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# Support Vector Machines for Regression and Novelty Detection

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#### **Abstract**

This paper evaluates three regression models. We first construct a linear ridge regression model that makes a prediction y on some matrix X. We then compare this model to an  $\epsilon$  SVR and a  $\nu$  SVR.

#### **Index Terms**

Ridge regression, SVM, SVR,  $\epsilon$  SVR,  $\nu$  SVR.

#### I. Introduction

**R** EGRESSION refers to the estimation of random variables using a set of observations. If we exist inside a space  $\mathbb{R}^n$ , we can approximate a distribution of points-provided our approximation follows the same distribution as the original. For example, if we have a distribution we know is gaussian, and the distribution follows what appears to be a linear path in  $\mathbb{R}^2$ , we can approximate that path with a linear function. If we consider a distribution with an infinitely large margin, we can apply the SVM classifier where the points act practically as the slack variables  $\xi_i$ . Vapnik defines this process using an alternate form of the representer theorem, sometimes referred to as the epsilon-insensitive loss function [1].

## II. THE DATA FOR OUR EXPERIMENTS

Our data is represented as a matrix in CSV format. The first column of the file represents our testing regressors without noise; likewise, the second column represents our training regressors—also without noise. Our matrices  $\underline{\underline{\mathbf{X}}}_{test}$  and  $\underline{\underline{\mathbf{X}}}_{train}$  come from columns 3-21 and 22-40 respectively. Our vectors  $\underline{\underline{\mathbf{y}}}_{test}$  and  $\underline{\underline{\mathbf{y}}}_{train}$  come from columns 41 and 42, respectively. For each experiment, we split the data 80:20 to validate that our predictions follow approximately the same distribution at the original.

### III. COMPARING REGRESSION ALGORITHMS

The following three experiments were conducted in Python. For each experiment, the program reads from a CSV using the Pandas library, performs regression, and examines the error by comparing the results of the predictions with the actual results from the CSV.

### A. Linear Ridge Regression

This first model uses the methods of *Tikhonov Regularization* to apply a regression using MMSE. First, we define a vector we call  $\gamma$  that contains 1000 logarithmically spaced elements, such that:

$$\gamma_i \in [-1, 2] \tag{1}$$

For each value of  $\gamma$ , we perform ridge regression and calculate the MMSE. A plot of the validation square error as a function of  $\gamma$  is shown below.

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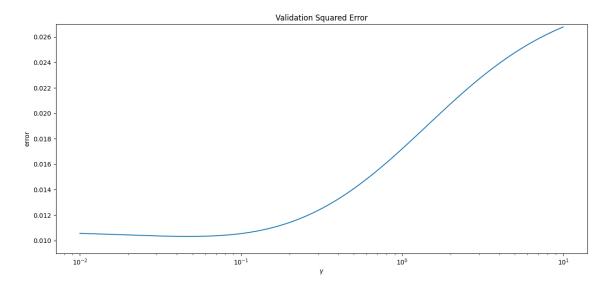


Fig. 1. Validation square error for linear ridge regression as a function of  $\gamma$ 

After choosing the optimal value for  $\gamma$  (in our case,  $\gamma \approx 0.047$ ), we apply ridge regression once more, and compare our predictions in this model with the actual value stored in the CSV.

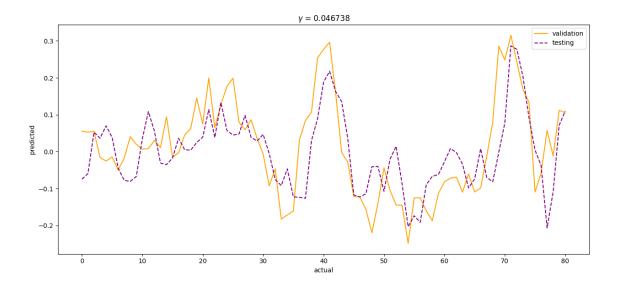


Fig. 2. Linear ridge regression distribution using  $\gamma$  = 0.46738

# B. The $\epsilon$ SVR

The  $\epsilon$  SVR is similar to linear ridge regression; however, this method uses cross-validation to determine an optimal value for  $\epsilon$ -a free parameter in the SVR that represents a positive error tolerance. For this model, we represent  $\epsilon$  and C as vectors of 100 logarithmically spaced elements such that:

$$\epsilon_i \in [-4, 1] \tag{2}$$

$$C_i \in [-2, 2] \tag{3}$$

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After calculating the mean square error for all the possible combinations of C and  $\epsilon$ , we pass the results into a contour plot to better visualize the gradient.

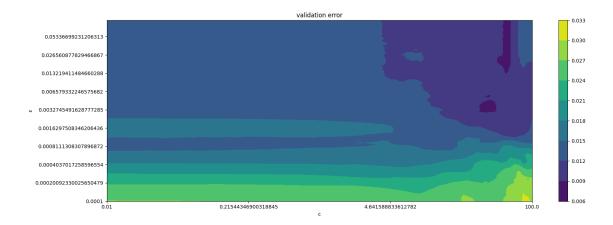


Fig. 3. MMSE for the  $\epsilon$  SVR

We find that our minimum value occurs when  $\epsilon \approx 0.498$  and  $C \approx 1.050$ . After setting these values as static and plotting the training data and the testing data, we see the following results.

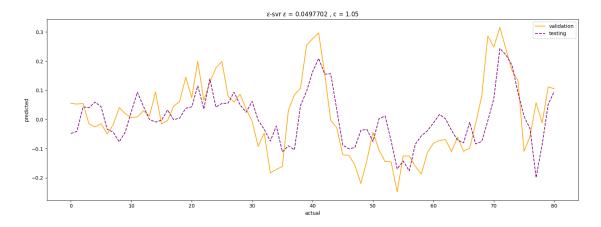


Fig. 4.  $\epsilon$  SVR distribution using  $\epsilon$  = 0.498 and C = 1.05

# C. The $\nu$ SVR

The  $\nu$ -SVR is similar to the  $\epsilon$ -SVR; however, the  $\nu$ -SVR tunes  $\epsilon$  by applying the method described by Smola and Scholkopf [1]. For the  $\nu$ -SVR, we represent C and  $\nu$  as 100 logarithmically spaced vectors, such that:

$$\nu_i \in [-3, 0] \tag{4}$$

$$C_i \in [-2, 2] \tag{5}$$

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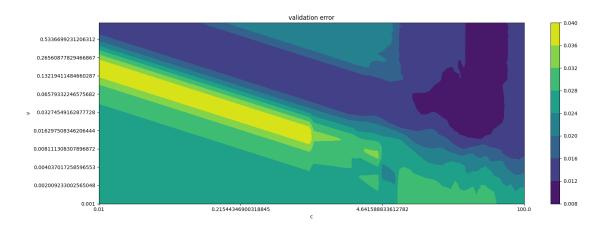


Fig. 5. MMSE for the  $\nu$  SVR

Similar to the method in the  $\epsilon$ -SVR, we use a contour plot of the values for  $\nu$  and C so that we can find their optimal values using inspection. For our training data, we find an optimal value where  $\nu \approx 0.572$  and  $C \approx 1.05$ . We plot the regressors again below.

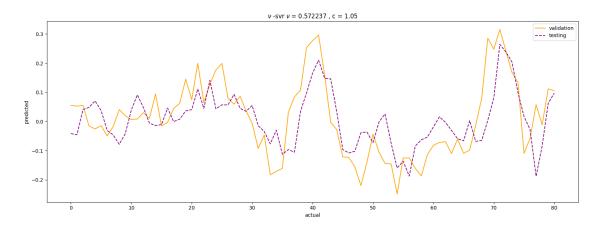


Fig. 6.  $\nu$  SVR distribution using  $\nu = 0.572$  and C = 1.05

## IV. CONCLUSION

From the following experiments we see that for this particular set of data, the three SVR algorithms produce similar regressors; however, the  $\nu$ -SVR and the  $\epsilon$ -SVR differ in the number of support vectors and complexity of the learning task.

## V. SOURCE CODE

https://github.com/keithhbova/support\_vector\_machines/

### REFERENCES

[1] A Tutorial on Support Vector Regression. Alex J. Smola and Bernhard Scholkopf [Online]. Available: https://alex.smola.org/papers/2003/SmoSch03b.pdf