

# Graphs in the **gRbase** package

Søren Højsgaard

January 28, 2013

## Contents

<b>1</b>	<b>Introduction</b>	<b>1</b>
<b>2</b>	<b>Graphs</b>	<b>2</b>
2.1	Undirected graphs . . . . .	2
2.2	Directed acyclic graphs (DAGs) . . . . .	4
2.3	Graph coercion . . . . .	5
2.4	Plotting graphs . . . . .	6
<b>3</b>	<b>Advanced graph operations</b>	<b>7</b>
3.1	Moralization . . . . .	7
3.2	Topological sort . . . . .	8
3.3	Getting cliques . . . . .	9
3.4	Maximum cardinality search . . . . .	10
3.5	Triangulation . . . . .	11
3.6	RIP ordering / junction tree . . . . .	12
3.7	Minimal triangulation and maximum prime subgraph decomposition . . . .	13
<b>4</b>	<b>Time and space considerations</b>	<b>13</b>
4.1	Time . . . . .	13
4.2	Space . . . . .	14
<b>5</b>	<b>Graph queries</b>	<b>15</b>

## 1 Introduction

For the R community, the packages **graph**, **RBGL**, **Rgraphviz** and **igraph** are extremely useful tools for graph operations, manipulation and layout. The **gRbase** package adds some additional tools to these fine packages. The most important tools are:

1. Undirected and directed acyclic graphs can be specified using formulae or an adjacency list using the functions `ug()` and `dag()`. This gives graphs represented in one of the following forms:<sup>1</sup>
  - A `graphNEL` object (the default),
  - A dense adjacency matrix. By this we mean a “standard” matrix in R.
  - A sparse adjacency matrix. By this we mean a `dgCMatrix` from the `Matrix` package.
2. Some graph algorithms are implemented in **gRbase**. These can be applied to graphs represented as `graphNELs` and matrices. The most important algorithms are:
  - `moralize()`, (moralize a directed acyclic graph)
  - `mcs()`, (maximum cardinality search for undirected graph)
  - `triangulate()`, (triangulate undirected graph)
  - `rip()`, (RIP ordering of cliques of triangulated undirected graph)
  - `getCliques()`, (get the (maximal) cliques of an undirected graph)
  - `minimalTriang()` (minimal triangulation of undirected graph)<sup>2</sup>
  - `mpd()` (maximal prime subgraph decomposition of undirected graph)<sup>3</sup>

The general scheme is the following: There is a `mcs()` function and `mcs()` methods for `graphNELs` and for the two matrix types. The workhorse is the function `mcsMAT()` and the various methods coerces the graph to a (sparse) matrix and invokes `mcsMAT()`.

## 2 Graphs

Undirected graphs can be created by the `ug()` function and directed acyclic graphs (DAGs) by the `dag()` function.

The graphs can be specified either using formulae or a list of vectors; see examples below.

### 2.1 Undirected graphs

An undirected graph is created by the `ug()` function.

---

<sup>1</sup>There is a fourth form: `igraph` objects. These, however, will probably not be supported in the future.

<sup>2</sup>Needs more work

<sup>3</sup>Needs more work

**As graphNEL:** The following specifications are equivalent (notice that “.” and “\*” can be used interchangeably):

```
R> ug11 <- ug(~a:b:c + c:d + d:e + a:e + f:g)
R> ug11 <- ug(~a*b*c + c*d + d*e + a*e + f*g)
R> ug12 <- ug(c("a","b","c"),c("c","d"),c("d","e"),c("a","e"),c("f","g"))
R> ug13 <- ug(~a:b:c, ~c:d, ~d:e + a:e + f:g)
R> ug13 <- ug(~a*b*c, ~c*d, ~d*e + a*e + f*g)
```

```
R> ug11
```

```
A graphNEL graph with undirected edges
Number of Nodes = 7
Number of Edges = 7
```

**As adjacency matrix:** A representation as an adjacency matrix can be obtained with one of the following equivalent specifications:

```
R> ug11m <- ug(~a*b*c + c*d + d*e + a*e + f*g, result="matrix")
R> ug12m <- ug(c("a","b","c"),c("c","d"),c("d","e"),c("a","e"),c("f","g"),
               result="matrix")
```

```
R> ug11m
```

```
  a b c d e f g
a 0 1 1 0 1 0 0
b 1 0 1 0 0 0 0
c 1 1 0 1 0 0 0
d 0 0 1 0 1 0 0
e 1 0 0 1 0 0 0
f 0 0 0 0 0 0 1
g 0 0 0 0 0 1 0
```

```
R> ug11M <- ug(~a*b*c + c*d + d*e + a*e + f*g, result="Matrix")
R> ug12M <- ug(c("a","b","c"),c("c","d"),c("d","e"),c("a","e"),c("f","g"),
               result="Matrix")
```

```
R> ug11M
```

```
7 x 7 sparse Matrix of class "dgCMatrix"
  a b c d e f g
a . 1 1 . 1 . .
b 1 . 1 . . . .
c 1 1 . 1 . . .
d . . 1 . 1 . .
e 1 . . 1 . . .
f . . . . . 1 .
g . . . . . 1 .
```

## 2.2 Directed acyclic graphs (DAGs)

A directed acyclic graph is created by the `dag()` function.

**As graphNEL:** The following specifications are equivalent (notice that “:” and “\*” can be used interchangeably):

```
R> dag11 <- dag(~a + b:a + c:a:b + d:c:e + e:a + g:f)
R> dag11 <- dag(~a + b*a + c*a*b + d*c*e + e*a + g*f)
R> dag12 <- dag("a", c("b","a"), c("c","a","b"), c("d","c","e"),
               c("e","a"),c("g","f"))
R> dag13 <- dag(~a, ~b:a, ~c:a:b, ~d:c:e, ~e:a, ~g:f)
R> dag13 <- dag(~a, ~b*a, ~c*a*b, ~d*c*e, ~e*a, ~g*f)
```

```
R> dag11
```

```
A graphNEL graph with directed edges
Number of Nodes = 7
Number of Edges = 7
```

Here `~a` means that “a” has no parents while `~d:b:c` means that “d” has parents “b” and “c”.

**As adjacency matrix:** A representation as an adjacency matrix can be obtained with

```
R> dag11m <- dag(~a + b:a + c:a:b + d:c:e + e:a + g:f, result="matrix")
R> dag12m <- dag("a", c("b","a"), c("c","a","b"), c("d","c","e"),
               c("e","a"),c("g","f"), result="matrix")
```

```
R> dag11m
```

```
   a b c d e g f
a 0 1 1 0 1 0 0
b 0 0 1 0 0 0 0
c 0 0 0 1 0 0 0
d 0 0 0 0 0 0 0
e 0 0 0 1 0 0 0
g 0 0 0 0 0 0 0
f 0 0 0 0 0 1 0
```

```
R> dag11M <- dag(~a + b:a + c:a:b + d:c:e + e:a + g:f, result="Matrix")
R> dag12M <- dag("a", c("b","a"), c("c","a","b"), c("d","c","e"),
               c("e","a"),c("g","f"), result="Matrix")
```

```
R> dag11M
```

```

7 x 7 sparse Matrix of class "dgCMatrix"
  a b c d e g f
a . 1 1 . 1 . .
b . . 1 . . . .
c . . . 1 . . .
d . . . . . . .
e . . . 1 . . .
g . . . . . . .
f . . . . . 1 .

```

## 2.3 Graph coercion

Graphs can be coerced between different representations using `as()`; for example

```
R> as(ug11,"matrix")
```

```

  a b c d e f g
a 0 1 1 0 1 0 0
b 1 0 1 0 0 0 0
c 1 1 0 1 0 0 0
d 0 0 1 0 1 0 0
e 1 0 0 1 0 0 0
f 0 0 0 0 0 0 1
g 0 0 0 0 0 1 0

```

```
R> as(as(ug11,"matrix"),"dgCMatrix")
```

```

7 x 7 sparse Matrix of class "dgCMatrix"
  a b c d e f g
a . 1 1 . 1 . .
b 1 . 1 . . . .
c 1 1 . 1 . . .
d . . 1 . 1 . .
e 1 . . 1 . . .
f . . . . . 1 .
g . . . . . 1 .

```

```
R> as(as(as(ug11,"matrix"),"dgCMatrix"),"graphNEL")
```

```

A graphNEL graph with undirected edges
Number of Nodes = 7
Number of Edges = 7

```

```
R> as(as(as(as(ug11,"matrix"),"Matrix"),"graphNEL"),"igraph")
```

```

IGRAPH UNW- 7 7 --
+ attr: name (v/c), label (v/c), weight (e/n)

```

**NOTICE!** There is one thing to notice when coercing a dense matrix to a sparse matrix. Consider this

```
R> m <- matrix(1:4,nrow=2)
R> as(m, "Matrix")

2 x 2 Matrix of class "dgeMatrix"
      [,1] [,2]
[1,]    1    3
[2,]    2    4

R> as(m, "dgCMatrix")

2 x 2 sparse Matrix of class "dgCMatrix"

[1,] 1 3
[2,] 2 4
```

In the first case, the matrix coercion method will, based on properties of the matrix, impose a specific type on the result. The graph algorithms in **gRbase** are based on the **dgCMatrix** form and hence the latter case is more safe. Coercion to a **dgCMatrix** can be made much faster with the following function from **gRbase**:

```
R> asdgCMatrix(m)

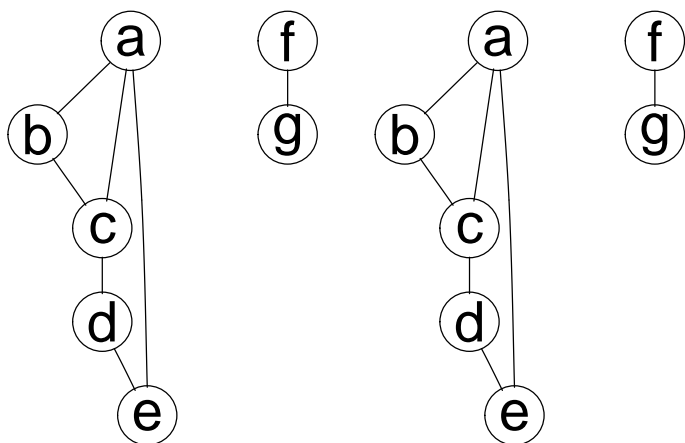
2 x 2 sparse Matrix of class "dgCMatrix"

[1,] 1 3
[2,] 2 4
```

## 2.4 Plotting graphs

Graphs represented as **graphNEL** objects are displayed with **plot()**. There is no **plot()** method for graphs represented as adjacency matrices, so here coercion is one option:

```
R> par(mfrow=c(1,2))
R> plot(ug11)
R> plot(as(ug11m, "graphNEL"))
```



An alternative for adjacency matrices is the `gplot()` function in the **sna** package:

```

R> par(mfrow=c(1,2))
R> library(sna)
R> gplot(ug11m, label=colnames(ug11m),gmode="graph")
R> gplot(dag11m, label=colnames(dag11m))

```

## 3 Advanced graph operations

### 3.1 Moralization

```

R> apropos("^moralize\\.")

```

```

[1] "moralize.Matrix"    "moralize.graphNEL" "moralize.igraph"
[4] "moralize.matrix"

```

A moralized directed acyclic graph is obtained with

```

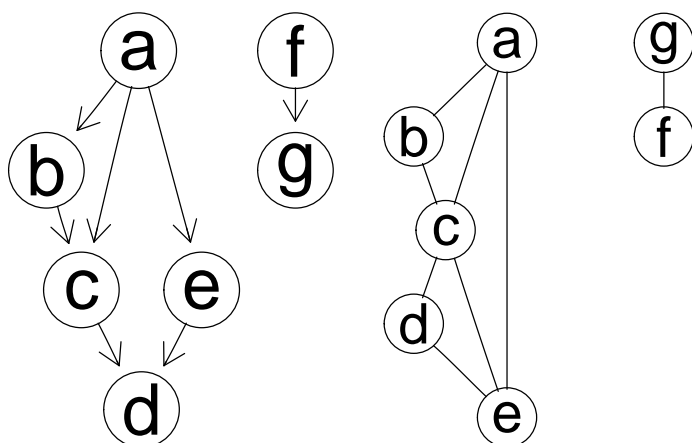
R> dag11.mor <- moralize(dag11)

```

```

R> par(mfrow=c(1,2))
R> plot(dag11)
R> plot(dag11.mor)

```



For the alternative representations

```
R> moralize(dag11m)
```

```

  a b c d e g f
a 0 1 1 0 1 0 0
b 1 0 1 0 0 0 0
c 1 1 0 1 1 0 0
d 0 0 1 0 1 0 0
e 1 0 1 1 0 0 0
g 0 0 0 0 0 0 1
f 0 0 0 0 0 1 0

```

```
R> moralize(dag11M)
```

```

7 x 7 sparse Matrix of class "dgCMatrix"
  a b c d e g f
a . 1 1 . 1 . .
b 1 . 1 . . . .
c 1 1 . 1 1 . .
d . . 1 . 1 . .
e 1 . 1 1 . . .
g . . . . . 1
f . . . . . 1 .

```

## 3.2 Topological sort

A topological ordering of a directed graph is a linear ordering of its vertices such that, for every edge  $(u \rightarrow v)$ ,  $u$  comes before  $v$  in the ordering. A topological ordering is possible if and only if the graph has no directed cycles, that is, if it is a directed acyclic graph (DAG). Any DAG has at least one topological ordering.

```
R> topoSort(dag11)
```

```
[1] "a" "b" "c" "e" "f" "d" "g"
```

```
R> topoSort(dag11m)
```

```
[1] "a" "b" "c" "e" "f" "d" "g"
```

```
R> topoSort(dag11M)
```

```
[1] "a" "b" "c" "e" "f" "d" "g"
```

The `dag()` function actually allows specification of a directed graph with cycles (a check can be imposed with `forceCheck=TRUE`). Below is a directed graph with a cycle

```
R> topoSort(dag(~a:b+b:c+c:a))
```

```
character(0)
```

### 3.3 Getting cliques

In graph theory, a clique is often a complete subset of a graph. A maximal clique is a clique which can not be enlarged. In statistics (and that is the convention we follow here) a clique is usually understood to be a maximal clique.

Finding the cliques of a general graph is an NP complete problem. Finding the cliques of a triangulated graph is linear in the number of cliques.

```
R> str(getCliques(ug11))
```

```
List of 5
```

```
$ : chr [1:3] "a" "b" "c"  
$ : chr [1:2] "a" "e"  
$ : chr [1:2] "d" "c"  
$ : chr [1:2] "d" "e"  
$ : chr [1:2] "f" "g"
```

```
R> str(getCliques(ug11m))
```

```
List of 5
```

```
$ : chr [1:3] "a" "b" "c"  
$ : chr [1:2] "a" "e"  
$ : chr [1:2] "d" "c"  
$ : chr [1:2] "d" "e"  
$ : chr [1:2] "f" "g"
```

```
R> str(getCliques(ug11M))
```

```
List of 5
```

```
$ : chr [1:3] "a" "b" "c"  
$ : chr [1:2] "a" "e"  
$ : chr [1:2] "d" "c"  
$ : chr [1:2] "d" "e"  
$ : chr [1:2] "f" "g"
```

### 3.4 Maximum cardinality search

```
R> apropos("^mcs\\.")
```

```
[1] "mcs.Matrix"    "mcs.graphNEL" "mcs.igraph"    "mcs.matrix"
```

Testing for whether a graph is triangulated is based on Maximum Cardinality Search. If `character(0)` is returned the graph is not triangulated. Otherwise a linear ordering of the nodes is returned.

```
R> mcs(ug11)
```

```
character(0)
```

```
R> mcs(ug11m)
```

```
character(0)
```

```
R> mcs(ug11M)
```

```
character(0)
```

```
R> mcs(dag11.mor)
```

```
[1] "a" "b" "c" "e" "d" "g" "f"
```

```
R> mcs(as(dag11.mor,"matrix"))
```

```
[1] "a" "b" "c" "e" "d" "g" "f"
```

```
R> mcs(as(dag11.mor,"Matrix"))
```

```
[1] "a" "b" "c" "e" "d" "g" "f"
```

```
R> mcs(dag11)
```

```
character(0)
```

### 3.5 Triangulation

```
R> apropos("^triangulate\\.")
```

```
[1] "triangulate.Matrix"    "triangulate.graphNEL" "triangulate.igraph"  
[4] "triangulate.matrix"
```

Triangulate an undirected graph by adding extra edges to the graph:

```
R> (tug11<-triangulate(ug11))
```

```
A graphNEL graph with undirected edges  
Number of Nodes = 7  
Number of Edges = 8
```

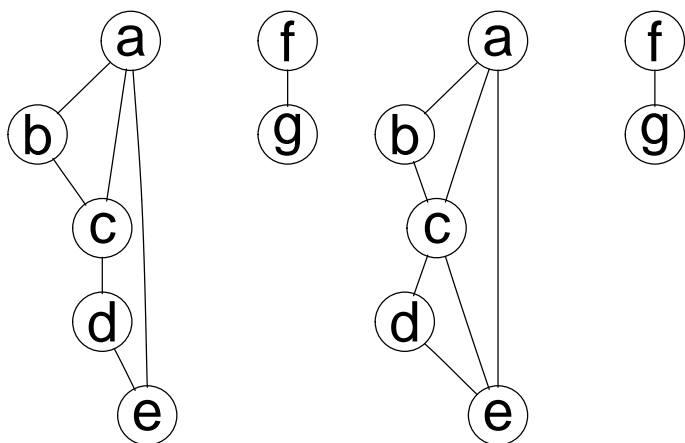
```
R> (tug11m<-triangulate(ug11m))
```

```
  a b c d e f g  
a 0 1 1 0 1 0 0  
b 1 0 1 0 0 0 0  
c 1 1 0 1 1 0 0  
d 0 0 1 0 1 0 0  
e 1 0 1 1 0 0 0  
f 0 0 0 0 0 0 1  
g 0 0 0 0 0 1 0
```

```
R> (tug11M<-triangulate(ug11M))
```

```
7 x 7 sparse Matrix of class "dgCMatrix"  
  a b c d e f g  
a . 1 1 . 1 . .  
b 1 . 1 . . . .  
c 1 1 . 1 1 . .  
d . . 1 . 1 . .  
e 1 . 1 1 . . .  
f . . . . . 1  
g . . . . . 1 .
```

```
R> par(mfrow=c(1,2))  
R> plot(ug11)  
R> plot(tug11)
```



### 3.6 RIP ordering / junction tree

```
R> apropos("^rip\\.")
```

```
[1] "rip.Matrix"    "rip.graphNEL" "rip.igraph"    "rip.matrix"
```

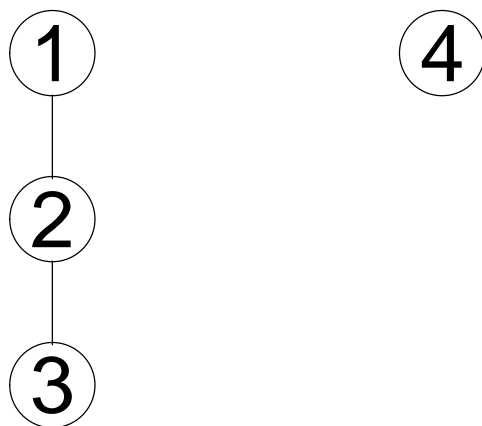
A RIP ordering of the cliques of a triangulated graph can be obtained as:

```
R> rr <- rip(tug11)
R> rr
```

```
cliques
 1 : c a b
 2 : e a c
 3 : d c e
 4 : g f
separators
 1 :
 2 : a c
 3 : c e
 4 :
parents
 1 : 0
 2 : 1
 3 : 2
 4 : 0
```

```
R> rr <- rip(tug11m)
R> rr <- rip(tug11M)
```

```
R> plot(rr)
```



### 3.7 Minimal triangulation and maximum prime subgraph decomposition

An undirected graph  $uG$  is triangulated (or chordal) if it has no cycles of length  $\geq 4$  without a chord which is equivalent to that the vertices can be given a perfect ordering. Any undirected graph can be triangulated by adding edges to the graph, so called fill-ins which gives the graph  $TuG$ . A triangulation  $TuG$  is minimal if no fill-ins can be removed without breaking the property that  $TuG$  is triangulated. A related concept is the minimum triangulation, which is the the graph with the smallest number of fill-ins. The minimum triangulation is unique. Finding the minimum triangulation is NP-hard.

## 4 Time and space considerations

### 4.1 Time

It is worth noticing that working with graphs representated as **graphNEL** objects is somewhat slower working with graphs represented as adjacency matrices. Consider finding the cliques of an undirected graph represented as a **graphNEL** object or as a matrix:

```
R> system.time({for (ii in 1:200) maxClique(ug11)}) ## in RBGL
```

```

user  system elapsed
0.10   0.00   0.11

```

```
R> system.time({for (ii in 1:200) maxCliqueMAT(ug11m)}) ## in gRbase
```

```

user  system elapsed
0.03   0.00   0.03

```

Working with sparse matrices rather than standard matrices slows indexing down:

```
R> system.time({for (ii in 1:2000) ug11m[2,]})
```

```
   user  system elapsed
    0      0      0
```

```
R> system.time({for (ii in 1:2000) ug11M[2,]})
```

```
   user  system elapsed
 0.69    0.00    0.69
```

However, **gRbase** has some functionality for indexing sparse matrices quickly:

```
R> system.time({for (ii in 1:2000) sp_getXj(ug11M,2)})
```

```
   user  system elapsed
 0.01    0.00    0.02
```

## 4.2 Space

The **graphNEL** representation is – at least – in principle more economic in terms of space requirements than the adjacency matrix representation (because the adjacency matrix representation uses a 0 to represent a “missing edge”. The sparse matrix representation is clearly only superior to the standard matrix representation if the graph is sparse:

```
R> V <- 1:100
R> M <- 1:10
R> ## Sparse graph
R> ##
R> g1 <- randomGraph(V, M, 0.05)
R> length(edgeList(g1))

[1] 128
```

```
R> c(NEL=object.size(g1),
    mat=object.size(as(g1, "matrix")),
    Mat=object.size(as.adjMAT(g1, "Matrix")))
```

```
   NEL    mat    Mat
142896  51648  16176
```

```
R> ## More dense graph
R> ##
R> g1 <- randomGraph(V, M, 0.5)
R> length(edgeList(g1))
```

```
[1] 4618
```

```
R> c(NEL=object.size(g1),  
    mat=object.size(as(g1, "matrix")),  
    Mat=object.size(as.adjMAT(g1, "Matrix")))
```

NEL	mat	Mat
3519160	51648	123936

## 5 Graph queries

The **graph** and **RBGL** packages implement various graph operations for **graphNEL** objects. See the documentation for these packages. The **gRbase** implements a few additional functions, see Section 1. An additional function in **gRbase** for graph operations is **querygraph()**. This function is intended as a wrapper for the various graph operations available in **gRbase**, **graph** and **RBGL**. There are two main virtues of **querygraph()**: 1) **querygraph()** operates on any of the three graph representations described above<sup>4</sup> and 2) **querygraph()** provides a unified interface to the graph operations. The general syntax is

```
R> args(querygraph)
```

```
function (object, op, set = NULL, set2 = NULL, set3 = NULL)  
NULL
```

---

<sup>4</sup>Actually not quite yet, but it will be so in the future.