

# Allometry\_testing

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## Notes about setup—there are two options:

1. Clone the pecan repo directly, then open this rmd in the repository. (Maybe best option because of dependancies)
2. You can install the allometry module from github with the code below (from allometry vignette):  
`library(devtools) # if not run, need to install the pecan modules install_github("PecanProject/pecan",subdir="base/logging")`  
`install_github("PecanProject/pecan",subdir="allometry")`

## Basic Pecan allometry tutorial

A lot of this text and code is just annotated from the allometry vignette. To view the different components of the tree for which you could develop allomtries + their abbreviations:

```
data("allom.components")
allom.components
```

##		description	abbreviation	ID
## 1	Complete tree (above + belowground)		BTT	1
## 2	Whole tree (aboveground)		BAT	2
## 3	Whole tree (above stump)			3
## 4	Stem (wood only)		BSW	4
## 5	Stem (bark only)		BSB	5
## 6	Stem (wood + bark)		BST	6
## 7	Stem top			7
## 8	Branches live		BBL	8
## 9	Branches live < 2.5 cm		BBL_1	9
## 10	Branches live 2.5-7.6 cm		BBL_2	10
## 11	Branches live > 7.6 cm		BBL_3	11
## 12	Branches dead		BBD	12
## 13	Branches total (live + dead)		BBT	13
## 14	Stem + branches (bark only)			14
## 15	Stem + branches (wood only)			15
## 16	Stem + branches (live)		BAP	16
## 17	Wood, bark, branches (live + dead; no twigs or foliage)		BAE	17
## 18	Foliage total		BFT	18
## 19	Foliage new		BFN	19
## 20	Foliage old		BFO	20
## 21	Twigs total		BBG	21
## 22	Twigs old		BBG_0	22
## 23	Foliage + twigs		BFG	23
## 24	Crown (branches + foliage + twigs)		BCT	24
## 25	Roots, coarse > 3 mm dia		BKL	25

## 26	Coarse stump roots	BSR 26
## 27	Coarse lateral roots	BLR 27
## 28	Fine roots	BFR 28
## 29	Roots total	BRT 29
## 30	Stump wood	30
## 31	Stump bark	31
## 32	Stump total	32
## 33	Stump + roots	33
## 34	Cones	34
## 35	Live crown (branches + foliage + twigs)	BCL 35
## 36	Dead crown (branches + foliage + twigs)	BCD 36
## 37	Small branches	BBS 37
## 38	Height	HT 40
## 39	Rooting Depth	Rd 41
## 40	Rooting Volume	Vol 42
## 41	Canopy Area	43

```
pfts = list(PIPO = data.frame(spcd=122,acronym='PIPO')) # list our "Pfts--plant functional types" of in
```

## AllomAve function in PEcAn allometry module

- Fits the bayesian allometric models for the pfts. ngibbs indicates the number of gibbs/mcmc samples
- looks like there is a burn in of ~200
- automatically outputs a pdf with the traceplots titled “allom.pipo.6.mcmc” and saves and .Rdata object.
- not totally clear to be what the different betas are

```
allom.stats = AllomAve(pfts,components =6, ngibbs=1000)
```

```
## [1] "writing output to"
## [2] "/home/miranda/agb_allometry/pecan/modules/allometry"
## $spcd
## [1] 122
##
## $acronym
## [1] PIPO
## Levels: PIPO
##
## [[3]]
## [1] 6
##
## [1] "Dropping allom rows: "
## |
```

|

```
allom.stats$PIPO
```

```
## [[1]]
## NULL
##
## [[2]]
## NULL
##
## [[3]]
## NULL
##
## [[4]]
## NULL
##
## [[5]]
## NULL
##
## [[6]]
##
## Iterations = 1:750
## Thinning interval = 1
## Number of chains = 3
## Sample size per chain = 750
##
## 1. Empirical mean and standard deviation for each variable,
##    plus standard error of the mean:
##
##      Mean      SD Naive SE Time-series SE
## b0.1  -0.9523  0.7621 0.016065      0.016659
## b0.2  -0.9313  1.3981 0.029475      0.029477
## b0.3  -2.5069  0.7266 0.015319      0.021378
## b0.4  -2.7044  0.8791 0.018534      0.025256
## b0.5  -2.7890  0.9093 0.019170      0.024355
## b0.6  -3.4818  0.5453 0.011495      0.016797
## b1.1   1.1348  0.4301 0.009068      0.009437
## b1.2   1.1938  0.7471 0.015749      0.015750
## b1.3   2.3865  0.2451 0.005167      0.007142
## b1.4   2.3977  0.3038 0.006405      0.008457
## b1.5   2.5033  0.3103 0.006543      0.008246
## b1.6   2.6447  0.1303 0.002747      0.004001
## mu0   -2.2168  0.9403 0.019822      0.024052
## mu1    2.0343  0.5206 0.010976      0.011944
## sigma  0.1315  0.0685 0.001444      0.001684
## tau11   4.4226  8.2397 0.173709      0.251329
## tau12  -2.4383  5.2708 0.111119      0.145956
## tau22   1.5184  3.6768 0.077513      0.093979
## D      66.6461 60.7116 1.279913      1.530546
## Bg0    -2.3370  0.6158 0.012983      0.012996
## Bg1     2.3338  0.1759 0.003709      0.003621
## Sg      0.4620  0.2381 0.005019      0.004873
## Dg     183.4849 54.2587 1.143873      1.130324
##
## 2. Quantiles for each variable:
##
```

##	2.5%	25%	50%	75%	97.5%
## b0.1	-2.69658	-1.37241	-0.8308	-0.41975	0.2091
## b0.2	-4.50763	-1.64672	-0.6186	0.08552	0.9032
## b0.3	-3.92746	-2.96619	-2.5104	-2.05727	-1.0503
## b0.4	-4.47173	-3.26932	-2.6821	-2.13836	-0.9816
## b0.5	-4.59567	-3.35675	-2.7766	-2.23158	-1.0209
## b0.6	-4.54504	-3.83625	-3.4744	-3.14461	-2.3972
## b1.1	0.46900	0.82182	1.0737	1.38727	2.0667
## b1.2	0.16537	0.63913	1.0529	1.59236	3.0019
## b1.3	1.90295	2.23364	2.3886	2.54015	2.8621
## b1.4	1.80130	2.19708	2.3923	2.59235	3.0033
## b1.5	1.88430	2.31358	2.5032	2.69454	3.0920
## b1.6	2.38859	2.56335	2.6429	2.72939	2.9038
## mu0	-4.06829	-2.75280	-2.2013	-1.68668	-0.3246
## mu1	0.94255	1.75547	2.0396	2.32275	3.0656
## sigma	0.05771	0.08685	0.1132	0.15488	0.3190
## tau11	0.46344	1.43589	2.5545	4.83135	18.3376
## tau12	-10.03039	-2.62083	-1.4225	-0.79702	-0.2328
## tau22	0.19125	0.53116	0.9020	1.61909	5.8851
## D	-13.74784	27.78100	54.9703	91.45138	218.8039
## Bg0	-3.46674	-2.77279	-2.3433	-1.90377	-1.1147
## Bg1	1.98478	2.21225	2.3353	2.45911	2.6583
## Sg	0.13629	0.28115	0.4161	0.61514	1.0171
## Dg	72.84246	146.65021	187.6497	226.26288	276.6505

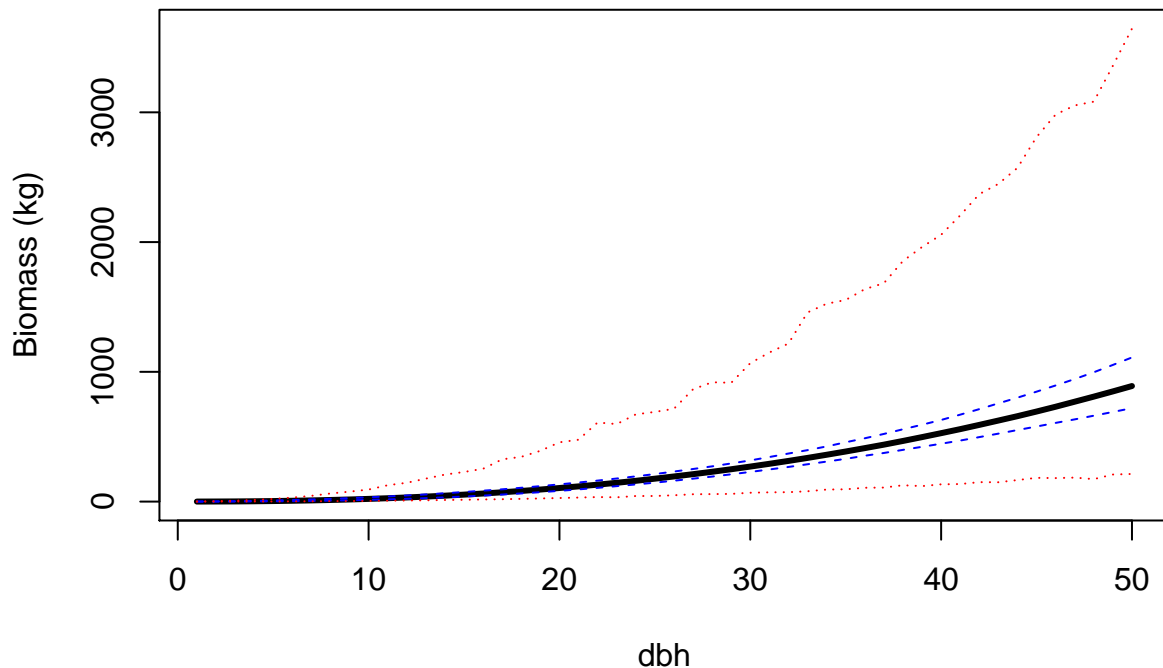
## Predict for individual trees:

- the allom.predict fuction will use the fit relationships from allomAve to predict the biomass component for the pft of interest (PIPO) over the DBH vector values:
- dbh = diameter to predict at
- component is which component of biomass to predict see defintation in the allom.components object
- unclear to me whether component = 6 is predicting Stem biomass, or is only suing the diameter of the stem in the allometry equation
- use = Bg

*##Predict for individual trees:*

```
allom.fit = load.allom(getwd()) # get the allometry listed in our working directory
dbh = 1:50 # vector of DBH values to predict over
```

```
pred = allom.predict(allom.fit, dbh = dbh, pft = "PIPO", component =6, use = "Bg", interval = "prediction")
conf = allom.predict(allom.fit,dbh = dbh,pft = "PIPO",component = 6,use = "Bg",interval = "confidence")
PI = apply(pred,2,quantile,c(0.025,0.5,0.975),na.rm=TRUE)
CI = apply(conf,2,quantile,c(0.025,0.5,0.975),na.rm=TRUE)
plot(dbh,CI[2,],type='l',lwd=3,ylim=range(PI),ylab="Biomass (kg)")
lines(dbh,CI[1,],lty=2,col="blue")
lines(dbh,CI[3,],lty=2,col="blue")
lines(dbh,PI[1,],lty=3,col="red")
lines(dbh,PI[3,],lty=3,col="red")
```



### other stuff I learned from digging:

- there is a file in the pecan repository that has a table with all the beta parameters from jenkins
- really we want to use the `query.allom.data` function to query the “field” data, but there is a database that this function connects to via “con=” argument...and I can’t figure out where this database is located/what the connection should be

```
jenkins.table <- read.csv(system.file("data/Table3_GTR-NE-319.v2.csv", package = "PEcAn.allometry"), header = TRUE)
#colnames(jenkins.table)[1] <- c("")
head(jenkins.table)
```

```
## Table.3.....Equations.and.parameters.for.diameter.based.biomass.equations
## 1
## 2
## 3 Species
## 4 0
## 5 0
## 6 0
## 7 0
## 8
## 9 X X.1 X.2 X.3 X.4 X.5
## 10
## 11
## 12 Common Name Component ID Equation Form ID a b c
## 13 eastern conifers 2 7 0.5 15000 2.7
## 14 softwoods (general) 3 1 -1.01 2.41 1
## 15 softwoods (general) 3 4 4.5966 -0.2364 0.00411
## 16 softwoods (general) 6 4 4.142 -0.227 0.003
## 17 X.6 X.7 X.8 X.9 X.10 X.11 X.12
## 18
## 19
## 20 d e Diameter Corrected for bias Bias correction (CF) r R2
## 21 364946 d.b.h. no 0 0 0.98
## 22 d.b.h. no 0 0 0.99
## 23 2 d.b.h. no 0 0 0.96
```

```

## 6      2      d.b.h.      no      0      0 0.97
##      X.13      X.14      X.15      X.16      X.17      X.18
## 1
## 2 MinDiameter MaxDiameter Sample size Stump height Top d.o.b. Units diameter
## 3      1.00      72.00      83      0      cm
## 4      0.80      34.10      108      6      cm
## 5      2.50      25.00      131      12      mm
## 6      12.50      55.00      131      12      0      mm
##      X.19      X.20      X.21      X.22      X.23
## 1
## 2 Units biomass Component Component sum Ratio equation Segmented equation
## 3      kg      na      a
## 4      kg      na
## 5      kg      na
## 6      kg      na
##      X.24      X.25
## 1
## 2 Equation number Source
## 3      1      140
## 4      1      51
## 5      1      107
## 6      1      107
##
## 1
## 2
## 3 assume 0-inch stump height, but data from other studies so stump heights are probably mixed; 43 pi
## 4      no stump height; tree data
## 5
## 6      12-inch stump including enti

```

```

# other functions
###AllomAve

###AllomUnitCoef # converts length units FROM cm TO specified units converts mass units TO kg FROM spec

###allom.BayesFit #Module to fit a common power-law allometric model to a mixture of raw data and allom

###allom.predict # Function for making tree-level Monte Carlo predictions from allometric equations est

###load.allom() # load .allom files

# This is proabably what we want
###query.allom.data# Module to grab allometric information from the raw data table Will grab both origi
#Tallied equation format based on Jenkins et al 2004 USFS General Technical Report

###read.allom.data # Extracts PFT- and component-specific data and allometeric equations from the speci

# note that query.allom.data is not an exported fuction from the stand alone module....

```

```
#query.allom.data("PIPO")
```