This document explains how to use the parameters in an object returned by \texttt{svm()} for own prediction functions.

1 Binary Classifier

For class prediction in the binary case, the class of a new data vector \( n \) is usually given by \textit{the sign} of

\[
\sum_i a_i y_i K(x_i, n) + \rho
\]

where \( x_i \) is the \( i \)-th support vector, \( y_i \) the corresponding label, \( a_i \) the corresponding coefficient, and \( K \) is the kernel (for example the linear one, i.e. \( K(u, v) = u^T v \)).

Now, the \texttt{libsvm} library interfaced by the \texttt{svm()} function actually returns \( a_i y_i \) as \( i \)-th coefficient and the negative \( \rho \), so in fact uses the formula:

\[
\sum_i \text{coef}_i K(x_i, n) - \rho
\]

where the training examples (=training data) are labeled \{1,-1\} (!). A simplified R function for prediction with linear kernel would be:

\[
\texttt{svmpred <- function (m, newdata, K=crossprod)}
\
\{
  \texttt{## this guy does the computation:}
  \texttt{pred.one <- function (x)}
  \texttt{    sign(sum(sapply(1:m$tot.nSV, function (j) K(m$SV[j,], x) * m$coefs[j] ) - m$rho )} \\
  \texttt{## this is just for convenience:}
  \texttt{if (is.vector(newdata))}
  \texttt{    newdata <- t(as.matrix(x))}
  \texttt{sapply (1:nrow(newdata),}
  \texttt{        function (i) pred.one(newdata[i,])})
\}
where `pred.one()` does the actual prediction for one new data vector, the remainder is just a convenience for prediction of multiple new examples. It is easy to extend this to other kernels, just replace $K()$ with the appropriate function (see the help page for the formulas used) and supply the additional constants.

As we will see in the next section, the multi-class prediction is more complicated, because the coefficients of the diverse binary SVMs are stored in a compressed format.

## 2 Multiclass-classifier

To handle $k$ classes, $k > 2$, `svm()` trains all binary subclassifiers (one-against-one-method) and then uses a voting mechanism to determine the actual class. Now, this means $k(k-1)/2$ classifiers, hence in principle $k(k-1)/2$ sets of SVs, coefficients and rhos. These are stored in a compressed format:

1. Only one SV is stored in case it were used by several classifiers. The `model$SV-matrix` is ordered by classes, and you find the starting indices by using `nSV` (number of SVs):

   ```r
   start <- c(1, cumsum(model$nSV))
   start <- start[-length(start)]
   ```

   `sum(nSV)` equals the total number of (distinct) SVs.

2. The coefficients of the SVs are stored in the `model$coefs`-matrix, grouped by classes. Because the separating hyperplanes found by the SVM algorithm has SVs on both sides, you will have two sets of coefficients per binary classifier, and e.g., for 3 classes, you could build a block-matrix like this for the classifiers $(i, j)$ ($i,j$=class numbers):

   $\begin{array}{ccc}
   & 0 & 1 & 2 \\
   0 & X & set(0,1) & set(0,2) \\
   1 & set(1,0) & X & set(1,2) \\
   2 & set(2,0) & set(2,1) & X \\
   \end{array}$

   where `set(i,j)` are the coefficients for the classifier $(i,j)$, lying on the side of class $j$. Because there are no entries for $(i,i)$, we can save the diagonal and shift up the lower triangular matrix to get

   $\begin{array}{ccc}
   & 0 & 1 & 2 \\
   0 & set(1,0) & set(0,1) & set(0,2) \\
   1 & set(2,0) & set(2,1) & set(1,2) \\
   \end{array}$

   Each set $(.,j)$ has length `nSV[j]`, so of course, there will be some filling 0s in some sets.

   `model$coefs` is the transposed of such a matrix, therefore for a data set with, say, 6 classes, you get 6-1=5 columns.

   The coefficients of $(i,j)$ start at `model$coefs[start[i],j]` and those of $(j,i)$ at `model$coefs[start[j],i-1]`.

3. The $k(k-1)/2$ rhos are just linearly stored in the vector `model$rho`.
The following code shows how to use this for prediction:

```r
## Linear Kernel function
K <- function(i,j) crossprod(i,j)
predsvm <- function(object, newdata)
{
  ## compute start-index
  start <- c(1, cumsum(object$nSV)+1)
  start <- start[-length(start)]

  ## compute kernel values
  kernel <- sapply (1:object$tot.nSV,
      function (x) K(object$SV[x,], newdata))

  ## compute raw prediction for classifier (i,j)
predone <- function (i,j)
  {
    ## ranges for class i and j:
    ri <- start[i] : (start[i] + object$nSV[i] - 1)
    rj <- start[j] : (start[j] + object$nSV[j] - 1)

    ## coefs for (i,j):
    coef1 <- object$coefs[ri, j-1]
    coef2 <- object$coefs[rj, i]

    ## return raw values:
    crossprod(coef1, kernel[ri]) + crossprod(coef2, kernel[rj])
  }

  ## compute votes for all classifiers
  votes <- rep(0,object$nclasses)
  c <- 0 # rho counter
  for (i in 1 : (object$nclasses - 1))
    for (j in (i + 1) : object$nclasses)
      if (predone(i,j) > object$rho[c <- c + 1])
        votes[i] <- votes[i] + 1
      else
        votes[j] <- votes[j] + 1

  ## return winner (index with max. votes)
  object$levels[which(votes %in% max(votes))[1]]
}
```

In case data were scaled prior fitting the model (note that this is the default for `svm()`), the new data needs to be scaled as well before applying the prediction functions, for example using the following code snippet (object is an object returned by `svm()`, newdata a data frame):

```r
if (any(object$scaled))
  newdata[,object$scaled] <-
    scale(newdata[,object$scaled, drop = FALSE],
      center = object$x.scale$scaled:center",
      scale = object$x.scale$scaled:scale"
)
```

For regression, the response needs to be scaled as well before training, and the predictions need to be scaled back accordingly.