# SUPPORT VECTOR CLASSIFIER VIA MATHEMATICA

Béla PALÁNCZ<sup>1</sup> and Lajos VÖLGYESI<sup>2,3</sup>

 <sup>1</sup> Department of Photogrammetry and Geoinformatics
 <sup>2</sup>Department of Geodesy and Surveying
 <sup>3</sup> Physical Geodesy and Geodynamic Research Group of the Hungarian Academy of Sciences Budapest University of Technology and Economics H–1521 Budapest, Hungary e-mail: palancz@epito.bme.hu

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### Abstract

In this case study a Support Vector Classifier function has been developed in *Mathematica*. Starting with a brief summary of support vector classification method, the step by step implementation of the classification algorithm in *Mathematica* is presented and explained. To check our function, two test problems, learning a chess board and classification of two intertwined spirals are solved. In addition, an application to filtering of airborne digital land image by pixel classification is demonstrated using a new SVM kernel family, the KMOD, a kernel with moderate decreasing.

Keywords: software Mathematica, kernel methods, pixel classification, remote sensing.

## Introduction

Kernel methods represent a relatively new family of algorithms that presents a series of useful features for pattern analysis in datasets. Kernel methods combine the simplicity and computational efficiency of linear algorithms, such as the perception algorithm or ridge regression, with flexibility of non-linear systems, such as for example neural networks, and rigour of statistical approaches such as regularization methods in multivariate statistics. As a result of the special way they represent functions, these algorithms typically reduce the learning step to convex optimization problem that can always be solved in polynomial time, avoiding the problem of local minima typical of neural networks, decision trees and other non-linear approaches [1].

### 1. Support Vector Classification

### 1.1. Binary classification

In the case of binary classification, we try to estimate a real-valued function  $f: X \subseteq \mathbb{R}^n \to \mathbb{R}$  using training data, that is *n*-dimensional patterns  $x_i$  and class labels  $y_i \in \{-1, 1\}$ 

$$((x_1, y_1), ..., (x_m, y_m)) \in \mathbb{R}^n \times \{-1, 1\}$$

such that f will correctly classify new examples (x, y) – that is, f(x) = y for samples(x, y), which were generated from the same underlying probability distribution P(x, y) as the training data. If we put no restriction on the class of functions that we choose our estimate f from, however, even a function that does well on the training data – for example by satisfying  $f(x_i) = y_i$  for i = 1...m – need not be generalized well to unseen examples. Suppose we know nothing additional about f (for example, about its smoothness), then the values on the training patterns carry no information whatsoever about values on novel patterns. Hence learning is impossible, and minimizing the training error does not imply any small expected test error.

Statistical learning theory, or Vapnik-Chervonenkis theory, shows that it is crucial to restrict the class of functions that the learning machine can implement to one with capacity that is suitable for the amount of available training data.

### 1.2. Optimal hyperplane classifier

To design learning algorithms, we thus must come up with a class of functions whose capacity can be computed. SV classifiers are based on the class of hyperplanes

$$\langle w, x \rangle + b = 0$$
  $w \in \mathbb{R}^n$ ,  $b \in \mathbb{R}$ 

corresponding to decision functions

$$f(x) = \operatorname{sign}(\langle w, x \rangle + b).$$

We can show that the optimal hyperplane, defined as the one with the maximal margin of separation between the two classes (see *Fig. 1*), has the lowest capacity, which ensure that the classifier learned from training samples will misclassify the less elements of the test samples originated from the same probability distribution.

### 1.3. Maximal margin classifier

The optimization problem to find the optimal w vector and the threshold b is the following, given a set of linearly separable training samples

$$S = ((x_1, y_1), ..., (x_m, y_m))$$

the hyperplane  $(w^*, b^*)$  that maximizes the geometric margin.

minimize<sub>w,b</sub> 
$$\langle w, w \rangle$$

subject to  $y_i \langle w, x_i \rangle + b \ge 1$ , i = 1, ..., m.

Then the geometric margin can be computed considering that

$$\langle w^*, x_1 \rangle + b^* = 1$$
  
 $\langle w^*, x_2 \rangle + b^* = -1$ 

then

$$\langle w^*, (x_1 - x_2) \rangle = 2$$

rescaling

$$\left\langle \frac{w^*}{\|w^*\|}, (x_1 - x_2) \right\rangle = \frac{2}{\|w^*\|}$$

therefore the margin is



*Fig. 1.* A separable classification problem. The optimal hyperplane is orthogonal to the shortest line connecting the convex hulls of the two classes, and intersects it half way. There is a weight vector w and a threshold b such that  $y_i (\langle w, x_i \rangle + b) > 0$ . Rescaling w and b such that the point(s) closest to the hyperplane satisfy  $|\langle w, x_i \rangle + b| = 1$ , we obtain a form (w, b) of the hyperplane with  $y_i (\langle w, x_i \rangle + b) \ge 1$ . Note that the margin, measured perpendicularly to the hyperplane, equals 1/||w||. To maximize the margin, we thus have to minimize ||w|| subject to  $y_i (\langle w, x_i \rangle + b) \ge 1$  [2].

The training patterns lying the closest to the hyperplane (see *Fig. 1* two balls and one diamond) are called support vectors, carrying all relevant information about the classification problem. The number of support vectors, SV are equal to or less than the number of the training patterns, m.

This minimization problem can be transformed into a dual maximization problem leading to a quadratic programming task, whose solution w has an expansion

$$w = \sum_{i=1}^{SV} v_i x_i$$

Consequently, the final decision function is

$$f(x) = \operatorname{sign}\left(\sum_{i=1}^{SV} v_i \langle x, x_i \rangle + b\right)$$

which only depends on dot products between patterns. This lets us generalize to the nonlinear case.

#### 1.4. Feature spaces and kernels

*Fig.* 2 shows the basic idea of SV machines, which is to map the data into some other dot space, called the feature space F via a nonlinear map,

$$\Phi: \mathbb{R}^n \to \mathbb{F}$$

and perform the above linear algorithm in F. This only requires the evaluation of dot products,

$$K(u, v) = \langle \Phi(u), \Phi(v) \rangle$$

Clearly, if F is high dimensional, the dot product on the right hand side will be very expensive to compute. In some cases, however, there is a simple kernel that can be evaluated efficiently. For instance, the polynomial kernel

$$K(u, v) = \langle u, v \rangle^{d}$$

can be shown to correspond to a map  $\Phi$  into the space spanned by all products of exactly d dimensions of  $\mathbb{R}^n$ . For d = 2 and  $u, v \in \mathbb{R}^2$ , for example, we have

$$\langle u, v \rangle^2 = \left( \left( \begin{pmatrix} u_1 \\ u_2 \end{pmatrix}, \begin{pmatrix} v_1 \\ v_2 \end{pmatrix} \right) \right)^2 = \left\langle \left( \begin{pmatrix} u_1^2 \\ \sqrt{2} \\ u_1 \\ u_2^2 \end{pmatrix}, \begin{pmatrix} v_1^2 \\ \sqrt{2} \\ v_1 \\ v_2 \end{pmatrix} \right) \right\rangle = \langle \Phi(u), \Phi(v) \rangle$$

defining  $\Phi(x) = (x_1^2, \sqrt{2} x_1 x_2, x_2^2).$ 



*Fig.* 2. The idea of SV machines: map the training data nonlinearly into a higher dimensional feature space via  $\Phi$ , and construct a separating hyperplane with maximum margin there. This yields a nonlinear decision boundary in input space. By the use of kernel function, it is possible to compute the separating hyperplane without explicitly carrying out the map into the feature space [3].

More generally, we can prove that for every kernel that gives rise to a positive matrix (kernel matrix)  $M_{ij} = K(x_i, x_j)$  we can construct a map such that  $K(u, v) = \langle \Phi(u), \Phi(v) \rangle$  holds.

#### 1.5. Optimization as a dual quadratic programming problem

Now the dual minimization problem of margin maximization is the following: Consider classifying a set of training samples

$$S = ((x_1, y_1), \dots (x_m, y_m))$$

using the feature space implicitly defined by the kernel K(x, z) and suppose the parameters  $\alpha^*$  solve the following quadratic optimization problem

minimize 
$$W(\alpha) = \sum_{i=1}^{m} \alpha_i - \frac{1}{2} \sum_{i,j=1}^{m} y_i y_j \alpha_i \alpha_j \left( K(x_i, x_j) + \frac{1}{c} \delta_{ij} \right)$$
  
subject to  $\sum_{i=1}^{m} y_i \alpha_i = 0$ ,  $\alpha_i \ge 0$ ,  $i = 1, \dots m$ .  
Let  $f(x) = \sum_{i=1}^{m} y_i \alpha_i^* K(x_i, x) + b^*$ , where  $b^*$  is chosen so that  $y_i f(x_i) = 1 - \frac{\alpha_i^*}{c}$  for any  $i$  with  $\alpha_i^* \ne 0$ .

Then the decision rule given by sign(f(x)) is equivalent to the hyperplane in the feature space implicitly defined by the kernel K(x, z), which solves the optimization problem, where the geometric margin is

$$\gamma = \left(\sum_{i \in sv} \alpha_i^* - \frac{1}{c} \langle \alpha^*, \; \alpha^* \rangle\right)^{-1/2}$$

where set *sv* corresponds to indexes *i*, for which  $\alpha_i^* \neq 0$ ,

$$sv = \{ i : \alpha_i^* \neq 0; i = 1, ..., m \}$$

Training samples,  $x_i$  for which  $i \in sv$  are called support vectors giving contribution to the definition of f(x).

### 2. Implementation of SVC in Mathematica

### 2.1. Steps of implementation

The dual optimization problem can be solved conveniently using *Mathematica*. In this section, the steps of the implementation of SVC algorithm are shown by solving XOR problem. The truth table of XOR, using bipolar values for the output, is

Table 1. Truth table of XOR problem

x1	x2	у
0	0	-1
0	1	1
1	0	1
1	1	-1

The input and output data lists are

xym={{0,0},{0,1},{1,0},{1,1}}; zm={-1,1,1,-1};

Let us employ Gaussian kernel with  $\beta$  gain

 $\beta$ =10.; K[u\_,v\_]:=Exp[- $\beta$  (u-v).(u-v)]

The number of the data pairs in the training set, m is

m=Length[zm] 4 Create the objective function  $W(\alpha)$  to be maximized, with regularization parameter,

c=5.;

First, we prepare a matrix M, which is an extended form of the kernel matrix,

M=(Table[N[K[xym[[i]],xym[[j]]]],
{i,1,m},{j,1,m}]+(1/c)IdentityMatrix[m]);

then the objective function can be expressed as

$$W = \sum_{i=1}^{m} \alpha_i - (1/2) \sum_{i=1}^{m} \sum_{j=1}^{m} (\operatorname{zm}[[i]]\operatorname{zm}[[j]] \alpha_i \alpha_j M[[i,j]]);$$

The constraints for the unknown variables are

g=Apply[And,Join[Table[ 
$$\alpha_i \geq 0$$
,{i,1,m}], { $\sum_{i=1}^m \text{ zm[[i]]}\alpha_i$ ==0}]];

However, the maximization problem is a convex quadratic problem, for practical reasons to maximize the objective function the built-in function NMaximize is applied. NMaximize implements several algorithms for finding constrained global optima. The methods are flexible enough to cope with functions that are not differentiable or continuous, and are not easily trapped by local optima. Possible settings for the Method option include "RandomSearch", "NelderMead", "DifferentialEvolution" and "SimulatedAnnealing".

Here we use DifferentialEvolution, which is a genetic algorithm that maintains a population of specimens,  $x_1, ..., x_n$ , represented as vectors of real numbers ('genes'). Every iteration, each  $x_i$  chooses random integers a, b, and c and constructs the mate  $y_i = x_i + \gamma (x_a + (x_b - x_c))$ , where  $\gamma$  is the value of ScalingFactor. Then  $x_i$  is mated with  $y_i$  according to the value of CrossProbability, giving us the child  $z_i$ . At this point  $x_i$  competes against  $z_i$  for the position of  $x_i$  in the population. The default value of SearchPoints is Automatic, which is Min[10\*d, 50], where d is the number of variables.

We need the list of unknown variables  $\alpha$ ,

vars=Table[ $\alpha_i$ ,{ i,1,m }];

Then the solution of the maximization problem is,

 $\begin{array}{l} \texttt{sol=NMaximize[{ W,g }, vars, Method } \rightarrow \texttt{DifferentialEvolution]} \\ \{1.66679, \{\alpha_1 \rightarrow 0.833396, \alpha_2 \rightarrow \\ 0.833396, \alpha_3 \rightarrow 0.833396, \alpha_4 \rightarrow 0.833396\} \}\end{array}$ 

The consistency of this solution can be checked by computing values of b for every data point. Theoretically, these values should be the same for any data points, however, in general, this is only approximately true.

 $\begin{aligned} &\texttt{bdata=Table[((1-\alpha_j/c)/zm[[j]]-} \\ &\sum_{i=1}^{m} \texttt{zm[[i]]}\alpha_i \texttt{K[xym[[i]],xym[[j]])/.sol[[2]],{j,1,m}]} \\ &\{-1.89729 \times 10^{-16}, 6.65294 \times 10^{-17}, 3.45251 \times 10^{-16}, 0.\} \end{aligned}$ 

The value of *b* can be chosen as the average of these values

```
b=Apply[Plus,bdata]/m 5.55126 \times 10^{-17}
```

Then the classifier function is,

$$f[w_]:=((\sum_{i=1}^{m} zm[[i]]\alpha_iK[w,xym[[i]]])+b)/.sol[[2]]$$

In symbolic form

Clear [x,y] f[{x,y}] 5.55126 × 10<sup>-17</sup> - 0.833396 e<sup>-10.((-1+x)<sup>2</sup>+(-1+y)<sup>2</sup>)</sup> + 0.833396 e<sup>-10.(x<sup>2</sup>+(-1+y)<sup>2</sup>)</sup> + 0.833396 e<sup>-10.((-1+x)<sup>2</sup>+y<sup>2</sup>)</sup> - 0.833396 e<sup>-10.(x<sup>2</sup>+y<sup>2</sup>)</sup>

Let us display the contour lines of the continuous classification function,

```
 << ExtendGraphics'LabelContour' 
 p=ContourPlot[f[{x,y}], {x, 0, 1}, {y,0,1}, 
 Contours <math>\rightarrow {-0.5,-0.05,0,0.05,0.5}, 
 ContourShading \rightarrow False, 
 PlotPoints \rightarrow 50,DisplayFunction \rightarrow Identity]; 
 p1=LabelContourLines[p,LabelPlacement \rightarrowAutomatic, 
 LabelFont \rightarrow {"Courier",10},DisplayFunction \rightarrow Identity]; 
 <<Graphics'MultipleListPlot' 
 p2=MultipleListPlot[{0,0},{1,1}},{0,1},{1,0}}, 
 SymbolStyle \rightarrow {Hue[.0],Hue[0.8]}, 
 SymbolStyle \rightarrow {PlotSymbol[Box,3], 
 PlotSymbol[Star,7]}, 
 Frame\rightarrowTrue,Axes\rightarrow False,AspectRatio\rightarrow1, 
 DisplayFunction \rightarrow Identity]; 
 Show[{p1,p2},DisplayFunction \rightarrow $DisplayFunction];
```



*Fig. 3.* The contour lines of the continuous classification function f(x1, x2) for XOR problem

The discrete classifier, the decision rule using signum function is Plot3D[Sign[f[{x,y}]], {x,0,1},{y,0,1},PlotRange $\rightarrow$ All];



*Fig.* 4. The decision rule, sign(f(x1, x2)) for XOR problem

Map[Sign[f[#]]&,xym]
{ -1, 1, 1, -1}
and
zm==%
True

#### 2.2. Mathematica modul for SVC

These steps can be collected in a module where the vector *xm* contains the input vectors (the training set), and the vector *ym* contains the corresponding scalar output values (the labels of the training set),

```
\begin{split} & \text{SupportVectorClassifier[xm_,ym_,K_,c_]:=Module[} \\ & \{\text{m,n,M,i,j,W,g,vars,sol,bdata,b}\}, \\ & \text{m=Length[ym];n=Length[xm[[1]]];} \\ & \text{M=Table[K[xm[[i]],xm[[j]]],{i,1,m}},{j,1,m}]+(1/c) \\ & \text{IdentityMatrix[m];} \\ & \text{W} = \sum_{i=1}^{m} \alpha_i - (1/2) \sum_{i=1}^{m} \sum_{j=1}^{m} (ym[[i]]ym[[j]] \alpha_i \alpha_j M[[i,j]]); \\ & \text{g=Apply[And,Join[Table[\alpha_i \ge 0,{i,1,m}],{\sum_{i=1}^{m} ym[[i]]\alpha_i == 0}]]; \\ & \text{vars=Table[\alpha_i,{i,1,m}];} \\ & \text{sol=NMaximize[{W,g}, vars,Method \rightarrow DifferentialEvolution][[2]]; \\ & \text{bdata=Table[((1-\alpha_j/c)/ym[[j]] - \sum_{i=1}^{m} ym[[i]]\alpha_i K[xm[[i]],xm[[j]])/.sol[[2]],{j,1,m}]; \\ & \text{bata=Table[((1-\alpha_j/c)/ym[[j]] - \sum_{i=1}^{m} ym[[i]]\alpha_i K[xm[[i]],xm[[j]])/.sol[[2]],{j,1,m}]; \\ & \text{bata=Table[((1-\alpha_j/c)/ym[[j]] - \sum_{i=1}^{m} ym[[i]]\alpha_i K[xm[[i]],xm[[j]])/.sol[[2]],{j,1,m}]; \\ & \text{bata=Table[} \} \\ & \text{vars=Iable[} \\ & \text{vars=Iable[} \} \\ & \text{vars=Iable[} \\ & \text{vars=Iab
```

The results of this module are the analytic form of the continuous classifier function and the values of  $\alpha_i$  's. Let us check the solution of the XOR problem

```
\begin{split} & \texttt{SupportVectorClassifier[xym,zm,K,c]} \\ & \{5.55112 \times 10^{-17} - 0.833396 \, \text{e}^{-10.((-1+x1)^2 + (-1+x2)^2)} \\ & + 0.833396 \, \text{e}^{-10.(x1^2 + (-1+x2)^2)} \\ & + 0.833396 \, \text{e}^{-10.((-1+x1)^2 + x2^2)} \\ & - 0.833396 \, \text{e}^{-10.(x1^2 + x2^2)}, \\ & \{0.833396, 0.833396, 0.833396, 0.833396\} \end{split}
```

```
%[[1]]==f[{x<sub>1</sub>,x<sub>2</sub> }]//Chop
True
```

### 3. Two test problems

### 3.1. Learning a chess board

Let us consider a  $2 \times 2$  chess board. The training points are generated by uniformly distributed random numbers from the interval [-1,1] × [-1, 1]. The chess board matrix

```
 M=\{\{1,-1\},\{-1,1\}\}; \\ p1=ListDensityPlot[M,MeshRange \rightarrow \{\{-1,1\},\{-1,1\}\}, \\ DisplayFunction \rightarrow Identity];
```

Creating the training set using 50 random samples,

```
xym={};zm={};
Do[x1=Random[Real,{-0.99,0.99}];
x2=Random[Real,{-0.99,0.99}];
If[x1 x2>0,z=1,z=-1];
AppendTo[xym,{x1,x2}];
AppendTo[zm,z],{k,1,50}];
```

Preparation data to display them

```
data=Transpose[Join[Transpose[xym],{zm}]];
data1=Map[Drop[#,-1]&,
Select[data,#[[3]]>0&]];
data2=Map[Drop[#,-1]&,
Select[data,#[[3]]<0&]];
p2=MultipleListPlot[data1,data2,
SymbolShape→{PlotSymbol[Triangle,5],
PlotSymbol[Box,2]},SymbolStyle→{Hue[.7],Hue[.0]},
Frame→True, AspectRatio→1,DisplayFunction→Identity];
```

Let us employ the same Gaussian kernel, but now with gain  $\beta = 15$ .

 $\beta$  =15;

employing parameter c = 100,

c=100;

The solution is

### F=SupportVectorClassifier[xym,zm,K,c];

This run could take some minutes.

 $\begin{array}{l} p3=ContourPlot[F[[1]], \{x_1, -1, 1\}, \{x_2, -1, 1\}, Contours \rightarrow \{0\}, \\ PlotPoints \rightarrow 100, ContourShading \rightarrow False, \\ DisplayFunction \rightarrow Identity]; \\ p4=DensityPlot[Sign[F[[1]]], \{x_1, -1, 1\}, \{x_2, -1, 1\}, \\ PlotPoints \rightarrow 100, Mesh \rightarrow False, DisplayFunction \rightarrow Identity]; \\ Show[GraphicsArray[{Show[{p2,p3}], p1, p4}], \\ DisplayFunction \rightarrow $DisplayFunction]; \end{array}$ 



*Fig. 5.* SVC result from 50 random samples: the functions f(x) with the random samples, the ideal chess board and sign(f(x))

The SV solution of a larger chess board problem can be found in [3].

### 3.2. Two intertwined spirals

The two intertwined spirals represent a challenging classification benchmark originated from the field of neural networks [4].

Parametric equations of the spirals are

x1[t\_]:=2 Cos[t]  $e^{t/10}$ y1[t\_]:=1.5 Sin[t]  $e^{t/10}$ x2[t\_]:=2.7 Cos[t]  $e^{t/10}$ y2[t\_]:=2.025 Sin[t]  $e^{t/10}$ 

Generating 26 discrete points for each spiral,

s1=Table[{x1[t],y1[t]},{t,  $\pi$ , 3.5 $\pi$ , 2.5 $\pi$ /25}]; s2=Table[{x2[t],y2[t]},{t,  $-\pi$  /2, 2.5 $\pi$  , 3.0 $\pi$  /25}];

then displaying these points,

```
\begin{split} & S1=ListPlot[s1,PlotStyle \rightarrow \{RGBColor[1,0,0],PointSize[0.02]\},\\ & AspectRatio \rightarrow 1,DisplayFunction \rightarrow Identity];\\ & S2=ListPlot[s2,PlotStyle \rightarrow \{RGBColor[0,0,1],PointSize[0.015]\},\\ & AspectRatio \rightarrow 1,DisplayFunction \rightarrow Identity];\\ & pspiral=Show[\{S1,S2\},DisplayFunction \rightarrow \$DisplayFunction]; \end{split}
```



Fig. 6. Two intertwined spirals represented by 26 points each

Creating the teaching set, putting these points into one list

# xym=Join[s1,s2];

Generating the labels of the samples,

```
zm=Join[Table[1.,{26}],Table[-1.,{26}]];
Length[zm]
52 Dimensions[xym]
{52,2}
```

Applying wavelet kernel [5] with parameter a = 1.8, in the case of dimension n = 2

n=2;a=1.8;  
K[u\_,v\_]:=
$$\prod_{i=1}^{n}$$
 (Cos[1.75(u[[i]]-v[[i]])/a] Exp[-(u[[i]]-v[[i]])<sup>2</sup>/2a<sup>2</sup>])

and with parameter c = 100

c=100.;

The solution is

```
F=SupportVectorClassifier[xym,zm,K,c];
```

This run could take some minutes.

```
\begin{array}{l} \texttt{psvc1=ContourPlot[F[[1]], \{x_1, -7, 7\}, \{x_2, -7, 7\},}\\ \texttt{Contours} \rightarrow \{0\},\\ \texttt{PlotPoints} \rightarrow 50, \texttt{AspectRatio} \rightarrow 1, \texttt{ContourShading} \rightarrow \texttt{False},\\ \texttt{DisplayFunction} \rightarrow \texttt{Identity]};\\ \texttt{Show[{pspiral,psvc1}, DisplayFunction} \rightarrow \texttt{SDisplayFunction]}; \end{array}
```



Fig. 7. Classification results of SVC with the nonlinear decision boundary

The continuous classification function and the decision rule can be displayed in 3D, too

```
\begin{array}{l} \texttt{P1=Plot3D[F[[1]],} \{x_1, -7, 7\}, \{x_2, -7, 7\}, \texttt{PlotRange} \rightarrow \texttt{All}, \\ \texttt{PlotPoints} \rightarrow \{30, 30\}, \\ \texttt{DisplayFunction} \rightarrow \texttt{Identity]}; \\ \texttt{P2=Plot3D[Sign[F[[1]]],} \{x_1, -7, 7\}, \{x_2, -7, 7\}, \texttt{PlotRange} \rightarrow \texttt{All}, \\ \texttt{PlotPoints} \rightarrow \{40, 40\}, \texttt{DisplayFunction} \rightarrow \texttt{Identity]}; \\ \texttt{Show[GraphicsArray[}\{\texttt{P1,P2}\}], \ \texttt{DisplayFunction} \Rightarrow \texttt{SDisplayFunction}; \\ \end{array}
```



Fig. 8. Classification function and the corresponding decision rule

## 4. Image Classification

Classification of digital images in order to separate different categories of land cover types like urban area, water, vegetation, agricultural area etc., and to carry out thematic change analysis are frequent tasks in geoscience and remote sensing. Many different methods are used, traditional maximum-likelihood, *k*-nearest neighbor, rule based (classification and regression tree), supervised and unsupervised neural network, fuzzy logic and neuro-fuzzy and even support vector classifiers with Gaussian-kernel.

SVCs were used for classification of land cover using polarimetric synthetic aperture radar (SAR) images, and for classification of clouds, snow and ice, however, with usual kernels like RBF (Radial Basis Function). In this illustrative example, we use KMOD type kernel for synthetic image data.

Let us load the Image Processing Application of Mathematica,

```
<<ImageProcessing'
img=ImageRead["F: \ ImageProcessing\Syntetic.JPG"];
```

Let us consider a relatively small image of synthetic data,

```
ImageDimensions[img]
{201, 301}
Show[Graphics[img]];
```

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Fig. 9. Digital image of synthetic data

4.1. Binary Classification

We should filter out pixels being different from yellowy type ones. For this task an SVM binary classifier can be developed. Pixels from the three different categories: yellowy, brownish and bluish spots are picked up randomly and their RGB values are stored in three different files. Unfortunately, this journal supports black and white images, only. To enjoy the colors, please, see [9] on Web.

We have ten yellowish pixels,

```
rgb1=
ReadList["F:\ImageProcessing\yellow.dat", {Number,Number,Number}];
Show[
Graphics[RasterArray[{RGBColor[#[[1]],#[[2]],#[[3]]]&/@rgb1}]]];
```



Fig. 10. Colors (grayscale) of the ten yellowish training data

ten brownish pixels,

```
rgb2=
ReadList["F:\ImageProcessing\dark.dat", {Number,Number,Number}];
Show [
Graphics[RasterArray[{RGBColor[#[[1]], #[[2]],#[[3]]]&/@rgb2}]]];
```

and ten bluish pixels,





```
rgb3=
ReadList["F:\ImageProcessing \light.dat",{Number,Number,Number}];
Show [
Graphics[RasterArray[{RGBColor[#[[1]],#[[2]],#[[3]]]&/@rgb3}]]];
```



Fig. 12. Colors (grayscale) of the ten bluish training data

These three sets are jointly building the input for the classifier,

### xym=Join[rgb1,rgb2,rgb3];

The first ten elements are labeled with 1, and the remaining elements with -1,

```
zm=Join[Table[1,{10}],Table[-1,{20}]];
```

Now, we shall employ a Kernel with MOderate Decreasing (KMOD) having two parameters,

 $\gamma = 0.5; \sigma = 3.;$ 

The motivation of the employment of this kernel can be explained as follows. In most commonly used kernels (eg. RBF), points very close to each other are strongly correlated whereas points far apart have uncorrelated images in the augmented space. The aim is to force the images of the original points to be linearly separable in the augmented space. In order to get such a behavior, a kernel must turn very close points from the original space into weakly correlated elements (as weak as possible) while still preventing the closeness information from vanishing. To achieve this tradeoff, we need the following couple features: a quick decrease in the neighborhood of zero and a moderate decrease towards infinity. The RBF kernel may satisfy correctly the first requirement but not the second, whereas the exponential RBF does not respond correctly to both of the requirements. Alternatively, the KMOD is proposed [7], whose analytic expression is

 $K[u_,v_]:=Exp[\gamma/(Norm[u-v]^2+\sigma^2)]-1$ 

In the case of n = 1

Plot3D[  $Exp[\gamma/(Abs[u-v]^2 + \sigma^2)] - 1, \{u, 0, 1\}, \{v, 0, 1\}];$ 



Fig. 13. Kernel with MOderate Decreasing (KMOD)

The value of the regularization parameter is,

## c=100;

Training the continuous classifier function,

### F1=SupportVectorClassifier[xym,zm,K,c];

we get its analytical form,

 $\begin{aligned} & \texttt{Short[F1[[1]],15]} \\ & -0.108208 - 91.611(-1 + e^{0.5/(9.+\text{Abs}[-0.352941 + x1]^2 + \text{Abs}[-0.639216 + x2]^2 + \text{Abs}[-0.596078 + x3]^2)}) \\ & -43.766(-1 + e^{0.5/(9.+\text{Abs}[-0.34902 + x1]^2 + \text{Abs}[-0.627451 + x2]^2 + \text{Abs}[-0.596078 + x3]^2)}) \\ & + << 30 >> + \\ & + 0.0(-1 + e^{0.5/(9.+\text{Abs}[-0.25098 + x1]^2 + \text{Abs}[-0.262745 + x2]^2 + \text{Abs}[-0.219608 + x3]^2)}) + \\ & + 0.0(-1 + e^{0.5/(9.+\text{Abs}[-0.235294 + x1]^2 + \text{Abs}[-0.247059 + x2]^2 + \text{Abs}[-0.211765 + x3]^2)}) \end{aligned}$ 

The RGB values of the original image vector

### xx=Flatten[RawImageData[img],1];

and its dimensions

```
nd=Dimensions[xx] \{60501, 3\}
```

Now, the discrete classifier is applied,

```
decision=
Map[Sign[F1[[1]]]/.{x_1 \rightarrow \#[[1]],x_2 \rightarrow \#[[2]],x_3 \rightarrow \#[[3]]}&,xx/255];
```

We shall use colors for yellowish and not yellowish spots, with RGB values patches and others, respectively,

patches={0.5,0.5,0};others={1,0.5,0.75};

The RGB value of pixels classified as yellowish (labeled with 1 by the classifier) will be overwritten by the RGB values of patches, and the others pixels (labeled with -1) by the RGB values of others,

```
Do[If[decision[[i]]==1,xx[[i]]=patches,xx[[i]]=others],
{i,1,nd[[1]]};
```

We partition the RGB image vector to form a matrix, and transform this image matrix into an image object,

```
U=ToRGBColor[Partition[xx,301]];
```

Here are the original and the filtered image,

# Show[GraphicsArray[{Graphics[img],Graphics[U]}]];



Fig. 14. The original and the filtered image

The results show the data generalization ability of SVC, which was trained with only 30 pixels, and successfully represents more than 60 thousands.

#### 4.2. Multi-class Classification

Support Vector Classifiers were originally designed for binary classification. How to effectively extend it for multi-class classification is still an on-going research issue. Currently there are two types of approaches for multi-class SVC. One is by constructing and combining several binary classifiers, while the other is by directly considering all data in one optimization formulation. Methods of this latter category are called all-together methods.

Here we employ the *one-against-all* method [8] belonging to the first category, the combination of several binary classifiers. It constructs k SVC models, which means k decision functions, where k is the number of classes. The *i*th SVC is trained with all of the examples in the *i*-th class with positive labels, and all other examples with negative labels. Thus given m training data  $(x_1, y_1), \ldots, (x_m, y_m)$ , where  $x_i \in \mathbb{R}^n$ ,  $i = 1, \ldots, m$ , and  $y_i \in \{1, k\}$ . We say  $x_i$  is in the class which has the largest value of the decision function. In our case  $k = 3, m = 201 \times 301 = 60501$  and n = 3.

Let us define the list of labels for the second class:

### zm=Join[Table[-1,{10}],Table[1,{10}],Table[-1,{10}]];

The second decision function is

### F2=SupportVectorClassifier[xym,zm,K,c];

in symbolic form

```
\begin{aligned} & \texttt{Short}[\texttt{F2}[\texttt{[1]}],\texttt{15}] \\ & -0.0635132 - 60.3727(-1 + \\ & e^{0.5/(9.+Abs[-0.352941+x1]^2+Abs[-0.639216+x2]^2+Abs[-0.596078+x3]^2)}) - \\ & -32.4338(-1 + e^{0.5/(9.+Abs[-0.34902+x1]^2+Abs[-0.627451+x2]^2+Abs[-0.596078+x3]^2)}) + \\ & + << 27 >> + \\ & +42.0541(-1 + e^{0.5/(9.+Abs[-0.25098+x1]^2+Abs[-0.262745+x2]^2+Abs[-0.219608+x3]^2)}) + \\ & + 37.0146(-1 + e^{0.5/(9.+Abs[-0.235294+x1]^2+Abs[-0.247059+x2]^2+Abs[-0.211765+x3]^2)}) \end{aligned}
```

The third class has the list of labels as follows

### zm=Join[Table[-1,{20}],Table[1,{10}]];

Then the third decision function is

#### F3=SupportVectorClassifier[xym,zm,K,c];

in symbolic form

```
\begin{aligned} & \texttt{Short[F3[[1]],15]} \\ & -0.874528 + 157635(-1 + \\ & + e^{0.5/(9.+Abs[-0.352941+x1]^2 + Abs[-0.639216+x2]^2 + Abs[-0.596078+x3]^2)}) + 78.926(-1 + \\ & + e^{0.5/(9.+Abs[-0.34902+x1]^2 + Abs[-0.627451+x2]^2 + Abs[-0.596078+x3]^2)}) - \\ & -51.2964(-1 + e^{0.5/(9.+Abs[-0.811765+<<1>)]^2 + Abs[<<1>>]^2 + Abs[-0.588235+x3]^2)}) + \\ & + 81.8819(-1 + e^{0.5/(9.+<<1>>2+<<1>>2+Abs[<<1>>]^2)}) + \\ & + <<32 >> \end{aligned}
```

Now we restore the RGB values of the original image vector which was overwritten during the binary classification

### xx=Flatten[RawImageData[img],1];

Then applying the three different continuous classifiers to the pixel vector of the image

```
decision1=

Map[F1[[1]]/.{x_1 \rightarrow \#[[1]], x_2 \rightarrow \#[[2]], x_3 \rightarrow \#[[3]]}&,xx/255];

decision2=

Map[F2[[1]]/.{x_1 \rightarrow \#[[1]], x_2 \rightarrow \#[[2]], x_3 \rightarrow \#[[3]]}&,xx/255];

decision3=

Map[F3[[1]]/.{x_1 \rightarrow \#[[1]], x_2 \rightarrow \#[[2]], x_3 \rightarrow \#[[3]]}&,xx/255];
```

We construct a list having elements as list referring to a pixel, and containing three values, the values resulted by the three different classifiers applied to the pixels

### d123=Transpose[{decision1,decision2,decision3}];

The pixel will be assigned to the class for which it has the largest value of the decision functions:

### d=Map[Position[#,Max[#]]&,d123]//Flatten;

Let us assign the following three colors to the different classes:

class1={1,0,0};class2={0,1,0};class3={0,0,1};

We paint each pixel with the color of the proper class,

```
Do[Which[d[[i]]==1,xx[[i]]=class1,d[[i]]==2,xx[[i]]=class2,
d[[i]]==3,xx[[i]]=class3],{i,1,nd[[1]]};
```

then partition the RGB image vector to form an image matrix, and transform this image matrix into an image object:

```
U=ToRGBColor[Partition[xx,301]];
```

Here are the original and the three-class classified image:

# Show[GraphicsArray[{Graphics[img],Graphics[U]}]];



Fig. 15. The original and the three-class classified image

## 5. Conclusions

Support vector classification method has been developed in software *Mathematica* and the method is ready to use for different technical applications. The step by step implementation of the support vector classification algorithm in *Mathematica* was presented and explained here.

Support vector classification method provides a very promising application possibility in photogrammetry and petrological microscopy. One of the most important applications of this method in remote sensing is the filtering of airborne digital land images by pixel classification.

The Mathematica notebook form of this paper is available on Web [9].

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