Wrapper function parlMICE

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For big datasets or high number of imputations, performing multiple imputation with function mice from package mice (Van Buuren and Groothuis-Oudshoorn 2011) might take a long time. As a solution, wrapper function parlMICE was created to enable the imputation procedure to be run in parallel. This is done by dividing the imputations over multiple cores (or CPUs), thereby speeding up the process.

First, this vignette will show the time gain with parlMICE for a small and (relatively) large dataset. Second, parlMICE's arguments will be disussed. For more information about running functions in parallel, we gladly refer to the parallel manual and the document *How-to-go parallel in R*, written by Max Gordon. We thank Max Gordon for his previous work on this subject.

Time gain with small datasets

In Figure 1, the processing time is compared between mice and parlMICE. The simulated dataset has 1000 cases and 4 variables with covariance structure

$$\Sigma = \begin{array}{cccc} V_1 & V_2 & V_3 & V_4 \\ V_1 & 1 & 0.5 & 0.5 & 0.5 \\ V_2 & 0.5 & 1 & 0.5 & 0.5 \\ V_3 & 0.5 & 0.5 & 1 & 0.5 \\ V_4 & 0.5 & 0.5 & 0.5 & 1 \end{array}$$

A MCAR missingness mechanism is imposed on the data. In total, 80 percent of the cases (rows) has missingness on one variable. All variables have missing values.

Multiple imputation is performed with the default arguments of mice. The procedure with parlMICE is repeated for 1, 2 and 3 cores. The results are shown in Figure 1.

From Figure 1 it becomes apparent that for a small number of imputations, the conventional mice function works faster than the wrapper function. This is the case for $m \leq 35$. From that point on, wrapper function parlMICE returns the imputations faster.

The right figure of Figure 1 shows the processing time for very high number of imputations. After the intersection, the lines follow the same trend: 3 cores perform the work faster than 2 cores, which work faster than 1 core. Naturally, using parlMICE with 1 core is similar to function mice, as the figure shows with the green and orange lines.

Time gain with large datasets

Another test was run with a (relatively) large dataset of 10000 cases and 8 variables. The covariance matrix was simply extended to the situation of 8 variables. In Figure 2, the results are shown.

Figure 2 shows that for all imputations, function parlMICE works faster than mice. Even for very small number of imputations, running mice in parallel with parlMICE saves a significant amount of time. This time gain can easily reach 70 percent for 10 imputations and more.

There is not a huge difference between using 2 and 3 cores with wrapper function parlMICE. For all number of imputations, the procedure runs faster with 3 cores, even though the imputations have to be divided over the cores. It might therefore be desirable to use always as many cores as possible (but always keep 1!).

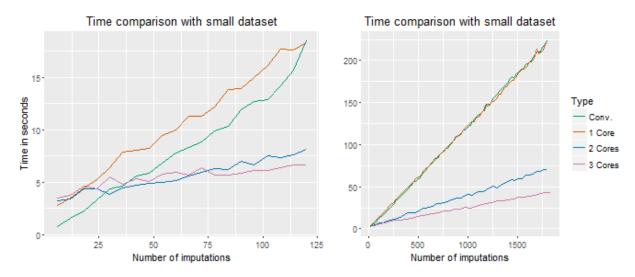


Figure 1: Processing time for small datasets. Multiple imputation are performed with mice (conventional) and wrapper function parlMICE (1, 2 and 3 cores respectively). The dataset has 1000 cases and 4 variables with a correlation of 0.5. 80 percent of the cases have one missing value based on MCAR missingness.

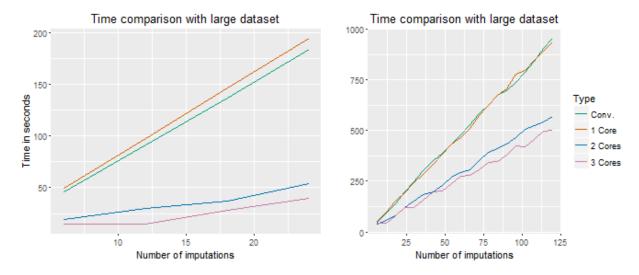


Figure 2: Processing time for large datasets. Multiple imputations are performed with mice (conventional) and wrapper function parlMICE (1, 2 and 3 cores respectively). The dataset has 10000 cases and 8 variables with a correlation of 0.5. 80 percent of the cases have one missing value based on MCAR missingness.

Default settings

We will now dig into the arguments of function parlMICE. Easy imputation of an incomplete dataset (say, nhanes) can be performed with parlMICE in the following way.

```
imp <- parlMICE(nhanes)
class(imp)</pre>
```

```
## [1] "mids"
```

The function returns a mids object as created by mice. In fact, parlMICE makes use of function ibind to combine the mids objects returned by the different cores. Therefore, the call of the mids object has slightly changed.

imp\$call

```
## [[1]]
## mice(data = data, m = n.imp.core, printFlag = FALSE)
##
## [[2]]
## ibind(x = imp, y = imps[[i]])
##
## [[3]]
## ibind(x = imp, y = imps[[i]])
```

All other parts of the mids object are standard.

Using mice arguments

Function parlMICE is able to deal with the conventional mice arguments. In order to change the imputation method from its default (predictive mean matching) to, for example, Bayesian linear regression, the method argument can be adjusted. For other possibilities with mice, we refer to the manual.

```
imp <- parlMICE(nhanes, method = "norm")
imp$method</pre>
```

```
## age bmi hyp chl
## "norm" "norm" "norm"
```

In mice, the number of imputations is specified with argument m. In parlMICE, it is possible to use this argument as well, but we advise to rather use the arguments n.core and n.imp.core. The next section discusses these arguments.

Arguments n.core, n.imp.core

With n.core, the number of cores (or CPUs) is given. n.imp.core defines the number of imputations per core. Subsequently, the total number of imputations equals n.core * n.imp.core. As a default, n.core is specified as the number of available, logical cores minus 1. The default number of imputations per core has been set to 2.

Subsequently, running the default parlMICE function on a computer with,

detectCores()

```
## [1] 4
```

available, logical cores, results to m = (4 - 1) * 2 = 6 imputations. We can check this by evaluating the m shown in the mids object.

```
imp$m
```

[1] 6

Argument seed

In simulation studies, it is often desired to set a seed to make the results reproducible. In contrast to mice, the seed value for parlMICE cannot be defined outside the function. This has to do with the different cores running at the same time (for more information we gladly refer to the Parallel manual).

A seed value can be used in parlMICE with argument seed:

```
imp <- parlMICE(nhanes, seed = 123)</pre>
```

Systems other than Windows

Function parlMICE calls for function parlapply from the Parallel package. Although other options are available, we have chosen for parlapply because it allows the use of multiple cores on all computers, including a Windows computer. For the cluster, we rely on the default option in the function makeCluster, which is "PSOCK".

On systems other than Windows, cluster type "FORK" might be faster (see the document written by Max Gordon). Therefore, we advise to change the type to "FORK" if your computer does not run on Windows.

```
imp <- parlMICE(nhanes, type = "FORK")</pre>
```

References

Gordon, M. (2015). How-to go parallel in R - basics + tips. Available at http://gforge.se/2015/02/how-to-go-parallel-in-r-basics-tips/

Manual base-package Parallel, available at https://stat.ethz.ch/R-manual/R-devel/library/parallel/doc/parallel.pdf

Manual package MICE, available at https://cran.r-project.org/web/packages/mice/mice.pdf

Van Buuren, S., and C.G.M. Groothuis-Oudshoorn. 2011. "Mice: Multivariate Imputation by Chained Equations in R." *Journal of Statistical Software* 45 (3).