# Lab 7. Evaluating the Recommender Systems

# Preparing the data to evaluate models

We need two sets - trainig set and testing set - to evaluate the model.

#### Splitting the data

I will split the data so that th training set consists of 80% of the data, and the test set - 20%.

```
percentage_training <- 0.8
```

For each user in the test set, we need to define how many items to use to generate recommendations. For this, I will first check the minimum number of items rated by users to be sure there will be no users with no items to test.

```
min(rowCounts(ratings movies))
## [1] 18
items to keep <- 15 #number of items to generate recommendations
rating threshold <- 3 # threshold with the minimum rating that is considered good
n eval <- 1 #number of times to run evaluation
eval sets <- evaluationScheme(data = ratings movies,</pre>
                              method = "split",
                              train = percentage training,
                              given = items to keep,
                              goodRating = rating threshold,
                              k = n eval)
eval sets
## Evaluation scheme with 15 items given
## Method: 'split' with 1 run(s).
## Training set proportion: 0.800
## Good ratings: >=3.000000
## Data set: 560 x 332 rating matrix of class 'realRatingMatrix' with 55298 ratings.
getData(eval sets, "train") # training set
## 448 x 332 rating matrix of class 'realRatingMatrix' with 44470 ratings.
getData(eval sets, "known") # set with the items used to build the recommendations
## 112 x 332 rating matrix of class 'realRatingMatrix' with 1680 ratings.
getData(eval sets, "unknown") # set with the items used to test the recommendations
## 112 x 332 rating matrix of class 'realRatingMatrix' with 9148 ratings.
qplot(rowCounts(getData(eval sets, "unknown"))) +
```

```
geom_histogram(binwidth = 10) +
ggtitle("unknown items by the users")
```

### Bootstrapping the data

Bootrstrapping is another approach to split the data. The same user can be sampled more than once and, if the training set has the same size as it did earlier, there will be more users in the test set.

#### Using k-fold to validate models

This approach is the most accurate one, although it's computationally heavier.

# Evavluating the ratings

I will use the k-fold approach for evaluation. In the following code, I re-define the evaluation sets, build IBCF model and a matrix with predicted ratings.

```
eval_sets <- evaluationScheme(data = ratings_movies,</pre>
                                method = "cross-validation",
                                k = n \text{ fold,}
                                given = items to keep,
                                goodRating = rating threshold)
model to evaluate <- "IBCF"</pre>
model parameters <- NULL</pre>
eval_recommender <- Recommender(data = getData(eval_sets, "train"),</pre>
                                  method = model to evaluate,
                                  parameter = model parameters)
items_to_recommend <- 10
eval prediction <- predict(object = eval_recommender,</pre>
                             newdata = getData(eval sets, "known"),
                             n = items_to_recommend,
                             type = "ratings")
qplot(rowCounts(eval prediction)) +
  geom\ histogram(binwidth = 10) +
  ggtitle("Distribution of movies per user")
```

Now, I compute the accuracy measures for each user. Most of the RMSEs are in the range of 0.8 to 1.4:

```
ggtitle("Distribution of the RMSE by user")
```

In order to have a performance index for the whole model, I specify *byUser* as FALSE and compute the average indices:

### Evaluating the recommendations

Another way to measure accuracies is by comparing the recommendations with the purchases having a positive rating.

```
results <- evaluate(x = eval sets,
                    method = model to evaluate,
                    n = seq(10, 100, 10))
## IBCF run fold/sample [model time/prediction time]
     1 [7.47sec/0.46sec]
     2 [7.09sec/0.48sec]
##
     3 [7.32sec/0.45sec]
##
     4 [8.34sec/0.47sec]
head(getConfusionMatrix(results)[[1]])
                       FP
                                FN
                                         TN precision
## 10 3.142857 6.857143 67.02857 239.9714 0.3142857 0.04466012 0.04466012
## 20 6.321429 13.678571 63.85000 233.1500 0.3160714 0.08891903 0.08891903
## 30 9.192857 20.807143 60.97857 226.0214 0.3064286 0.13193468 0.13193468
  40 11.921429 28.078571 58.25000 218.7500 0.2980357 0.17033031 0.17033031
  50 14.678571 35.321429 55.49286 211.5071 0.2935714 0.20963292 0.20963292
  60 17.321429 42.678571 52.85000 204.1500 0.2886905 0.24824396 0.24824396
             FPR
## 10 0.02719801
## 20 0.05426466
## 30 0.08295660
## 40 0.11213301
## 50 0.14119344
```

```
## 60 0.17070228
```

In order to have a look at all the splits at the same time, I sum up the indices:

Finally, I plot the ROC and the precision/recall curves:

```
plot(results, annotate = TRUE, main = "ROC curve")
```

```
plot(results, "prec/rec", annotate = TRUE, main = "Precision-recall")
```

# Comparing models

In order to compare different models, I define them as a following list: - Item-based collaborative filtering, using the Cosine as the distance function - Item-based collaborative filtering, using the Pearson correlation as the distance function - User-based collaborative filtering, using the Cosine as the distance function - User-based collaborative filtering, using the Pearson correlation as the distance function - Random recommendations to have a base line

)

Now, I define a different numbers of recommended movies and run and evaluate the models:

```
n recommendations <- c(1, 5, seq(10, 100, 10))
list results <- evaluate(x = eval sets,</pre>
                         method = models to evaluate,
                         n = n recommendations)
  IBCF run fold/sample [model time/prediction time]
     1 [7.23sec/0.2sec]
     2 [9.79sec/0.39sec]
     3 [8sec/0.51sec]
     4 [7.17sec/0.38sec]
##
## IBCF run fold/sample [model time/prediction time]
     1 [17.35sec/0.37sec]
##
     2 [17.72sec/0.28sec]
##
     3 [16.46sec/0.37sec]
##
     4 [18.5sec/0.34sec]
## UBCF run fold/sample [model time/prediction time]
     1 [0.02sec/8.1sec]
##
     2 [0sec/9.29sec]
##
     3 [0.01sec/8.15sec]
##
     4 [0.01sec/6.54sec]
##
## UBCF run fold/sample [model time/prediction time]
     1 [0.02sec/17.52sec]
     2 [0sec/17.32sec]
     3 [0.02sec/15.24sec]
     4 [0sec/9.46sec]
##
## RANDOM run fold/sample [model time/prediction time]
     1 [0.01sec/0.21sec]
##
     2 [0sec/0.26sec]
##
##
     3 [0sec/0.19sec]
     4 [0sec/0.33sec]
##
sapply(list results, class) == "evaluationResults"
## IBCF_cos IBCF_cor UBCF_cos UBCF_cor
                                        random
      TRUE
                TRUE
                        TRUE
                                 TRUE
                                            TRUE
avg matrices <- lapply(list_results, avg)</pre>
head(avg matrices$IBCF cos[, 5:8])
##
     precision
                    recall
                                    TPR
                                                FPR
```

```
## 1 0.3989979 0.005637331 0.005637331 0.00239457

## 5 0.3221017 0.020800675 0.020800675 0.01344330

## 10 0.3195863 0.043228764 0.043228764 0.02720080

## 20 0.3140114 0.085113290 0.085113290 0.05508793

## 30 0.3085196 0.127515994 0.127515994 0.08360939

## 40 0.3071197 0.171622620 0.171622620 0.11176470
```

# Identifying the most suitable model

I compare the models by building a chart displaying their ROC curves and Precision/recall curves.

```
plot(list_results, annotate = 1, legend = "topleft")
title("ROC curve")
```

```
plot(list_results, "prec/rec", annotate = 1, legend = "bottomright")
title("Precision-recall")
```

# Optimizing a numeric parameter

IBCF takes account of the k-closest items. I will explore more values, ranging between 5 and 40:

```
vector_k <- c(5, 10, 20, 30, 40)
models_to_evaluate <- lapply(vector_k, function(k) {
    list(name = "IBCF",
        param = list(method = "cosine", k = k))
})
names(models_to_evaluate) <- paste0("IBCF_k_", vector_k)</pre>
```

Now I will build and evaluate the same models:

```
##
     4 [10.01sec/0.47sec]
## IBCF run fold/sample [model time/prediction time]
     1 [9.34sec/0.28sec]
     2 [9.31sec/0.39sec]
##
     3 [10.56sec/0.37sec]
##
     4 [9.18sec/0.3sec]
##
   IBCF run fold/sample [model time/prediction time]
       [4.7sec/0.17sec]
##
     2 [5sec/0.34sec]
##
     3 [5sec/0.19sec]
##
     4 [4.35sec/0.23sec]
   IBCF run fold/sample [model time/prediction time]
     1 [5.72sec/0.2sec]
##
     2 [8.15sec/0.39sec]
##
     3 [7.86sec/0.2sec]
##
     4 [8.45sec/0.47sec]
##
  IBCF run fold/sample [model time/prediction time]
     1 [7.02sec/0.37sec]
##
     2 [8sec/0.25sec]
##
     3 [9.03sec/0.32sec]
##
     4 [7.7sec/0.41sec]
##
plot(list results, annotate = 1, legend = "topleft")
title("ROC curve")
```

```
plot(list_results, "prec/rec", annotate = 1, legend = "bottomright")
title("Precision-recall")
```

Based on the ROC curve's plot, the k having the biggest AUC is 10. Another good candidate is 5, but it can never have a high TPR. This means that, even if we set a very high n value, the algorithm won't be able to recommend a big percentage of items that the user liked. The IBCF with k = 5 recommends only a few items similar to the purchases. Therefore, it can't be used to recommend many items.

Based on the precision/recall plot, k should be set to 10 to achieve the highest recall. If we are more interested in the precision, we set k to 5.