bnclassify: Learning Bayesian Network Classifiers

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Abstract

The **bnclassify** package provides state-of-the art algorithms for learning Bayesian network classifiers from data. For structure learning it provides variants of the greedy hill-climbing search, a well-known adaptation of the Chow-Liu algorithm and averaged one-dependence estimators. It provides Bayesian and maximum likelihood parameter estimation, as well as three naive-Bayes-specific methods based on discriminative score optimization and Bayesian model averaging. The implementation is efficient enough to allow for time-consuming discriminative scores on medium-sized data sets. **bnclassify** provides utilities for model evaluation, such as cross-validated accuracy and penalized log-likelihood scores, and analysis of the underlying networks, including network plotting via the **Rgraphviz** package. It is extensively tested, with over 200 automated tests that give a code coverage of 94%. Here we present the main functionalities, illustrate them with a number of data sets, and comment on related software.

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1 Introduction

Bayesian network classifiers [Bielza and Larrañaga, 2014, Friedman et al., 1997] are competitive performance classifiers [e.g., Zaidi et al., 2013] with the added benefit of interpretability. Their simplest member, the naive Bayes (NB) [Minsky, 1961], is well-known [Hand and Yu, 2001]. More elaborate models exist, taking advantage of the Bayesian network [Pearl, 1988, Koller and Friedman, 2009] formalism for representing complex probability distributions. The tree augmented naive Bayes [Friedman et al., 1997] and the averaged one-dependence estimators (AODE) [Webb et al., 2005] are among the most prominent.

A Bayesian network classifier is simply a Bayesian network applied to classification, that is, the prediction of the probability $P(c \mid \mathbf{x})$ of some discrete (class) variable C given some features X. The bilearn [Scutari and Ness, 2018, Scutari, 2010] package already provides state-of-the art algorithms for learning Bayesian networks from data. Yet, learning classifiers is specific, as the implicit goal is to estimate $P(c \mid \mathbf{x})$ rather than the joint probability $P(\mathbf{x}, c)$. Thus, specific search algorithms, network scores, parameter estimation and inference methods have been devised for this setting. In particular, many search algorithms consider a restricted space of structures such as that of augmented naive Bayes [Friedman et al., 1997] models. Unlike with general Bayesian networks, it makes sense to omit a feature X_i from the model as long as the estimation of $P(c \mid \mathbf{x})$ is no better than that of $P(c \mid \mathbf{x} \setminus x_i)$. Discriminative scores, related to the estimation of $P(c \mid \mathbf{x})$ rather than $P(c, \mathbf{x})$, are used to learn both structure [Keogh and Pazzani, 2002, Grossman and Domingos, 2004, Pernkopf and Bilmes, 2010, Carvalho et al., 2011 and parameters [Zaidi et al., 2013, 2017]. Some of the prominent classifiers [Webb et al., 2005] are ensembles of networks, and there are even heuristics applied at inference time, such as the lazy elimination technique [Zheng and Webb, 2006]. Many of these methods [e.g., Dash and Cooper, 2002, Zaidi et al., 2013, Keogh and Pazzani, 2002, Pazzani, 1996] are at best available in standalone implementations published alongside the original papers.

The **bnclassify** package implements state-of-the-art algorithms for learning structure and parameters. The implementation is efficient enough to allow for time-consuming discriminative scores on relatively large data sets. It provides utility functions for prediction and inference, model evaluation with network scores and cross-validated estimation of predictive performance, and model analysis, such as querying structure type or plotting via the **Rgraphviz** package [Hansen et al., 2017]. It integrates with the **caret** [Kuhn et al., 2017, Kuhn, 2008] and **mlr** [Bischl et al., 2017, 2015] packages for straightforward use in machine learning pipelines. Currently it supports only discrete variables. The functionalities are illustrated in an introductory vignette, while an additional vignette provides details on the implemented methods. It includes over 200 unit and integration tests that give a code coverage of 94 percent¹.

The rest of this paper is structured as follows. Section 2 provides background on Bayesian network classifiers. Section 3 describes the implemented functionalities. Section 4 illustrates usage with a synthetic data set. Section 5 discusses implementation while Section 6 briefly surveys related software. Finally, Section 7 concludes and outlines future work.

2 Background

2.1 Bayesian network classifiers

A Bayesian network classifier is a Bayesian network used for predicting a discrete class variable C. It assigns \mathbf{x} , an observation of n predictor variables (features) $\mathbf{X} = (X_1, \dots, X_n)$, to the most probable class:

$$c^* = \underset{c}{\operatorname{arg\,max}} P(c \mid \mathbf{x}) = \underset{c}{\operatorname{arg\,max}} P(\mathbf{x}, c).$$

The classifier factorizes $P(\mathbf{x},c)$ according to a Bayesian network $\mathcal{B} = \langle \mathcal{G}, \boldsymbol{\theta} \rangle$. \mathcal{G} is a directed

¹See https://codecov.io/github/bmihaljevic/bnclassify?branch=master

acyclic graph with a node per each of the variables in (\mathbf{X}, C) , encoding conditional independences: a variable X is independent of its nondescendants in \mathcal{G} given the values $\mathbf{pa}(x)$ of its parents. \mathcal{G} thus factorizes the joint into local conditional distributions over subsets of variables:

$$P(\mathbf{x}, c) = P(c \mid \mathbf{pa}(c)) \prod_{i=1}^{n} P(x_i \mid \mathbf{pa}(x_i)).$$

The parameters $\boldsymbol{\theta}$ specify the local conditional distributions.

2.2 Learning structure

We learn \mathcal{B} from a data set $\mathcal{D} = \{(\mathbf{x}^1, c^1), \dots, (\mathbf{x}^N, c^N)\}$ of N observations of \mathbf{X} and C. There are two main approaches to learning the structure \mathcal{G} from \mathcal{D} : a) testing for conditional independence among triplets of variables and b) searching a space of possible structures in order to optimize a network quality score. Under assumptions such as a limited number of parents per variable, approach a) can produce the correct network in polynomial time [Cheng et al., 2002, Tsamardinos et al., 2003]. On the other hand, finding the optimal structure even with at most two parents per variable is NP-hard [Chickering et al., 2004]. Thus, heuristic search algorithms, such as greedy hill-climbing, are commonly used [see e.g., Koller and Friedman, 2009]. Ways to reduce model complexity, in order to avoid overfitting the training data \mathcal{D} , include searching in restricted structure spaces and penalizing it with adequate scores.

Common scores in structure learning are the penalized log-likelihood scores, such as the Akaike information criterion (AIC) [Akaike, 1974] and Bayesian information criterion (BIC) [Schwarz, 1978]. They measure the model's fitting of the empirical distribution $\widehat{P}(c, \mathbf{x})$ adding a penalty term that is a function of structure complexity. They are decomposable with respect to \mathcal{G} , allowing for efficient search algorithms. Yet, with limited N and a large n, discriminative scores based on $P(c \mid \mathbf{x})$, such as conditional log-likelihood and classification accuracy, are more suitable to the classification task [Friedman et al., 1997]. These, however, are not decomposable according to \mathcal{G} . While one can add a complexity penalty to discriminative scores [e.g., Grossman and Domingos, 2004], they are instead often cross-validated to induce preference towards structures that generalize better, making their computation even more time demanding.

For Bayesian network classifiers, a common [see Bielza and Larrañaga, 2014] structure space is that of augmented naive Bayes [Friedman et al., 1997] models (see Figure 1), factorizing $P(\mathbf{X}, C)$ as

$$P(\mathbf{X}, C) = P(C) \prod_{i=1}^{n} P(X_i \mid \mathbf{Pa}(X_i)), \tag{1}$$

with $C \in \mathbf{Pa}(X_i)$ for all X_i and $\mathbf{Pa}(C) = \emptyset$. Models of different complexity arise by extending or shrinking the parent sets $\mathbf{Pa}(X_i)$, ranging from the NB [Minsky, 1961] with $\mathbf{Pa}(X_i) = \{C\}$ for all X_i , to those with a limited-size $\mathbf{Pa}(X_i)$ [Friedman et al., 1997, Sahami, 1996], to those with unbounded $\mathbf{Pa}(X_i)$ [Pernkopf and O'Leary, 2003]. While the NB can only represent linearly separable classes [Jaeger, 2003], more complex models are more expressive [Varando et al., 2015]. Simpler models, with sparser $\mathbf{Pa}(X_i)$, may perform better with less training data, due to their lower variance, yet worse with more data as the bias due to wrong independence assumptions will tend to dominate the error.

The algorithms that produce the above structures are generally instances of greedy hill-climbing [Keogh and Pazzani, 2002, Sahami, 1996], with arc inclusion and removal as their search operators. Some [e.g., Pazzani, 1996] add node inclusion or removal, thus embedding feature selection [Guyon and Elisseeff, 2003] within structure learning. Alternatives include the adaptation [Friedman et al., 1997] of the Chow-Liu [Chow and Liu, 1968] algorithm to find the optimal one-dependence estimator (ODE) with respect to decomposable penalized log-likelihood scores in time quadratic in n. Some structures, such as NB or AODE, are fixed and thus require no search.

2.3 Learning parameters

Given \mathcal{G} , learning $\boldsymbol{\theta}$ in order to best approximate the underlying $P(C, \mathbf{X})$ is straightforward. For discrete variables X_i and $\mathbf{Pa}(X_i)$, Bayesian estimation can be obtained in closed form by assuming a Dirichlet prior over $\boldsymbol{\theta}$. With all Dirichlet hyper-parameters equal to α ,

$$\theta_{ijk} = \frac{N_{ijk} + \alpha}{N_{\cdot j} + r_i \alpha},\tag{2}$$

where N_{ijk} is the number of instances in \mathcal{D} such that $X_i = k$ and $\mathbf{pa}(x_i) = j$, corresponding to the j-th possible instantiation of $\mathbf{pa}(x_i)$, $N_{\cdot j}$ is the number of instances in which $\mathbf{pa}(x_i) = j$, while r_i is the cardinality of X_i . $\alpha = 0$ in Equation 2 yields the maximum likelihood estimate of θ_{ijk} . With incomplete data, the parameters of local distributions are no longer independent and we cannot separately maximize the likelihood for each X_i as in Equation 2. Optimizing the likelihood requires a time-consuming algorithm like expectation maximization [Dempster et al., 1977] which does not guarantee convergence to the global optimum.

While the NB can separate any two linearly separable classes given the appropriate θ , learning by approximating $P(C, \mathbf{X})$ cannot render the optimal θ in some cases [Jaeger, 2003]. Multiple methods [Hall, 2007, Zaidi et al., 2013, 2017] learn a weight $w_i \in [0, 1]$ for each feature and then update θ as

$$\theta_{ijk}^{weighted} = \frac{(\theta_{ijk})^{w_i}}{\sum_{k=1}^{r_i} (\theta_{ijk})^{w_i}}.$$

A $w_i < 1$ reduces the effect of X_i on the class posterior, with $w_i = 0$ omitting X_i from the model, making weighting more general than feature selection. The weights can be found by maximizing a discriminative score [Zaidi et al., 2013] or computing the usefulness of a feature in a decision tree [Hall, 2007]. Mainly applied to naive Bayes models, a generalization for augmented naive Bayes classifiers has been recently developed [Zaidi et al., 2017].

Another parameter estimation method for the naive Bayes is by means of Bayesian model averaging over the 2^n possible naive Bayes structures with up to n features [Dash and Cooper, 2002]. It is computed in time linear in n and provides the posterior probability of an arc from C to X_i .

2.4 Inference

Computing $P(c \mid \mathbf{x})$ for a fully observed \mathbf{x} means multiplying the corresponding $\boldsymbol{\theta}$. With an incomplete \mathbf{x} , however, exact inference requires summing over parameters of the local distributions and is NP-hard in the general case [Cooper, 1990], yet can be tractable with limited-complexity structures. The AODE ensemble computes $P(c \mid \mathbf{x})$ as the average of the $P_i(c \mid \mathbf{x})$ of the n base models. A special case is the lazy elimination [Zheng and Webb, 2006] heuristic which omits x_i from Equation 1 if $P(x_i \mid x_j) = 1$ for some x_j .

3 Functionalities

There are four groups of functionalities:

- 1. Learning network structure
- 2. Learning network parameters
- 3. Evaluating the model
- 4. Predicting with the model

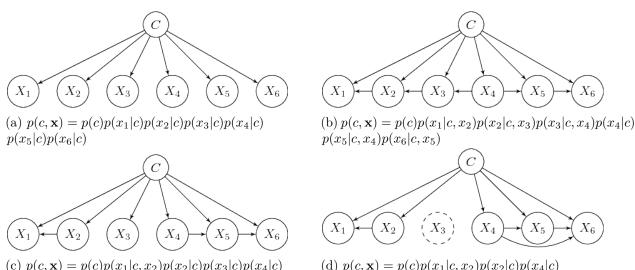
The first step is to learn structure and the second, usually, to learn parameters. The obtained models can be evaluated, used for prediction or analyzed. The following provides a brief overview. For details on some of the underlying methods please see the "methods" vignette.

3.1 Structures

The learning algorithms produce the following network structures:

- Naive Bayes (NB) (Figure 1a) [Minsky, 1961]
- k-dependence Bayesian classifier (k-DB) [Sahami, 1996, Pernkopf and Bilmes, 2010]
 - One-dependence estimators (ODE)
 - * Tree-augmented naive Bayes (TAN) (Figure 1b) [Friedman et al., 1997]
 - * Forest-augmented naive Bayes (FAN) (Figure 1c)
- Semi-naive Bayes (SNB)(Figure 1d) [Pazzani, 1996]
- Averaged one-dependence estimators (AODE) [Webb et al., 2005]

Figure 1 shows some of these structures and their factorizations of $P(c, \mathbf{x})$. We use k-DB in the sense meant by Pernkopf and Bilmes [2010] rather than that by Sahami [1996], as we impose no minimum on the number of augmenting arcs. SNB is the only structure whose complexity is not a priori bounded: the feature subgraph might be complete in the extreme case.



(c) $p(c, \mathbf{x}) = p(c)p(x_1|c, x_2)p(x_2|c)p(x_3|c)p(x_4|c)$ $p(x_5|c, x_4)p(x_6|c, x_5)$

(d) $p(c, \mathbf{x}) = p(c)p(x_1|c, x_2)p(x_2|c)p(x_4|c)$ $p(x_5|c, x_4)p(x_6|c, x_4, x_5)$

Figure 1: Augmented naive Bayes models produced by the **bnclassify** package. (a) NB; (b) TAN (c) FAN (d) SNB. k-DB and AODE not shown. The NB assumes that the features are independent given the class. ODE allows each predictor to depend on at most one other predictor: the TAN is a special case with exactly n-1 augmenting arcs (i.e., inter-feature arcs) while a FAN may have less than n-1. The k-DB allows for up to k parent features per feature X_i , with NB and ODE as its special cases with k=0 and k=1, respectively. The SNB does not restrict the number of parents but requires that connected feature subgraphs be complete (connected, after removing C, subgraphs in (d): $\{X_1, X_2\}$, and $\{X_4, X_5, X_6\}$), also allowing the removal of features (X_3) omitted in (d). The AODE is not a single structure but an ensemble of n ODE models in which one feature is the parent of all others (a super-parent).

3.2 Algorithms

Each structure learning algorithm is implemented by a single R function. Table 1 lists these algorithms along with the corresponding structures that they produce, the scores they can be combined with, and their R functions. Below we provide their abbreviations, references, brief comments and illustrate function calls.

3.2.1 Fixed structure

We implement two algorithms:

- NB
- AODE

The NB and AODE structures are fixed given the number of variables, and thus no search is required to estimate them from data. For example, we can get a NB structure with

n <- nb('class', dataset = car)</pre>

where class is the name of the class variable C and car the dataset containing observations of C and \mathbf{X} .

3.2.2 Optimal ODEs with decomposable scores

We implement one algorithm:

• Chow-Liu for ODEs (CL-ODE; Friedman et al. [1997])

Maximizing log-likelihood will always render a TAN while maximizing penalized log-likelihood may render a FAN since including some arcs can degrade such a score. With incomplete data our implementation does not guarantee the optimal ODE as that would require computing maximum likelihood parameters. The parameters of the tan_cl function are the network score to use and, optionally, the root for features' subgraph:

n <- tan_cl('class', car, score = 'AIC', root = 'buying')</pre>

3.2.3 Greedy hill-climbing with global scores

bnclassify implements five algorithms:

- Hill-climbing tree augmented naive Bayes (HC-TAN) [Keogh and Pazzani, 2002]
- Hill-climbing super-parent tree augmented naive Bayes (HC-SP-TAN) [Keogh and Pazzani, 2002]
- Backward sequential elimination and joining (BSEJ) [Pazzani, 1996]
- Forward sequential selection and joining (FSSJ) [Pazzani, 1996]
- Hill-climbing k-dependence Bayesian classifier (k-DB) [Pernkopf and Bilmes, 2010]

These algorithms use the cross-validated estimate of predictive accuracy as score. Only the FSSJ and BSEJ provide feature selection. The parameters of the corresponding functions include the number of cross-validation folds k and the minimal absolute score improvement epsilon required for continuing the search:

 $fssj \leftarrow fssj('class', car, k = 5, epsilon = 0)$

Table 1: Implemented structure learning algorithms.

Structure	Search algorithm	Score	Feature selection	Function
NB	-	-	-	nb
TAN/FAN	CL-ODE	log-lik, AIC, BIC	-	tan_cl
TAN	TAN-HC	accuracy	-	tan_hc
TAN	TAN-HCSP	accuracy	-	tan_hcsp
SNB	FSSJ	accuracy	forward	fssj
SNB	BSEJ	accuracy	backward	bsej
AODE	-	-	-	aode
kDB	kDB	accuracy	-	kdb

3.3 Parameters

bnclassify only handles discrete features. With fully observed data, it estimates the parameters with maximum likelihood or Bayesian estimation, according to Equation 2, with a single α for

all local distributions. With incomplete data it uses available case analysis and substitutes $N_{\cdot j}$. in Equation 2 with $N_{ij} = \sum_{k=1}^{r_i} N_{ijk}$, i.e., with the count of instances in which $\mathbf{Pa}(X_i) = j$ and X_i is observed.

We implement two methods for weighted naive Bayes parameter estimation:

- Weighting attributes to alleviate naive Bayes' independence assumption (WANBIA) [Zaidi et al., 2013]
- Attribute-weighted naive Bayes (AWNB) [Hall, 2007]

And one method for estimation by means of Bayesian model averaging over all NB structures with up to n features:

• Model averaged naive Bayes (MANB) [Dash and Cooper, 2002]

It makes little sense to apply WANBIA, MANB and AWNB to structures other than NB. WANBIA, for example, learns the weights by optimizing the conditional log-likelihood of the NB. Parameter learning is done with the 1p function. For example,

a <- lp(n, smooth = 1, manb_prior = 0.5)

computes Bayesian parameter estimates with $\alpha=1$ (the smooth argument) for all local distributions, and updates them with the MANB estimation obtained with a 0.5 prior probability for each class-to-feature arc.

3.4 Utilities

Single-structure-learning functions, as opposed to those that learn an ensemble of structures, return an S3 object of class bnc_dag. The following functions can be invoked on such objects:

- Plot the network: plot
- Query model type: is tan, is ode, is nb, is aode, ...
- Query model properties: narcs, families, features, ...
- Convert to **gRain** and **bnlearn** objects: as_grain

Ensembles are of type bnc_aode and only print and model type queries can be applied on such objects. Fitting the parameters (by calling lp) of a bnc_dag produces a bnc_bn object. In addition to all bnc_dag functions, the following are meaningful:

- Predict class labels and class posterior probability: predict
- Predict joint distribution: compute_joint
- Network scores: AIC, BIC, logLik, clogLik
- Cross-validated accuracy: cv
- Query model properties: nparams
- Parameter weights: manb_arc_posterior, weights

The above functions for bnc_bn can also be applied to an ensemble with fitted parameters.

3.5 Documentation

This vignette provides an overview of the package and background on the implemented methods. Calling <code>?bnclassify</code> gives a more concise overview of the functionalities, with pointers to relevant functions and their documentation. The "usage" vignette presents more detailed usage examples and shows how to combine the functions. The "methods" vignette provides details on the underlying methods and documents implementation specifics, especially where they differ from or are undocumented in the original paper.

4 Usage example

The available functionalities can be split into four groups:

- 1. Learning network structure and parameters
- 2. Analyzing the model
- 3. Evaluating the model
- 4. Predicting with the model

We illustrate these functionalities with the synthetic car data set with six features. We begin with a simple example for each functionality group and then elaborate on the options in the following sections. We first load the package and the dataset,

```
library(bnclassify)
data(car)
then learn a naive Bayes structure and its parameters,
nb <- nb('class', car)</pre>
nb \leftarrow lp(nb, car, smooth = 0.01)
get the number of arcs in the network,
narcs(nb)
[1] 6
get the 10-fold cross-validation estimate of accuracy,
cv(nb, car, k = 10)
[1] 0.861684
and finally classify the entire data set
p <- predict(nb, car)</pre>
head(p)
[1] unacc unacc unacc unacc unacc unacc
Levels: unacc acc good vgood
```

4.1 Learning

The functions for structure learning, shown in Table 1, correspond to the different algorithms. They all receive the name of the class variable and the data set as their first two arguments, which are then followed by optional arguments. The following runs the CL-ODE algorithm with the AIC score, followed by the FSSJ algorithm to learn another model:

```
ode_cl_aic <- tan_cl('class', car, score = 'aic')
set.seed(3)
fssj <- fssj('class', car, k = 5, epsilon = 0)</pre>
```

The bnc function is a shorthand for learning structure and parameters in a single step,

```
ode_cl_aic <- bnc('tan_cl', 'class', car, smooth = 1, dag_args = list(score = 'aic'))</pre>
```

where the first argument is the name of the structure learning function while and optional arguments go in dag_args.

4.2 Analyzing

Printing the model, such as the above ode_cl_aic object, provides basic information about it. ode_cl_aic

```
Bayesian network classifier
```

num. arcs: 9
free parameters: 131

learning algorithm: tan_cl

While plotting the network is especially useful for small networks, printing the structure in the **deal** [Bottcher and Dethlefsen, 2013] and **bnlearn** format may be more useful for larger ones:

```
ms <- modelstring(ode_cl_aic)
strwrap(ms, width = 60)
[1] "[class] [buying|class] [doors|class] [persons|class]"
[2] "[maint|buying:class] [safety|persons:class]"
[3] "[lug_boot|safety:class]"</pre>
```

We can query the type of structure; params lets us access the conditional probability tables (CPTs); while features lists the features:

```
is_ode(ode_cl_aic)
[1] TRUE
params(nb)$buying
       class
buying
                                            good
               unacc
                               acc
                                                        vgood
  low
        0.2132243562 0.2317727320 0.6664252607 0.5997847478
 med
        0.2214885458 \ 0.2994740131 \ 0.3332850521 \ 0.3999077491
  high 0.2677680077 0.2812467451 0.0001448436 0.0001537515
  vhigh 0.2975190903 0.1875065097 0.0001448436 0.0001537515
length(features(fssj))
[1] 5
```

For example, fssj has selected five out of six features.

manb_arc_posterior provides the MANB posterior probabilities for arcs from the class to each of the features:

With the posterior probability of 0% for the arc from class to doors, and 100% for all others, MANB renders doors independent from the class while leaving the other features' parameters unaltered. We can see this by printing out the CPTs:

params(manb)\$doors

```
class
doors unacc acc good vgood
2 0.25 0.25 0.25 0.25
3 0.25 0.25 0.25 0.25
4 0.25 0.25 0.25 0.25
5more 0.25 0.25 0.25 0.25
all.equal(params(manb)$buying, params(nb)$buying)
[1] TRUE
```

For more functions for querying a structure with parameters (bnc_bn) see ?inspect_bnc_bn. For a structure without parameters (bnc_dag), see ?inspect_bnc_dag.

4.3 Evaluating

Multiple scores can be computed:

```
logLik(ode_cl_aic, car)
'log Lik.' -13307.59 (df=131)
```

```
AIC(ode_cl_aic, car)
[1] -13438.59
```

The cv function estimates the predictive accuracy of one or more models with stratified cross-validation. In the following we assess the above models produced by NB and CL-ODE algorithms:

Above, k is the desired number of folds, and dag = TRUE evaluates structure and parameter learning, while dag = FALSE keeps the structure fixed and evaluates just the parameter learning. The output gives 86% and 93% accuracy estimates for NB and CL-ODE, respectively. mlr and caret packages provide additional options for evaluating predictive performance, such as different metrics, and bnclassify is integrated with both (see the "usage" vignette).

4.4 Predicting

As shown above, we can predict class labels with predict. We can also get the class posterior probabilities:

5 Implementation

With complete data, **bnclassify** implements prediction for augmented naive Bayes models as well as for ensembles of such models. It multiplies the corresponding θ in logarithmic space, applying the log-sum-exp trick before normalizing, to reduce the chance of underflow. On instances with missing entries, it uses the **gRain** package [Højsgaard, 2016, Højsgaard, 2012] to perform exact inference, which is notably slower. Network plotting is implemented by the **Rgraphviz** package. Some functions are implemented in C++ with **Rcpp** for efficiency. The package is extensively tested, with over 200 unit and integrated tests that give a 94% code coverage.

6 Related software

NB, TAN, and AODE are available in general-purpose tools such as **bnlearn** and Weka, while WANBIA² and MANB³ are only available in stand-alone software, published along with the original publications. We are not aware of available implementations of the remaining methods implemented in **bnclassify**.

bnlearn implements algorithms for learning general purpose Bayesian networks. Among them, algorithms for Markov blanket learning by testing for independences, such as IAMB [Tsamardinos and Aliferis, 2003] and GS [Margaritis and Thrun, 2000], can be very useful for classification

²https://sourceforge.net/projects/rawnaivebayes

³http://www.dbmi.pitt.edu/content/manb

as they can look for the Markov blanket of the class variable. **bnlearn** combines the search algorithms, such as greedy hill-climbing and tabu search [Glover and Laguna, 2013], only with generative scores such as penalized log-likelihood. Among classification models, it implements the discrete NB and CL-ODE. It does not handle incomplete data and provides cross-validation and prediction only for the NB and TAN models, but not for the unrestricted Bayesian networks.

Version 3.8 of Weka [Hall et al., 2009, Bouckaert, 2004] provides variants of the AODE [Webb et al., 2005] as well as the CL-ODE and NB. It implements five additional search algorithms, such as K2 [Cooper and Herskovits, 1992], tabu search and simulated annealing [Kirkpatrick et al., 1983], combining them only with generative scores. Except for the NB, Weka only handles discrete data and uses simple imputation (replacing with the mode or mean) to handle incomplete data. It provides two constraint-based algorithms, but performs conditional independence tests in an ad-hoc way [Bouckaert, 2004]. Weka provides Bayesian model averaging for parameter estimation [Bouckaert, 1995].

jBNC⁴ (version 1.2.2) is a Java library which learns ODE classifiers from Sacha et al. [2002] by optimizing penalized log-likelihood or the cross-validated estimate of accuracy. The CGBayes (version 7.14.14) package [McGeachie et al., 2014] for MATLAB implements conditional Gaussian networks [Lauritzen and Wermuth, 1989]. It provides four structure learning algorithms, including a variant of K2 and a greedy hill-climber, all optimizing the marginal likelihood of the data given the network.

7 Conclusion

bnclassify implements multiple state-of-the art algorithms for learning Bayesian network classifiers. It also provides features such as model analysis and evaluation. It is reasonably efficient and can handle large data sets. We hope that **bnclassify** will be useful to practitioners as well as to researchers wishing to compare their methods to existing ones.

Future work includes handling real-valued feature via conditional Gaussian models. Straightforward extensions include adding flexibility to the hill-climbing algorithm, such as restarts to avoid local minima.

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