Machine learning 2

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notes

```
print("scatter plot für koordinaten und box plot für varianz in long und lat") # why did we want this?

## [1] "scatter plot für koordinaten und box plot für varianz in long und lat"

print("boxplot für distances für jedes modell -> um diese vergleichen zu können")

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print("für finales ergebniss, ein karte, welche ? ist und soll koorinaten anzeigt und mit linien verbin

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```

librarys

```
\#install.packages("installr")
#library("installr")
#install.Rtools()
#install.packages("ggmap")
#install.packages("maptools")
#install.packages("maps")
#install.packages("glmnet")
#install.packages("ISLR")
#TODO the one below necessary?
\#install.packages("rgl")
#install.packages("dplyr")
#install.packages("keras")
#install.packages("tensorflow")
library("glmnet")
## Lade nötiges Paket: Matrix
## Loaded glmnet 4.1-3
library("ggplot2")
library("ggmap")
## Google's Terms of Service: https://cloud.google.com/maps-platform/terms/.
```

Please cite ggmap if you use it! See citation("ggmap") for details.

```
library("maptools")
## Lade nötiges Paket: sp
## Checking rgeos availability: FALSE
## Please note that 'maptools' will be retired by the end of 2023,
## plan transition at your earliest convenience;
## some functionality will be moved to 'sp'.
##
        Note: when rgeos is not available, polygon geometry
                                                                 computations in maptools depend on gpcl
##
        which has a restricted licence. It is disabled by default;
        to enable gpclib, type gpclibPermit()
library("maps")
library("dplyr")
##
## Attache Paket: 'dplyr'
## Die folgenden Objekte sind maskiert von 'package:stats':
##
##
       filter, lag
## Die folgenden Objekte sind maskiert von 'package:base':
##
       intersect, setdiff, setequal, union
library(ISLR)
options(rgl.printRglwidget = TRUE)
library(rgl)
library(keras)
library(tensorflow)
```

Comparison of machine learning algorithms / Introduction / Theory

This report attempts to compare two regression/classification algorithms on its behavior on a specific data set. What should be found out?

Data set

Here should be a bit of a short summary of the data set with some key characteristics of the data set.

What is it about? Which variables are included? What type of variables? What is missing? What about missing values?

The data used in this project is provided by the authors of the paper "Predicting the Geographical Origin of Music" [Fang Zhou, Claire Q, Ross D. King (2015) how crated and used this set as basis for this paper. The key information embedded in this data set are attributes of music pieces labeled by there point of origin.

The authors collected 1,142 pieces of music from a personal CD-collection. The target values represented by longitude and latitude are the main country/area of residence of the artists that produced the music. Over all the authors collected music from 73 countries/areas not including western style music, since this category is called to have a global influence and therefor unfitted for predicting a specific country/region.

The attributes to describe the music pieces are automatically created by a software called MARSYAS (Tzanetakis 2007), a software created to extract audio features from wave files. With MARSYAS the authors could

convert every music peace to a set of 116 features called 'chromatic features.' The features are numerical and the authors claim to have normalized them into a gaussian normal distribution.

```
distances <- function(predicted, actual_value) {</pre>
  dif <- predicted-actual_value</pre>
  dif <- dif * (40030/360) # scaling coordinates to km by the factor circumference (km) / 360°
 mse \leftarrow sqrt(dif[,1]^2 + dif[,2]^2)
 return(mse)
data <- read.csv("Data/default_plus_chromatic_features_1059_tracks.txt", header=FALSE)</pre>
data <- as.data.frame(data)</pre>
colnames(data)[117:118] <- c("Latitude", "Longitude")</pre>
# Maybe some more preprocessing could be done here.
anyNA(data) # testing if there is at least a single NA -> but in this dataset there isn't
## [1] FALSE
# Maybe some more pre-processing could be done here.
anyDuplicated(data) # testing for duplicates -> 0 found
## [1] 0
# feature scaling
# separate labels from features
#label_column_names <- c("Longitude", "Latitude")
#data_labels <- data[label_column_names]</pre>
#data_features <- subset(data, select = -label_column_names)
# scale features
#data features scaled <- as.data.frame(scale(data features))</pre>
# add labels to the now scaled features
#data <- cbind(data features scaled, data labels)
```

some insights into the data

```
# check if data is already standardized

# get columns without the target cols ("long" and "lat")
data_without_target_cols <- subset(data, select=-c(Latitude,Longitude))

# for each column in the data get sd and median
sd_per_col <- apply(data_without_target_cols, 2, sd) # the two stands for columns, if we would have use
sd_per_col_df <- data.frame(sd_per_col)

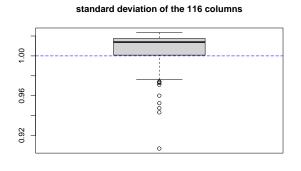
mean_per_col <- apply(data_without_target_cols, 2, mean)
mean_per_col_df <- data.frame(mean_per_col)

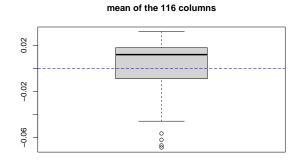
sd_and_mean_per_col_df <- merge(sd_per_col_df, mean_per_col_df, by="row.names", all=TRUE)

#par(mar = c(4, 4, .1, .1)) # to make the two plots show side by side and not above each other

boxplot(sd_per_col, data=sd_and_mean_per_col_df, xlab="standard deviation (blue line at 1)", y_lab="val-abline(h=1, col = "blue", lty=5)</pre>
```

boxplot(mean_per_col, data=sd_and_mean_per_col_df, xlab="mean (blue line at 0)", y_lab="value", main="mabline(h=0, col = "blue", lty=5)



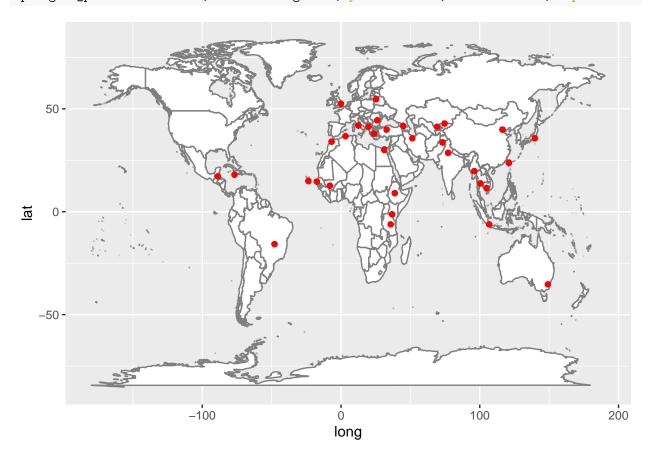


standard deviation (blue line at 1)

mean (blue line at 0)

```
# basic world map with music origins
mapWorld <- borders("world", colour="gray50", fill="white")
mp <- ggplot() + mapWorld

mp + geom_point(data = data, aes(x = Longitude, y = Latitude), color = "red", alpha = 0.5)</pre>
```



conspicuousness

When looking at the map, it seems like there are way fewer unique data points than expected. The paper states that there are 1,142 pieces from 73 countries/areas, but counting the point on the map just returnes 33 data points. The following code investigates this difference.

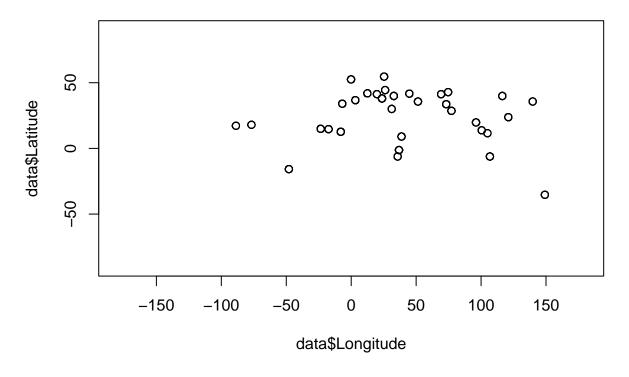
```
# investigate why there are fewer points on the map than regions on the map
# data is the full data set

# group data by unique combinations of long and lat and safe in data frame
# and count occurrences of each unique combination
# the unique combinations of long and lat represent regions
occurences_per_region <- data.frame(data %>% count(Longitude, Latitude, sort=TRUE))

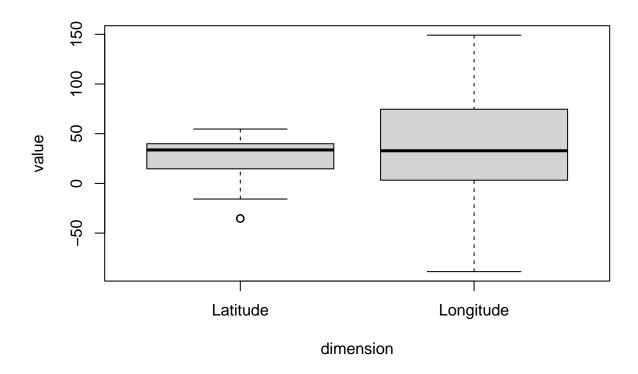
nrow(occurences_per_region) # returns 33 -> 33 regions in the data set and not 71 like proposed in the
## [1] 33
sum(occurences_per_region[, 'n']) # returns 1059 -> therefore there are 1095 tracks in the data set
## [1] 1059
```

The results suggest that there are actually only 33 unique combinations of latitude and longitude in the data set. Therefore the pieces can only be categorized into 33 different categories. The reason for this discrepancy to the suggested number in the paper (73) might be that some regions have been aggregated prior to uploading the data or some row are missing. The second statement is also supported, by the fact that there are not 1,142 pieces in the data set as described in the paper, but only 1059 pieces. Whatever this (small) difference does not influence the tasked tackled/perused by this report.

```
plot(data$Longitude, data$Latitude, xlim=c(-180, 180), ylim=c(-90, 90))
```



```
df_lat <- data.frame(value = data$Latitude, dimension = "Latitude")
df_long <- data.frame(value = data$Longitude, dimension = "Longitude")
boxplot_df <- rbind(df_lat, df_long)
boxplot(value~dimension, data=boxplot_df)</pre>
```



```
set.seed(1)
n <-dim(data)[1]
train <- sample(1:n, 0.8*n)
test <- (1:n)[-train]</pre>
```

Method

Short summary about the algorithms. Which are used? What do we do? Classification or Regression?

Baseline - Linear Regression

First, we take a baseline to get a basic understanding on how well our chosen algorithms perform. Therefore, we decided to use a linear regression. First we created a model which includes all variables. The lm() command cannot compute a model for both output variables at the same time. So it creates two separate linear models, one for each output variable:

```
model.lm.all <- lm(cbind(Longitude, Latitude)~., data=data[train,])
pred.all <- predict(model.lm.all, newdata=data[test,])</pre>
```

To calculate a good meaningful measurement for the goodness of fit for the predictions the distance from the true location is calculated. The distances are calculated as the euclidean distances of the Longitude and Latitude between the predictions and the true location. They are measured in [km]. These will be used for all the comparisons of the algorithms between each other but also with the baseline and also with the literature (we need here another citation to the original paper)

The predictions are on average quite far from the true destination:

```
## [1] 5850.727
```

The impact of the variables was analysed. It seemed that not all of them have a significant influence on the models. So a second model was developed, just using the variable which have a significant influence on the full model, either on the Latitude or on the Longitude.

The predictions for the smaller models are on average a bit better:

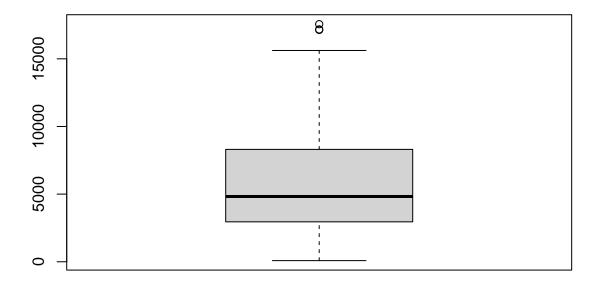
```
## [1] 5635.057
```

An ANOVA was calculated to check if there is a significant difference between the two models.

```
anova(model.lm.all, model.lm.sig)
```

```
## Analysis of Variance Table
##
## Model 1: cbind(Longitude, Latitude) ~ V1 + V2 + V3 + V4 + V5 + V6 + V7 +
##
       V8 + V9 + V10 + V11 + V12 + V13 + V14 + V15 + V16 + V17 +
##
       V18 + V19 + V20 + V21 + V22 + V23 + V24 + V25 + V26 + V27 +
       V28 + V29 + V30 + V31 + V32 + V33 + V34 + V35 + V36 + V37 +
##
       V38 + V39 + V40 + V41 + V42 + V43 + V44 + V45 + V46 + V47 +
##
       V48 + V49 + V50 + V51 + V52 + V53 + V54 + V55 + V56 + V57 +
##
##
       V58 + V59 + V60 + V61 + V62 + V63 + V64 + V65 + V66 + V67 +
##
       V68 + V69 + V70 + V71 + V72 + V73 + V74 + V75 + V76 + V77 +
       V78 + V79 + V80 + V81 + V82 + V83 + V84 + V85 + V86 + V87 +
##
##
       V88 + V89 + V90 + V91 + V92 + V93 + V94 + V95 + V96 + V97 +
##
       V98 + V99 + V100 + V101 + V102 + V103 + V104 + V105 + V106 +
##
       V107 + V108 + V109 + V110 + V111 + V112 + V113 + V114 + V115 +
##
       V116
  Model 2: cbind(Longitude, Latitude) ~ V4 + V9 + V16 + V30 + V32 + V33 +
##
       V37 + V38 + V61 + V90 + V91 + V92 + V95 + V96 + V104 + V5 +
##
       V6 + V8 + V9 + V11 + V15 + V34 + V39 + V63 + V94 + V97
##
     Res.Df Df Gen.var.
                         Pillai approx F num Df den Df
##
                                                           Pr(>F)
## 1
        774
                 666.34
## 2
        821 47
                 688.29 0.17239
                                  1.5533
                                             94
                                                   1548 0.0007525 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
```

The model tells that there is a difference and, as seen before, the smaller model performs better.



Algorithm 1 - Ridge Regression

Short introduction of the first algorithm. What does it do? What are the strengths? What are weaknesses? How is it implemented, including major code snippets.

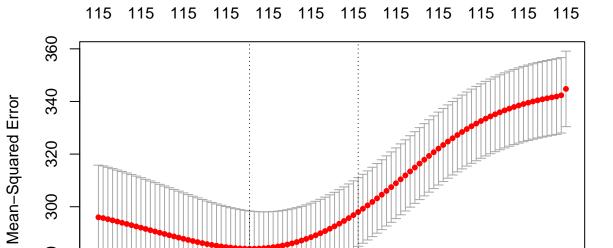
The first algorithm we tried was ridge regression. This algorithm is similar to a linear regression but while the linear regression tries to minimize the difference between the weighted input variables and the output data, the ridge regression adds a regularisation term on the input variables.

[Here has to be added the formula of the ridge regression]

In fact, this is a possibility to fight over-fitting. If lambda is big the model tends to just take the b0 into account. So the predicted value is the mean of the output variable. If lambda is small, then the model tends to be normal non-regularised model, hence the one from the linear regression.

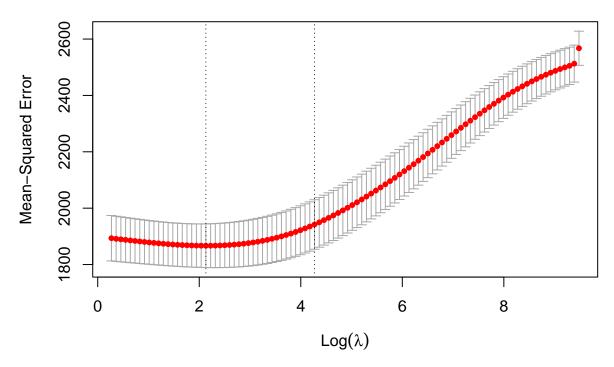
The command glmnet is used to perform the ridge regression. Cross-validation is performed to find the optimal lambda values. In general, it would be possible with glmnet to just calculate one single model for both output variables. But unfortunately, this option is not available when doing the cross-validation. So, again two independent cross-validations are done to get two values for lambda, one for each output variable.

[1] "data.frame"



 $\text{Log}(\lambda)$





It can be seen that the distributions for lambda are quite similar for both models. Although the MSE for both models differ the optimal lambda is quite similar:

[1] 10.83272

[1] 8.419139

The model performs better with ridge regression as with the baseline. In this model two lambdas are used.

[1] 4492.287

A second result is calculated with just one lambda. This lambda is calculated as the mean of the before calculated values of lambda. The result for the adapted version is just slightly worse than with the model with two independent lambdas:

[1] 4496.877

Algorithm 2

Short introduction of the second algorithm. What does it do? What are the strengths? What are weaknesses? How is it implemented, including major code snippets.

For the second algorithm we'll be using Artificial Neural Networks (NNs) as they can be used for regression problems as well. To use NNs for our predictions we will, as with all machine learning models, fit the data to the model. This is done by minimising the loss function, which for NN-regression is the mean of squared errors MSE over the training set. In the case of fitting a neural network, it is much faster if the data are scaled/normalised first and also gives more appropriate results if the features that are used have different scales and ranges. But this is already the case here.

The NN takes all features from the training data and outputs two values for the Latitude and Longitude, so it is a multi-output regression. Inbetween Input and Output Layer are a few more layers defined which

all use ReLU as activation function. For compiling and fitting the model we again use the Mean Squared Error (MSE) and Adam as the optimizer. We let the model train for 100 epochs with a batch size of 128 and validation split of 0.2.

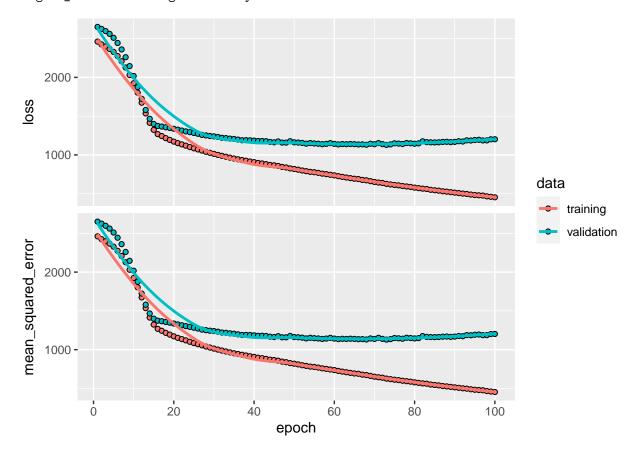
The results are visualized below.

Loaded Tensorflow version 2.7.0

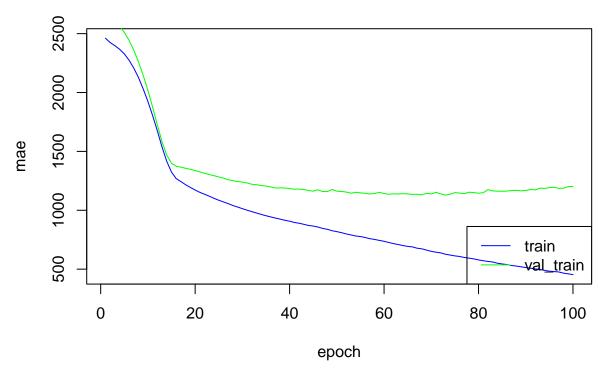
Model: "sequential"

##			
## ##	Layer (type)	Output Shape	Param #
##	dense_3 (Dense)	(None, 64)	7488
## ##	dense_2 (Dense)	(None, 16)	1040
## ##	dense_1 (Dense)	(None, 8)	136
## ##	dense (Dense)	(None, 2)	18
##	=======================================		========
##	Total params: 8,682		
##	Trainable params: 8,682		
##	Non-trainable params: 0		
##			

`geom_smooth()` using formula 'y ~ x'



Model's Mean Squared Error

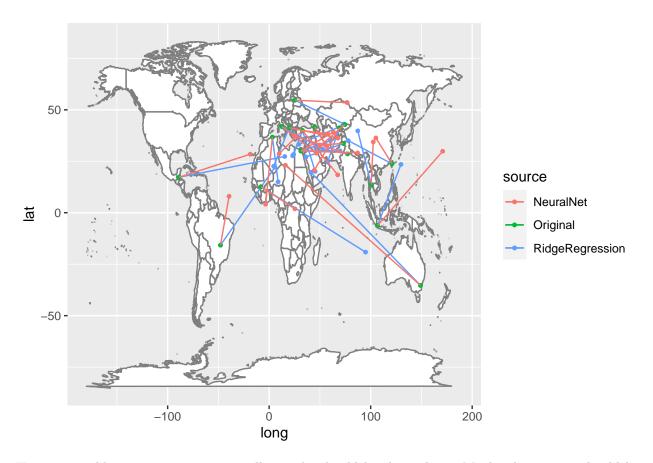


For evaluation we predict the Longitude and Latitude for the same test data as for the other regression cases and also use the custom distances metric. We get the following result:

[1] 4421.732

Results

- ## [1] 212 4
- **##** [1] 212 4
- ## [1] 212 4
- ## [1] 424



Here some tables, summaries or especially graphs should be shown here. Maybe this section should be separated into two to show the algorithms for themselves

target_values <- data.frame (Latitude = c(-15.75, 14.91, 12.65), Longitude = c(-47.95, -23.51, -8.00)) target_values index < -1: $nrow(target_values)target_values$ source <- "Original"

neural_net_results <- data.frame (Latitude = c(2.274030, 23.708105, 54.149323), Longitude = c(-23.0598640, 41.1416664, 14.7137728)) neural_net_results index < -1: $nrow(neural_net_results)neural_net_results$ ource <- "NeuralNet"

ridge_regression_results <- data.frame (Latitude = c(14.827996, 27.492521, 45.357388), Longitude = c(-8.6322704, 39.3965714, 32.7124573)) ridge_regression_results index < -1: $nrow(ridge_regression_results)ridge_regression_results$ <- "RidgeRegression"

aggregated_results_ridge_regression <- rbind(target_values, ridge_regression_results) aggregated_results_neural_net <- rbind(target_values, neural_net_results)

mapWorld <- borders("world," colour="gray50," fill="white") mp_2 <- ggplot() + mapWorld + geom_point(size = 2, data = aggregated_results_ridge_regression, aes(Longitude, Latitude, group=index, color=source)) + geom_line(data = aggregated_results_ridge_regression, aes(Longitude, Latitude, group=index, color="RidgeRegression")) + geom_point(size = 2, data = aggregated_results_neural_net, aes(Longitude, Latitude, group=index, color=source)) + geom_line(data = aggregated_results_neural_net, aes(Longitude, Latitude, group=index, color="NeuralNet")) mp_2

Discussion

Here follows the discussion of the results. What are the major findings? How did the algorithms perform? Which one was better overall? Is it always better or were the findings which were better by the other one?

Which one should be implemented? How could the algorithm be tweaked to perform even better? Where were the problems during implementation? Where are the limits for the algorithms?

How precise do we predict the cities? How far is the difference in kilometers? The authors of the paper where the dataset comes from have a mean great circle error of 3113km? Are we above or below and by how much?

Conclusion

At final some conclusions about the key findings and which algorithm should be used. What was the goal? Were and how were they achieved?

Fang Zhou, Claire Q, Ross D. King. 2015. "Predicting the Geographical Origin of Music." Paper. https://ieeexplore.ieee.org/document/7023456.

Tzanetakis, George. 2007. "Marsyas: A Case Study in Implementing Music Information Retrieval Systems." In *Intelligent Music Information Systems: Tools and Methodologies*, edited by Shepherd Shen and Liu Cui. Information Science Reference. http://marsyas.sness.net/pdfs/0000/0007/imis_bookchapter.pdf.