 1 )
```



## Introduction to Keras

- Keras is a deep-learning framework for Python that provides a convenient way to define and train almost any kind of deep-learning model
- Keras handles three backend engines that can be plugged seamlessly into Keras:
   TensorFlow, Theano, and Microsoft Cognitive Toolkit



## Development of deep learning models in Keras

#### Define the model

Create a Sequential model and add configured layers

#### 2. Compile the model

Specify the loss function and optimizers call the compile() function on the model

#### 3. Fit the model

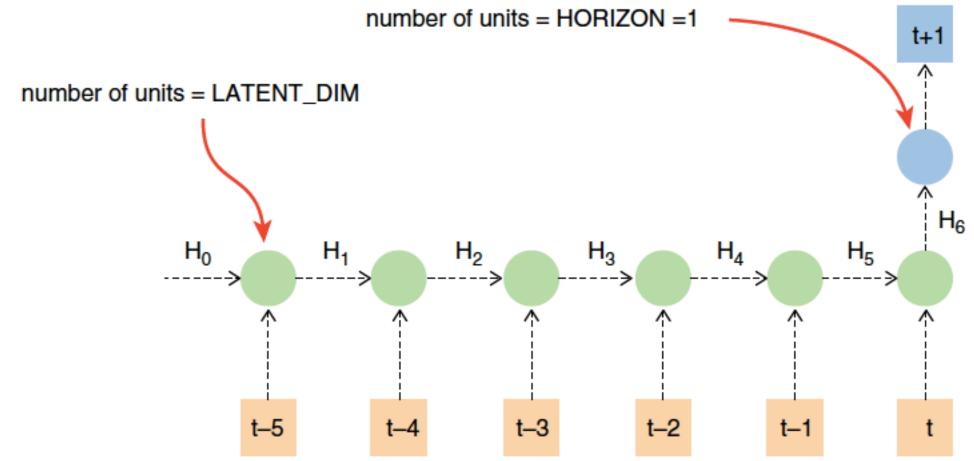
Train the model on a sample of data by calling the fit() function

#### 4. Make predictions

Use the model to generate predictions on new data by calling the predict() and evaluate() functions

## **Univariate Models**

In univariate models we have a single series of observations, and a model is required to learn from the series of past observations to predict the next value in the sequence



## Implementing a GRU as a single-step model

#### There must be defined:

- the latent dimension (LATENT\_DIM): is number of units in the RNN layer
- the batch size (BATCH\_SIZE): is the number of samples per mini-batch
- the **epochs** (EPOCHS): is the maximum number of times the training algorithm will cycle through all samples

# Implementing a GRU as a single-step model

```
HORIZON = 1
LATENT DIM = 5
BATCH SIZE = 32
EPOCHS = 10
model = Sequential()
model.add(GRU(LATENT DIM, input shape=(T, 1)))
model.add(Dense(HORIZON))
model.compile(optimizer='RMSprop', loss='mse')
model.summary()
GRU earlystop = EarlyStopping(monitor='val loss',
                               min delta=0,
                               patience=5)
```

## Implementing a GRU as a single-step model

```
model history = model.fit(X train,
                           y train,
                           batch size=BATCH SIZE,
                           epochs=EPOCHS,
                           validation data=(X valid, y valid),
                           callbacks=[GRU earlystop],
                           verbose=1)
y pred = model.predict(X test)
y pred
```

## Multivariate Models

In multivariate time series there is more than one observation for each time step

