

Package ‘BayesTree’

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Title Bayesian Methods for Tree Based Models

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Description Implementation of BART:Bayesian Additive Regression Trees, Chipman, George,
McCulloch (2009)

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Description

BART is a Bayesian “sum-of-trees” model.

For numeric response y , we have $y = f(x) + \epsilon$, where $\epsilon \sim N(0, \sigma^2)$.

For a binary response y , $P(Y = 1|x) = F(f(x))$, where F denotes the standard normal cdf (probit link).

In both cases, f is the sum of many tree models. The goal is to have very flexible inference for the unknown function f .

In the spirit of “ensemble models”, each tree is constrained by a prior to be a weak learner so that it contributes a small amount to the overall fit.

Usage

```
bart (
  x.train, y.train, x.test=matrix(0.0,0,0),
  sigest=NA, sigdf=3, sigquant=.90,
  k=2.0,
  power=2.0, base=.95,
  binaryOffset=0,
  ntree=200,
  ndpost=1000, nskip=100,
  printevery=100, keepevery=1, keeptrainfits=TRUE,
  usequants=FALSE, numcut=100, printcutoffs=0,
  verbose=TRUE)
## S3 method for class 'bart':
plot(
  x,
  plquants=c(.05,.95), cols =c('blue','black'),
  ...)
```

Arguments

<code>x.train</code>	<p>Explanatory variables for training (in sample) data. May be a matrix or a data frame, with (as usual) rows corresponding to observations and columns to variables. If a variable is a factor in a data frame, it is replaced with dummies. Note that q dummies are created if $q > 2$ and one dummy is created if $q = 2$, where q is the number of levels of the factor. <code>makeind</code> is used to generate the dummies. <code>bart</code> will generate draws of $f(x)$ for each x which is a row of <code>x.train</code>.</p>
<code>y.train</code>	<p>Dependent variable for training (in sample) data. If y is numeric a continuous response model is fit (normal errors). If y is a factor (or just has values 0 and 1) then a binary response model with a probit link is fit.</p>

<code>x.test</code>	Explanatory variables for test (out of sample) data. Should have same structure as <code>x.train</code> . <code>bart</code> will generate draws of $f(x)$ for each x which is a row of <code>x.test</code> .
<code>sigest</code>	The prior for the error variance (σ^2) is inverted chi-squared (the standard conditionally conjugate prior). The prior is specified by choosing the degrees of freedom, a rough estimate of the corresponding standard deviation and a quantile to put this rough estimate at. If <code>sigest=NA</code> then the rough estimate will be the usual least squares estimator. Otherwise the supplied value will be used. Not used if y is binary.
<code>sigdf</code>	Degrees of freedom for error variance prior. Not used if y is binary.
<code>sigquant</code>	The quantile of the prior that the rough estimate (see <code>sigest</code>) is placed at. The closer the quantile is to 1, the more aggressive the fit will be as you are putting more prior weight on error standard deviations (σ) less than the rough estimate. Not used if y is binary.
<code>k</code>	For numeric y , k is the number of prior standard deviations $E(Y x) = f(x)$ is away from ± 0.5 . The response (<code>y.train</code>) is internally scaled to range from -0.5 to 0.5 . For binary y , k is the number of prior standard deviations $f(x)$ is away from ± 0.5 . In both cases, the bigger k is, the more conservative the fitting will be.
<code>power</code>	Power parameter for tree prior.
<code>base</code>	Base parameter for tree prior.
<code>binaryOffset</code>	Used for binary y . The model is $P(Y = 1 x) = F(f(x) + \text{binaryOffset})$. The idea is that f is shrunk towards 0, so the offset allows you to shrink towards a probability other than 0.5 .
<code>ntree</code>	The number of trees in the sum.
<code>ndpost</code>	The number of posterior draws after burn in, <code>ndpost/keepevery</code> will actually be returned.
<code>nskip</code>	Number of MCMC iterations to be treated as burn in.
<code>printevery</code>	As the MCMC runs, a message is printed every <code>printevery</code> draws.
<code>keepevery</code>	Every <code>keepevery</code> draw is kept to be returned to the user. A “draw” will consist of values of the error standard deviation (σ) and $f^*(x)$ at $x = \text{rows}$ from the <code>train</code> (optionally) and <code>test</code> data, where f^* denotes the current draw of f .
<code>keeptrainfits</code>	If true the draws of $f(x)$ for $x = \text{rows of } x.train$ are returned.
<code>usequants</code>	Decision rules in the tree are of the form $x \leq c$ vs. $x > c$ for each variable corresponding to a column of <code>x.train</code> . <code>usequants</code> determines how the set of possible c is determined. If <code>usequants</code> is true, then the c are a subset of the values $(xs[i]+xs[i+1])/2$ where xs is unique sorted values obtained from the corresponding column of <code>x.train</code> . If <code>usequants</code> is false, the cutoffs are equally spaced across the range of values taken on by the corresponding column of <code>x.train</code> .
<code>numcut</code>	The number of possible values of c (see <code>usequants</code>). If a single number is given, this is used for all variables. Otherwise a vector with length equal to <code>ncol(x.train)</code> is required, where the i^{th} element gives the number of c used for the i^{th} variable

	in <code>x.train</code> . If <code>usequants</code> is false, <code>numcut</code> equally spaced cutoffs are used covering the range of values in the corresponding column of <code>x.train</code> . If <code>usequants</code> is true, then <code>min(numcut, the number of unique values in the corresponding columns of x.train - 1)</code> <code>c</code> values are used.
<code>printcutoffs</code>	The number of cutoff rules <code>c</code> to printed to screen before the MCMC is run. Give a single integer, the same value will be used for all variables. If 0, nothing is printed.
<code>verbose</code>	Logical, if FALSE suppress printing.
<code>x</code>	Value returned by <code>bart</code> which contains the information to be plotted.
<code>plquants</code>	In the plots, beliefs about $f(x)$ are indicated by plotting the posterior median and a lower and upper quantile. <code>plquants</code> is a double vector of length two giving the lower and upper quantiles.
<code>cols</code>	Vector of two colors. First color is used to plot the median of $f(x)$ and the second color is used to plot the lower and upper quantiles.
<code>...</code>	Additional arguments passed on to plot.

Details

BART is an Bayesian MCMC method. At each MCMC iteration, we produce a draw from the joint posterior $(f, \sigma)|(x, y)$ in the numeric y case and just f in the binary y case.

Thus, unlike a lot of other modelling methods in R, we do not produce a single model object from which fits and summaries may be extracted. The output consists of values $f^*(x)$ (and σ^* in the numeric case) where $*$ denotes a particular draw. The x is either a row from the training data (`x.train`) or the test data (`x.test`).

Value

The `plot` method sets `mfrow` to `c(1,2)` and makes two plots.

The first plot is the sequence of kept draws of σ including the burn-in draws. Initially these draws will decline as BART finds fit and then level off when the MCMC has burnt in.

The second plot has y on the horizontal axis and posterior intervals for the corresponding $f(x)$ on the vertical axis.

`bart` returns a list assigned class 'bart'. In the numeric y case, the list has components:

<code>yhat.train</code>	A matrix with <code>(ndpost/keepevery)</code> rows and <code>nrow(x.train)</code> columns. Each row corresponds to a draw f^* from the posterior of f and each column corresponds to a row of <code>x.train</code> . The (i, j) value is $f^*(x)$ for the i^{th} kept draw of f and the j^{th} row of <code>x.train</code> . Burn-in is dropped.
<code>yhat.test</code>	Same as <code>yhat.train</code> but now the <code>x</code> 's are the rows of the test data.
<code>yhat.train.mean</code>	train data fits = mean of <code>yhat.train</code> columns.
<code>yhat.test.mean</code>	test data fits = mean of <code>yhat.test</code> columns.
<code>sigma</code>	post burn in draws of <code>sigma</code> , length = <code>ndpost/keepevery</code> .

<code>first.sigma</code>	burn-in draws of sigma.
<code>varcount</code>	a matrix with (ndpost/keepevery) rows and nrow(x.train) columns. Each row is for a draw. For each variable (corresponding to the columns), the total count of the number of times that variable is used in a tree decision rule (over all trees) is given.
<code>sigest</code>	The rough error standard deviation (σ) used in the prior.
<code>y</code>	The input dependent vector of values for the dependent variable. This is used in plot.bart.

In the binary y case, the returned list has the components `yhat.train`, `yhat.test`, and `varcount` as above. In addition the list has a `binaryOffset` component giving the value used.

Note that in the binary y case `yhat.train` and `yhat.test` are $f(x) + \text{binaryOffset}$. If you want draws of the probability $P(Y = 1|x)$ you need to apply the normal cdf (`pnorm`) to these values.

Note

There was a bug in BayesTree_0.1-0 (now fixed of course).

If the number of test observations was less than the number of trees (200 is the default), the `yhat.test` and `yhat.test.mean` components were suspect.

Author(s)

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References

- Chipman, H., George, E., and McCulloch, R. (2009) BART: Bayesian Additive Regression Trees.
- Chipman, H., George, E., and McCulloch R. (2006) Bayesian Ensemble Learning. Advances in Neural Information Processing Systems 19, Scholkopf, Platt and Hoffman, Eds., MIT Press, Cambridge, MA, 265-272.
- both of the above at: <http://www.rob-mcculloch.org>
- Friedman, J.H. (1991) Multivariate adaptive regression splines. *The Annals of Statistics*, **19**, 1–67.

See Also

[pdbart](#)

Examples

```
##simulate data (example from Friedman MARS paper)
f = function(x){
  10*sin(pi*x[,1]*x[,2]) + 20*(x[,3]-.5)^2+10*x[,4]+5*x[,5]
}
sigma = 1.0 #y = f(x) + sigma*z , z~N(0,1)
n = 100     #number of observations
set.seed(99)
x=matrix(runif(n*10),n,10) #10 variables, only first 5 matter
```

```

Ey = f(x)
y=Ey+sigma*rnorm(n)
lmFit = lm(y~.,data.frame(x,y)) #compare lm fit to BART later
##run BART
set.seed(99)
bartFit = bart(x,y)
plot(bartFit) # plot bart fit
##compare BART fit to linear matter and truth = Ey
fitmat = cbind(y,Ey,lmFit$fitted,bartFit$yhat.train.mean)
colnames(fitmat) = c('y','Ey','lm','bart')
print(cor(fitmat))

```

makeind

Build x matrix from x data frame (convert factors to dummies)

Description

Converts factors to dummies.

Note that with more than one level, BART needs a dummy for each level of a factor (unlike in linear regression where one of the dummies is dropped).

Usage

```

makeind(
  x,
  all=TRUE)

```

Arguments

<code>x</code>	Data frame of explanatory variables.
<code>all</code>	If <code>all=TRUE</code> , a factor with <code>p</code> levels will be replaced by all <code>p</code> dummies. If <code>all=FALSE</code> , the <code>p</code> th dummy is dropped.

Details

Uses function `class.ind` from the `nnet` library. Note that if you have train and test data frames, it may be best to `rbind` the two together, apply `makeind` to the result, and then pull them back apart.

Value

A matrix.
Numerical variables come first, and then the appended dummies.

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See Also[bart](#)**Examples**

```
x1 = 1:10
x2 = as.factor(c(rep(1,5),rep(2,5)))
x3 = as.factor(c(1,1,1,2,2,2,4,4,4,4))
levels(x3) = c('rob','hugh','ed')
x = data.frame(x1,x2,x3)

junk = makeind(x)
```

pdbart

*Partial Dependence Plots for BART***Description**

Run `bart` at test observations constructed so that a plot can be created displaying the effect of a single variable (`pdbart`) or pair of variables (`pd2bart`). Note the y is a binary with $P(Y = 1|x) = F(f(x))$ with F the standard normal cdf, then the plots are all on the f scale.

Usage

```
pdbart (
  x.train, y.train,
  xind=1:ncol(x.train), levs=NULL, levquants=c(.05, (1:9)/10, .95),
  pl=TRUE, plquants=c(.05,.95), ...)
## S3 method for class 'pdbart':
plot (
  x,
  xind = 1:length(x$fd),
  plquants =c(.05,.95), cols=c('black','blue'), ...)
pd2bart (
  x.train, y.train,
  xind=1:2, levs=NULL, levquants=c(.05, (1:9)/10, .95),
  pl=TRUE, plquants=c(.05,.95), ...)
## S3 method for class 'pd2bart':
plot (
  x,
  plquants =c(.05,.95), contour.color='white',
  justmedian=TRUE, ...)
```

Arguments

<code>x.train</code>	<p>Explanatory variables for training (in sample) data. Must be a matrix (typeof double) with (as usual) rows corresponding to observations and columns to variables. Note that for a categorical variable you need to use dummies and if there are more than two categories, you need to put all the dummies in (unlike linear regression).</p>
<code>y.train</code>	<p>Dependent variable for training (in sample) data. Must be a vector (typeof double) with length equal to the number of observations (equal to the number of rows of <code>x.train</code>).</p>
<code>xind</code>	<p>Integer vector indicating which variables are to be plotted. In <code>pdbart</code>, variables (columns of <code>x.train</code>) for which plot is to be constructed. In <code>plot.pdbart</code>, indices in list returned by <code>pdbart</code> for which plot is to be constructed. In <code>pd2bart</code>, integer vector of length 2, indicating the pair of variables (columns of <code>x.train</code>) to plot.</p>
<code>levs</code>	<p>Gives the values of a variable at which the plot is to be constructed. List, where i^{th} component gives the values for i^{th} variable. In <code>pdbart</code>, should have same length as <code>xind</code>. In <code>pd2bart</code>, should have length 2. See also argument <code>levquants</code>.</p>
<code>levquants</code>	<p>If <code>levs</code> in NULL, the values of each variable used in the plot is set to the quantiles (in <code>x.train</code>) indicated by <code>levquants</code>. Double vector.</p>
<code>pl</code>	<p>For <code>pdbart</code> and <code>pd2bart</code>, if true, plot is made (by calling <code>plot.*</code>).</p>
<code>plquants</code>	<p>In the plots, beliefs about $f(x)$ are indicated by plotting the posterior median and a lower and upper quantile. <code>plquants</code> is a double vector of length two giving the lower and upper quantiles.</p>
<code>...</code>	<p>Additional arguments. In <code>pdbart</code>,<code>pd2bart</code>, passed on to <code>bart</code>. In <code>plot.pdbart</code>, passed on to <code>plot</code>. In <code>plot.pd2bart</code>, passed on to <code>image</code></p>
<code>x</code>	<p>For <code>plot.*</code>, object returned from <code>pdbart</code> or <code>pd2bart</code>.</p>
<code>cols</code>	<p>Vector of two colors. First color is for median of f, second color is for the upper and lower quantiles.</p>
<code>contour.color</code>	<p>Color for contours plotted on top of the image.</p>
<code>justmedian</code>	<p>Boolean, if true just one plot is created for the median of $f(x)$ draws. If false, three plots are created one for the median and two additional ones for the lower and upper quantiles. In this case, <code>mfrow</code> is set to <code>c(1,3)</code>.</p>

Details

We divide the predictor vector x into a subgroup of interest, x_s and the complement $x_c = x \setminus x_s$. A prediction $f(x)$ can then be written as $f(x_s, x_c)$. To estimate the effect of x_s on the prediction,

Friedman suggests the partial dependence function

$$f_s(x_s) = \frac{1}{n} \sum_{i=1}^n f(x_s, x_{ic})$$

where x_{ic} is the i^{th} observation of x_c in the data. Note that (x_s, x_{ic}) will generally not be one of the observed data points. Using BART it is straightforward to then estimate and even obtain uncertainty bounds for $f_s(x_s)$. A draw of $f_s^*(x_s)$ from the induced BART posterior on $f_s(x_s)$ is obtained by simply computing $f_s^*(x_s)$ as a byproduct of each MCMC draw f^* . The median (or average) of these MCMC draws $f_s^*(x_s)$ then yields an estimate of $f_s(x_s)$, and lower and upper quantiles can be used to obtain intervals for $f_s(x_s)$.

In `pdbart` x_s consists of a single variable in x and in `pd2bart` it is a pair of variables.

This is a computationally intensive procedure. For example, in `pdbart`, to compute the partial dependence plot for 5 x_s values, we need to compute $f(x_s, x_c)$ for all possible (x_s, x_{ic}) and there would be $5n$ of these where n is the sample size. All of that computation would be done for each kept BART draw. For this reason running BART with `keepevery` larger than 1 (eg. 10) makes the procedure much faster.

Value

The plot methods produce the plots and don't return anything.

`pdbart` and `pd2bart` return lists with components given below. The list returned by `pdbart` is assigned class 'pdbart' and the list returned by `pd2bart` is assigned class 'pd2bart'.

<code>fd</code>	<p>A matrix whose (i, j) value is the i^{th} draw of $f_s(x_s)$ for the j^{th} value of x_s. "fd" is for "function draws".</p> <p>For <code>pdbart</code> <code>fd</code> is actually a list whose k^{th} component is the matrix described above corresponding to the k^{th} variable chosen by argument <code>xind</code>.</p> <p>The number of columns in each matrix will equal the number of values given in the corresponding component of argument <code>levs</code> (or number of values in <code>levquants</code>).</p> <p>For <code>pd2bart</code>, <code>fd</code> is a single matrix. The columns correspond to all possible pairs of values for the pair of variables indicated by <code>xind</code>. That is, all possible (x_i, x_j) where x_i is a value in the <code>levs</code> component corresponding to the first x and x_j is a value in the <code>levs</code> components corresponding to the second one. The first x changes first.</p>
<code>levs</code>	<p>The list of levels used, each component corresponding to a variable. If argument <code>levs</code> was supplied it is unchanged. Otherwise, the levels in <code>levs</code> are as constructed using argument <code>levquants</code>.</p>
<code>xlbs</code>	<p>vector of character strings which are the plotting labels used for the variables.</p>

The remaining components returned in the list are the same as in the value of `bart`. They are simply passed on from the BART run used to create the partial dependence plot. The function `plot.bart` can be applied to the object returned by `pdbart` or `pd2bart` to examine the BART run.

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References

Chipman, H., George, E., and McCulloch, R. (2006) BART: Bayesian Additive Regression Trees.
 Chipman, H., George, E., and McCulloch R. (2006) Bayesian Ensemble Learning.
 both of the above at: <http://faculty.chicagogsb.edu/robert.mcculloch/research/rob-mcculloch-cv.html>
 Friedman, J.H. (2001) Greedy function approximation: A gradient boosting machine. *The Annals of Statistics*, **29**, 1189–1232.

Examples

```
##simulate data
f = function(x) { return(.5*x[,1] + 2*x[,2]*x[,3]) }
sigma=.2 # y = f(x) + sigma*z
n=100 #number of observations
set.seed(27)
x = matrix(2*runif(n*3)-1,ncol=3) ; colnames(x) = c('rob','hugh','ed')
Ey = f(x)
y = Ey + sigma*rnorm(n)
lmFit = lm(y~.,data.frame(x,y)) #compare lm fit to BART later
par(mfrow=c(1,3)) #first two for pdbart, third for pd2bart
##pdbart: one dimensional partial dependence plot
set.seed(99)
pdb1 = pdbart(x,y,xind=c(1,2),
  levs=list(seq(-1,1,.2),seq(-1,1,.2)),pl=FALSE,
  keepevery=10,ntree=100)
plot(pdb1,ylim=c(-.6,.6))
##pd2bart: two dimensional partial dependence plot
set.seed(99)
pdb2 = pd2bart(x,y,xind=c(2,3),
  levquants=c(.05,.1,.25,.5,.75,.9,.95),pl=FALSE,
  ntree=100,keepevery=10,verbose=FALSE)
plot(pdb2)
##compare BART fit to linear model and truth = Ey
fitmat = cbind(y,Ey,lmFit$fitted,pdb1$yhat.train.mean)
colnames(fitmat) = c('y','Ey','lm','bart')
print(cor(fitmat))
## plot.bart(pdb1) displays the BART run used to get the plot.
```

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