

TDDE15_Lab4_group

Alice Velander(alive213)

Oskar Hidén (oskhi827)

David Björelind (davbj395)

Samuel Persson (sampe028)

10/20/2020

Part 1 - Implementing GP Regression

```
#install.packages('kernlab')
library(kernlab)

# Computing the whole covariance matrix K from the kernel.
cov_function = function(X, Xstar){
  return(kernelMatrix(kernel = SEkernel, x=X, y=Xstar))
}

# Algorithm 2.1 sunns in  $n^3/6$  instead of  $O(n^3)$ . GP can be estimated faster,  $O(n)$  with eg. KISS-GP
posteriorGP = function(X_input, y_targets, k_cov_function, sigmaNoise=1, XStar){
  A = k_cov_function(X_input, X_input)
  A = A + diag(length(X_input))*sigmaNoise^2
  L = t(chol(A)) # chol Returns t(L)
  L_y = solve(L, y_targets)
  alpha = solve(t(L), L_y)

  k_star = k_cov_function(X_input, XStar)
  f_star = t(k_star)%*%alpha

  v = solve(L,k_star)
  V_f_star = k_cov_function(XStar, XStar) - t(v)%*%v
  return(list("mean"=f_star, "cov" = V_f_star))
}

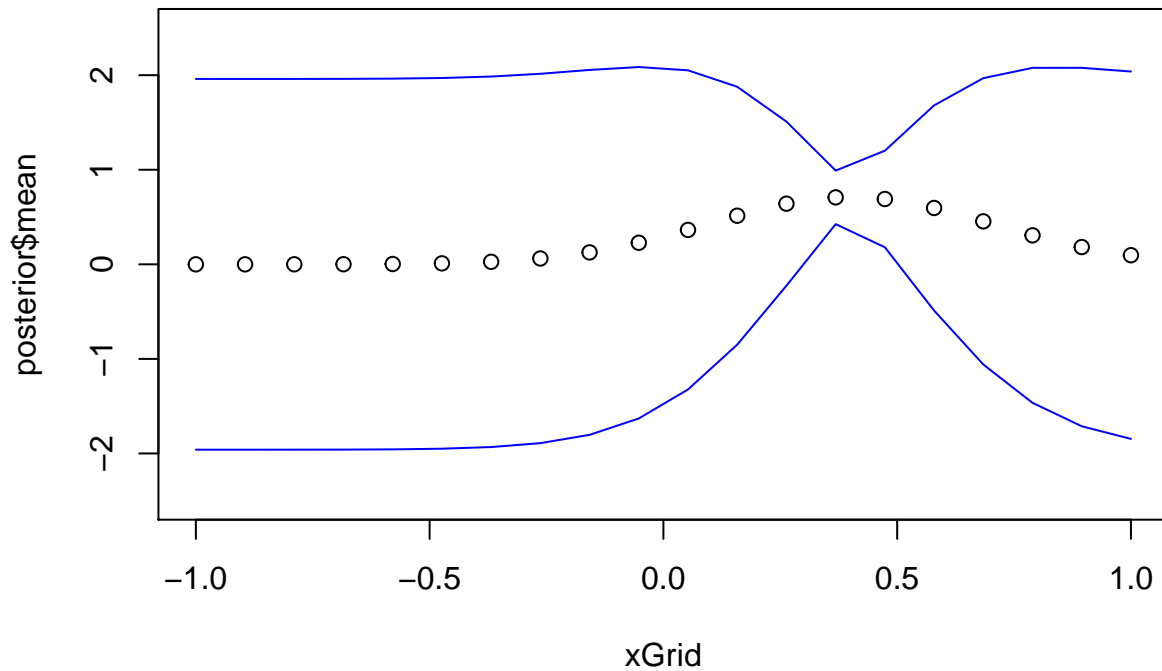
sigma_f = 1
ell = 0.3
SEkernel <- rbfdot(sigma = 1/(2*ell^2)) # Reparametrize the rbfdot (which is the SE kernel) in kernlab.
sigma_n = 0.1

# 2 - Update prior with a single observation
x=0.4
y=0.719
xGrid <- seq(-1,1,length=20) # x_star

posterior = posteriorGP(x, y, cov_function, sigma_n, xGrid)

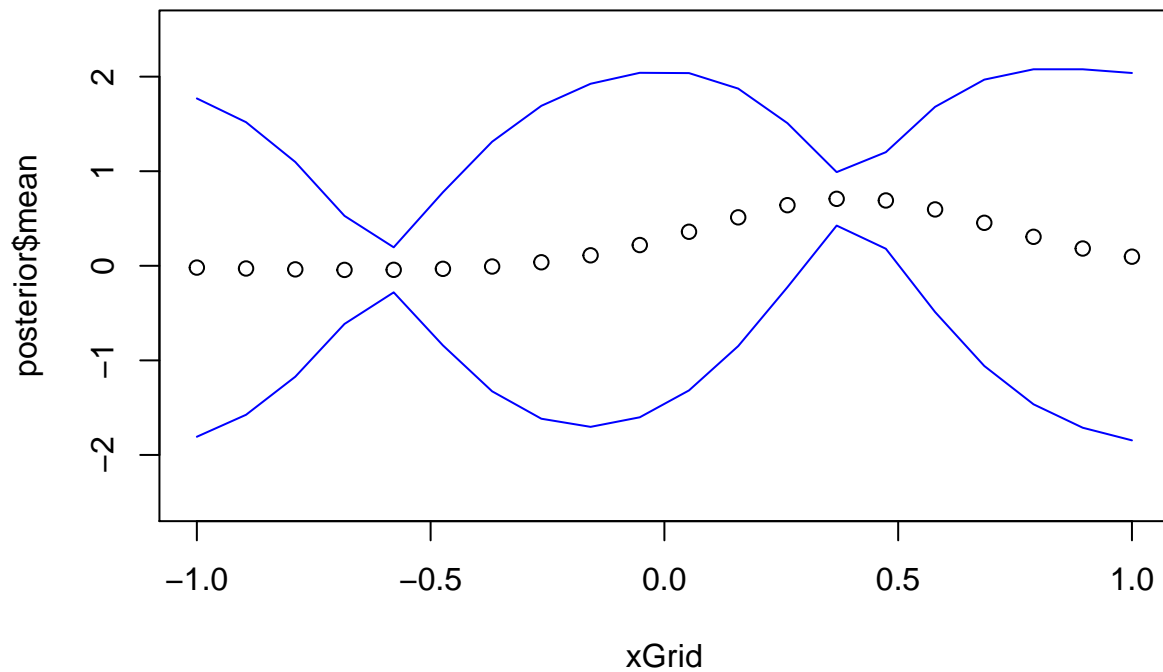
plot(xGrid, posterior$mean, ylim = c(-2.5,2.5)) # posterior mean
```

```
std_dev = sqrt(diag(posterior$cov))
points(xGrid, posterior$mean + 1.96*std_dev, col="blue", type="l")
points(xGrid, posterior$mean - 1.96*std_dev, col="blue", type="l")
```



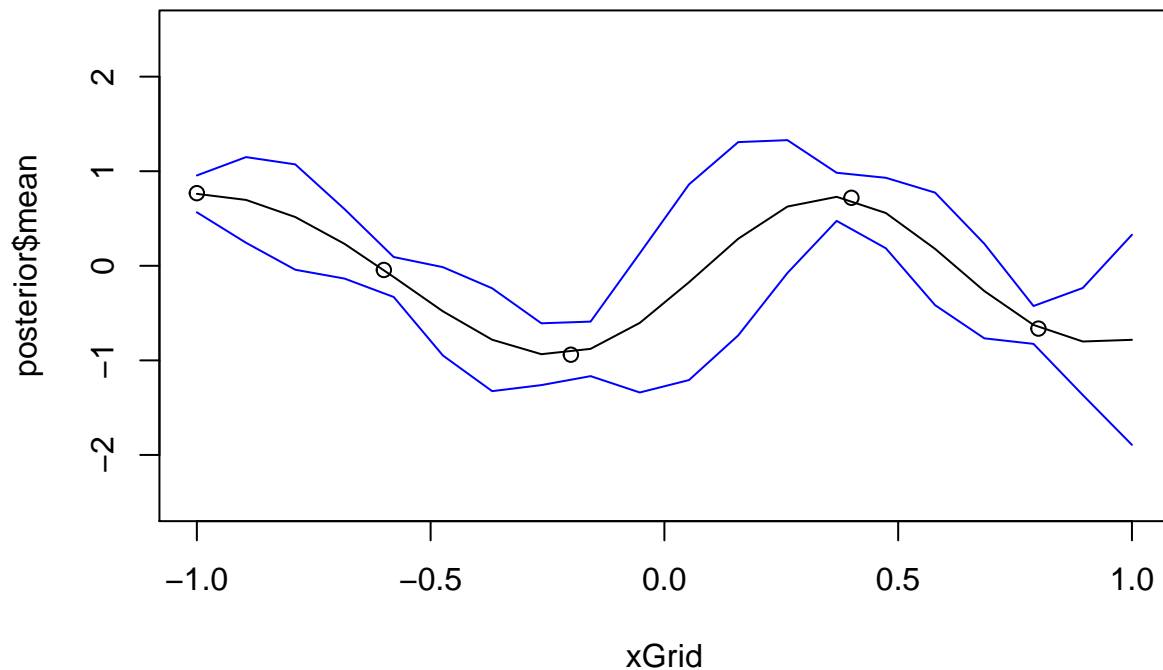
```
# 3 - Update posterior with another point. (by recalculating posterior with Algorithm 2.1)
x = c(0.4, -0.6)
y = c(0.719, -0.044)
posterior = posteriorGP(x, y, cov_function, sigma_n, xGrid)

plot(xGrid, posterior$mean, ylim = c(-2.5, 2.5)) # posterior mean
std_dev = sqrt(diag(posterior$cov))
points(xGrid, posterior$mean + 1.96*std_dev, col="blue", type="l")
points(xGrid, posterior$mean - 1.96*std_dev, col="blue", type="l")
```



```
# 4 - Calculate the posterior with five observations.
x = c(-1.0, -0.6, -0.2, 0.4, 0.8)
y = c(0.768, -0.044, -0.940, 0.719, -0.664)
posterior = posteriorGP(x, y, cov_function, sigma_n, xGrid)

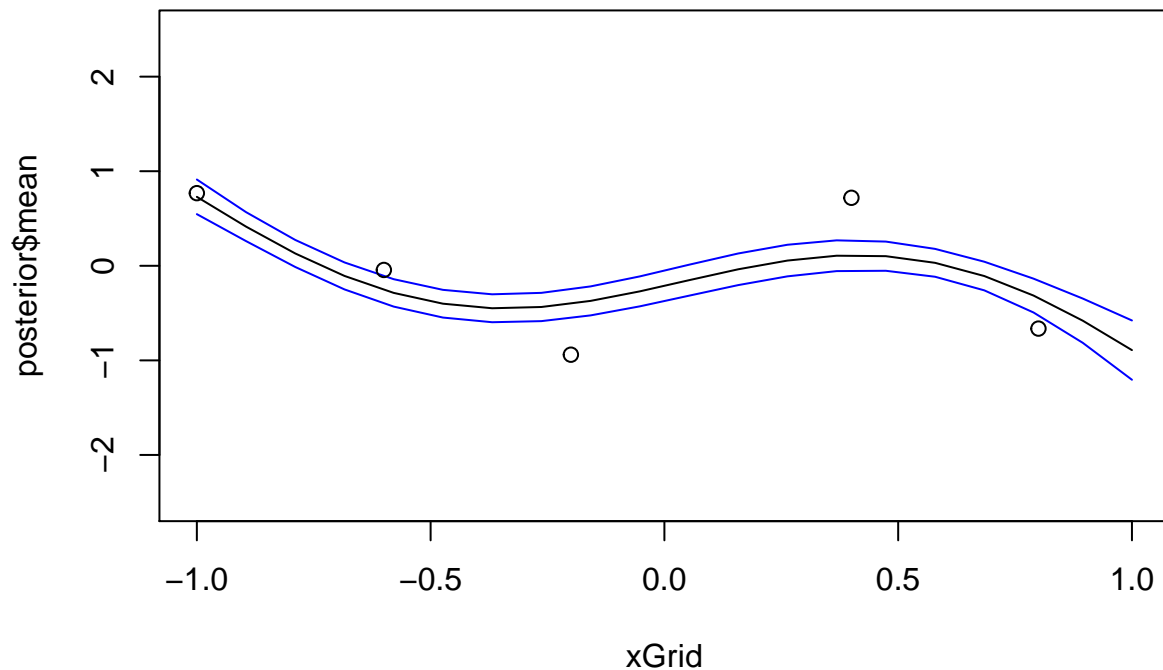
plot(xGrid, posterior$mean, ylim = c(-2.5, 2.5), type="l") # posterior mean
std_dev = sqrt(diag(posterior$cov))
points(xGrid, posterior$mean + 1.96*std_dev, col="blue", type="l")
points(xGrid, posterior$mean - 1.96*std_dev, col="blue", type="l")
points(x, y)
```



```
# 5 - Redo (4) with the same sigma_f but with a higher ell.
sigma_f = 1
ell = 1
SEkernel <- rbfdot(sigma = 1/(2*ell^2))

x = c(-1.0, -0.6, -0.2, 0.4, 0.8)
y = c(0.768, -0.044, -0.940, 0.719, -0.664)
posterior = posteriorGP(x, y, cov_function, sigma_n, xGrid)

plot(xGrid, posterior$mean, ylim = c(-2.5,2.5), type="l") # posterior mean
std_dev = sqrt(diag(posterior$cov))
points(xGrid, posterior$mean + 1.96*std_dev, col="blue", type="l")
points(xGrid, posterior$mean - 1.96*std_dev, col="blue", type="l")
points(x,y)
```



The two plots has the same `sigma_f` but different `ell`-values. The last plot has a smoother mean prediction, with a higher `ell` value. A higher `ell` gives a higher covariance for two points that are further apart. When we used more training data, we get a narrower confidence intervall. We are also more confident in our posterior distribution around our training points.

Part 2 - GP Regression with kernlab

```
temperature = read.csv("https://github.com/STIMALiU/AdvMLCourse/raw/master/GaussianProcess/Code/TempTull.
time = (1:2190)
nr_year = 2190/365
day = rep((1:365), nr_year)

reduce_data = function(array , nth){
  array[seq(1, length(array), nth)]
}

temp = reduce_data(temperature$temp,5)
time = reduce_data(time, 5)
day = reduce_data(day, 5)

set_se_kernel = function(ell, sigma_f){
  se_kernel = function(x , y){
    r_square = sum((x-y)*(x-y)) # Euclidian distance^2
    return(sigma_f^2*exp(-r_square/(2*ell^2)))
  }
}
```

```

}
class(se_kernel) <- "kernel"
return(se_kernel)
}

```

```
se_kernel = set_se_kernel(ell=1, sigma_f=1)
```

```

# Evaluate kernel in x=1 and x'=2
se_kernel(1,2)

```

```
## [1] 0.6065307
```

```

k = function(x, x_star){
  return(kernelMatrix(kernel = se_kernel, x=x, y=x_star))
}

```

```

# Evaluate the covariance matrix, k
x = c(1, 3, 4)
x_star = c(2, 3, 4)
k(x, x_star)

```

```

## An object of class "kernelMatrix"
##          [,1]      [,2]      [,3]
## [1,] 0.6065307 0.1353353 0.0111090
## [2,] 0.6065307 1.0000000 0.6065307
## [3,] 0.1353353 0.6065307 1.0000000

```

```

# 2 - Find f* with function: gausspr(kernlab)
sigma_f = 20
ell = 0.2
se_kernel = set_se_kernel(ell=ell, sigma_f=sigma_f)

```

```

# Find sigma_n
quadratic_reg = lm(temp ~ time + I(time^2))
sigma_n = sd(quadratic_reg$residuals)

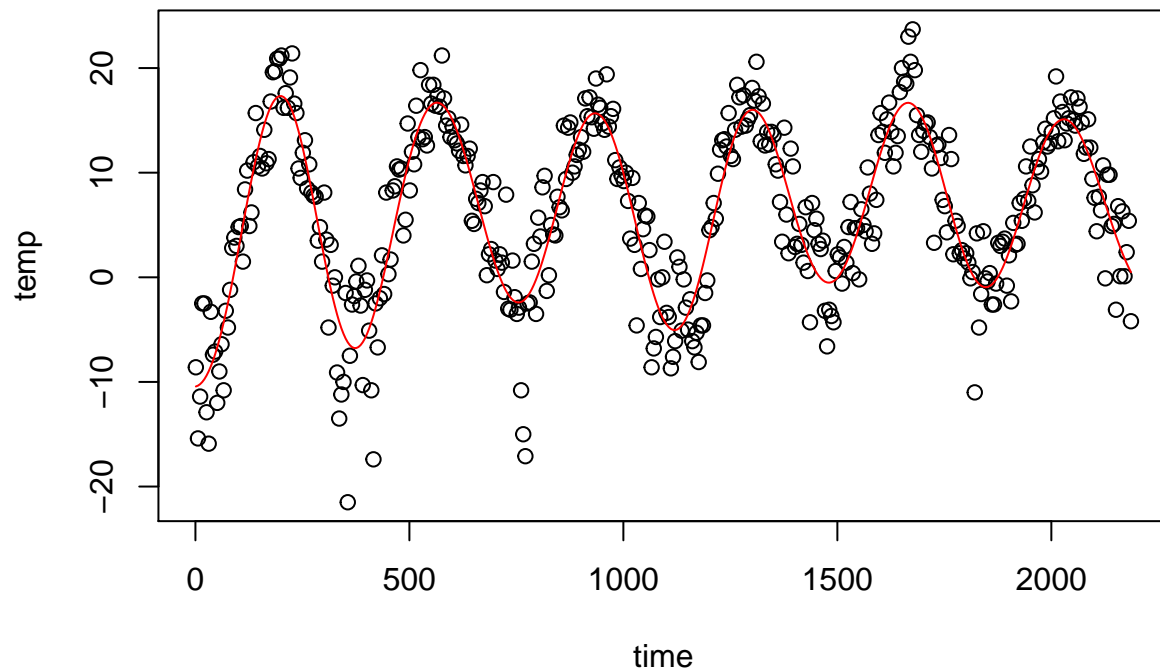
```

```
gp_model = gausspr(x = time, y = temp, type="regression", kernel=se_kernel, var = sigma_n^2, variance.m
```

```

mean_pred = predict(gp_model, time)
plot(time, temp)
points(time, mean_pred, type = "l", col="red")

```



```
#lines(time, mean_pred + 1.96 * predict(gp_model, time, type="sdeviation")) # ---- wrong due to scaling
```

```
# 3 - Scale and calculate with Algorithm 2.1
```

```
mean_temp = mean(temp)
```

```
sd_temp = sd(temp)
```

```
time_scale = scale(time)
```

```
temp_scale = scale(temp)
```

```
# Find sigma_n
```

```
quadratic_reg = lm(temp_scale ~ time_scale + I(time_scale^2))
```

```
sigma_n = sd(quadratic_reg$residuals)
```

```
post_gp = posteriorGP(time_scale, temp_scale, k, sigma_n, time_scale)
```

```
st_dev_new = sqrt(diag(post_gp$cov))*sd_temp
```

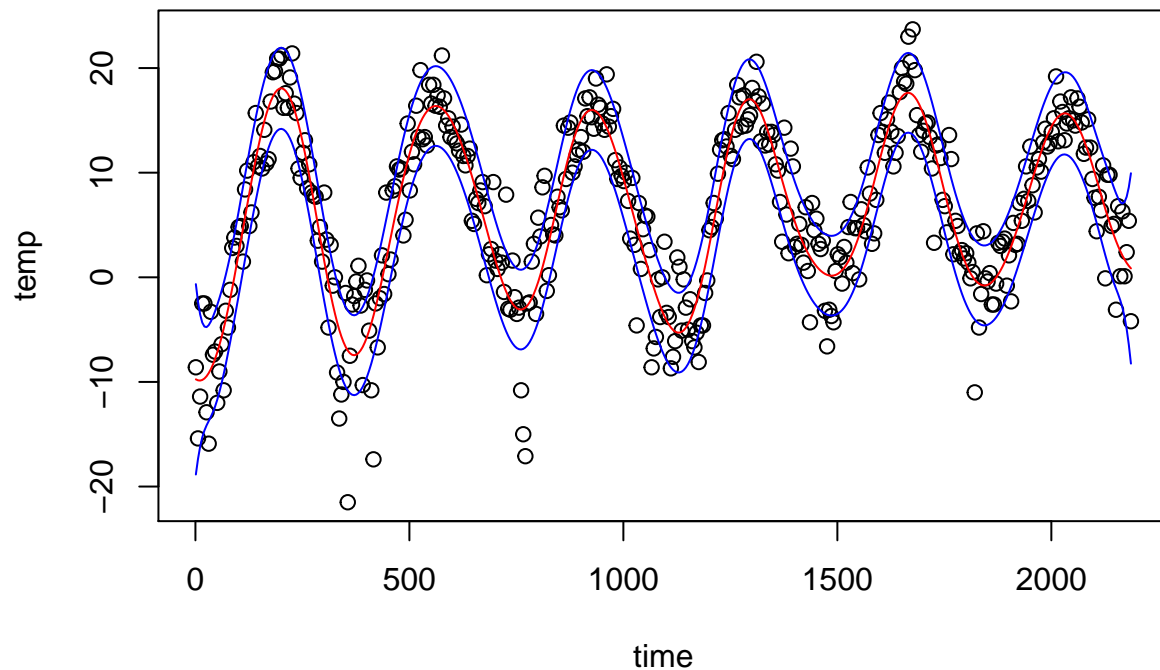
```
mean_pred_new = post_gp$mean*sd_temp+mean_temp
```

```
plot(time, temp)
```

```
lines(time, mean_pred_new, col="red")
```

```
lines(time, mean_pred_new+1.96*st_dev_new, col="blue")
```

```
lines(time, mean_pred_new-1.96*st_dev_new, col="blue")
```

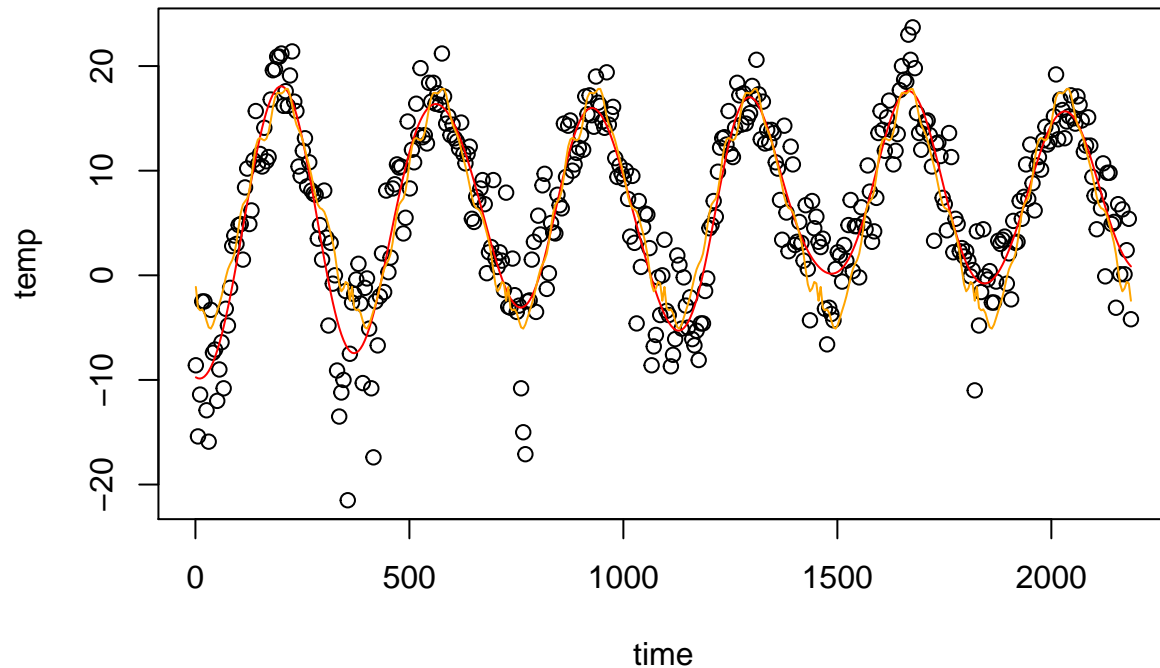


```
# 4 - use of Day
day_scale = scale(day)

#Find sigma_n
quadratic_reg = lm(temp_scale ~ day_scale + I(day_scale^2))
sigma_n = sd(quadratic_reg$residuals)

post_gp = posteriorGP(day_scale, temp_scale, k, sigma_n, day_scale)
st_dev_day = sqrt(diag(post_gp$cov))*sd_temp
mean_pred_day = post_gp$mean*sd_temp+mean_temp

# old plot
plot(time, temp)
lines(time, mean_pred_new, col = "red")
#lines(time, mean_pred_new+1.96*st_dev_new, col="blue")
#lines(time, mean_pred_new-1.96*st_dev_new, col="blue")
# new lines from day
lines(time, mean_pred_day, col = "orange")
```

```
#lines(time, mean_pred_day+1.96*st_dev_day, col="green")
#lines(time, mean_pred_day-1.96*st_dev_day, col="green")
```

When predicting using time and square exponential kernel we have a low correlation between the same date between two consecutive years. But when using day and square exponential kernel we assume the same correlation between the same date, no matter which year that date is from. And we get a prediction that does not take the long term change (ex. warmer winters) into account. Day is not that smooth when we change year, because day=1 and day=365 will not get a strong correlation, even though they are close.

```
# 5 - Generalization periodic kernel
sigma_f = 20
ell_1 = 1
ell_2 = 10
d = 365/sd(time)

set_gen_kernel = function(ell_1, ell_2, sigma_f, d){
  gen_kernel = function(x, y){
    r_square = sum((x-y)*(x-y)) #euclidian distance ^2
    return(sigma_f^2 * exp(-(2*sin(pi*sqrt(r_square)/d)^2)/(ell_1^2)) * exp(-r_square/(2*ell_2^2)))
  }
  class(gen_kernel) <- "kernel"
  return(gen_kernel)
}

k_gen = function(x, x_star){
```

```

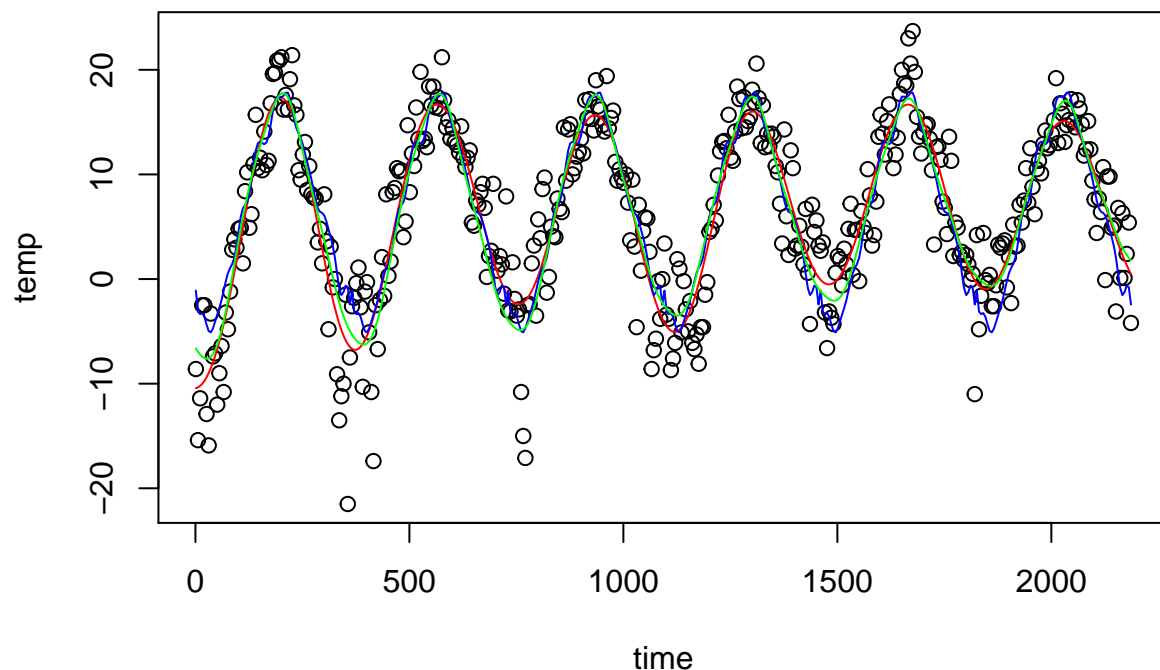
    return(kernelMatrix(kernel = gen_kernel, x=x, y=x_star))
}

# Set sigma_n
quadratic_reg = lm(temp ~ time + I(time^2))
sigma_n = sd(quadratic_reg$residuals)

gen_kernel = set_gen_kernel(ell_1, ell_2, sigma_f, d)
gp_model = gausspr(x = time, y = temp, type="regression", kernel=gen_kernel, var = sigma_n^2)

mean_pred_gen = predict(gp_model, time)
plot(time, temp)
points(time, mean_pred, type="l", col="red")
points(time, mean_pred_day, type = "l", col="blue")
points(time, mean_pred_gen, type = "l", col="green")

```



The periodic kernel will take both time and periodisation into account. Two equal dates will have a higher covariance if they are from two consecutive years. This means that we manage to capture the long term trend better, because of the second “exp()” in the kernel.

Part 3 - GP Classification with kernlab

```

#install.packages("AtmRay") # To make 2D grid like in Matlab's meshgrid.
library(AtmRay)

```

```

data <- read.csv("https://github.com/STIMaLiU/AdvMLCourse/raw/master/GaussianProcess/Code/banknoteFraud")
names(data) <- c("varWave", "skewWave", "kurtWave", "entropyWave", "fraud")
data[,5] <- as.factor(data[,5])

set.seed(111)
SelectTraining <- sample(1:dim(data)[1], size = 1000, replace = FALSE)
train = data[SelectTraining,]

# 1 - predicition with two covariats.
gp_fraud = gausspr(fraud ~ varWave + skewWave, data=train, type="classification")

```

```
## Using automatic sigma estimation (sigest) for RBF or laplace kernel
```

```

fraud_pred = predict(gp_fraud, train[,1:4])
# Confusion matrix
table = table(fraud_pred, train[,5])
missclass = 1 - sum(diag(table))/sum(table)
# Create the basis of the plot
x1 = seq(min(train[,1]), max(train[,1]), length=100)
x2 = seq(min(train[,2]), max(train[,2]), length=100)
grid_points = meshgrid(x1, x2)
grid_points = cbind(c(grid_points$x), c(grid_points$y))

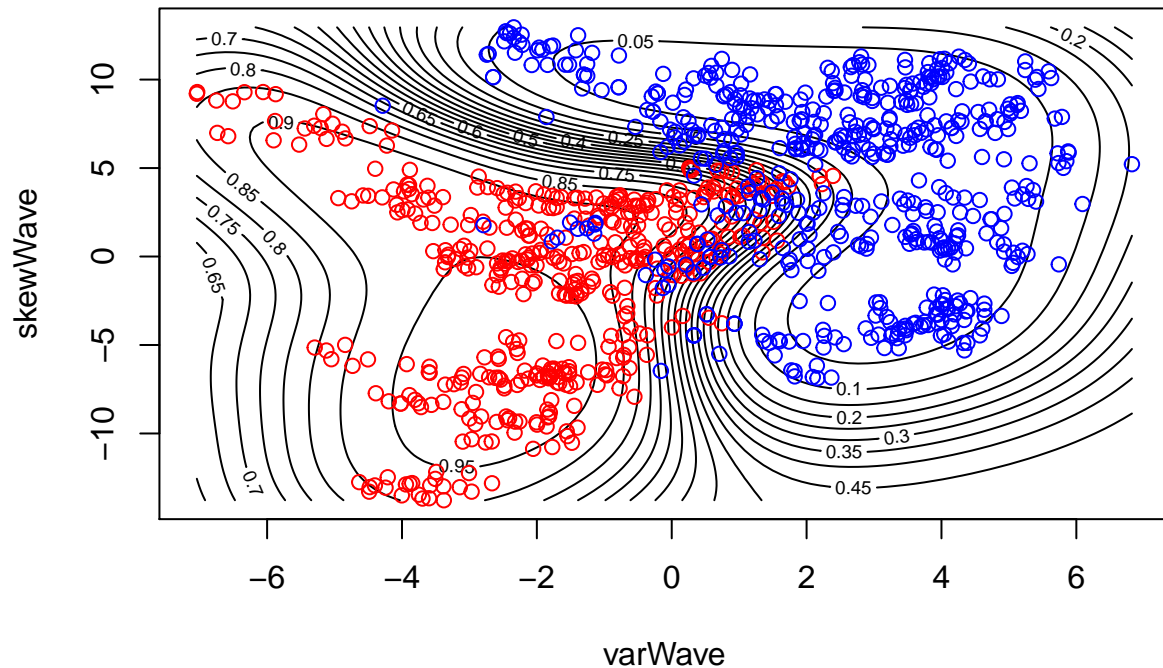
grid_points = data.frame(grid_points)
names(grid_points) = names(train)[1:2]

prob_pred = predict(gp_fraud, grid_points, type="probabilities")

# Plot the plot
contour(x1,x2,matrix(prob_pred[,2],100,byrow = TRUE), 20, xlab = "varWave", ylab = "skewWave", main = '1')
points(train[train[,5]==1,1],train[train[,5]==1,2],col="red")
points(train[train[,5]==0,1],train[train[,5]==0,2],col="blue")

```

Fraud is red



```
cat("Confusion matrix: ")
```

```
## Confusion matrix:
```

```
table
```

```
##
## fraud_pred  0   1
##            0 503  18
##            1  41 438
```

```
cat("missclassification rate: ", missclass)
```

```
## missclassification rate:  0.059
```

```
# 2 apply model on test data.
test = data[-SelectTraining,]
test_pred = predict(gp_fraud, test[,1:4])
table = table(test_pred, test[,5])
missclass = 1 - sum(diag(table))/sum(table)

cat("Confusion matrix:")
```

```
## Confusion matrix:
```

```
table
```

```
##  
## test_pred    0    1  
##           0 199    9  
##           1  19 145
```

```
cat("missclassification rate: ", missclass)
```

```
## missclassification rate:  0.07526882
```

```
# 3 training model with all 4 covariats
```

```
gp_fraud = gausspr(fraud ~ varWave + skewWave + kurtWave + entropyWave, data=train, type="classification")
```

```
## Using automatic sigma estimation (sigest) for RBF or laplace kernel
```

```
test_pred = predict(gp_fraud, test[,1:4])  
table = table(test_pred, test[,5])  
missclass = 1 - sum(diag(table))/sum(table)
```

```
cat("Confusion matrix:")
```

```
## Confusion matrix:
```

```
table
```

```
##  
## test_pred    0    1  
##           0 216    0  
##           1   2 154
```

```
cat("missclassification rate: ", missclass)
```

```
## missclassification rate:  0.005376344
```

The model with all four covariats is performing better while predicting fraud in the test data. One reason could be that the model now uses more input data (more information), while predicting the data. This is not true for all input data, because there is a risk of overfitting the model to the training data.