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- (a) I apply SVM model to this classification problem.
 - 1) Input: Since each image is a 5x5 grid of black or white cells, I represent it as a 5x5 matrix with elements O (white) and I (black). However, the input of SVM must be a vector, so I flatten the matrix into a 1x25 vector.
 - 2) Output: According to the problem, I need to classify an image into class A or class B. So, the output is a value that I can use to classify the image.
 - 3) Dotaset: The dataset for training is all the images of class A and class B. The size of it is 10, including 5 images of A and 5 images of B. Since SVM is a supervised learning method and this is a binary classification problem, I label each image of class A as "1" and each image of B as "-1". Therefore, the inputs of training process is a list of 10 tuples. The first element of each tuple is a 1×25 vector representation of an image and the second element is the label (101-1) of the image. The Test data set is 5 unlabeled images. Each image is to be classified after the training process.
 - 4) Loss function: I choose the quadratic loss function.

 $L_{045} = \sum_{i=1}^{2} (y_i - \phi(\underline{x}^i))^2$

The reason I choose it is I want to consider the cost of classification errors instead of the size of margin. In other words, for class A(y=1), I want $\phi(x) > 0$ and for class B (y=-1), I want $\phi(x) < 0$. Thus,

I use $(y_i - \phi(x^i))^2$ to measure the error.

But how to represent $\phi(\underline{x}^i)$? I use $\hat{\phi}(\underline{x}^i) = d_1 k(\underline{x}^i, \underline{x}^i) + \cdots + d_n k(\underline{x}^i, \underline{x}^n)$ to represent it. The reason is as follows.

Assume $\phi(x) \in V$ and $\mathcal{T} = \{d_1 k(x, x') + d_2 k(x, x^2) + \cdots + d_n k(x, x^n)\}$ is a subspace of V. $\phi(x) \in \mathcal{T}$, so $\phi(x)$ can be represented as:

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$$\widehat{\phi}(\underline{x}) = \lambda_1 K(\underline{x}, \underline{x}') + \lambda_2 k(\underline{x}, \underline{x}^2) + \cdots + \lambda_n k(\underline{x}, \underline{x}^n)$$

It's easy to know that when $\beta(x)$ is the map of $\beta(x)$ the Ω , $\beta(x)$ is the best representation of $\beta(x)$.

$$\langle (X) - \beta(X) \rangle + \langle (X, X_i) \rangle = 0$$

$$\Rightarrow \langle \phi(\underline{x}), k(\underline{x}, \underline{x}^{i}) \rangle = \langle \phi(\underline{x}), k(\underline{x}, \underline{x}^{i}) \rangle$$

$$\Rightarrow \phi(\underline{X}^{i}) = \widehat{\phi}(\underline{X}^{i}) = \sum_{j=1}^{n} d_{j} K(\underline{X}^{i}, \underline{X}^{j})$$

Then the goal is to minimize the loss function as follows.

Notice that the optimization is without constraints.

5) Kernel Function: The Kernel I choose is Gaussian Kernel.

$$K(\underline{X}^{i},\underline{X}^{j}) = \exp(-\frac{1}{h} \cdot ||\underline{X}^{i} - \underline{X}^{j}||^{2})$$

The reason is that theoretically, it can be used to fit any non-linear problem.

b) Result: After training the model, I can simply substitute X^i into $\beta(X^i)$ and see whether $\beta(X^i)$ is bigger than 0. If $\beta(X^i) > 0$, then classify X^i as class A, else classify it as class B. Name the unlabled images in the original order as C_1 , C_2 , C_3 , C_4 , C_5 . The result is $C_1 \rightarrow B$, $C_4 \rightarrow A$, $C_5 \rightarrow B$. The result is reasonable because I find that most of black cells of class A appear

in the upper left part and most of black cells of class B appear in the lower right part.

(b) In order to avoid overfitting, I add a regularization into the objective function. So now the objective function is as follows.

 $\min_{d_{i,...,dn}} \sum_{j=1}^{n} (y_{i} - \sum_{j=1}^{n} d_{j} K(\underline{x}^{i}, \underline{x}^{j}))^{2} + \lambda \cdot \| \vec{p}(x) \|^{2}$

By using the regularization term, the model is more likely to generate simple parameters. The reason I choose Lz regularization is that it has derivatives everywhere, which is a good mathematical property. The other way to avoid overfitting is to assign a high value to h in the kernel function.

(c) I also apply K-Nearest Neighbor model to this problem. The input is same as previous model. However, the output is no longer a value. It is a list containing labels of K nearest neighbors. In this model, we don't need a loss function or training process. Just compute the distances between the unlabeled point and any other labeled points. Pick K neareast neighbors and count the number of each class, then find which one is bigger and finally decide the classification.

The result is: $C_1 \rightarrow A$, $C_2 \rightarrow A$, $C_3 \rightarrow B$, $C_4 \rightarrow A$, $C_5 \rightarrow A$ when K is S. Obviously C_1 and C_5 are different from the prediction of SVM. However, I think the result makes sense because black cells of C_1 and C_5 are kind of evenly distributed on the map. The difference I think is from the uncertainty of values of K. I changed K to K, then it has the exactly same result of SVM.