Assignment 2: Bayesian networks

Graphical and Hidden Markov Models

Ignacio Almodóvar Cárdenas & Javier Muñoz Flores

20-04-2022

Explain Dataset

We have chosen a dataset located in kaggle. It contains information about the classification of certain drug types based on different features such as the age, the sex, the blood pressure levels, the cholesterol levels and the sodium-to-potassium ratio.

The dataset contains different types of variables, including continuous and categorical. These are:

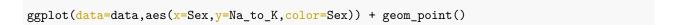
- Age: Continuous integer variable showing the age of each observation.
- Sex: Categorical variable with two levels (F,M) for indicating Female and Male respectively.
- BP: Categorical variable referring to patient's blood pressure. The levels are NORMAL and ABNOR-MAL. This last level includes both LOW and HIGH blood pressure conditions.
- Cholesterol: That is again a categorical variable including the cholesterol levels of the patient. It includes both NORMAL and ABNORMAL levels, again the last one refers to both HIGH and LOW.
- Na to K: Continuous variable showing the Sodium-potassium balance
- Drug: This will be our variable to predict. It contains two different levels (DrugX, DrugY). Depending on the characteristics of each patient it will require one drug or the other one.

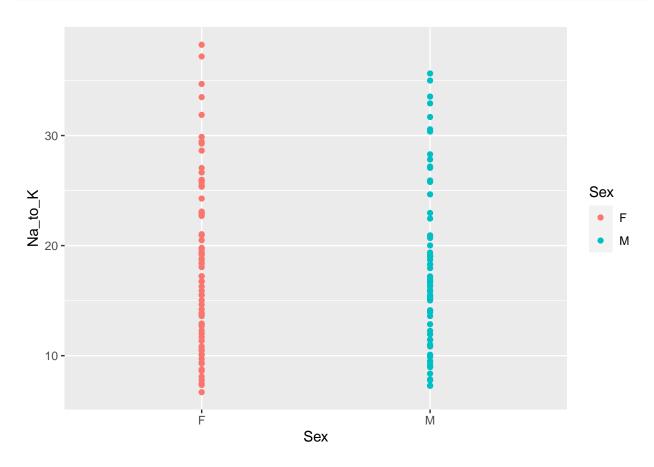
```
load(file = "data.RData")
data$Age %<>% as.numeric()
data$Sex %<>% as.factor()
data$BP %<>% as.factor()
data$Drug %<>% as.factor()
data$Cholesterol %<>% as.factor()
summary(data)
```

```
##
                     Sex
                                    BP
                                           Cholesterol
                                                            Na_to_K
                                                                              Drug
         Age
                                                        Min.
##
    Min.
           :15.00
                     F:74
                            ABNORMAL:86
                                           HIGH: 67
                                                                : 6.683
                                                                          drugX:54
##
    1st Qu.:30.00
                     M:71
                            NORMAL:59
                                           NORMAL:78
                                                        1st Qu.:11.939
                                                                          DrugY:91
   Median :43.00
                                                        Median :16.753
##
   Mean
           :43.85
                                                        Mean
                                                                :18.009
##
    3rd Qu.:58.00
                                                        3rd Qu.:22.905
##
    Max.
           :74.00
                                                        Max.
                                                                :38.247
```

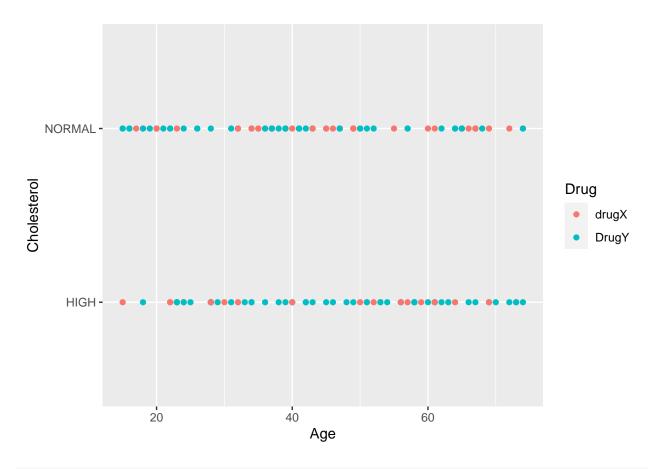
As it can be seen, the dataset is more or less well balanced for all the categorical variables except for the Drug one, where the number of levels have a significant difference. However, as this is just an experiment to deal with Bayesian networks and we do not really care about the results obtained, the data contained should be enough to make reasonable predictions.

Define a possible network structure and give a brief explanation as to why this structure could make sense. Include a graph of the proposed model.

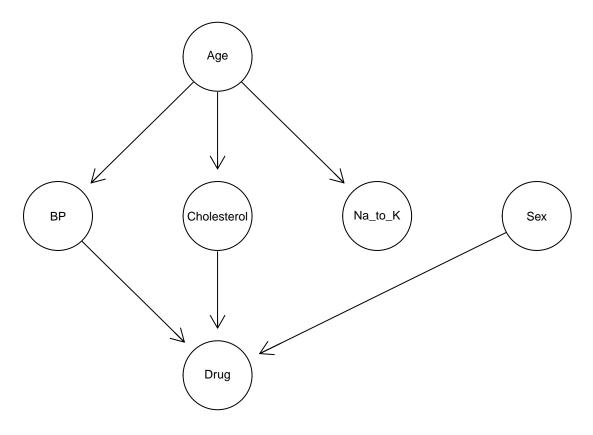




ggplot(data=data,aes(x=Age,y=Cholesterol,color=Drug)) + geom_point()

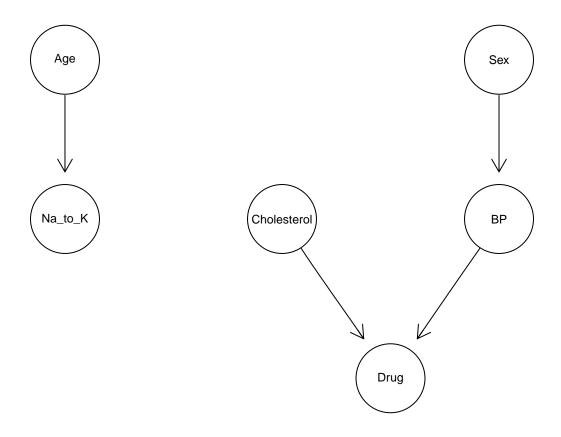


graph <- model2network("[Age][BP|Age][Cholesterol|Age][Na_to_K|Age][Sex][Drug|BP:Cholesterol:Sex]")
graphviz.plot(graph)</pre>



Tenemos dosa variables conitinuas. Sin embargo, para poder modela nuestra red como una bayesian network, tenemos que asumir condiciones gausianas como que los nodos categoricos no pueden tener padres conitnuois. Por lo tanto, quitaremos las relaciones supuestas inicialmente para poder plantear el problema correctamente.

graph_2=model2network("[Age][Sex][Cholesterol][Na_to_K|Age][BP|Sex][Drug|BP:Cholesterol]")
graphviz.plot(graph_2)



Fit the model and use it to make an out of sample prediction

```
bayesfit <- bn.fit(graph_2,data)
bayesfit</pre>
```

```
##
     Bayesian network parameters
##
##
##
     Parameters of node Age (Gaussian distribution)
##
## Conditional density: Age
## Coefficients:
##
   (Intercept)
##
      43.84828
## Standard deviation of the residuals: 16.75532
##
##
     Parameters of node BP (multinomial distribution)
##
## Conditional probability table:
##
##
             Sex
## BP
                                 М
     ABNORMAL 0.5945946 0.5915493
##
```

```
##
     NORMAL
              0.4054054 0.4084507
##
##
     Parameters of node Cholesterol (multinomial distribution)
##
## Conditional probability table:
        HIGH
              NORMAL
##
## 0.462069 0.537931
##
##
    Parameters of node Drug (multinomial distribution)
##
## Conditional probability table:
##
   , , Cholesterol = HIGH
##
##
##
          BP
## Drug
            ABNORMAL
                        NORMAL
     drugX 0.0000000 0.5405405
##
##
     DrugY 1.0000000 0.4594595
##
##
   , , Cholesterol = NORMAL
##
##
          BP
            ABNORMAL
                        NORMAL
## Drug
     drugX 0.3214286 0.7272727
##
     DrugY 0.6785714 0.2727273
##
##
##
     Parameters of node Na_to_K (Gaussian distribution)
##
##
## Conditional density: Na_to_K | Age
## Coefficients:
## (Intercept)
                        Age
## 19.00257413 -0.02267038
## Standard deviation of the residuals: 7.572041
##
    Parameters of node Sex (multinomial distribution)
##
##
## Conditional probability table:
##
            F
## 0.5103448 0.4896552
bayesfit$E
```

NULL

Now we are going to predict which drug will be taken a normal person, meaning a person with normal blood pressure and normal cholesterol.

```
v \leftarrow bayesfit \ Try \ this \ with \ bayesfit \ or \ expertfit \ instead. v[1,2,2]
```

[1] 0.7272727

Now we are going to calculate the probability of drugX in general

P(drugX) = P(drugX|Cholesterol, BP) = P(drugX, Cholesterol = HIGH) * P(Cholesterol = HIGH) + P(drugX, Cholesterol = Normeal) * P(cholesterol = NORMAL)

d) Try to fit the graph structure of the model using one or more of the different approaches we saw in class. Do the fits make sense? If not you could use whitelists or blacklists to make certain links impossible.

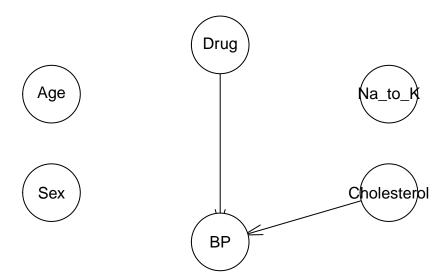
In the previous sections, we have assumed that the structure of the network were the correct one basing on different logical approaches. However, it exists different procedures to obtain the graph structure in a more unbiased way.

We will try to implement the three approaches saw in class:

- Constraint based: with the aimb method.
- Score based: with the homethod. We will use both AIC and BIC score functions to evaluate the fit.
- **Hybrid**: with the mmhc method.

First, we perform the first approach mentioned. This technique can be slow as it has to analyse each pair of nodes in the network. However, since our network is not too very large, the computational cost is not very high.

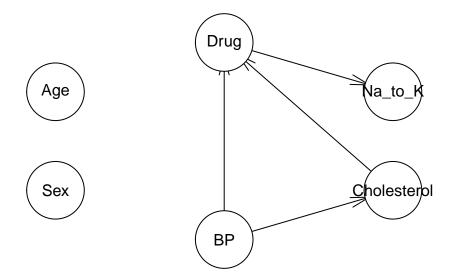
structure_iamb <- iamb(data)
plot(structure_iamb)</pre>



The structure obtained by this technique suggested that three out of the six nodes have not dependencies with others. This composition does not agree with our first approach, since it does not take into account some relationships we have assumed that were important and it eliminates so many dependencies.

Let's analyse the structure suggested by the **score based** methods. First, we look at the output basing on BIC.

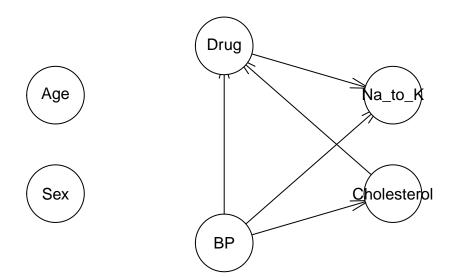
```
structure_bic <- hc(data, score = "bic-cg")
plot(structure_bic)</pre>
```



As we can see in the plot, this method includes the Cholesterol-Drug and $Drug-Na_to_K$ dependencies. While the first of the two relationships could makes sense, the second one does not make it. It could make sense an opposite direction of the arrow between $Drug-Na_to_K$, but in Conditional Gaussian Bayesian networks like this, categorical nodes have no continuous node parents, so it cannot be possible.

Let's see the structure proposed using AIC method.

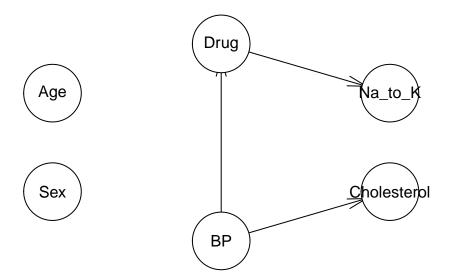
```
structure_aic <- hc(data, score = "aic-cg")
plot(structure_aic)</pre>
```



We can see that the dependency $BP-Na_to_K$ have been added, which is reasonable. Nevertheless, the dependency $Drug-Na_to_K$ stays.

Finally, we will see the struture suggested by the last of the approaches mentioned: \mathbf{hybrid} \mathbf{method} , a combination of the two previous ones.

structure_mmhc <- mmhc(data)
plot(structure_mmhc)</pre>



Again, it considers $Drug-Na_to_K$ dependency and it does not include important relationships which we have found through the score based methods.

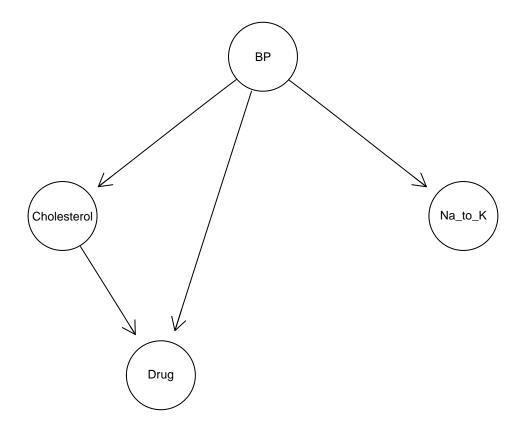
e) Which of the proposed graphical structures is the best?

In view of the results obtained in the previous section, where we have used algorithms which have helped us to learn the conditional independence structure, we can draw the following conclusions:

- We should eliminate the $Drug-Na_to_K$ since we consider it does not make sense with the analysis of or network.
- The nodes Age and Sex should be eliminated since it does not appear in any of the four structures proposed by the methods.
- The structure suggested by the score based methods adds important relationships which should be considered from a health point of view.
- The structure proposed by the AIC method should be selected as the best one instead of the one suggested by the BIC method because it considers *BP-Na_to_K* dependency. According to some research, the sodium (Na)/potassium (K) ratio could be associated with blood pressure (BP), so this relationship should be added to our network (Reference).

Thus, we delete the $Drug-Na_to_K$ dependency as well as the nodes Age and Sex and we draw the most optimal graphical structure.

final_structure <- model2network("[BP] [Na_to_K|BP] [Cholesterol|BP] [Drug|BP:Cholesterol]")
graphviz.plot(final_structure)</pre>



e) Provide a brief summary of your results.

To sum up, we conclude this work enumerating the most important insights obtained:

- The first approach for modelling the graph structure, i.e. the structure assumed of ourselves, has not been matched with the ones proposed by the algorithms. However, the dependency *Cholesterol-Drug*, which stays in almost all the structures proposed, seems to be the most reliable.
- It has been very important the assumptions in a network with both categorical and continuous variables like ours, i.e. a Conditional Gaussian Bayesian network. It is because we have taken into account that categorical nodes have no continuous node parents and it has not allowed to draw some dependencies that at first we could have considered to include.
- The probability that a random patient whose health we know nothing about, is prescribed to take the drugY is around 63%. Nevertheless, if we know that a patient has normal levels of blood pressure and cholesterol, it is more likely that he should take the drugY.
- The algorithms that provide a network structure agree that the variables Age and Sex should not be included in the network.