# Introduction to Machine Learning. Lec.5 Support Vector Regression

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## Regression is...

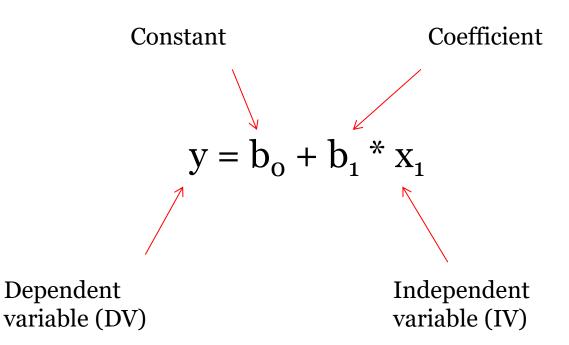
• a technique for determining the statistical **relationship between** two or more variables where a change in a dependent variable is associated with, and depends on, a change in one or more independent variables.

http://www.businessdictionary.com/definition/regression.html

## Types of regression models

- Simple Linear Regression
- Multiple Linear Regression
- Polynomial Regression
- Support Vector Regression (SVR)
- Decision Tree Regression
- Random Forest Regression

## SLR. Formula



## MLR. Formula

$$y = b_0 + b_1 x_1 + b_2 x_2 + b_3 x_3 + ... + b_n x_n$$

## PR. Formula

$$y = b_0 + b_1 x_1 + b_2 x_1^2 + b_3 x_1^3 + ... + b_n x_1^n$$

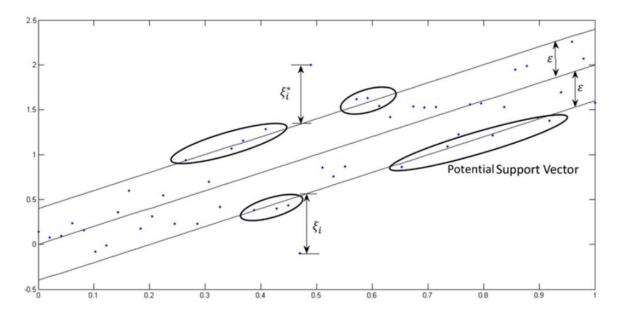
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- As in LR models (i.g. SLR), SVR has kind of a best fitting line. But other than this, it has parallels (supporting vectors).
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- SVR has a different goal compared to LR. In LR we are trying to minimize the error between the prediction and data. In SVR our goal is to make sure that errors **do not exceed the threshold**
- This threshold is controlled by hyper parameter Epsilon ( $\varepsilon$ , also called as width of a street)

## SVR. Regression vs. Classification

• The regression problem is a generalization of the classification problem, in which the model returns a continuous-valued output, as opposed to an output from a finite set. In other words, a regression model estimates a continuous-valued multivariate function.

## **SVM**

- SVMs solve binary classification problems by formulating them as convex optimization problems
- The optimization problem entails finding the maximum margin separating the hyperplane, while correctly classifying as many training points as possible
- SVMs represent this optimal hyperplane with support vectors

## SVM -> SVR

- SVM generalization to SVR is accomplished by introducing an  $\varepsilon$ -insensitive region around the function, called the  $\varepsilon$ -tube
- This tube reformulates the optimization problem to find the tube that best approximates the continuous-valued function, while balancing model complexity and prediction error

## SVM -> SVR

More specifically, SVR is formulated

- as an optimization problem by first defining a convex  $\varepsilon$  -insensitive loss function to be minimized
- and finding the flattest tube that contains most of the training instances.

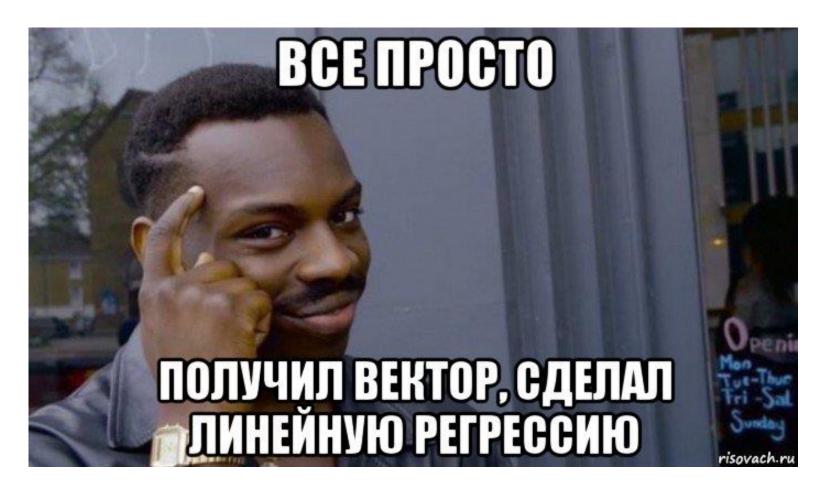
## What SVR does?

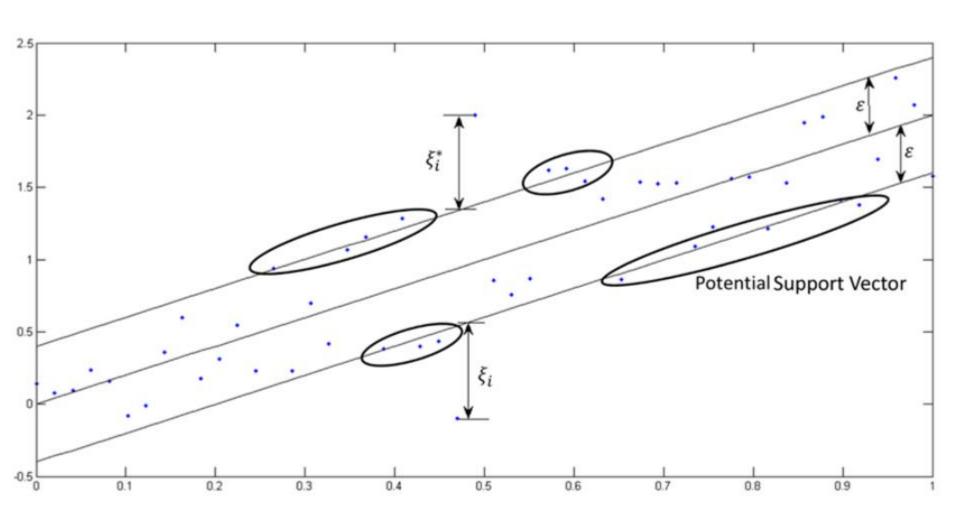
- SVR performs linear regression in a higher dimensional space
- We can think of SVR as if each data point in the training set represents it's own dimension.
- When you evaluate your kernel between a test point and a point in the training set the resulting value gives you the coordinate of your test point in that dimension

## What SVR does?

- The vector we get when we evaluate the test point for all points in the training set,  $\vec{k}$  is the representation of the test point in the higher dimensional space
- Once you have that vector you then use it to perform a linear regression

## What SVR does?





## General rule

- It requires a training set :  $\tau = \{\overrightarrow{X}, \overrightarrow{Y}\}$  which covers the domain of interest and is accompanied by solutions on that domain
- The work of the SVM is to approximate the function we used to generate the training set

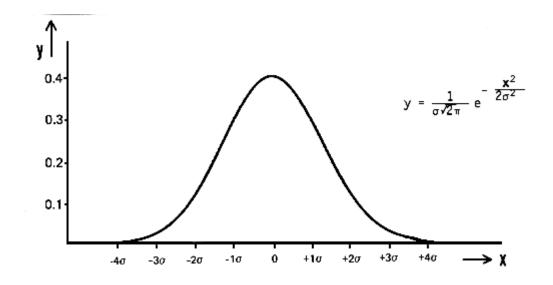
$$F(X) = \vec{Y}$$

# The algorithm of SVR. Step 1

- 1. Collect a training set  $\tau = \{\overrightarrow{X}, \overrightarrow{Y}\}$
- 2. Choose a kernel and it's parameters as well as any regularization needed
- 3. Form the correlation matrix,  $\vec{K}$
- 4. Train your machine, exactly or approximately, to get contraction coefficients  $\vec{\alpha} = \{\alpha_i\}$
- 5. Use those coefficients, create your estimator  $f(\vec{X}, \vec{\alpha}, x^*) = y^*$

# The algorithm of SVR. Step 2

- 1. Choose a kernel: Gaussian
- 2. Choose regularization: Noise



## Correlation matrix

$$K_{i,j} = \exp\left(\sum_{k} \theta_{k} |x_{k}^{i} - x_{k}^{j}|^{2}\right) + \epsilon \delta_{i,j}$$

At this time we have a main part of the algorithm

$$\overline{K}\vec{\alpha} = \vec{y}$$

- $\vec{y}$  is the vector of values corresponding the your training set
- $\overline{K}$  is you correlation matrix
- $\vec{\alpha}$  is a set of unknowns we need to solve for.

$$\vec{\alpha} = \overline{K}^{-1} \vec{y}$$

- Once  $\vec{\alpha}$  parameters are known form the estimator
- We use the coefficients we found during the optimization step and the kernel we started with
- To estimate the value  $y^*$  for a test point,  $\overrightarrow{x^*}$  compute the correlation vector  $\overrightarrow{k}$
- $y^* = \vec{\alpha} \cdot \vec{k}$

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- Finally, we compute the element of k

$$k_i = \exp\left(\sum_k \theta_k |x_k^i - x_k^{\star}|^2\right)$$