# Simulation of compound hierarchical models

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## 1 Introduction

Hierarchical probability models are widely used for data classified in a treelike 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

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

where  $X_t$  represents actual data. The random variables  $\Theta$  and  $\Lambda$  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

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

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):

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

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

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  $C_i$ ;
- 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  $i = 1, \dots, I, \ j = 1, \dots, J_i, \ t = 1, \dots, n_{ij}$  and with  $N_{ijt} | \Lambda_{ij}, \Phi_i \sim \operatorname{Poisson}(w_{ijt}\Lambda_{ij}) \qquad C_{ijtu} | \Theta_{ij}, \Psi_i \sim \operatorname{Lognormal}(\Theta_{ij}, 1)$   $\Lambda_{ij} | \Phi_i \sim \operatorname{Gamma}(\Phi_i, 1) \qquad \Theta_{ij} | \Psi_i \sim N(\Psi_i, 1)$   $\Phi_i \sim \operatorname{Exponential}(2) \qquad \Psi_i \sim N(2, 0.1).$ 

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, O Numeric, O Numeric, O NA
[3,] Numeric, O Numeric, O Numeric, O 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, O Numeric, O Numeric, O Numeric, O Numeric, O
[7,] Numeric, 3 Numeric, 4 Numeric, 2 Numeric, 2 Numeric, 0
> pf$classification
     cohort contract
[1,]
          1
[2,]
          1
[3,]
          1
                   3
[4,]
          1
                    4
          2
[5,]
                   1
          2
                    2
[6,]
          2
                    3
[7,]
```

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 31.37 7.521 11.383 0.000
[2,]
                      0.00 0.000 0.000 0.000
                                                   NA
[3,]
                      0.00 72.706 0.000 23.981
                                                   NA
[4,]
                4 0.00 98.130 50.622 55.705
                                                   NA
[5,]
                1 0.00 11.793 2.253
                                        2.397
                                                10.48
[6,]
         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
```

2 14.94 20.11 19.95 5.662

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:

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#### > frequency(pf)

[2,]

> frequency(pf, by = "cohort")

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.1	l1
[1,]	NA	. NA	NA	NA	NA		۱A
[2,]	NA	. NA	NA	NA	NA		۱A
[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	
Γ7.]	14.99	25.11	40.15	17.44	4.426	10.1	L6

\$split
NULL

> severity(pf, splitcol = 1)

#### \$main

Jilla I I	!						
	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,]
                         7.974
                                  23.40
                    1
[2,]
           1
                    2
                            NA
                                     NA
                                               NA
[3,]
          1
                    3
                            NA
                                     NA
                                              NA
          1
                                              NA
[4,]
                            NA
                                     NA
           2
                     1
                            NA
[5,]
                                     NA
                                               NA
           2
                     2
                            NA
[6,]
                                      NA
                                               NA
[7,]
                        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 0.8361 2.115 1.2699 1.1555
          1
                   2 1.7042 1.709 0.7493 1.0892
[2,]
                                                      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
                            2.320 1.6208 2.0114 1.2583
[7,]
                   3 1.6995
```

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
     37.53 3.556
                   8.9638 0.000
[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
                                 10.95
[5,]
      0.00 6.183 0.9818
                          2.263
      0.00 0.000 0.0000 0.000
[6,]
                                   0.00
     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, ..., 5 and with weights  $w_{it}$  simulated from

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

Strictly speaking, this is not a hierarchical model since the random variables  $\alpha$  and  $\beta$  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))</pre>
```

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  $\Lambda_i$ . 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
          2
                 0
                        0
                               0
                                      0
                                              0
[2,]
[3,]
          3
                 0
                        1
                               0
```

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).

П

## References

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