

Project 1

February 28, 2024

Contents

Problem 3: Learning Bayesian networks from protein data	1
Preparation to run the code	1
Loading data	1
Variables and observations	1
Visualisation of transformed data	4
Defining functions	4
Default parameters	ļ
Different parameters	ļ
Retraining the best BN	(
Render this .rmd into a pdf	7

Problem 3: Learning Bayesian networks from protein data

Preparation to run the code

Setting seed for reproducibility and loading packages. Some packages are additionally loaded inside the functions that are run in parallel.

```
library("GGally")
library("BiDAG")
library("igraph")
library("parallel")
library("doParallel")
```

Loading data

```
data <- read.table("2005_sachs_2_cd3cd28icam2_log_std.csv", sep=",", header=TRUE)</pre>
```

Variables and observations

```
num_variables <- ncol(data)
num_observations <- nrow(data)

print(paste("Number of variables:", num_variables))

## [1] "Number of variables: 11"

print(paste("Number of observations:", num_observations))</pre>
```

```
## [1] "Number of observations: 902"
```

Visualisation of transformed data

```
png(file="plot.png", width=650, height=1000)
ggpairs(data, progress=FALSE)
dev.off()
## pdf
## 2
```

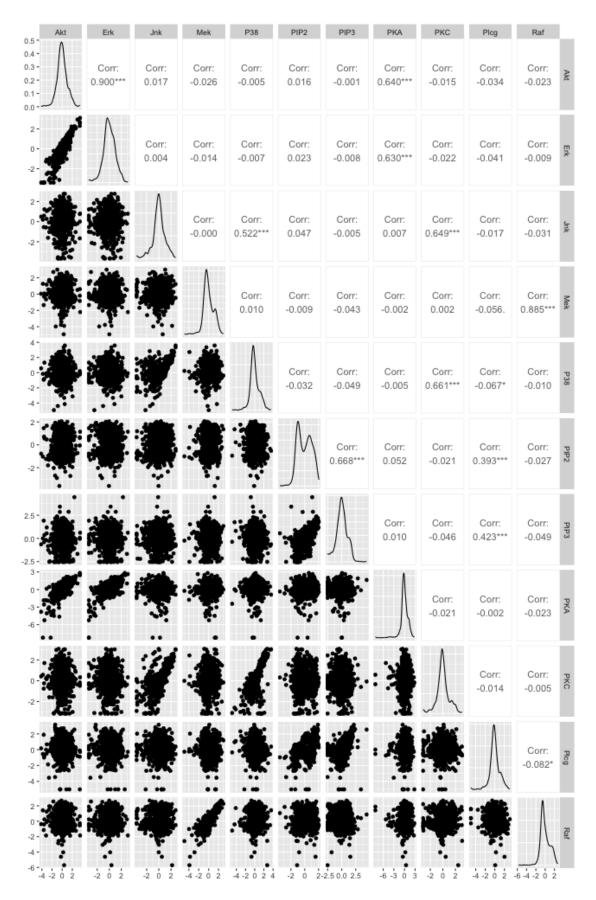


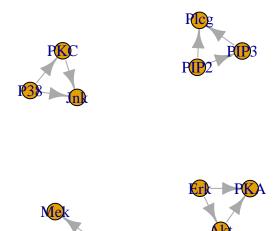
Figure 1: Plot of transformed data

Defining functions

```
splitting_data <- function(data) {</pre>
    # Shuffling data
    indices <- 1:nrow(data)</pre>
    indices <- sample(length(indices))</pre>
    # Splitting data
    train_size <- ceiling(nrow(data)*0.8)</pre>
    train_indices <- indices[1:train_size]</pre>
    test_indices <- indices[(train_size+1):(length(indices))]</pre>
    # Checking if there is no overlap
    if(length(unique(c(train_indices, test_indices))) != length(c(train_indices, test_indices))) {
        print("Overlap!")
    }
    train_data <- data[row.names(data) %in% train_indices, ]</pre>
    test_data <- data[row.names(data) %in% test_indices, ]</pre>
    split_data <- list("train_data"=train_data, "test_data"=test_data)</pre>
    return(split_data)
}
training_BN <- function(data, bgepar) {</pre>
    library("BiDAG")
    init_score_par <- scoreparameters("bge", data$train_data, bgepar)</pre>
    learnt_BN <- iterativeMCMC(init_score_par, verbose=FALSE)</pre>
    return(learnt_BN)
}
testing_BN <- function(data, BN, bgepar) {</pre>
    library("BiDAG")
    test_score_par <- scoreparameters("bge", data$test_data, bgepar)</pre>
    test_score <- scoreagainstDAG(test_score_par, BN$DAG)</pre>
    return(mean(test_score))
}
plot_DAG <- function(BN) {</pre>
    library("igraph")
    g <- graph_from_adjacency_matrix(BN$DAG)</pre>
    plot(g)
}
get_number_of_edges <- function(BN) {</pre>
    library("igraph")
    return(length(E(graph_from_adjacency_matrix(BN$DAG))))
}
```

Default parameters

```
bgepar <- list(am=1, aw=NULL, edgepf=1)
split_data <- splitting_data(data)
learnt_BN <- training_BN(split_data, bgepar)
mean_test_score <- testing_BN(split_data, learnt_BN, bgepar)
print(paste0("Average BGe score on testing data: ", mean_test_score))
## [1] "Average BGe score on testing data: -12.4139751176187"
plot_DAG(learnt_BN)</pre>
```



Different parameters

```
# Set the number of cores to use
num_cores <- detectCores()

# Register parallel backend
cl <- makeCluster(num_cores)
registerDoParallel(cl)

ams <- c(10^(-5), 10^(-3), 10^(-1), 10, 10^2)

res <- foreach(am = ams, .combine=c) %dopar% {
    set.seed(42)
    bgepar <- list(am=am, aw=NULL, edgepf=1)</pre>
```

```
numbers_of_edges <- c()</pre>
    mean_test_scores <- c()</pre>
    for(i in 1:10) {
        RNGkind("L'Ecuver-CMRG")
        split_data <- splitting_data(data)</pre>
        learnt_BN <- training_BN(split_data, bgepar)</pre>
        number_of_edges <- get_number_of_edges(learnt_BN)</pre>
        numbers of edges <- append(numbers of edges, number of edges)
        mean_test_score <- testing_BN(split_data, learnt_BN, bgepar)</pre>
        mean_test_scores <- append(mean_test_scores, mean_test_score)</pre>
    }
    return(c(am, mean(numbers_of_edges), mean(mean_test_scores)))
stopCluster(cl)
res_m <- t(matrix(data=res, nrow=5, ncol=3, byrow=TRUE))</pre>
rownames(res_m) <- c("Parameter am", "Average number of edges", "Average BGe score of the test data")
print(res_m)
##
                                             [,1]
                                                        [,2]
                                                                   [,3]
                                                                             [,4]
                                                    0.00100
## Parameter am
                                          0.00001
                                                               0.10000 10.00000 100.00000
## Average number of edges
                                          7.00000
                                                    7.00000 9.10000 13.40000 16.70000
\#\# Average BGe score of the test data -12.73416 -12.73416 -12.70972 -12.80457 -13.58552
```

Retraining the best BN

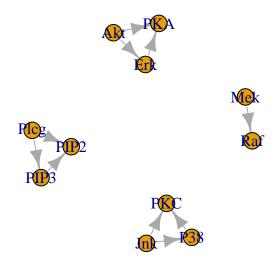
The best BN (based on the highest BGe score) was the one with $\alpha_m = 10^{-1}$.

We observe that the average number of edges increases with increasing α_m .

```
set.seed(42)
bgepar <- list(am=10^(-1), aw=NULL, edgepf=1)
split_data <- splitting_data(data)
learnt_BN <- training_BN(split_data, bgepar)

mean_test_score <- testing_BN(split_data, learnt_BN, bgepar)
print(paste0("Average BGe score on testing data: ", mean_test_score))

## [1] "Average BGe score on testing data: -12.1209831813562"
plot_DAG(learnt_BN)</pre>
```



Render this .rmd into a pdf

```
library(rmarkdown)
render("1.Rmd", pdf_document(TRUE), "1.pdf")
```