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#Daniel Tompsett January 2018
#Install the required software
library(devtools)
install_github("moreno-betancur/mice",
ref="f0e838f5a7fdcfcb9bfa27207e903d13d79cefaa")
library(foreign)
library(mice)
#To use NARFCS with the "mice" function refer to instructions in
#https://github.com/moreno-betancur/NARFCS.
#
#PREAMBLE:
#Before each algorithm, create the arguments for use in the function
"mice"
#Except for the argument parmSens, this will change as part of the
algorithm,
#   data=The data to impute
#   m=Number of Multiply imputed datasets
#   maxit=Number of Cycles
#   method=Imputation Methods
#   predictorMatrix=Predictor Matrix, involving all estimable terms
#   predictorSens=Predictor Matrix for sensitivity parameters
#   seed=Random seed
#   parmSens= These contain CSPs, and are left UNSPECIFIED here
#
#   The arguments of the calibration functions are provided using a
#   structure following that of the 'parmSens' argument, which are used
#   in the procedure to fix a range of possible values for the entries
#   of 'parmSens', that is the CSPs. Be sure to understand the
structure
#   of this argument before attempting calibration
#
#   The algorithms here only support MSP models with 1 sensitivity
#   parameter per model. The code would need to
#   be modified for models with multiple MSPs at once.
#####
###
#####Algorithm 1: Elicited ranges of
MSPs#####
#####
###
#Input:
#   MSPmodels=List of MSP models in form list(...,...)
#               Written in the form X~Y (no quotations)
#               THE FIRST ENTRY AFTER ~ MUST BE THAT FOR WHICH THE
#               THE COEFFICIENT IS THE MSP.
#
#   MSPfamily=Vector of MSPmodels regression families in form c(...,...)
#               arguments passed to family argument of glm.mids
#               "gaussian" for normal linear regression
#               "biomial" for logistic regression
#
#   Substantive=Your substantive model of interest
#               Written in the form X~Y (no quotations)
#               THE FIRST VARIABLE ENTRY AFTER ~ MUST BE THAT FOR WHICH

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#           THE COEFFICIENT IS THE MAIN TARGET PARAMETER.
#
#   Subfamily= Substantive model's regression family, written as c("")
#
#   Intercept=If effect of interest is the intercept, set to TRUE
#             Set as FALSE by default
#
#   l=Vector of lower bounds of test ranges for the CSPs
#   u=Vector of upper bounds of test ranges for the CSPs
#   step= Vector of intervals in which test points are taken
#
#   parms=List with the same structure as one would use if one were
#         specifying the 'parmSens' argument, with ALL entries set as
NA
#
#   NOTE:
#   The arguments parms, l, u and step fill in the entries of
#'parmSens'
#   with values over fixed ranges of the CSPs: parms gives the
structure
#   of 'parmSens', and l, u and step fill in the entries of 'parmSens'
#   with values between l and u in intervals of length step. Entries
#   in 'parmSens' corresponding to complete values are set to "".
#
#   Hence l, u and step must have an entry for each
#   element of parmSens in the order in which they appear. Set
#   l,u and step for entries corresponding to complete variables to NA.

calibrationr.fn<-
function(MSPmodels,MSPfamily,Substantive,Subfamily,Intercept=FALSE,l,u,step,parms){
MSP<-as.list(NULL)
substantive<-as.list(NULL)
range<-as.list(NULL)
for(k in 1:length(l)){
if (is.na(l[k])==TRUE){
range[[k]]<-list("")
}
else{
range[[k]]<-seq(l[k],u[k],step[k])
}
}
points<-expand.grid(range)
for(j in 1:nrow(points)){
parmsnew<-relist(as.matrix(points[j,]),parms)
parmSens<-parmsnew
set.seed(seed)
narfcs<-mice(data=data,m=m,maxit=maxit,method=method,
predictorMatrix=predictorMatrix,predictorSens=predictorSens,
parmSens=parmSens,print=F)
MSPlist<-as.vector(NULL)
for(n in 1:length(MSPmodels)){
MSPlm<-glm.mids(MSPmodels[[n]],narfcs,family=get(MSPfamily[n]))

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MSPlist[n]<-pool (MSPlm) $qbar[2]
}
MSP[[j]]<-MSPlist
lms<-glm.mids (Substantive,narfcs, family=get (Subfamily))

if (Intercept==TRUE){
substantive[[j]]<-summary(pool (lms)) [1,c (1,5:7)]}
else{substantive[[j]]<-summary(pool (lms)) [2,c (1,5:7)]
}
}

Msp=do.call (rbind,MSP)
colnames (Msp)<-rep ("c",ncol (Msp))
for(i in 1:ncol (Msp)){
colnames (Msp) [i]<-paste ("MSP",i)
}

Csps<-points
colnames (Csps)<-rep ("c",ncol (Csps))
for(i in 1:ncol (Csps)){
colnames (Csps) [i]<-paste ("CSP",i)
}

analysis=do.call (rbind,substantive)
results=cbind (Msp,analysis)

return (list (Csps=Csps,Msp=Msp,
analysis=analysis,
results=results)
)
}
#OUTPUT:
#      CSPs=Full list of tested sets of CSPs
#      Given in the same order as they appear in parmSens
#      Msp=Full list of estimated MSPs
#      Given in the same order they appear in MSPmodels
#      analysis=List of substantive effects of interest
#      With associated p value and 95% confidence interval
#      Results=The Msp and substantive analysis appended together
#####
###
#####Algorithm 2: Elicited Values All Variables at
Once#####
#####
###
#Input:
#      MSPmodels, MSPfamily, l, u, step and parms, All same as in
algorithm 1
#      Two new arguments
#      elicited=Vector of elicited values of the MSPs
#      tol=Vector of tolerances for each MSP for the algorithm
#      elicited and tol are written as c(... )in same order as MSPmodels
calibrationAAO.fn<-
function (MSPmodels,MSPfamily,elicited,l,u,step,tol,parms){

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MSP<-as.list(NULL)
range<-as.list(NULL)
for(k in 1:length(l)){
  if (is.na(l[k])==TRUE){
    range[[k]]<-list("")
  }
  else{
    range[[k]]<-seq(l[k],u[k],step[k])
  }
}
points<-expand.grid(range)
for(j in 1:nrow(points)){
  parmsnew<-relist(as.matrix(points[j,]),parms)
  parmSens<-parmsnew
  set.seed(seed)
  narfcs<-mice(data=data,m=m,maxit=maxit,method=method,
  predictorMatrix=predictorMatrix,predictorSens=predictorSens,
  parmSens=parmSens,print=F)
  MSPlist<-as.vector(NULL)
  for(n in 1:length(MSPmodels)){
    MSPlm<-glm.mids(MSPmodels[[n]],narfcs,family=MSPfamily[n])
    MSPlist[n]<-pool(MSPlm)$qbar[2]
  }
  MSP[[j]]<-MSPlist
}
dif<-as.vector(NULL)
for(p in 1:nrow(points)){
  dif[[p]]<-tol-abs(MSP[[p]]-elicited)
}
Msps=do.call(rbind,MSP)
colnames(Msps)<-rep("c",ncol(Msps))
for(i in 1:ncol(Msps)){
  colnames(Msps)[i]<-paste("MSP",i)
}

colnames(points)<-rep("c",ncol(points))
for(i in 1:ncol(points)){
  colnames(points)[i]<-paste("CSP",i)
}
Csps<-points

con<-lapply(dif,function(x)all(x>0))

MSPcal<-MSP[c(which(con==TRUE)) ]
calibratedMSP=do.call(rbind,MSPcal)
colnames(calibratedMSP)<-rep("c",ncol(calibratedMSP))
for(i in 1:ncol(calibratedMSP)){
  colnames(calibratedMSP)[i]<-paste("MSP",i)
}

CSPcal<-points[c(which(con==TRUE)), ]
colnames(CSPcal)<-rep("c",ncol(CSPcal))
for(i in 1:ncol(CSPcal)){

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colnames(CSPcal)[i]<-paste("CSP",i)
}

return(list(calibratedMSP=calibratedMSP,calibratedCSP=CSPcal,
Csps=Csps,Msp=Msps))
}
#OUTPUT:
# calibratedCSP= Sets of calibrated CSPs which satisfy the tolerances
# calibratedMsp= Estimated MSPs of caliibratedCSP
# Csps=Full list of tested sets of CSPs
# Msp=Full list of estimated MSPs
#####
###
#####Algorithm 3: One variable at a
time#####
#####
###
#INPUT:
# MSPmodels, MSPfamily, l, u, step, elicited and tol
# have same meaning as in algorithm 1. Since there is only 1 variable
# no need to write them in as list(...) or c(...).
#
# parms=List with the same structure as one would use if one were
# specifying the 'parmSens' argument.
#
# SPECIFIED DIFFERENTLY TO ALGORITHMS 1 and 2
# write as one would write 'parmSens', with the CSP you wish
# to vary set to NA. Set other CSPs to their chosen fixed
# values, and complete variables to "".

calibrationOAAT.fn<-
function(MSPmodel,MSPfamily,elicited,l,u,step,tol,parms){
MSPPr<-as.list(NULL)
CSPPr<-as.list(NULL)
range<-seq(l,u,step)
i<-1
test1<-100
test2<-0
while(abs(test1-test2)>tol){
dif<-as.vector(NULL)
for(j in 1:length(range)){
parmSens<-parms
for(l in 1:length(parmSens)){
parmSens[[l]][is.na(parmSens[[l]])]<-range[j]
}
set.seed(seed)
narfcs<-mice(data=data,m=m,maxit=maxit,method=method,
predictorMatrix=predictorMatrix,predictorSens=predictorSens,
parmSens=parmSens,print=F)
MSPlm<-glm.mids(MSPmodels,narfcs,family=MSPfamily)

dif[j]<-pool(MSPlm)$qbar[2]
}
}

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diff<-dif-elicited
indexp<-which(diff==min(diff[which(diff>=0)]))
indexn<-which(diff==max(diff[which(diff<=0)]))
MSPr[[i]]<-c(dif[indexn],dif[indexp])
CSPr[[i]]<-c(range[indexn],range[indexp])
test1<-dif[indexn]
test2<-dif[indexp]
lower<-min(range[indexn],range[indexp])
upper<-max(range[indexn],range[indexp])
range<-seq(lower,upper,(step/10^i))
i<-i+1
}
return(list=c(Msp=MSPr[i-1],Csp=CSPr[i-1]))
}
#Output:
#      Csp=The two closest calibrated values of the CSP within tolerance
#      Msp=The estimated MSP values for Csp

#####
#####
#####Example#####
#####
#####
#GENERATE EXAMPLE DATASET:
#      2 continuous variables with missing data, Y1 and Y2
#      Corresponding missingness indicators M1 and M2
#      Complete continuous variable X
set.seed(45678)
c1<-c(20,10,8)
c2<-c(10,15,11)
c3<-c(8,11,14)
sigma<-rbind(c1,c2,c3)
M1<-rbinom(1000,1,0.5)
M2<-rbinom(1000,1,0.5)
mu1=50-(1*M1+2*M2)
mu2=50-(4*M1+5*M2)
mu3=rep(50,1000)
library(MASS)
data<-matrix(,nrow=1000,ncol=3)
for(j in 1:1000){
data[j,]<-mvrnorm(1,mu=c(mu1[j],mu2[j],mu3[j]),Sigma=sigma)
}
Y1<-data[,1]
Y2<-data[,2]
X<-data[,3]
data<-data.frame(Y1,Y2,X,M1,M2)
data$Y1[data$M1=="1"]<-NA
data$Y2[data$M2=="1"]<-NA
attach(data)

#ARGUMENTS FOR MICE:
m=10
maxit=10

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method=c("normSens","normSens","","","")
#NARFCS models are
#   Y1~Y2+X+M2+M1
#   Y2~Y1+X+M1+M2
predictorMatrix<- diag(0, ncol(data))
rownames(predictorMatrix) <- names(data)
colnames(predictorMatrix) <- names(data)
predictorMatrix["Y1",c("Y2","X","M2")] <- 1
predictorMatrix["Y2",c("Y1","X","M1")] <- 1
predictorSens<- diag(0, ncol(data))
rownames(predictorSens) <- names(data)
colnames(predictorSens) <- names(data)
seed=123456786

#####ALGORITHM 1:
MSPmodels<-list(Y1~M1,Y2~M2)
MSPfamily=c("gaussian","gaussian")
Substantive<-Y1~X
Subfamily=c("gaussian")
#We will fix ranges for both the two CSPs at c(-10,0) taking points 1
unit apart
l<-c(-10,-10,NA,NA,NA)
u<-c(0,0,NA,NA,NA)
step<-c(1,1,NA,NA,NA)

#Entries for X, M1 and M2 set to NA as they are complete

parms<-list(list(NA),list(NA),list(NA),list(NA),list(NA))

#RUN ALGORITHM:
ans<-
calibrationr.fn(MSPmodels,MSPfamily,Substantive,Subfamily,Intercept=FALSE
,l,u,step,parms)

#####ALGORITHM 2:
#Arguments the same as with algorithm 2, except substantive model
arguments
elicited<-c(-10,-10)
tol<-c(0.6,0.6)

#RUN ALGORITHM:
ans<-calibrationAAO.fn(MSPmodels,MSPfamily,elicited,l,u,step,tol,parms)

#####ALGORITHM 3:
#Calibrate CSP for Y1 with CSP for Y2 at -10
MSPmodels<-Y1~M1
MSPfamily<-"gaussian" #QUOTATIONS VITAL
elicited<--10 #Elicited MSP is -10
l<--20
u<-0
step<-5
tol<-0.1
parms=list(list(NA),list(-10),list(""),list(""),list(""))

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# CSP for Y1 is set to NA, CSP for Y2 to -10.
# X, M1 and M1 are complete, entry is therefore list("")

#RUN ALGORITHM:
ans<-calibrationOAT.fcn(MSPmodels,MSPfamily,elicited,l,u,step,tol,parms)

#####PERFORMING SENSITIVITY ANALYSIS For Algorithm
1#####
#We give a short example of doing the analysis from the results of
algorithm 1
#Assume the range of interest is (-10,-5) for the Msp for Y1
#and(-8,-4) for the MSP of Y2.

####Graphical Methods

#Plot the Msps against the analysis estimates
#INSTALL PACKAGE 'scatterplot3d'
library(scatterplot3d)

#3d plot
scatterplot3d(ans$results[, "MSP 1"],ans$results[, "MSP
2"],ans$results[, "est"],type="l",
angle=30)

#The flat plane clearly indicates the 'one at a time' assumption holds
#Graph individually and mark areas of interest for MSP of Y1
#Graph for MSP of Y1

plot(ans$results[, "MSP 1"],ans$results[, "est"],type="l")
abline(v=-10,col="red")
abline(v=-5,col="red")

#As effect of interest is Y1~X, Msp for Y2 is less important

#NOTE:
#To apply a tipping point analysis, plot against ans$results[, "Pr(>|t|)"]

####By Tabulation
#Obtain the set of results such that the MSPs lie within their elicited
#ranges

ans$results[ans$results[, "MSP 1"]>=-10 &
ans$results[, "MSP 1"]<=-5 &
ans$results[, "MSP 2"]>=-8 &
ans$results[, "MSP 2"]<=-4 ,]

# It may be worth looking at values that lie slightly outside the
elicited ranges as well

```