Using BART with a virtual species!

Getting Started

So you're interested in using embarcadero to do species distribution modeling with Bayesian additive regression trees! That's great. BARTs are a powerful way to do machine learning and, while not a new method per se, they are very new for SDMs.

Most of the core functionality of embarcadero is actually a wrapper for dbarts, which runs the actual BART fitting process. This vignete will show you

- 1. How to run BARTs
- 2. Variable importance measures
- 3. Automated variable selection
- 4. Partial dependence plots
- 5. Visualizing the posterior distribution

There's also just going to be some general comments on the process of using BARTs, the challenges to working with them, and some things that are hopefully coming next.

If you want to install, do it using devtools for now:

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#> Loading required package: raster
#> Loading required package: sp
#>
#> Attaching package: 'raster'
#> The following object is masked from 'package:dplyr':
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       select
#>
#> The following object is masked from 'package:tidyr':
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       extract
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#> Loading required package: ROCR
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#> The following object is masked from 'package:raster':
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       area
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       get_legend
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       rotate
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       between, first, last
#> The following object is masked from 'package:purrr':
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       transpose
```

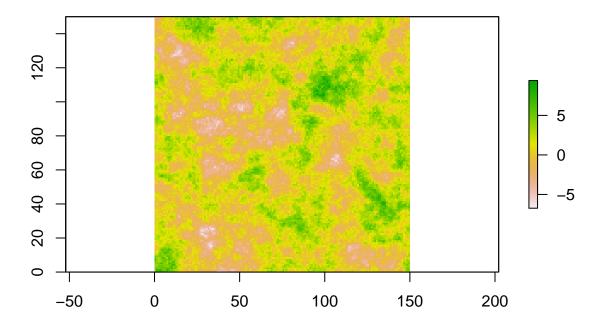
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#>
#> Attaching package: 'embarcadero'
#> The following object is masked from 'package:purrr':
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Doors are closing; please stand clear of the doors.

Creating the virtual species

First, let's create an imaginary landscape. We do this using the NLMR package:

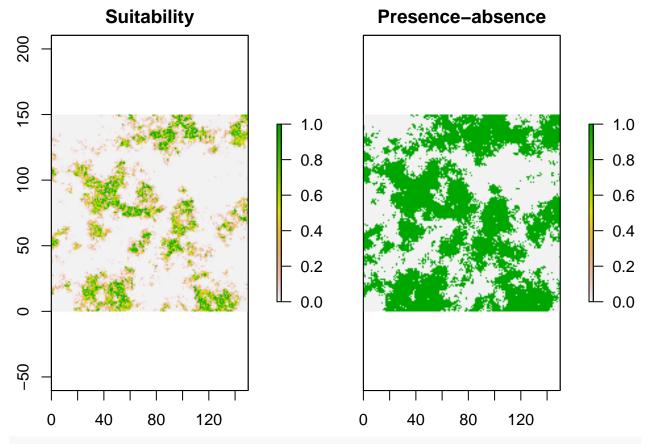
An imaginary variable

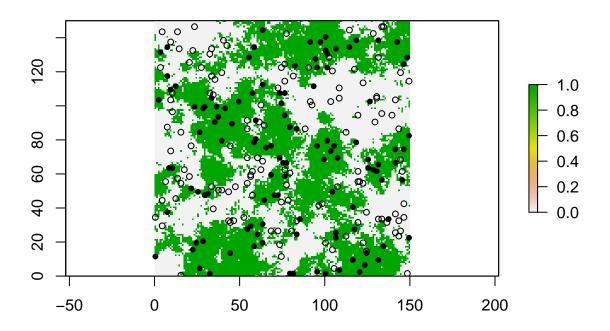


Next, let's make a species using the 'virtualspecies' package. Our imaginary species will only responds to variables 1-4, making variables 5-8 uninformative predictors (hopefully our model will drop them):

```
# Generate the species' climatic niche
random.sp <- generateRandomSp(climate[[1:4]],</pre>
```

```
# ^ These are the informative predictors
                              approach="pca",
                              relations='gaussian',
                              species.prevalence=0.5,
                              realistic.sp = TRUE,
                              PA.method='threshold')
#> - Perfoming the pca
   - Defining the response of the species along PCA axes
   - Calculating suitability values
      The final environmental suitability was rescaled between 0 and 1.
#>
#>
                     To disable, set argument rescale = FALSE
#>
   - Converting into Presence - Absence
#>
      Threshold conversion finished:
#> - cutoff = 0.00103681027010919
#> - species prevalence =0.5
```

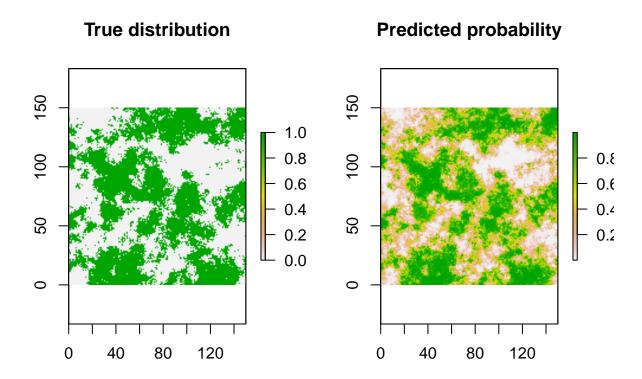




Alright. Now that we have the dataset, let's get to modeling!

We could easily throw all our data in one model, run it on defaults, make a map, and never think about it again. There's no laws against it.

```
#> use quantiles for rule cut points: false
#> data:
#> number of training observations: 250
#> number of test observations: 0
#> number of explanatory variables: 8
#>
\#> Cutoff rules c in x<=c vs x>c
#> Number of cutoffs: (var: number of possible c):
#> (1: 100) (2: 100) (3: 100) (4: 100) (5: 100)
#> (6: 100) (7: 100) (8: 100)
#> offsets:
#> reg : 0.00 0.00 0.00 0.00 0.00
#> Running mcmc loop:
#> iteration: 100 (of 1000)
#> iteration: 200 (of 1000)
#> iteration: 300 (of 1000)
#> iteration: 400 (of 1000)
#> iteration: 500 (of 1000)
#> iteration: 600 (of 1000)
#> iteration: 700 (of 1000)
#> iteration: 800 (of 1000)
#> iteration: 900 (of 1000)
#> iteration: 1000 (of 1000)
#> total seconds in loop: 3.233851
#>
#> Tree sizes, last iteration:
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#> Variable Usage, last iteration (var:count):
#> (1: 33) (2: 27) (3: 29) (4: 42) (5: 22)
#> (6: 40) (7: 37) (8: 33)
#> DONE BART
# Predict the species distribution!
map <- predict(sdm, climate, splitby=5)</pre>
#> Estimated time to total prediction (mins):
#> 1.111144
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```



This is.... alright? We could probably do better. One of the easiest ways is to cut some variables we don't think are performing well. In BART, pushing the models towards a smaller number of trees forces the variables to compete a bit, and preferentially upweights the better ones. We can plot that!

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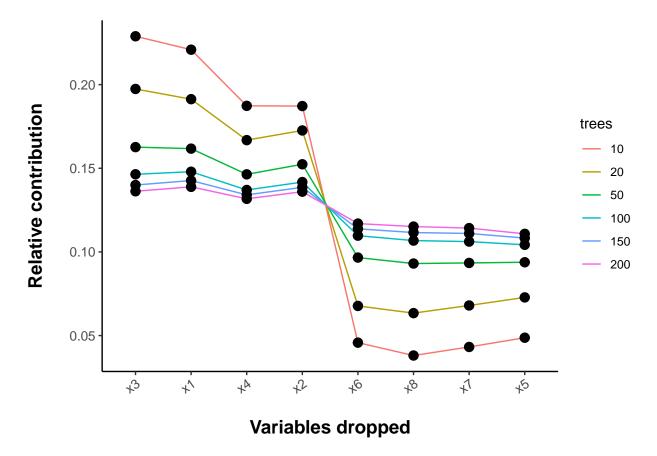
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OK. This would suggest that some of the variables, like x5 and x7, are less important while x3 and x4 are more important. But that doesn't give us a concrete list of what to drop. Luckily, embarcadero has a stepwise variable set reduction function. Let's run that, and then retrain the model.

```
# Stepwise variable set reduction
step.model <- variable.step(x.data=occ.df[,2:9],</pre>
                                 y.data=occ.df[,'Observed'])
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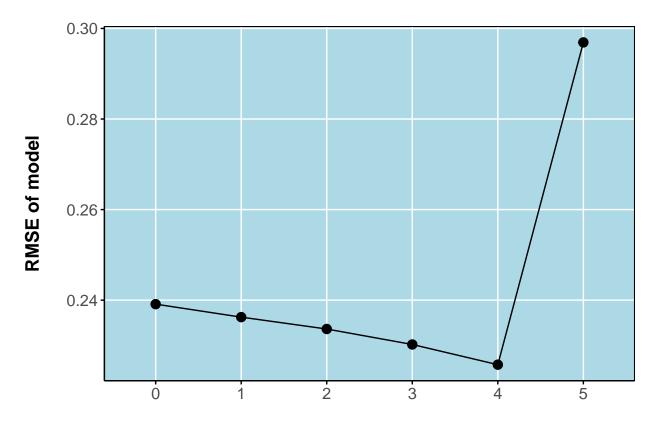
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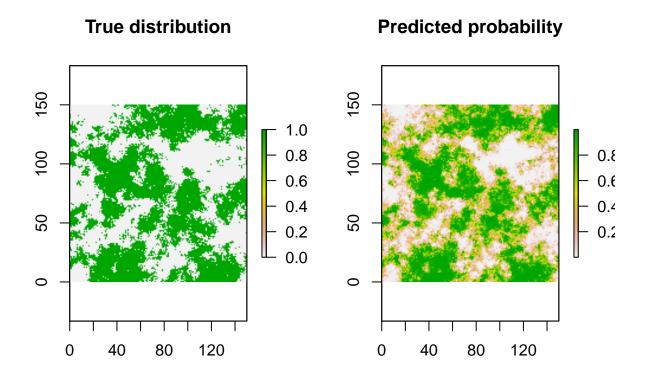
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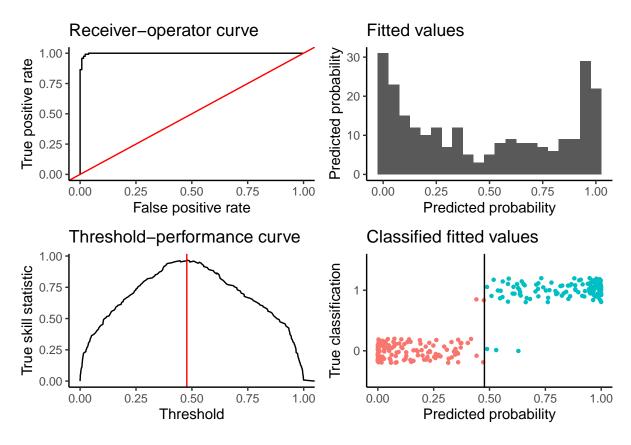
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#> [1] -----
#> [1] Final recommended variable list
#> [1] x1 x2 x3 x4
step.model
#> [1] "x1" "x2" "x3" "x4"
# Retrain the model
sdm <- bart(x.train=occ.df[, step.model], y.train=occ.df[,'Observed'],</pre>
           keeptrees = TRUE)
#>
#> Running BART with binary y
#> number of trees: 200
#> number of chains: 1, number of threads 1
#> Prior:
#> k: 2.000000
   power and base for tree prior: 2.000000 0.950000
#> use quantiles for rule cut points: false
#> number of training observations: 250
#> number of test observations: 0
#> number of explanatory variables: 4
#>
#> Cutoff rules c in x<=c vs x>c
#> Number of cutoffs: (var: number of possible c):
#> (1: 100) (2: 100) (3: 100) (4: 100)
```

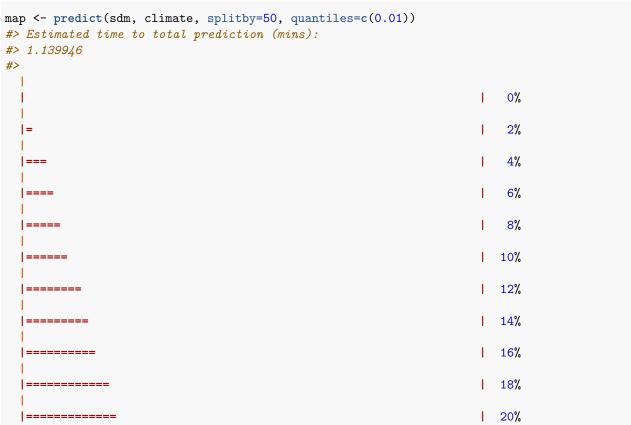
```
#>
#> offsets:
#> reg : 0.00 0.00 0.00 0.00 0.00
#> Running mcmc loop:
#> iteration: 100 (of 1000)
#> iteration: 200 (of 1000)
#> iteration: 300 (of 1000)
#> iteration: 400 (of 1000)
#> iteration: 500 (of 1000)
#> iteration: 600 (of 1000)
#> iteration: 700 (of 1000)
#> iteration: 800 (of 1000)
#> iteration: 900 (of 1000)
#> iteration: 1000 (of 1000)
#> total seconds in loop: 2.641285
#>
#> Tree sizes, last iteration:
#> [1] 2 3 2 3 2 4 3 2 2 2 2 2 2 3 3 2 3 2
#> 3 4 2 3 2 2 2 3 1 2 3 2 2 2 3 2 2 2 2 2
#> 3 2 2 2 4 2 2 2 2 2 2 2 2 2 3 2 2 3 3 2
#> 2 3 2 1 2 3 1 2 2 4 3 2 2 4 3 2 3 4 2 3
#> 2 2 2 2 3 2 3 1 4 2 3 2 2 3 1 3 3 3 2 2
#> 3 2 4 2 2 4 3 3 3 2 2 2 4 2 2 4 4 3 3 2
#> 2 2 3 3 3 2 1 2 2 3 3 2 3 3 3 2 3 2 2 2
#> 2 4 2 2 2 3 4 2 2 2 2 3 2 3 5 2 2 3 2 2
#> 2 3 3 2 3 2 2 2 2 3 3 3 2 4 3 3 5 3 2 3
#> 2 2
#>
#> Variable Usage, last iteration (var:count):
#> (1: 64) (2: 88) (3: 78) (4: 66)
#> DONE BART
# Predict the species distribution!
map <- predict(sdm, climate)</pre>
# How's it look?
par(mfrow=c(1,2))
plot(random.sp$pa.raster, main='True distribution')
plot(map, main='Predicted probability')
```



OK. We have a good-looking model. What's inside?

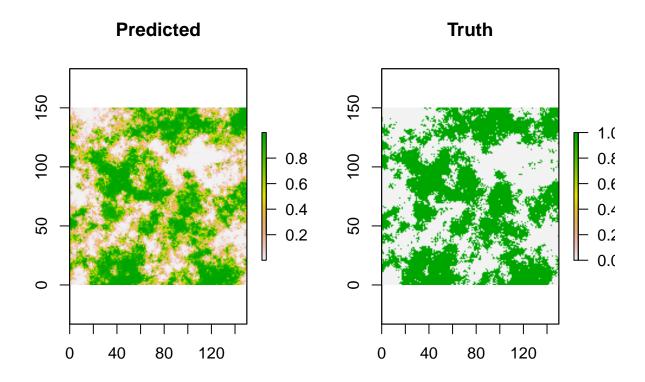
```
# How good is it?
summary(sdm)
#> Call: bart occ.df[, step.model] occ.df[, "Observed"] TRUE
#>
#> Predictor list:
#> x1 x2 x3 x4
#>
#> Area under the receiver-operator curve
#> AUC = 0.9983308
#>
#> Recommended threshold (maximizes true skill statistic)
#> Cutoff = 0.4771741
#> TSS = 0.9687982
#> Resulting type I error rate: 0.008474576
#> Resulting type II error rate: 0.02272727
```





	I	22%
	1	24%
	1	26%
	1	28%
	1	30%
	1	32%
	1	34%
	1	36%
	1	38%
	1	40%
 ===================================	1	42%
 ===================================	1	44%
 ===================================	1	46%
 ===================================	1	48%
	1	50%
	1	52%
	1	54%
	1	56%
	1	58%
	1	60%
	1	62%
	1	64%
	1	66%
 	1	68%
 	1	70%
	1	72%

```
| 74%
                                     76%
                                     78%
                                     80%
 _____
                                     82%
 84%
                                     86%
                                     88%
                                     90%
                                     92%
                                     94%
                                     96%
# Check
par(mfrow=c(1,2))
plot(map[[1]], main='Predicted')
plot(random.sp$pa.raster, main='Truth')
```



#plot(values(map) ~ values(random.sp\$pa.raster))