Simulation of compound hierarchical models

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1 Introduction

Hierarchical probability models are widely used for data classified in a tree-like structure and in Bayesian inference. The main characteristic of such models is to have the probability law at some level in the classification structure be conditional on the outcome in previous levels. For example, adopting a bottom to top description of the model, a simple hierarchical model could be written as

$$X_t | \Lambda, \Theta \sim \text{Poisson}(\Lambda)$$
$$\Lambda | \Theta \sim \text{Gamma}(3, \Theta) \qquad (1)$$
$$\Theta \sim \text{Gamma}(2, 2),$$

where X_t represents actual data. The random variables Θ and Λ are generally seen as uncertainty, or risk, parameters in the actuarial literature; in the sequel, we refer to them as mixing parameters.

The example above is merely a multi-level mixture of models, something that is simple to simulate "by hand". The following R expression will yield n variates of the random variable X_t:

```
> rpois(n, rgamma(n, 3, rgamma(n, 2, 2)))
```

However, for categorical data common in actuarial applications there will usually be many categories — or *nodes* — at each level. Simulation is then

complicated by the need to always use the correct parameters for each variate. Furthermore, actuaries often need to simulate both the frequency and the severity of claims for compound models of the form

$$S = C_1 + \cdots + C_N, \tag{2}$$

where $C_1, C_2, \ldots$ are mutually independent and identically distributed random variables each independent of N.

The package provides function `simul` to simulate data from compound models like (2) where both the frequency and the severity components can have a hierarchical structure. The function also supports weights (or volumes) in the model. We give here a brief description of the function and its usage; see Goulet and Pouliot (2008) for details about the supported models and more thorough examples.

2 Description of hierarchical models

We need a method to describe hierarchical models in R that will meet the following criteria:

1. simple and intuitive to go from the mathematical formulation of the model to the R formulation and back;

2. allows for any number of levels and nodes;

3. at any level, allows for any use of parameters higher in the hierarchical structure.

A hierarchical model is completely specified by the number of nodes at each level (I, $J_1, \ldots, J_I$ and $n_{11}, \ldots, n_{IJ}$, above) and by the probability laws at each level. The number of nodes is passed to `simul` by means of a named list where each element is a vector of the number of nodes at a given level. Vectors are recycled when the number of nodes is the same throughout a level. Probability models are expressed in a semi-symbolic fashion using an object of mode "expression". Each element of the object must be named — with names matching those of the number of nodes list — and should be a complete call to an existing random number generation function, but with the number of variates omitted. Hierarchical models are achieved by replacing one or more parameters of a distribution at a given level by any combination of the names of the levels above. If no mixing is to take place at a level, the model for this level can be NULL.

Example 1. Consider the following expanded version of model (1):

$$X_{ijt}|\Lambda_{ij}, \Theta_i \sim \text{Poisson}(\Lambda_{ij}), \qquad t = 1, \ldots, n_{ij}$$
$$\Lambda_{ij}|\Theta_i \sim \text{Gamma}(3, \Theta_i), \qquad j = 1, \ldots, J_i$$
$$\Theta_i \sim \text{Gamma}(2, 2), \qquad i = 1, \ldots, I,$$

with $I = 3$, $J_1 = 4$, $J_2 = 5$, $J_3 = 6$ and $n_{ij} \equiv n = 10$. Then the number of nodes is specified by

```
list(Theta = 3, Lambda = c(4, 5, 6), Data = 10)
```

and the probability model is expressed as

```
expression(Theta = rgamma(2, 2),
           Lambda = rgamma(3, Theta),
           Data = rpois(Lambda))
```

□

Storing the probability model requires an expression object in order to avoid evaluation of the incomplete calls to the random number generation functions. Function `simul` builds and executes the calls to the random generation functions from the top of the hierarchical model to the bottom. At each level, the function 1) infers the number of variates to generate from the number of nodes list, and 2) appropriately recycles the mixing parameters simulated previously.

The actual names in the list and the expression object can be anything; they merely serve to identify the mixing parameters. Furthermore, any random generation function can be used. The only constraint is that the name of the number of variates argument is `n`.

Function `simul` also supports usage of weights in models. These usually modify the frequency parameters to take into account the "size" of an entity. Weights are used in simulation wherever the name `weights` appears in a model.

3   Usage

Function `simul` can simulate data for structures where both the frequency model and the severity model are hierarchical. It has four main arguments: 1) `nodes` for the number of nodes list; 2) `model.freq` for the frequency model; 3) `model.sev` for the severity model; 4) `weights` for the vector of weights in lexicographic order, that is all weights of entity 1, then all weights of entity 2, and so on.

The function returns the variates in a list of class `"portfolio"` with a `dim` attribute of length two. The list contains all the individual claim amounts for each entity. Since every element can be a vector, the object can be seen as a three-dimension array with a third dimension of potentially varying length. The function also returns a matrix of integers giving the classification indexes of each entity in the portfolio.

The package also defines methods for four generic functions to easily access key quantities for each entity of the simulated portfolio:

1. a method of `aggregate` to compute the aggregate claim amounts S;

2. a method of `frequency` to compute the number of claims N;

3. a method of `severity` (a generic function introduced by the package) to return the individual claim amounts C_j;

4. a method of `weights` to extract the weights matrix.

In addition, all methods have a `classification` and a `prefix` argument. When the first is `FALSE`, the classification index columns are omitted from the result. The second argument overrides the default column name prefix; see the `simul.summaries` help page for details.

The following example illustrates these concepts in detail.

Example 2. Consider the following compound hierarchical model:

$$S_{ijt} = C_{ijt1} + \cdots + C_{ijtN_{ijt}},$$

for $i = 1, \ldots, I$, $j = 1, \ldots, J_i$, $t = 1, \ldots, n_{ij}$ and with

$$N_{ijt}|\Lambda_{ij}, \Phi_i \sim \text{Poisson}(w_{ijt}\Lambda_{ij}) \qquad C_{ijtu}|\Theta_{ij}, \Psi_i \sim \text{Lognormal}(\Theta_{ij}, 1)$$
$$\Lambda_{ij}|\Phi_i \sim \text{Gamma}(\Phi_i, 1) \qquad \Theta_{ij}|\Psi_i \sim N(\Psi_i, 1)$$
$$\Phi_i \sim \text{Exponential}(2) \qquad \Psi_i \sim N(2, 0.1).$$

Using as convention to number the data level 0, the above is a two-level compound hierarchical model.

Assuming that $I = 2$, $J_1 = 4$, $J_2 = 3$, $n_{11} = \cdots = n_{14} = 4$ and $n_{21} = n_{22} = n_{23} = 5$ and that weights are simply simulated from a uniform distribution on $(0.5, 2.5)$, then simulation of a data set with `simul` is achieved with:

```
> nodes <- list(cohort = 2,
+               contract = c(4, 3),
+               year = c(4, 4, 4, 4, 5, 5, 5))
> mf <- expression(cohort = rexp(2),
+                  contract = rgamma(cohort, 1),
+                  year = rpois(weights * contract))
> ms <- expression(cohort = rnorm(2, sqrt(0.1)),
+                  contract = rnorm(cohort, 1),
+                  year = rlnorm(contract, 1))
> wijt <- runif(31, 0.5, 2.5)
> pf <- simul(nodes = nodes, model.freq = mf, model.sev = ms, weights = wijt)
```

Object `pf` is a list of class `"portfolio"` containing, among other things, the aforementioned two-dimension list as element `data` and the classification matrix (subscripts i and j) as element `classification`:

```
> class(pf)

[1] "portfolio"

> pf$data
```

```
        year.1     year.2     year.3     year.4     year.5
[1,]  Numeric,2  Numeric,2  11.38      Numeric,0  NA
[2,]  Numeric,0  Numeric,0  Numeric,0  Numeric,0  NA
[3,]  Numeric,0  Numeric,3  Numeric,0  Numeric,2  NA
[4,]  Numeric,0  98.13      50.62      55.7       NA
[5,]  Numeric,0  11.79      2.253      2.397      Numeric,2
[6,]  Numeric,0  Numeric,0  Numeric,0  Numeric,0  Numeric,0
[7,]  Numeric,3  Numeric,4  Numeric,2  Numeric,2  Numeric,0
```

```
> pf$classification
```

```
      cohort contract
[1,]       1        1
[2,]       1        2
[3,]       1        3
[4,]       1        4
[5,]       2        1
[6,]       2        2
[7,]       2        3
```

The output of pf$data is not much readable. Printing the results of simul like this would bring many users to wonder what Numeric,n means. It is actually R's way to specify that a given element in the list is a numeric vector of length n — the third dimension mentioned above. To ease reading, the print method for objects of class "portfolio" only prints the simulation model and the number of claims in each node:

```
> pf
```

```
Portfolio of claim amounts

  Frequency model
    cohort    ~ rexp(2)
    contract ~ rgamma(cohort, 1)
    year      ~ rpois(weights * contract)
  Severity model
    cohort    ~ rnorm(2, sqrt(0.1))
    contract ~ rnorm(cohort, 1)
    year      ~ rlnorm(contract, 1)

  Number of claims per node:

      cohort contract year.1 year.2 year.3 year.4 year.5
[1,]       1        1      2      2      1      0     NA
[2,]       1        2      0      0      0      0     NA
[3,]       1        3      0      3      0      2     NA
[4,]       1        4      0      1      1      1     NA
[5,]       2        1      0      1      1      1      2
[6,]       2        2      0      0      0      0      0
[7,]       2        3      3      4      2      2      0
```

By default, the method of `aggregate` returns the values of S_{ijt} in a regular matrix (subscripts i and j in the rows, subscript t in the columns). The method has a by argument to get statistics for other groupings and a `FUN` argument to get statistics other than the sum:

```
> aggregate(pf)

     cohort contract year.1 year.2 year.3 year.4 year.5
[1,]      1        1  31.37  7.521 11.383  0.000     NA
[2,]      1        2   0.00  0.000  0.000  0.000     NA
[3,]      1        3   0.00 72.706  0.000 23.981     NA
[4,]      1        4   0.00 98.130 50.622 55.705     NA
[5,]      2        1   0.00 11.793  2.253  2.397  10.48
[6,]      2        2   0.00  0.000  0.000  0.000   0.00
[7,]      2        3  44.81 88.737 57.593 14.589   0.00

> aggregate(pf, by = c("cohort", "year"), FUN = mean)

     cohort year.1 year.2 year.3 year.4 year.5
[1,]      1  15.69  29.73  31.00 26.562     NA
[2,]      2  14.94  20.11  19.95  5.662  5.238
```

The method of `frequency` returns the values of N_{ijt}. It is mostly a wrapper for the `aggregate` method with the default sum statistic replaced by `length`. Hence, arguments by and `FUN` remain available:

```
> frequency(pf)

     cohort contract year.1 year.2 year.3 year.4 year.5
[1,]      1        1      2      2      1      0     NA
[2,]      1        2      0      0      0      0     NA
[3,]      1        3      0      3      0      2     NA
[4,]      1        4      0      1      1      1     NA
[5,]      2        1      0      1      1      1      2
[6,]      2        2      0      0      0      0      0
[7,]      2        3      3      4      2      2      0

> frequency(pf, by = "cohort")

     cohort freq
[1,]      1   13
[2,]      2   16
```

The method of `severity` returns the individual variates C_{ijtu} in a matrix similar to those above, but with a number of columns equal to the maximum number of observations per entity,

$$\max_{i,j} \sum_{t=1}^{n_{ij}} N_{ijt}.$$

Thus, the original period of observation (subscript t) and the identifier of the severity within the period (subscript u) are lost and each variate now constitute a "period" of observation. For this reason, the method provides an argument splitcol in case one would like to extract separately the individual severities of one or more periods:

```
> severity(pf)

$main
     cohort contract claim.1 claim.2 claim.3 claim.4 claim.5
[1,]      1        1   7.974  23.401   3.153   4.368  11.383
[2,]      1        2      NA      NA      NA      NA      NA
[3,]      1        3   3.817  41.979  26.910   4.903  19.078
[4,]      1        4  98.130  50.622  55.705      NA      NA
[5,]      2        1  11.793   2.253   2.397   9.472   1.004
[6,]      2        2      NA      NA      NA      NA      NA
[7,]      2        3  14.322  11.522  18.966  33.108  15.532
     claim.6 claim.7 claim.8 claim.9 claim.10 claim.11
[1,]      NA      NA      NA      NA       NA       NA
[2,]      NA      NA      NA      NA       NA       NA
[3,]      NA      NA      NA      NA       NA       NA
[4,]      NA      NA      NA      NA       NA       NA
[5,]      NA      NA      NA      NA       NA       NA
[6,]      NA      NA      NA      NA       NA       NA
[7,]   14.99   25.11   40.15   17.44    4.426    10.16

$split
NULL

> severity(pf, splitcol = 1)

$main
     cohort contract claim.1 claim.2 claim.3 claim.4 claim.5
[1,]      1        1   3.153   4.368  11.383      NA      NA
[2,]      1        2      NA      NA      NA      NA      NA
[3,]      1        3   3.817  41.979  26.910   4.903  19.078
[4,]      1        4  98.130  50.622  55.705      NA      NA
[5,]      2        1  11.793   2.253   2.397   9.472   1.004
[6,]      2        2      NA      NA      NA      NA      NA
[7,]      2        3  33.108  15.532  14.990  25.107  40.150
     claim.6 claim.7 claim.8
[1,]      NA      NA      NA
[2,]      NA      NA      NA
[3,]      NA      NA      NA
[4,]      NA      NA      NA
[5,]      NA      NA      NA
[6,]      NA      NA      NA
[7,]   17.44   4.426   10.16

$split
```

```
      cohort contract claim.1 claim.2 claim.3
[1,]       1        1   7.974   23.40      NA
[2,]       1        2      NA      NA      NA
[3,]       1        3      NA      NA      NA
[4,]       1        4      NA      NA      NA
[5,]       2        1      NA      NA      NA
[6,]       2        2      NA      NA      NA
[7,]       2        3  14.322   11.52   18.97
```

Finally, the weights matrix corresponding to the data in object `pf` is

```
> weights(pf)
```

```
      cohort contract year.1 year.2 year.3 year.4 year.5
[1,]       1        1 0.8361  2.115 1.2699 1.1555     NA
[2,]       1        2 1.7042  1.709 0.7493 1.0892     NA
[3,]       1        3 1.6552  1.762 1.5240 1.5100     NA
[4,]       1        4 1.5681  1.614 2.2358 2.1594     NA
[5,]       2        1 0.7229  1.907 2.2950 1.0595 0.9564
[6,]       2        2 0.5307  0.758 0.6868 0.9738 2.0823
[7,]       2        3 1.6995  2.320 1.6208 2.0114 1.2583
```

Combined with the argument `classification = FALSE`, the above methods can be used to easily compute loss ratios:

```
> aggregate(pf, classif = FALSE) / weights(pf, classif = FALSE)
```

```
      year.1 year.2  year.3 year.4 year.5
[1,]  37.53  3.556  8.9638  0.000     NA
[2,]   0.00  0.000  0.0000  0.000     NA
[3,]   0.00 41.264  0.0000 15.881     NA
[4,]   0.00 60.781 22.6412 25.796     NA
[5,]   0.00  6.183  0.9818  2.263  10.95
[6,]   0.00  0.000  0.0000  0.000   0.00
[7,]  26.37 38.244 35.5328  7.253   0.00
```

□

Example 3. Scollnik (2001) considers the following model for the simulation of claims frequency data in a Markov Chain Monte Carlo (MCMC) context:

$$S_{it}|\Lambda_i, \alpha, \beta \sim \text{Poisson}(w_{ij}\Lambda_i)$$
$$\Lambda_i|\alpha, \beta \sim \text{Gamma}(\alpha, \beta)$$
$$\alpha \sim \text{Gamma}(5,5)$$
$$\beta \sim \text{Gamma}(25,1)$$

for $i = 1, 2, 3$, $j = 1, \ldots, 5$ and with weights w_{it} simulated from

$$w_{it}|a_i, b_i \sim \text{Gamma}(a_i, b_i)$$
$$a_i \sim U(0, 100)$$
$$b_i \sim U(0, 100).$$

8

Strictly speaking, this is not a hierarchical model since the random variables α and β are parallel rather than nested. Nevertheless, with some minor manual intervention, function `simul` can simulate data from this model.

First, one simulates the weights (in lexicographic order) with

```
> wit <- rgamma(15, rep(runif(3, 0, 100), each = 5),
+               rep(runif(3, 0, 100), each = 5))
```

Second, one calls `simul` to simulate the frequency data. The key here consists in manually inserting the simulation of the shape and rate parameters of the gamma distribution in the model for Λ_i. Finally, wrapping the call to `simul` in `frequency` will immediately yield the matrix of observations:

```
> frequency(simul(list(entity = 3, year = 5),
+       expression(entity = rgamma(rgamma(1, 5, 5), rgamma(1, 25, 1)),
+           year = rpois(weights * entity)),
+       weights = wit))

     entity year.1 year.2 year.3 year.4 year.5
[1,]      1      0      0      0      0      0
[2,]      2      0      0      0      0      0
[3,]      3      0      1      0      1      1
```

☐

One will find more examples of `simul` usage in the `simulation` demo file. Function `simul` was used to simulate the data in Forgues et al. (2006).

References

A. Forgues, V. Goulet, and J. Lu. Credibility for severity revisited. *North American Actuarial Journal*, 10(1):49–62, 2006.

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