

# SC1\_Proj

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## Dataset

The dataset is TODO: describe

```
data <- read_csv("../data/data.csv")
glimpse(data)
```

```
## Observations: 569
## Variables: 33
## $ id <dbl> 842302, 842517, 84300903, 84348301, 8435840...
## $ diagnosis <chr> "M", "M", "M", "M", "M", "M", "M", "M", "M"...
## $ radius_mean <dbl> 17.990, 20.570, 19.690, 11.420, 20.290, 12....
## $ texture_mean <dbl> 10.38, 17.77, 21.25, 20.38, 14.34, 15.70, 1...
## $ perimeter_mean <dbl> 122.80, 132.90, 130.00, 77.58, 135.10, 82.5...
## $ area_mean <dbl> 1001.0, 1326.0, 1203.0, 386.1, 1297.0, 477....
## $ smoothness_mean <dbl> 0.11840, 0.08474, 0.10960, 0.14250, 0.10030...
## $ compactness_mean <dbl> 0.27760, 0.07864, 0.15990, 0.28390, 0.13280...
## $ concavity_mean <dbl> 0.30010, 0.08690, 0.19740, 0.24140, 0.19800...
## $ `concave points_mean` <dbl> 0.14710, 0.07017, 0.12790, 0.10520, 0.10430...
## $ symmetry_mean <dbl> 0.2419, 0.1812, 0.2069, 0.2597, 0.1809, 0.2...
## $ fractal_dimension_mean <dbl> 0.07871, 0.05667, 0.05999, 0.09744, 0.05883...
## $ radius_se <dbl> 1.0950, 0.5435, 0.7456, 0.4956, 0.7572, 0.3...
## $ texture_se <dbl> 0.9053, 0.7339, 0.7869, 1.1560, 0.7813, 0.8...
## $ perimeter_se <dbl> 8.589, 3.398, 4.585, 3.445, 5.438, 2.217, 3...
## $ area_se <dbl> 153.40, 74.08, 94.03, 27.23, 94.44, 27.19, ...
## $ smoothness_se <dbl> 0.006399, 0.005225, 0.006150, 0.009110, 0.0...
## $ compactness_se <dbl> 0.049040, 0.013080, 0.040060, 0.074580, 0.0...
## $ concavity_se <dbl> 0.05373, 0.01860, 0.03832, 0.05661, 0.05688...
## $ `concave points_se` <dbl> 0.015870, 0.013400, 0.020580, 0.018670, 0.0...
## $ symmetry_se <dbl> 0.03003, 0.01389, 0.02250, 0.05963, 0.01756...
## $ fractal_dimension_se <dbl> 0.006193, 0.003532, 0.004571, 0.009208, 0.0...
## $ radius_worst <dbl> 25.38, 24.99, 23.57, 14.91, 22.54, 15.47, 2...
## $ texture_worst <dbl> 17.33, 23.41, 25.53, 26.50, 16.67, 23.75, 2...
## $ perimeter_worst <dbl> 184.60, 158.80, 152.50, 98.87, 152.20, 103....
## $ area_worst <dbl> 2019.0, 1956.0, 1709.0, 567.7, 1575.0, 741....
## $ smoothness_worst <dbl> 0.1622, 0.1238, 0.1444, 0.2098, 0.1374, 0.1...
## $ compactness_worst <dbl> 0.6656, 0.1866, 0.4245, 0.8663, 0.2050, 0.5...
## $ concavity_worst <dbl> 0.71190, 0.24160, 0.45040, 0.68690, 0.40000...
## $ `concave points_worst` <dbl> 0.26540, 0.18600, 0.24300, 0.25750, 0.16250...
## $ symmetry_worst <dbl> 0.4601, 0.2750, 0.3613, 0.6638, 0.2364, 0.3...
## $ fractal_dimension_worst <dbl> 0.11890, 0.08902, 0.08758, 0.17300, 0.07678...
## $ X33 <chr> NA, NA, NA, NA, NA, NA, NA, NA, NA, NA, NA, NA,...
```

```
colnames(data)[3:32] <- c('radius_m','texture_m', 'perim_m','area_m','smooth_m','compact_m','concav_m',
```

## Dataset Preprocessing Visualisation and Exploration

```
colSums(is.na(data))
```

```
##          id      diagnosis      radius_m      texture_m      perim_m      area_m
##          0          0          0          0          0          0
##      smooth_m      compact_m      concav_m      concav_pt_m      symmetry_m      frac_dim_m
##          0          0          0          0          0          0
##      radius_se      texture_se      perim_se      area_se      smooth_se      compact_se
##          0          0          0          0          0          0
##      concav_se      concav_pt_se      symmetry_se      frac_dim_se      radius_w      texture_w
##          0          0          0          0          0          0
##      perim_w      area_w      smooth_w      compact_w      concav_w      concav_pt_w
##          0          0          0          0          0          0
##      symmetry_w      frac_dim_w      X33
##          0          0          569
```

```
data %<>% mutate_at(vars(diagnosis), factor)
```

```
train <- data %>% sample_frac(0.8)
test <- anti_join(data,train, by='id')
```

```
# need ids for later
id_train <- train$id
id_test <- test$id
```

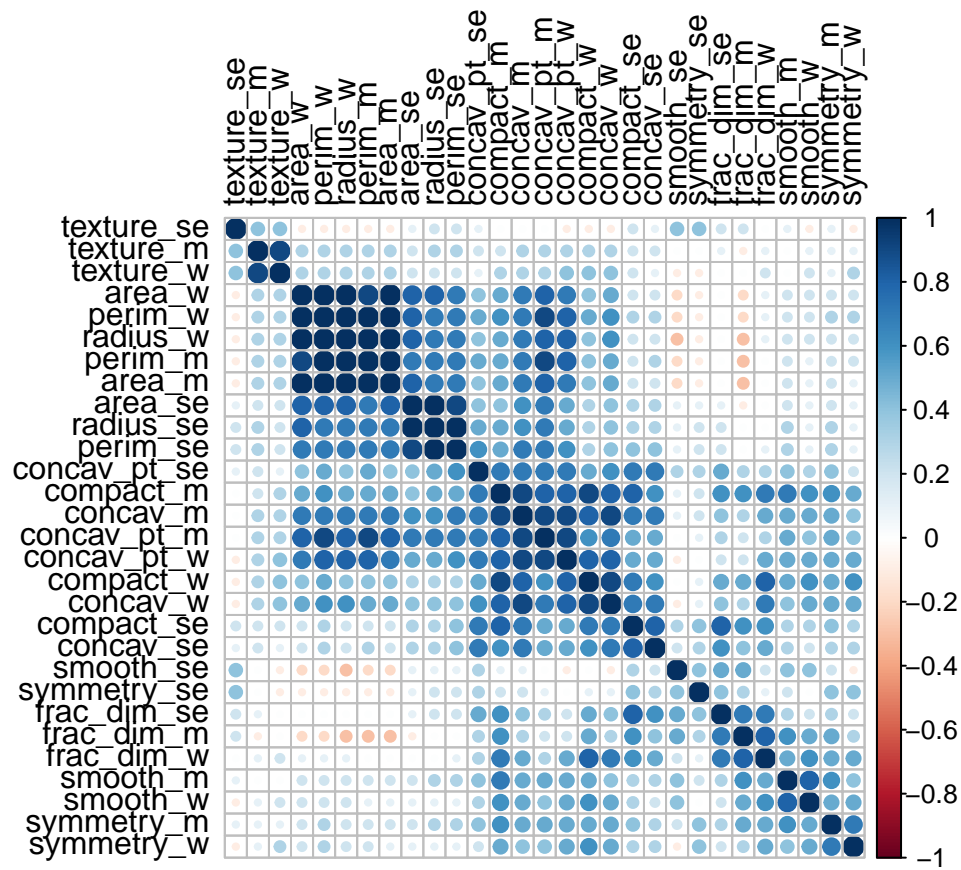
```
data %<>%
  dplyr::select(-c(id, X33))
train %<>%
  dplyr::select(-c(id, X33))
test %<>%
  dplyr::select(-c(id, X33))
```

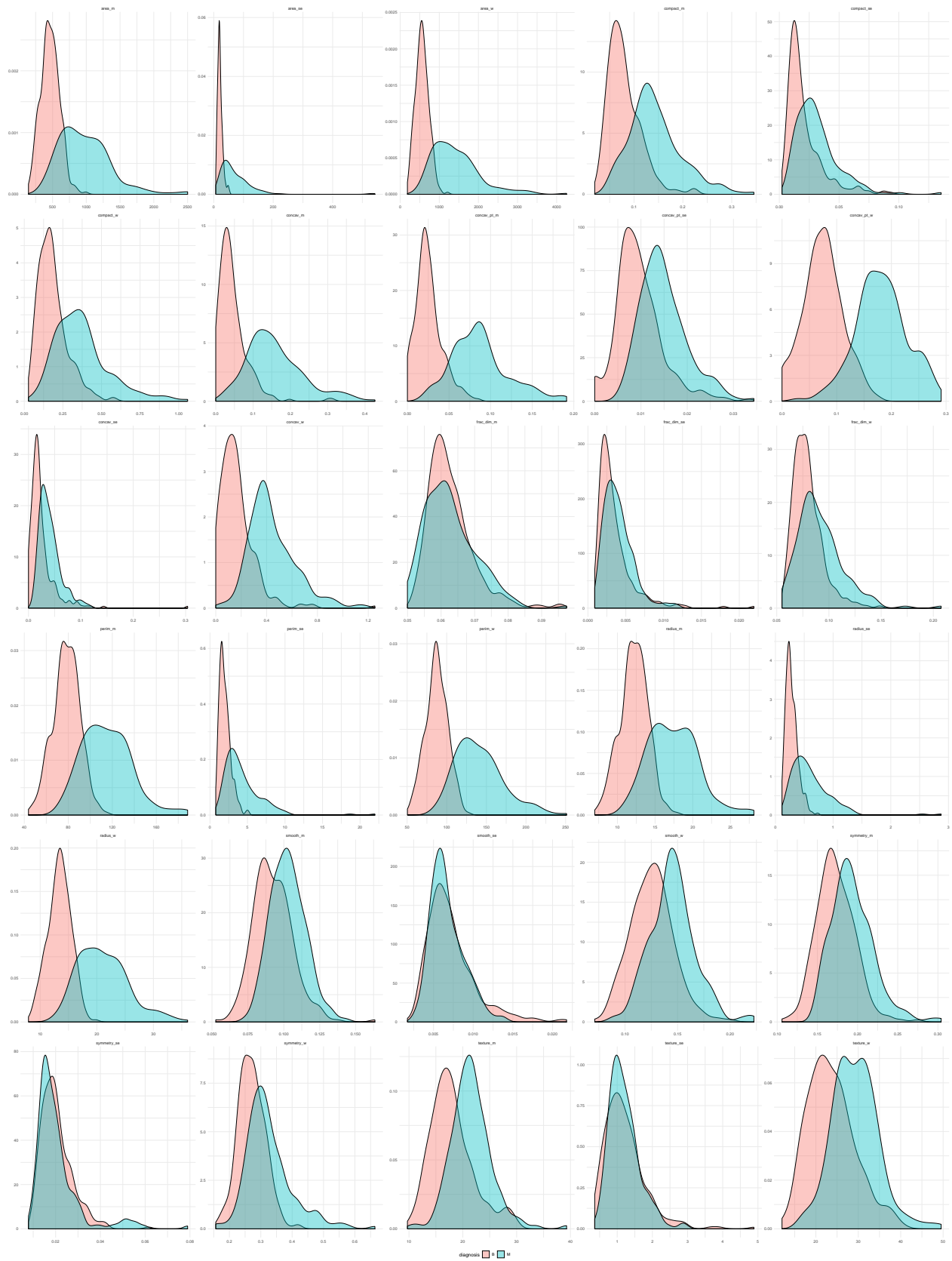
```
sum(is.na(data))
```

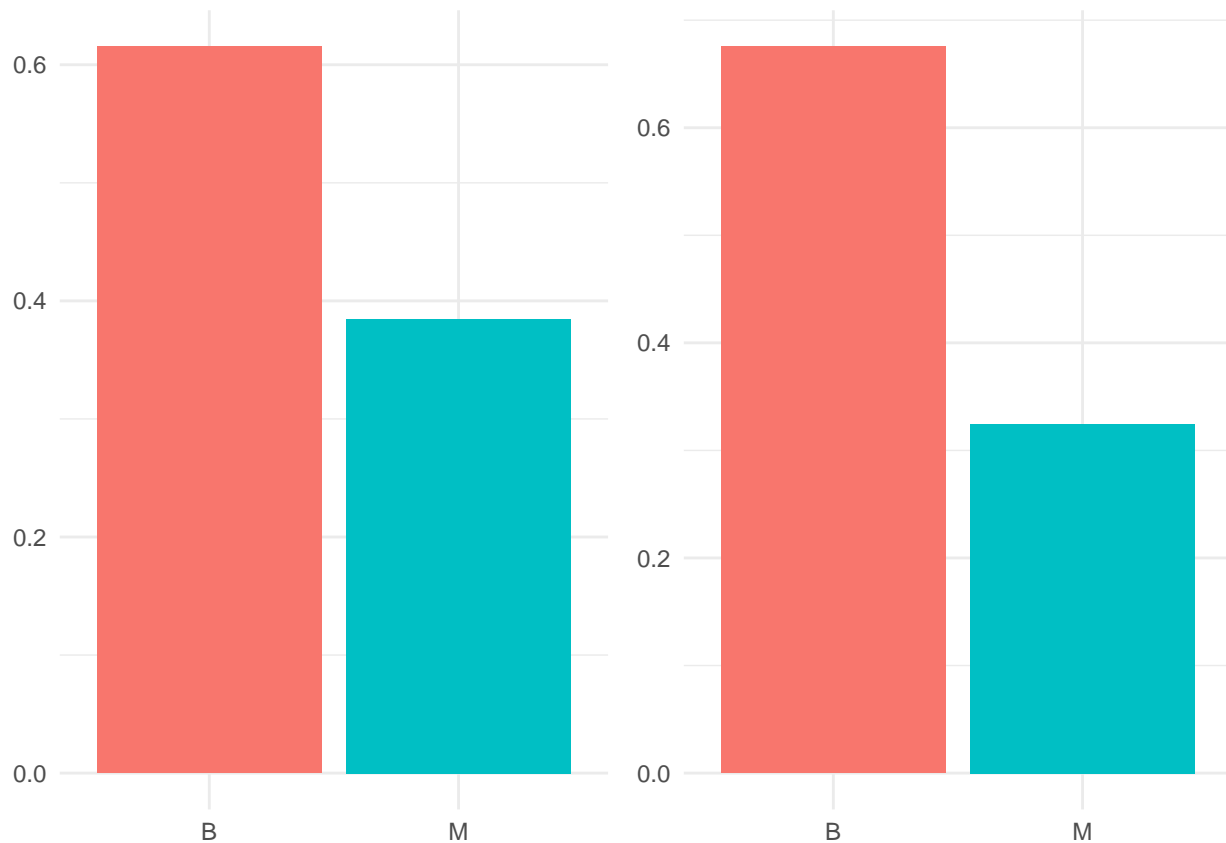
```
## [1] 0
```

```
training_data <- train[2:dim(train)[2]]
training_classes <- train[1]
```

```
test_data <- test[2:dim(test)[2]]
test_classes <- test[1]
```







```
confusion_plot <- function(actual,predicted){
  confusion_matrix <- as.data.frame(table(actual,predicted))
  g <-ggplot(confusion_matrix,aes(x=actual,y=predicted))+
    geom_tile(aes(fill=Freq))+
    geom_text(aes(label=sprintf("%1.0f", Freq)),color="white",fontface="bold")+
    labs(x="Actual class",y="Predicted class")+
    theme_minimal()
  return(g)
}
```

## Dimensionality Reduction and Feature Selection

### PCA

Code

```
normalise_z <- function(X){
  mean_cols <- colMeans(X)
  sd_cols <- apply(X, 2, sd)
  mean_normalised_X <- t(apply(X, 1, function(x){x - mean_cols}))
  normalised_X <- t(apply(mean_normalised_X, 1, function(x){x / sd_cols}))
  return(normalised_X)
}

pca <- function(X, number_components_keep) {
  normalised_X <- normalise_z(X)
```

```

corr_mat <- t(normalised_X) %*% normalised_X

eigenvectors <- eigen(corr_mat, symmetric=TRUE)$vectors

reduced_data <- X %*% eigenvectors[,1:number_components_keep]
relevant_eigs <- eigenvectors[,1:number_components_keep]
returnds <- list(reduced_data, relevant_eigs)
names(returnds) <- c("reduced_data", "reduction_matrix")
return(returnds)
}

```

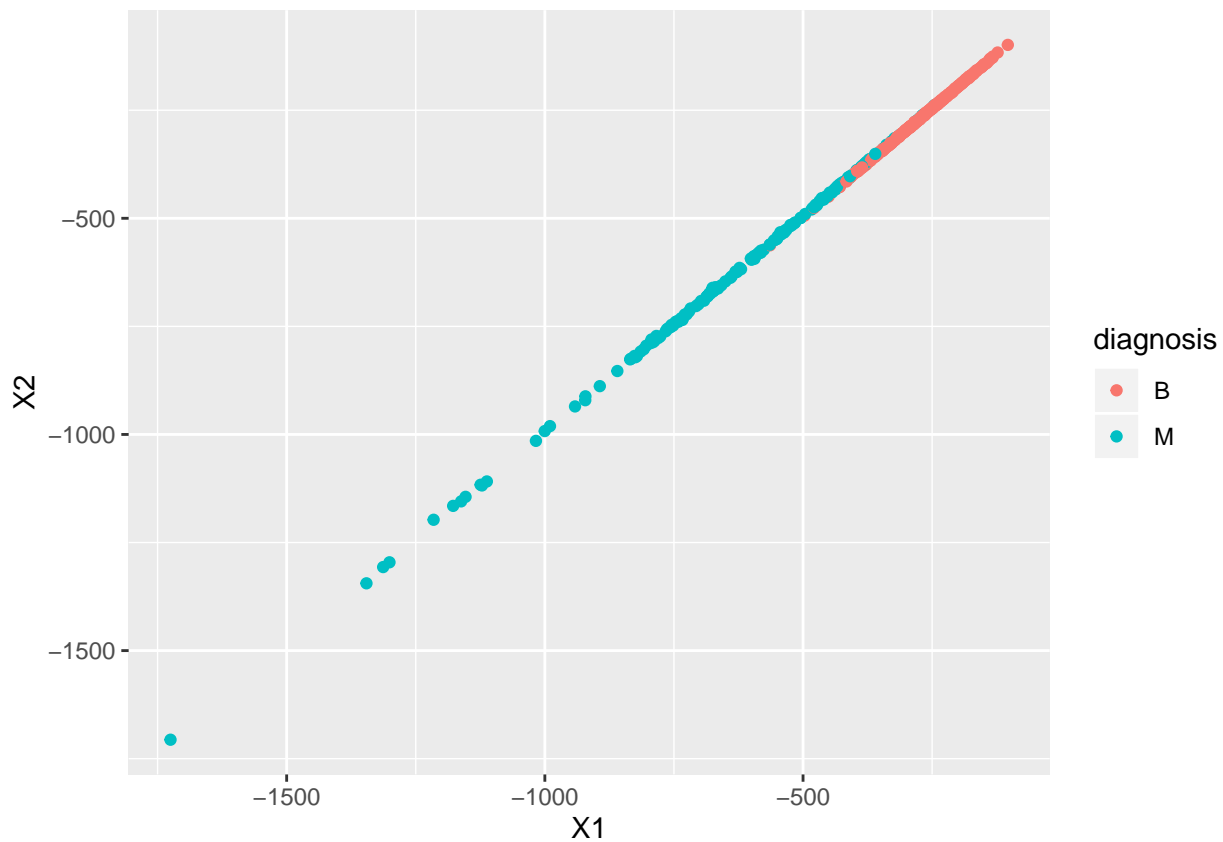
Apply to dataset

```

pca_result <- pca(as.matrix(training_data), 2)
pca_reduced_training_data <- data.frame(cbind(pca_result$reduced_data, training_classes))

ggplot(data=pca_reduced_training_data, aes(x=X1, y=X2)) + geom_point(aes(colour=diagnosis))

```



## tSNE

TODO: try different perplexity parameters

```
{r} #reduced_training_data <- tsne::tsne(training_data) # #reduced_train
<- data.frame(cbind(reduced_training_data, training_classes))
#ggplot(data=reduced_training_data, aes(x=X1, y=X2)) + geom_point(aes(co
#
```

- Correlation Feature Selection
- LDA

## Classification

To solve the problem of finding a SVM like classifier for non-separable data we must permit a certain number of points to violate the boundaries set however this number and the amount they violate the constraints by must be as small as possible. To formulate this we introduce a variable  $\epsilon_i$  for each data point into the objective functions and the constraints leading to the optimisation problem:

$$\begin{aligned} \min_{w, \epsilon_i} \quad & \frac{1}{2} w^T w + C \sum_{i=0}^n \epsilon_i \\ \text{such that } & w \cdot x_i + b + \epsilon_i > 1 \text{ if } y_i = 1 \\ & \text{and } w \cdot x_i + b + \epsilon_i < -1 \text{ if } y_i = -1 \end{aligned}$$

Note that we have swapped the sign of the  $b$  term in the equation for the hyperplane because I implemented it this way before realising they were different and am lazy.

As the above problem is convex (as it is quadratic) and Slater's condition holds then strong duality holds and we can take the Lagrangian of the optimisation problem and consider the result of the KKT conditions. By doing so we can reformulate the optimisation problem as the dual problem:

$$\begin{aligned} \min_{\lambda} \quad & \frac{\bar{\lambda} X X^T \bar{\lambda}^T}{4} + \lambda^T \mathbf{1} \\ \text{such that } & 0 \leq \lambda_i \leq C \\ & \text{and } \sum_i^n \lambda_i y_i = 0 \end{aligned}$$

where

$$X = \begin{pmatrix} x_1 \\ \vdots \\ x_n \end{pmatrix} \text{ and } \bar{\lambda} = [\lambda_1 \cdot y_1, \dots, \lambda_n \cdot y_n] \text{ and } \mathbf{1} = [1, \dots, 1] \in \mathbb{R}^n$$

As before we have to massage this optimisation problem into one that can be solved using `solve.QP`. In this formulation

$$d = \mathbf{1}$$

and

$$D = \begin{pmatrix} y_1 & 0 & \dots & 0 \\ 0 & y_2 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & y_n \end{pmatrix} X X^T \begin{pmatrix} y_1 & 0 & \dots & 0 \\ 0 & y_2 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & y_n \end{pmatrix}$$

$A$  and  $b_0$  require slightly more manipulation this time around with

$$A = \begin{pmatrix} y_1 & y_2 & \dots & y_n \\ & I & & \\ & & -I & \end{pmatrix}$$

and

$$b_0 = \begin{pmatrix} 0 \\ \mathbf{0} \\ -C \end{pmatrix}$$

where

$$\mathbf{0} = [0, \dots, 0]^T \in \mathbb{R}^n$$

and

$$C = [C, \dots, C]^T \in \mathbb{R}^n$$

The code for this applied to the non-separable data can be found below.

```
C <- 1

X <- as.matrix(combined_class)[,1:2]
y <- as.matrix(combined_class)[,3]
Dmat2 <- diag(y) * X %*% t(X) %*% diag(y)
diag(Dmat2) <- diag(Dmat2) + 1e-11
dv2 <- rep(1, 30)

A2 <- rbind( y,diag(30))
A2 <- rbind(A2, -1*diag(30))

bv2 <- c(c(0), rep(0, 30), rep(-C, 30) )
model <- solve.QP(Dmat2, dv2, t(A2), bv2, meq = 1)
```

In order to recover  $w$  and  $b$  from  $\lambda$  we use the relationship

$$w = \sum_{i=0}^{n-1} \lambda_i x_i^T y_i$$

and

$$b = \text{mean}\left(\sum_{i=0}^k y_i - w \cdot x_i\right) \cdot \forall i. 0 < \lambda_i < C$$

Which can be made as functions in R as so:

```
calculate_b <- function(w, X, y, a, C) {
  ks <- sapply(a, function(x){return(x > 0 && x < C)})
  indices <- which(ks)
  sum_bs <- 0
  for(i in indices) {
    sum_bs <- sum_bs + (y[i] - w %*% X[i,])
  }
  return(sum_bs / length(indices))
}
```



```

}

recover_w <- function(a, y, X){
  colSums(diag(a) %*% diag(y) %*% X)
}

```

We can see the results of using the dual regression below

## SVM

```

soft_margin_svm_plotter <- function(w, b) {
  plotter <- function(x) {
    return(1/w[2] * -(b + (w[1]*as.numeric(x))))
  }
  return(plotter)
}

```

```

factor_to_label <- function(x) {
  if(as.character(x) == "M") {
    return(1)
  }
  else {
    return(-1)
  }
}

```

```

label_to_factor <- function(x) {
  if(x == 1) {
    return(as.factor("M"))
  }
  else{
    return(as.factor("B"))
  }
}

numeric_test_labels <- apply(test_classes, 1, factor_to_label)
numeric_training_labels <- apply(training_classes, 1, factor_to_label)

```

Use PCA then do SVM

```

model <- svm(X=pca_result$reduced_data,
             classes=numeric_training_labels,
             C=100000, margin_type='soft',
             kernel_function = linear_kernel,
             feature_map = linear_basis_function)

reduced_prediction_fn <- model$prediction_function

pca_reduced_prediction_fn <- function(x) {
  p <- x %*% pca_result$reduction_matrix
  reduced_prediction_fn(t(p))
}

```

```

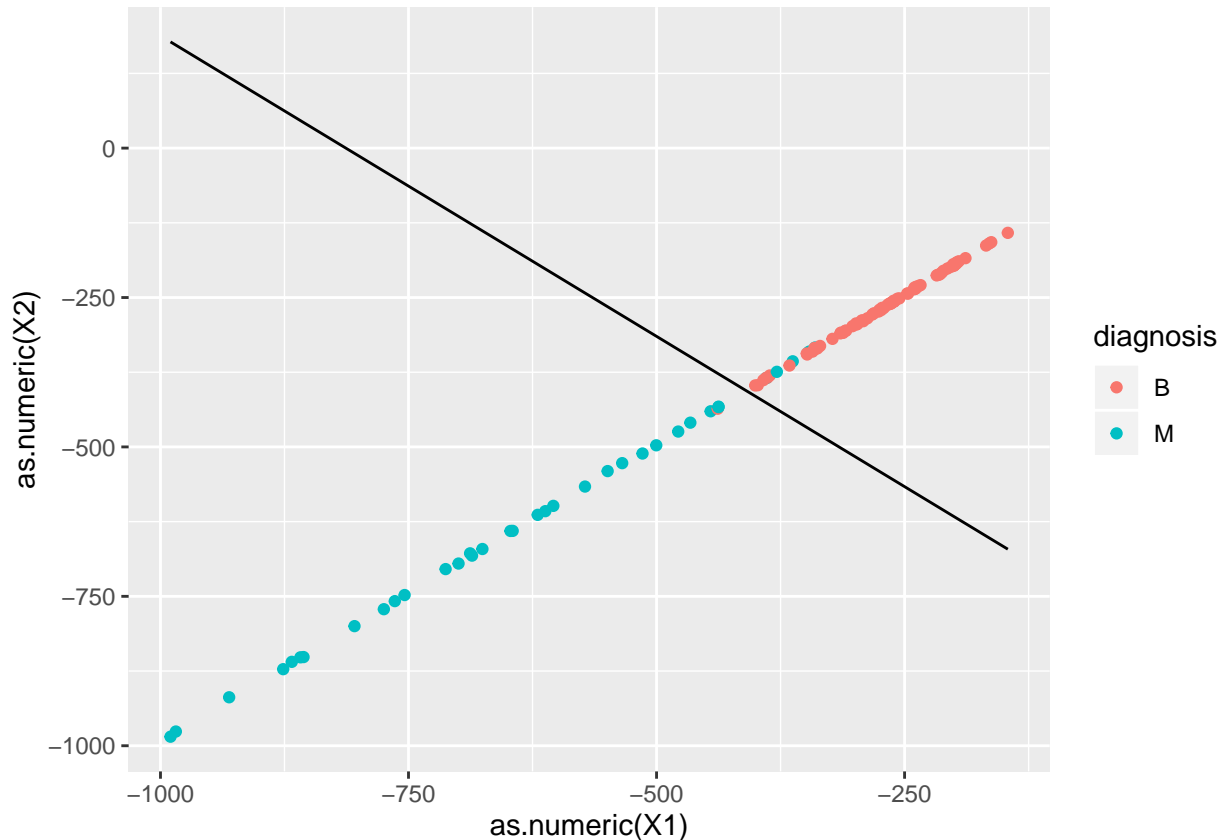
predictions_svm <- apply(as.matrix(test_data),1, pca_reduced_prediction_fn)
accuracy_calc(numeric_test_labels, predictions_svm)

## [1] 93.85965

svm_plotter <- soft_margin_svm_plotter(model$params$w, model$params$b)
embedded_test_data <- data.frame(cbind(as.matrix(test_data) %*% pca_result$reduction_matrix), test_class)

ggplot(embedded_test_data, aes(x=as.numeric(X1), y=as.numeric(X2))) +
  geom_point(aes(colour=diagnosis)) +
  stat_function(fun=svm_plotter)

```



## Naive Bayes

### Mathematical setting

Let  $y$  be the class label that we want to assign to an observation  $\mathbf{x} = (x_1, \dots, x_d)$ , where  $x_1, \dots, x_d$  are the features. The probability of an observation having label  $y$  is given by Bayes rule,

$$\begin{aligned}
 P(y|x_1, \dots, x_d) &= \frac{P(x_1, \dots, x_d|y_k)P(y)}{P(x_1, \dots, x_d)} \\
 &\propto P(x_1, \dots, x_d|y_k)P(y).
 \end{aligned}$$

The prior class probability  $P(y)$  can be easily obtained by the proportion of observation that are in the given class.

The main assumption is that every feature is conditionally independent given the class label  $y$ . The reason why this classifier is called *naive* is that very often this assumption is not actually realistic.

This assumption simplifies the posterior to

$$P(y|x_1, \dots, x_d) \propto P(y) \prod_{i=1}^d P(x_i|y).$$

There are various types of Naive Bayes classifiers based on the type of features. In our case, since we have continuous variables we assume that all features are normally distributed. Therefore, the conditional probabilities can be calculated as

$$P(x_i|y) = \frac{1}{\sqrt{2\pi\sigma_y^2}} \exp\left(-\frac{(x_i - \mu_y)^2}{2\sigma_y^2}\right)$$

Finally, to assign the class to an observation we use the Maximum A Posteriori decision rule. For every observation, we pick the class the has the highest probability

$$y = \underset{y}{\operatorname{argmax}} P(y) \prod_{i=1}^d P(x_i|y).$$

## Implementation

*Here are some code snippets just to illustrate how these theoretical aspects are implemented. The full code can be found in the package.*

The observations are stored as rows in  $X$  and the corresponding class labels are entire rows in the column matrix  $y$ .

First we calculate the prior class probabilities based on the number of observations in each class.

```
n <- dim(X)[1]
d <- dim(X)[2]
classes <- sort(unique(y)[, 1])
k <- length(classes)

prior <- rep(0, k)
for (i in 1:k) {
  prior[i] <- sum(y == classes[i]) / n
}
```

Then we create an array of the mean and sd of the data split by classes and features.

```
summaries <- array(rep(1, d * k * 2), dim = c(k, d, 2))
for (i in 1:k) {
  X_k <- X[which(y == (i - 1)), ]
  summaries[i, , 1] <- apply(X_k, 2, mean)
  summaries[i, , 2] <- apply(X_k, 2, sd)
}
```

Finally, the predictions are obtained by taking the largest posterior class probability. Note that in order to avoid underflow, we take the maximum of the *log* posterior class probabilities.

```
probs <- matrix(rep(0, n * k), nrow = n)
for (obs in 1:n) {
  for (class in 1:k) {
    class_prob <- log(prior[class])
```

```

    for (feat in 1:d) {
      mu <- summaries[class, feat, 1]
      sd <- summaries[class, feat, 2]
      cond <- dnorm(x_new[obs, feat], mu, sd, log = TRUE)
      class_prob <- class_prob + cond
    }
    probs[obs, class] <- class_prob
  }
}

pred <- apply(probs, 1, which.max)

```

## Fit model to dataset

```

install_github("andreabecsek/NaiveBayes")
library(NaiveBayes)

```

```

levels(training_classes$diagnosis) <- c(0,1)
training_classes %<>% as.matrix
mode(training_classes) <- 'numeric'

levels(test_classes$diagnosis) <-c(0,1)
test_classes %<>% as.matrix
mode(test_classes) <- 'numeric'

```

Fit the Naive Bayes model to the data, calculate predictions and check the accuracy using.

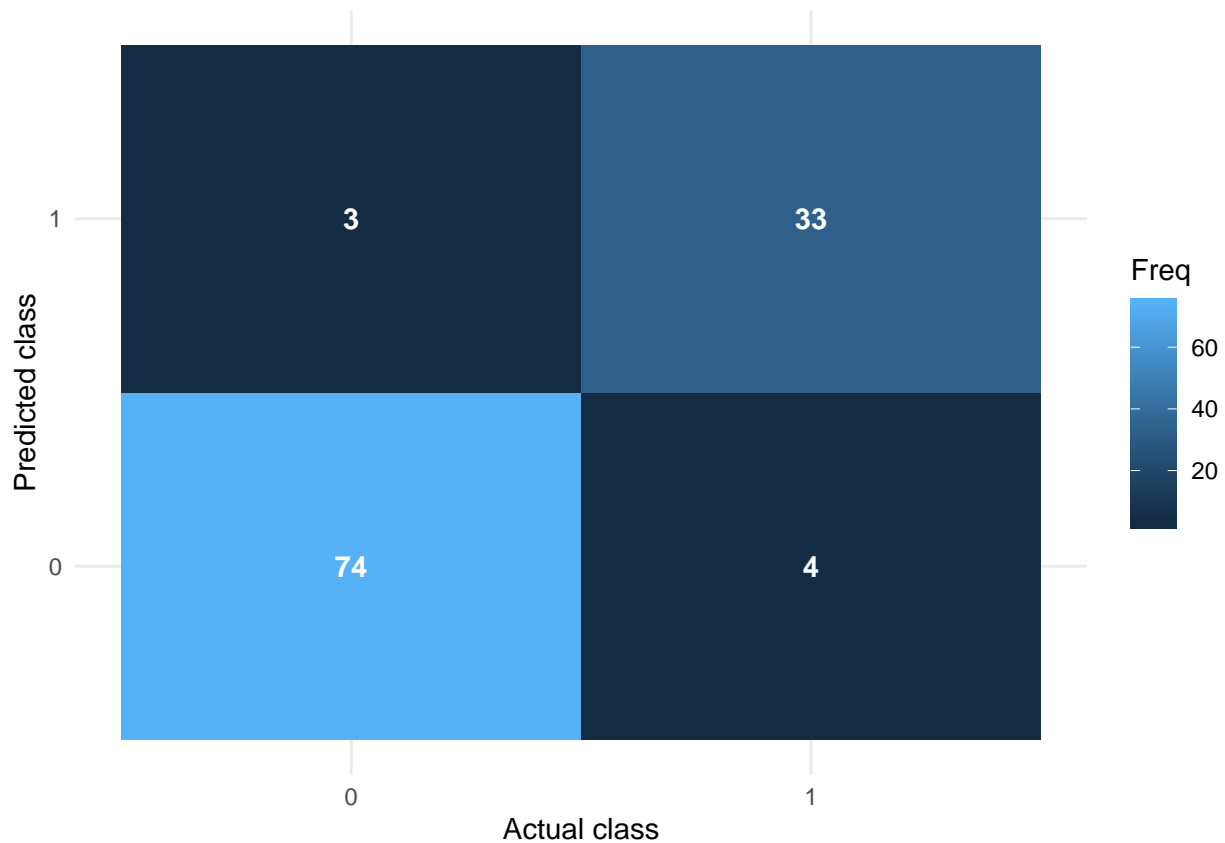
```

model_naive <- naive_bayes(training_data,training_classes)

predictions_naive <- predict(model_naive,as.matrix(test_data))

confusion_plot(test_classes,predictions_naive)

```



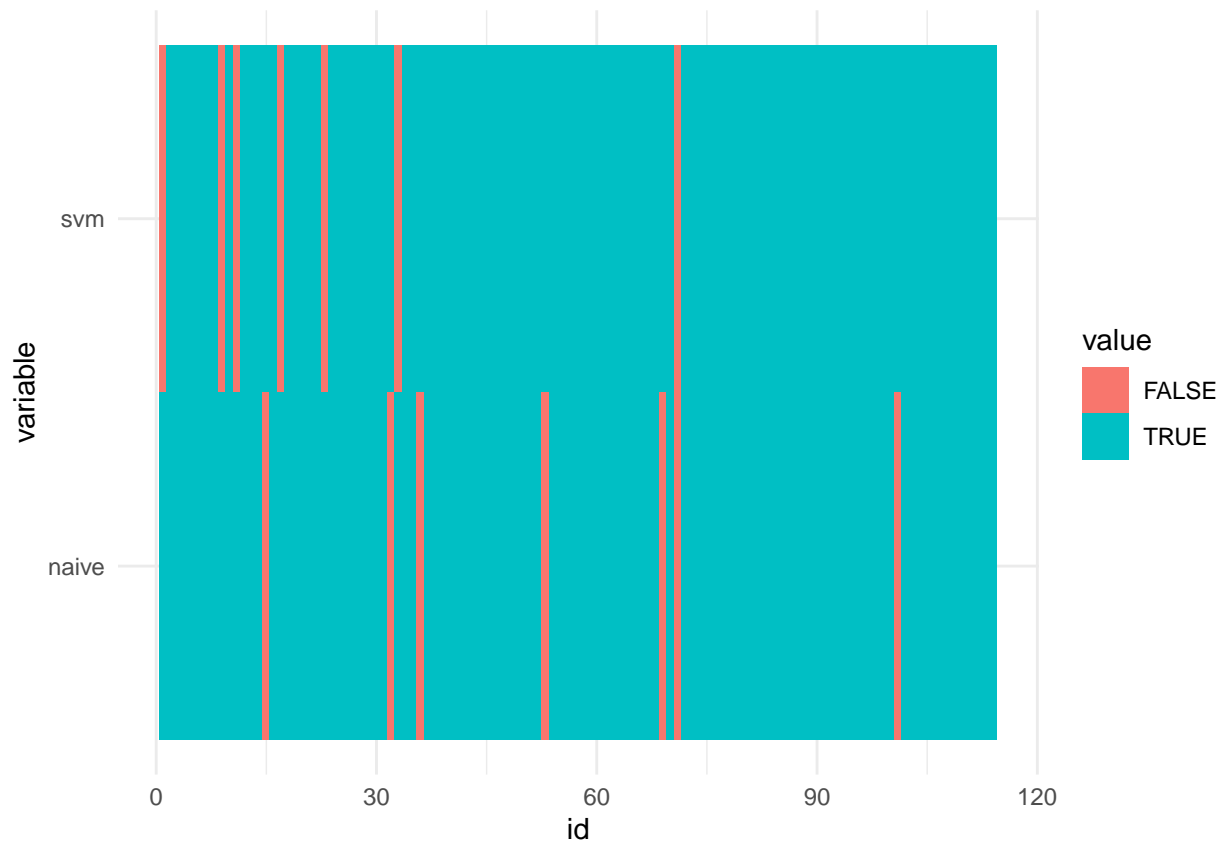
Merge all predictions

```
id <- seq(length(id_test))
all_predictions <- cbind(id,numeric_test_labels,predictions_naive,predictions_svm)
colnames(all_predictions) <- c('id','actual','naive','svm')
all_predictions[all_predictions==1] <-0
all_predictions %<>% as.data.frame()
```

```
errors <- all_predictions %>%
  mutate(naive=naive==actual) %>%
  mutate(svm=svm==actual) %>%
  dplyr::select(-actual)
```

```
a <- errors %>%
  melt(id='id')

ggplot(a,aes(x=id,y=variable,fill=value))+
  geom_raster()+
  theme_minimal()
```



TODO: analyse results

- Naive Bayes
- Logistic Regression

## Conclusion

- Evaluation of results
- Discuss outliers

TODO: create outlier plot