Multivariate Multiple Imputation

Utrecht University Winter School: Missing Data in R



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Outline

Flavors of MI

Fully Conditional Specification

Joint Modeling

Mixed Data Types



Joint Modeling vs. Fully Conditional Specification

When imputing with *Joint Modeling* (JM) approaches, the missing data are replaced by samples from the joint posterior predictive distribution.

• To impute *X*, *Y*, and *Z*, we draw:

$$X, Y, Z \sim P(X, Y, Z|\theta)$$

With *Fully Conditional Specification* (FCS), the missing data are replaced with samples from the conditional posterior predictive distribution of each incomplete variable.

• To impute *X*, *Y*, and *Z*, we draw:

$$X \sim P(X|Y, Z, \theta_X)$$

$$Y \sim P(Y|X, Z, \theta_Y)$$

$$Z \sim P(Z|Y, X, \theta_Z)$$



Joint Modeling: Strengths

When correctly implemented, JM approaches are guaranteed to produce *Bayesianly proper* imputations.

- A sufficient condition for properness is that the imputations are randomly sampled from the correctly specified joint posterior predictive distribution of the missing data.
 - This is the defining characteristic of JM methods.

When using the correct distribution, imputations produced by JM methods will be the best possible imputations.

- Unbiased parameter estimates
- Well-calibrated sampling variability



Joint Modeling: Weaknesses

JM approaches don't scale well.

 The computational burden increases with the number of incomplete variables.

JM approaches are only applicable when the joint distribution of all incomplete variables follows a known form.

 Mixes of continuous and categorical variables are difficult to accommodate.

Software Implementations

In R, MI via JM is available from several packages.

- Amelia (Honaker, King, & Blackwell, 2011)
 - Bootstrapped EM algorithm
- norm (Schafer, 2013)
 - · Classic data augmentation.
- mice (Van Buuren & Groothuis-Oudshoorn, 2011)
 - Data augmentation for block updating.

JM imputation is also available in SAS, Stata, SPSS, and Mplus.



Fully Conditional Specification: Strengths

FCS scales much better than JM.

• FCS only samples from a series of univariate distributions, not large joint distributions.

FCS approaches can create imputations for variables that don't have a sensible joint distribution.

• FCS can easily treat mixes of continuous and categorical variables.



Fully Conditional Specification: Weaknesses

FCS will usually be slower than JM.

• Each variable gets its own fully parameterized distribution, even if that granularity is unnecessary.

When the incomplete variables don't have a known joint distribution, FCS doesn't have theoretical support.

- There is, however, a large degree of empirical support for the tenability of the FCS approach.
- In practice, we usually choose FCS since real data rarely arise from a known joint distribution.

Software Implementations

The **mice** package is the most popular R implementation of FCS. (Van Buuren & Groothuis-Oudshoorn, 2011).

- Mature implementation
- · Well integrated into the larger R ecosystem
- Very active development

The mi package (Su, Yajima, Gelman, & Hill, 2011) offers another option.

- More focus on diagnostics
- Object oriented flavor
- Not very actively developed

FCS imputation is also available in SAS, SPSS, Stata, and Mplus.

Aside: Gibbs Sampling

Up to this point, most of the models we've explored could be approximated by sampling directly from their posterior distributions.

• This won't be true with arbitrary, multivariate missing data.

To make inference regarding a multivariate distribution with multiple, interrelated, unknown parameters, we can use *Gibbs sampling*.

• Sample from the conditional distribution of each parameter, conditioning on the current best guesses of all other parameters.

Aside: Gibbs Sampling

Suppose the following:

- 1. I want to make some inference about the tri-variate mean of $X,Y,Z=\mu_X,\mu_Y,\mu_Z\sim P(\mu|\theta)$
- 2. $P(\mu|\theta)$ is super hairy and difficult to sample
- 3. I can easily sample from the conditional distributions: $P(\mu_X|\hat{\mu}_Y,\hat{\mu}_Z,\theta)$, $P(\mu_Y|\hat{\mu}_X,\hat{\mu}_Z,\theta)$, and $P(\mu_Z|\hat{\mu}_X,\hat{\mu}_Y,\theta)$.

Then, I can approximate the full joint distribution $P(\mu|\theta)$ by sequentially sampling from $P(\mu_X|\hat{\mu}_Y,\hat{\mu}_Z,\theta)$, $P(\mu_Y|\hat{\mu}_X,\hat{\mu}_Z,\theta)$, and $P(\mu_Z|\hat{\mu}_X,\hat{\mu}_Y,\theta)$.

Aside: Gibbs Sampling

Starting with initial guesses of μ_Y , $\hat{\mu}_Y^{(0)}$, and μ_Z , $\hat{\mu}_Z^{(0)}$, and assuming θ is known, Gibbs sampling proceeds as follows:

$$\begin{split} \hat{\mu}_{X}^{(1)} &\sim P(\mu_{X}|\hat{\mu}_{Y}^{(0)},\hat{\mu}_{Z}^{(0)},\theta) \\ \hat{\mu}_{Y}^{(1)} &\sim P(\mu_{Y}|\hat{\mu}_{X}^{(1)},\hat{\mu}_{Z}^{(0)},\theta) \\ \hat{\mu}_{Z}^{(1)} &\sim P(\mu_{Z}|\hat{\mu}_{Y}^{(1)},\hat{\mu}_{X}^{(1)},\theta) \\ \hat{\mu}_{X}^{(2)} &\sim P(\mu_{X}|\hat{\mu}_{Y}^{(1)},\hat{\mu}_{X}^{(1)},\theta) \\ \hat{\mu}_{Y}^{(2)} &\sim P(\mu_{Y}|\hat{\mu}_{X}^{(2)},\hat{\mu}_{Z}^{(1)},\theta) \\ \hat{\mu}_{Y}^{(2)} &\sim P(\mu_{Z}|\hat{\mu}_{Y}^{(2)},\hat{\mu}_{X}^{(2)},\theta) \\ \vdots &\vdots \end{split}$$



Why do we care?

Multivariate MI employs the same logic as Gibbs sampling.

- The imputations are created by conditioning on the current estimates of the imputation model parameters.
- The imputation model parameters are updated by conditioning on the most recent imputations.
- With FCS, each variable is imputed by conditioning on the most recent imputations of all other variables.

Fully Conditional Specification



Procedure: Fully Conditional Specification

- 1. Fill the missing data with reasonable guesses.
- 2. For each incomplete variable, do a single iteration of univariate Bayesian MI (e.g., as seen in the last set of slides).
 - After each variable on the data set is so treated, we've completed one iteration.
- 3. Repeat Step 2 many times.
- After the imputation model parameters stabilize, save M imputed data sets.

Example: Data Generation

First we'll simulate some synthetic data.

```
## Simulate some data:
simData <-
   simCovData(nObs = 1000, sigma = 0.25, nVars = 4)
head(simData, 10)
           x1
                      x2
                             x3
                                             ×4
   0.06313632 -1.4057704 -0.01709217 1.47929405
  -1.31592547 1.2970920 -0.83500777 -0.44528158
  -0.30997023 0.9782580 0.02731853 0.35507390
4
   0.06927787 0.1836032 0.68794409 0.08049987
  -0.99354894 -0.3038956 0.80918329 -1.72143555
6
   0.36828016 -0.9423245 1.05155348 -0.11078496
   1.33333163 2.3089780 1.47203000 0.85877495
 -0.02759718 0.1714383 0.18927909 0.28627771
9
   2.37929433 2.5080935
                          2.18344726 1.56980951
10
   0.31841502
              0.7886025
                          0.73658136 0.39445970
```

Example: Data Generation

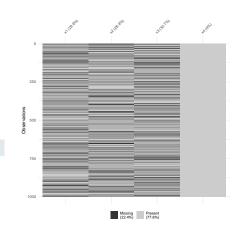
Next, we impose some missing values on the simulated data.

Check the results.

```
head(missData, 10)
           x1
                  x2
                        x3
                                          x4
   0.06313632
                    NA -0.01709217 1.47929405
           NA 1.2970920
                                NA -0.44528158
3
  -0.30997023
                    NA 0.02731853 0.35507390
           NΑ
                    NA 0.68794409 0.08049987
5
           NA
                    NA 0.80918329 -1.72143555
6
   0.36828016
                    NΑ
                                NA -0.11078496
   1.33333163 2.3089780
                                NA
                                    0.85877495
8
           NA
                    NA 0.18927909 0.28627771
   2.37929433 2.5080935
                                NA 1.56980951
   0.31841502 0.7886025 0.73658136 0.39445970
```

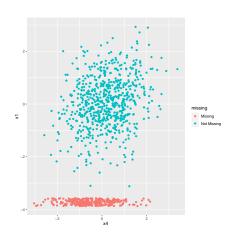
Use the naniar::vis_missing() function to visualize the pattern of missing values.

vis_miss(missData)



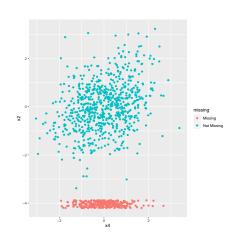
Use naniar::geom_miss_point() to visualize the response pattern.

```
ggplot(missData, aes(x4, x1)) +
   geom_miss_point()
```



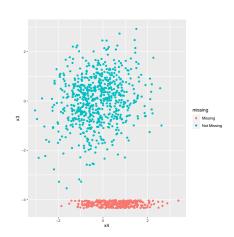
Use naniar::geom_miss_point() to visualize the response pattern.

```
ggplot(missData, aes(x4, x2)) +
   geom_miss_point()
```



Use naniar::geom_miss_point() to visualize the response pattern.

```
ggplot(missData, aes(x4, x3)) +
   geom_miss_point()
```



```
## Define iteration numbers:
nImps <- 100
nBurn <- 500
nSams <- nBurn + nImps
## Summarize missingness:
rMat <- !is.na(missData)
nObs <- colSums(rMat)
nMis <- colSums(!rMat)
## Fill the missingness with initial (bad) quesses:
mean0 <- colMeans(missData, na.rm = TRUE)
sigma0 <- cov(missData, use = "pairwise")</pre>
draws0 <- rmvnorm(nrow(missData), mean0, sigma0)</pre>
impData
        <- missData
impData[!rMat] <- draws0[!rMat]</pre>
```

Define an elementary imputation function:

```
eif <- function(data, rVec, v) {
    ## Get the expected model parameters:
    fit <- lm(paste(v, "~ ."), data = data[rVec, ])
    beta <- coef(fit)
    s2 <- summarv(fit)$sigma^2</pre>
    ## Partition data:
    vars <- setdiff(colnames(data), v)</pre>
    data <- as.matrix(data)
    xObs <- cbind(1, data[rVec, vars])
    xMis <- cbind(1, data[!rVec, vars])
    ## Sample sigma:
    sigmaSam <- rinvchisq(1, df = fit$df, scale = s2)
    ## Sample beta:
    betaVar <- sigmaSam * solve(crossprod(xObs))</pre>
    betaSam <- rmvnorm(1, mean = beta, sigma = betaVar)
    ## Return a randomly sampled imputation:
    xMis %*% t(betaSam) + rnorm(nrow(xMis), 0, sgrt(sigmaSam))
```

Apply the elementary imputation function to each incomplete variable:

Analyze the multiply imputed datasets:

Compare approaches:

	Est	SE	t	р	FMI
(Intercept)	-0.009	0.038	-0.238	1.188	0.406
x2	0.225	0.041	5.519	< 0.001	0.483
x3	0.161	0.049	3.283	0.001	0.637

Manual Version

	Est	SE	t	р	FMI
(Intercept)	-0.012	0.036	-0.342	1.267	0.351
x2	0.225	0.04	5.647	< 0.001	0.463
х3	0.167	0.045	3.681	< 0.001	0.58

MICE Version

Joint Modeling



Aside: Definition of Regression Parameters

So far, we've been using the least-squares estimates of α , β , and σ^2 to parameterize our posterior distributions.

• We can also define the parameters in terms of sufficient statistics.

Given μ and Σ , we can define all of our regression moments as:

$$\beta = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$$

$$= \mathsf{Cov}(\mathbf{X})^{-1} \mathsf{Cov}(\mathbf{X}, \mathbf{Y})$$

$$\alpha = \mu_{\mathbf{Y}} - \beta^T \mu_{\mathbf{X}}$$

$$\Sigma_{\varepsilon} = \Sigma_{\mathbf{Y}} - \beta^T \Sigma_{\mathbf{X}} \beta$$

These definitions are crucial for JM approaches.

 Within the subset of data define by a given response pattern, the outcome variables will be entirely missing.

Multivariate Bayesian Regression

Previously, we saw examples of univariate Bayesian regression which used the following model:

$$\begin{split} \beta &\sim \text{MVN}\left(\hat{\beta}_{ls}, \ \sigma^2(\mathbf{X}^T\mathbf{X})^{-1}\right) \\ \sigma^2 &\sim \text{Inv-}\chi^2\left(N-P, MSE\right) \end{split}$$

We can directly extend the above to the multivariate case:

$$\begin{split} & \boldsymbol{\Sigma^{(i)}} \sim \text{Inv-W}\left(N-\mathbf{1}, (N-\mathbf{1})\boldsymbol{\Sigma^{(i-1)}}\right) \\ & \boldsymbol{\mu^{(i)}} \sim \text{MVN}\left(\boldsymbol{\mu^{(i-1)}}, N^{-1}\boldsymbol{\Sigma^{(i)}}\right) \end{split}$$

We get α , β , and Σ_{ε} via the calculations on the preceding slide

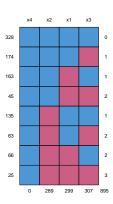
Procedure: Joint Modeling

In JM imputation, we estimate the imputation model via the Tanner and Wong (1987) *data augmentation* algorithm.

- 1. Partition the incomplete data by response pattern.
 - Produce *S* subsets wherein each row shares the same response pattern.
- **2.** Provide initial guesses for μ and Σ .
- 3. Within each subset, use the current guesses of μ and Σ to generate imputations via multivariate Bayesian regression.
- 4. Use the filled-in data matrix to updated the sufficient statistics.
- 5. Repeat Steps 3 and 4 many times.
- After the imputation model parameters have stabilized, save M imputed data sets produced in Step 3.

Use mice::md.pattern() to visualize the response patterns.

pats <- md.pattern(missData)</pre>



```
iStep <- function(data, pats, ind, pars) {
    ## Loop over non-trivial response patterns:
    for(i in 1:nrow(pats)) {
        ## Define the current response pattern:
       p1 <- pats[i, ]
        ## Replace missing data with imputations:
        data[ind == i, !p1] <- getImps(data = data[ind == i, ],</pre>
                                              = pars$mu,
                                        mıı
                                        sigma = pars$sigma,
                                              = p1)
                                        p1
    ## Return the imputed data:
    data
```

```
getImps <- function(data, mu, sigma, p1) {</pre>
    ## Partition the parameter matrices:
   mY <- matrix(mu[!p1])
   mX <- matrix(mu[p1])</pre>
    sY <- sigma[!p1, !p1]
    sX <- sigma[p1, p1]
    cXY <- sigma[p1, !p1]
    ## Compute the imputation model parameters:
    beta <- solve(sX) %*% cXY
    alpha <- mY - t(beta) %*% mX
    sE <- sY - t(beta) %*% sX %*% beta
    ## Pull out predictors:
    X <- as.matrix(data[ , p1])</pre>
    ## Generate and return the imputations:
    n \leftarrow nrow(X)
   matrix(1, n) %*% t(alpha) + X %*% beta + rmvnorm(n, sigma = sE)
```

```
pStep <- function(data) {
    ## Update the complete-data sufficient statistics:
    n <- nrow(data)
    m <- colMeans(data)
    s <- (n - 1) * cov(data)

## Sample the covariance matrix and mean vector:
    sigma <- riwish((n - 1), s)
    mu <- rmvnorm(1, m, (sigma / n))

## Return the updated parameters:
    list(mu = mu, sigma = sigma)
}</pre>
```

Now that we've defined the necessary functions, do the imputation:

```
## Some preliminaries:
impData <- missData
nIter <- 50
nImps <- 100
## Summarize response patterns:
rMat <- !is.na(impData)</pre>
pats <- uniquecombs(rMat, ordered = TRUE)</pre>
ind <- attr(pats, "index")</pre>
## Exclude the fully observed pattern:
if(tail(pats, 1) %>% all())
    pats <- pats[-nrow(pats), ]</pre>
## Get starting values for the parameters:
theta0 <- list(mu = colMeans(impData, na.rm = TRUE),
                sigma = cov(impData, use = "pairwise")
```

```
## Iterate over I- and P-Steps to generate imputations:
impList1 <- list()</pre>
for(m in 1:nImps) {
    ## Initialize the parameter vector:
    theta1 <- theta0
    for(rp in 1:nIter) {
        ## Do one I-Step:
        impData <- iStep(data = impData,</pre>
                           pats = pats,
                           ind = ind,
                           pars = theta1)
        ## Do one P-Step:
        theta1 <- pStep(impData)</pre>
    ## Save the final imputed dataset:
    impList1[[m]] <- impData</pre>
```

Do the same type of imputation with **norm**:

```
missData <- as.matrix(missData)</pre>
## Pre-process the data and get starting values via EM:
meta <- prelim.norm(missData)</pre>
theta0 <- em.norm(meta, showits = FALSE)
rngseed(235711)
impList2 <- list()</pre>
for(m in 1 : nImps) {
    ## Estimate the imputation model via data augmentation:
    theta1 <- da.norm(s = meta, start = theta0, steps = nIter)
    ## Impute missing values via a final I-Step of DA:
    impList2[[m]] <- imp.norm(s = meta, theta = theta1, x = missData)</pre>
```

Analyze the multiply imputed data:

Compare approaches:

	Est	SE	t	р	FMI
(Intercept)	-0.01	0.036	-0.276	1.217	0.355
x2	0.229	0.042	5.464	< 0.001	0.506
x3	0.17	0.044	3.878	< 0.001	0.542

Manual Version

	Est	SE	t	p	FMI
(Intercept)	-0.012	0.036	-0.337	1.264	0.359
x2	0.231	0.041	5.652	< 0.001	0.482
x3	0.162	0.046	3.504	0.001	0.591

NORM Version

Mixed Data Types



FCS for Mixed Data Types

FCS imputation can easily accommodate incomplete data that contain both continuous and categorical/non-normal variables.

- Replace the normal-theory elementary imputation model described above with an appropriate model for the distribution of each incomplete variable
 - Logistic regression (various flavors)
 - Donor-based methods
 - Tree-based methods

The FCS framework can essentially accommodate any data for which you can define an appropriate supervised model.

• Many useful methods are already implemented in the mice package.

JM for Mixed Data Types

When applying JM to incomplete data with mixed variable types, we have to general options.

- Impute under the multivariate normal model and round, coarsen, or truncate the continuous imputations to "match" the original data.
 - This was the old-school recommendation from the days before FCS (e.g., Allison, 2002; Schafer, 1997).
 - The **Amelia** package implements this approach.
 - This approach tends to perform poorly in methodological evaluations (e.g., Lang & Wu, 2017; Wu, Jia, & Enders, 2015).
- 2. Impute under an appropriate joint model for the data.
 - This approach is only available when a suitable joint model exists.
 - The mix package (Schafer, 2017) implements this approach for the general location model (Little & Schluchter, 1985).
 - This approach also doesn't do very well, in practice (Lang & Wu, 2017).

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