

COMPSCI 589

Lecture 10: Support Vector and Neural Network Regression

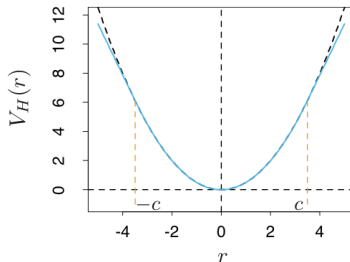
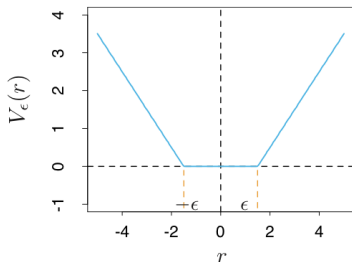
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Support Vector Regression

- Support Vector Regression (SVR) is the generalization of SVMs to the case of regression.
- As with SVMs, SVR is a linear regression model trained using a different objective function. In this case, the *epsilon insensitive loss*.



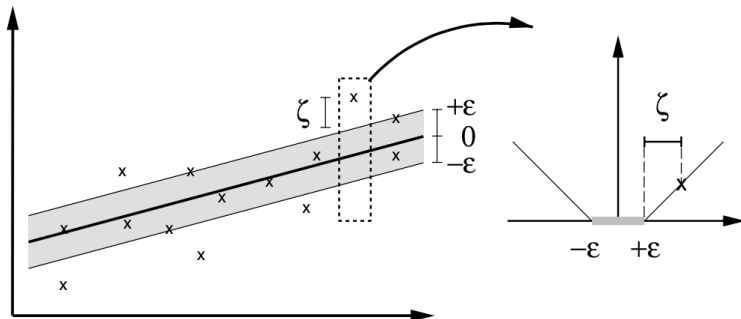
Support Vector Regression

$$f_{SVR}(\mathbf{x}) = \left(\sum_{d=1}^D w_d x_d \right) + b = \mathbf{x}\mathbf{w} + b$$

$$\mathbf{w}^*, b^* = \arg \min_{\mathbf{w}, b} C \sum_{i=1}^N V_{\epsilon}(y_i - \mathbf{x}_i \mathbf{w} - b) + \|\mathbf{w}\|_2^2$$

$$V_{\epsilon}(r) = \begin{cases} 0 & \dots \text{ if } |r| < \epsilon \\ |r| - \epsilon & \dots \text{ otherwise} \end{cases}$$

Support Vector Regression



Kernelization

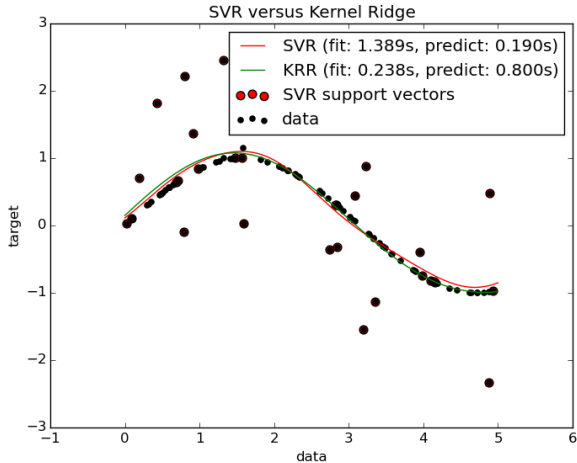
Using the same representer theorem used in classification, it can be shown that

$$f_{SVR}(\mathbf{x}) = \mathbf{x}\mathbf{w}^* + b^* = \sum_{i=1}^N \alpha_i \langle \mathbf{x}, \mathbf{x}_i \rangle + \sum_{i=1}^N \alpha_i \langle \mathbf{1}, \mathbf{x}_i \rangle$$

This can again be generalized using kernels to allow for non-linear models:

$$f_{SVR}(\mathbf{x}) = \mathbf{x}\mathbf{w}^* + b^* = \sum_{i=1}^N \alpha_i K(\mathbf{x}, \mathbf{x}_i)$$

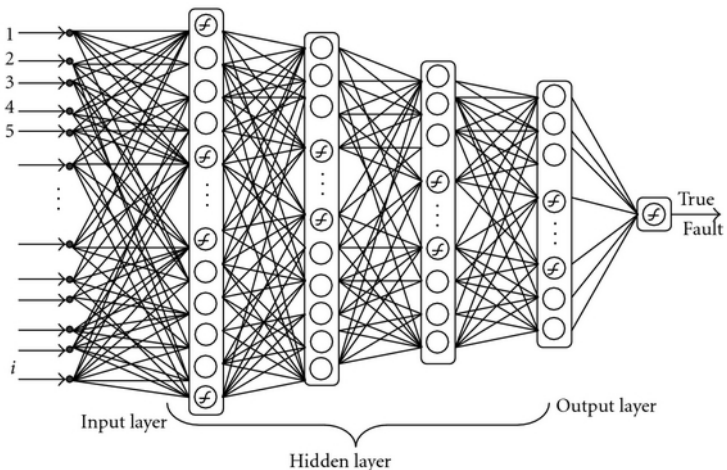
SVR vs KRR



Trade-Offs

- SVR is more robust to outliers than OLS (minimizing the MSE) due to the loss being linear in the tails instead of quadratic. This is related to a long line of work on robust regression.
- Kernel SVR is low bias and has good capacity control, but use of cross validation to select regularization hyperparameters is critical.
- The learning problem is convex for any choice of regularization parameters and thus has a unique global optimum.
- The kernel matrix computation is quadratic in the data dimension, but the model has a support vector property.
- You need to know what kernel to use or you need to use some form of validation to select from among several alternatives.

Multi-Layer Perceptron



Neural Network Regression

To convert an MLP from classification to regression, we only need to change the output activation function from logistic to linear.

- The hidden layer non-linearities are smooth functions:

$$h_k^1 = \frac{1}{1 + \exp(-(\sum_d w_{dk}^1 x_d + b_{dk}^1))}$$
$$h_k^i = \frac{1}{1 + \exp(-(\sum_l w_{lk}^i h_l^{(i-1)} + b_{lk}^i))} \text{ for } i = 2, \dots, L$$

- The output layer activation function is a linear function:

$$\hat{y} = \sum_l w_l^o h_l^L + b^o$$

Learning

Let θ be the complete collection of parameters defining a neural network model. Our goal is to find the value of θ that minimizes the MSE on the training data set $\mathcal{D} = \{y_i, \mathbf{x}_i\}_{i=1:N}$

$$\mathcal{L}_{MSE}(\mathcal{D}|\theta) = \frac{1}{N} \sum_{n=1}^N (y_n - \hat{y}_n)^2$$

Gradients

We need the gradient with respect to each of the parameters. Let's begin with w_l^o :

$$\frac{\partial \mathcal{L}_{MSE}(\mathcal{D}|\theta)}{\partial w_l^o} = \frac{\partial}{\partial w_l^o} \frac{1}{N} \sum_{n=1}^N (y_n - \hat{y}_n)^2 = 0 \quad (1)$$

$$= \frac{2}{N} \sum_{n=1}^N (y_n - \hat{y}_n) \frac{\partial \hat{y}_n}{\partial w_l^o} \quad (2)$$

$$= \frac{2}{N} \sum_{n=1}^N (y_n - \hat{y}_n) h_l^L \quad (3)$$

Gradients

It's also useful to define the derivatives wrt the hidden units for a single data case:

$$\epsilon_k^L = \frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial h_k^L} = \frac{\partial}{\partial h_k^L} (y - \hat{y})^2 \quad (4)$$

$$= 2(y - \hat{y}) \frac{\partial \hat{y}}{\partial h_k^L} \quad (5)$$

$$= 2(y - \hat{y}) w_k^o \quad (6)$$

In general, we can define: $\epsilon_k^j = \frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial h_k^j}$

Gradients

Suppose we're trying to compute the derivative with respect to the weight w_{kl}^j for some layer j and assume we have ϵ_l^j computed for all hidden units l in layer j .

$$\frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial w_{kl}^j} = \frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial h_l^j} \frac{\partial h_l^j}{\partial w_{kl}^j} \quad (7)$$

$$= \epsilon_l^j h_l^j (1 - h_l^j) h_k^{j-1} \quad (8)$$

The total derivative is then given by:

$$\frac{\partial \mathcal{L}_{MSE}(\mathcal{D}|\theta)}{\partial w_{kl}^j} = \frac{1}{N} \sum_{n=1}^N \frac{\partial \mathcal{L}_{MSE}(y_n, \mathbf{x}_n|\theta)}{\partial w_{kl}^j}$$

Gradients

Suppose we're trying to compute the error with respect to hidden unit k in layer $j - 1$ and assume we have ϵ_l^j computed for all hidden units l in layer j .

$$\frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial h_k^{j-1}} = \sum_l \frac{\partial \mathcal{L}_{MSE}(y, \mathbf{x}|\theta)}{\partial h_l^j} \frac{\partial h_l^j}{\partial h_k^{j-1}} \quad (9)$$

$$= \sum_l \epsilon_l^j h_l^j (1 - h_l^j) w_{kl}^{j-1} \quad (10)$$

Backpropagation

- The Backpropagation algorithm works by making a forward pass through the network for each data case and storing all the hidden unit values.
- The algorithm then computes the error at the output and makes a backward pass through the network computing the derivatives with respect to the parameters as well as the contribution of each hidden unit to the error. These are the ϵ_k^j values.
- The complete computation is just an application of the chain rule with caching of intermediate terms in the neural network graph structure.

Trade-Offs

- Neural network regression has low bias, but high variance.
- The objective function has local optima and requires iterative numerical optimization using backpropagation to compute the gradients, which can be slow.
- Making predictions with trained models can be very fast.
- Capacity control in these models can be crucial. The capacity parameters are the depth of the network and the size of each layer.
- These models can also be trained using ℓ_2 or ℓ_1 regularization or the more recent dropout scheme as an alternative to controlling network structure.