

# Package ‘gRain’

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**Version** 1.3.14

**Title** Graphical Independence Networks

**Author** Søren Højsgaard <sorenh@math.aau.dk>

**Maintainer** Søren Højsgaard <sorenh@math.aau.dk>

**Description** Probability propagation in graphical independence networks, also known as Bayesian networks or probabilistic expert systems. Documentation of the package is provided in vignettes included in the package and in the paper by Højsgaard (2012, <[doi:10.18637/jss.v046.i10](https://doi.org/10.18637/jss.v046.i10)>). See 'citation("gRain")' for details.

**License** GPL (>= 2)

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**Imports** igraph, stats4, broom, magrittr, Rcpp (>= 0.11.1)

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chest

*Chest clinic example*

**Description**

Conditional probability tables for the chest clinic example.

**Usage**

data(chest\_cpt)

**Format**

An object of class list of length 8.

**Examples**

```
## 'data' generated with the following code fragment
yn <- c("yes", "no")
a <- cptable(~asia, values=c(1,99),levels=yn)
t.a <- cptable(~tub|asia, values=c(5,95,1,99),levels=yn)
s <- cptable(~smoke, values=c(5,5), levels=yn)
l.s <- cptable(~lung|smoke, values=c(1,9,1,99), levels=yn)
b.s <- cptable(~bronc|smoke, values=c(6,4,3,7), levels=yn)
e.lt <- cptable(~either|lung:tub, values=c(1,0,1,0,1,0,0,1), levels=yn)
x.e <- cptable(~xray|either, values=c(98,2,5,95), levels=yn)
```

```
d.be <- ctable(~dysp|bronc:either, values=c(9,1,7,3,8,2,1,9), levels=yn)

grain(compileCPT(a, t.a, s, l.s, b.s, e.lt, x.e, d.be))

# 'data' generated from
# chest_cpt <- list(a, t.a, s, l.s, b.s, e.lt, x.e, d.be)

data(chest_cpt)
```

---

components\_extract      *Extract conditional probabilities and clique potentials from data.*

---

### Description

Extract list of conditional probability tables and list of clique potentials from data.

### Usage

```
extractCPT(data_, graph, smooth = 0)

extractPOT(data_, graph, smooth = 0)

extractMARG(data_, graph, smooth = 0)

marg2pot(mg)

pot2marg(pt)
```

### Arguments

data_	A named array or a dataframe.
graph	An igraph object or a list or formula which can be turned into a igraph object by calling ug or dag. For extractCPT, graph must be/define a DAG while for extractPOT, graph must be/define undirected triangulated graph.
smooth	See 'details' below.
mg	An object of class marg_rep
pt	An object of class pot_representation

### Details

If smooth is non-zero then smooth is added to all cell counts before normalization takes place.

### Value

- extractCPT: A list of conditional probability tables.
- extractPOT: A list of clique potentials.
- extractMARG: A list of clique marginals.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[compileCPT](#), [compilePOT](#), [grain](#)

**Examples**

```
## Extract cpts / clique potentials from data and graph
# specification and create network. There are different ways:

data(lizard, package="gRbase")

# DAG: height <- species -> diam
daG <- dag(~species + height:species + diam:species, result="igraph")

# UG : [height:species][diam:species]
uG <- ug(~height:species + diam:species, result="igraph")

pt <- extractPOT(lizard, ~height:species + diam:species)
cp <- extractCPT(lizard, ~species + height:species + diam:species)

pt
cp

# Both specify the same probability distribution
tabListMult(pt) %>% as.data.frame.table
tabListMult(cp) %>% as.data.frame.table

## Not run:
# Bayesian networks can be created as
bn.uG <- grain(pt)
bn.daG <- grain(cp)

# The steps above are wrapped into a convenience method which
# builds a network from at graph and data.
bn.uG <- grain(uG, data=lizard)
bn.daG <- grain(daG, data=lizard)

## End(Not run)
```

---

components\_gather      *Compile conditional probability tables / cliques potentials.*

---

### Description

Compile conditional probability tables / cliques potentials as a preprocessing step for creating a graphical independence network

### Usage

```
compileCPT(x, ..., forceCheck = TRUE)
```

```
compilePOT(x, ..., forceCheck = TRUE)
```

```
parse_cpt(xi)
```

### Arguments

x	To compileCPT x is a list of conditional probability tables; to compilePOT, x is a list of clique potentials.
...	Additional arguments; currently not used.
forceCheck	Controls if consistency checks of the probability tables should be made.
xi	cpt in some representation

### Details

\* ``compileCPT`` is relevant for turning a collection of `cptable`'s into an object from which a network can be built. For example, when specification of a `cpt` is made with `cptable` then the levels of the node is given but not the levels of the parents. ``compileCPT`` checks that the levels of variables in the `cpt`'s are consistent and also that the specifications define a dag.

\* ``compilePOT`` is not of direct relevance for the user for the moment. However, the elements of the input should be arrays which define a chordal undirected graph and the arrays should, if multiplied, form a valid probability density.

### Value

A list with a class attribute.

### Author(s)

Søren Højsgaard, <sorenh@math.aau.dk>

## References

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

## See Also

[extractCPT](#), [extractPOT](#), [extractMARG](#)

## Examples

```
data(chest_cpt)
x <- compileCPT(chest_cpt)
class(x)
grain(x)
```

---

cpt

*Create conditional probability table CPT.*

---

## Description

Create conditional probability table CPT.

## Usage

```
cpt(names, levels, values, smooth = 0)
```

## Arguments

names	Names of variables defining table; either a character vector or a right hand sided formula.
levels	1. a list with specification of the levels of the factors in names or 2) a vector with number of levels of the factors in names. See 'examples' below.
values	values to go into the array.
smooth	Should values be smoothed, see 'Details' below.

## Value

An array.

**Examples**

```

universe <- list(gender=c('male', 'female'),
                 answer=c('yes', 'no'),
                 rain=c('yes', 'no'))
t1 <- cpt(c("gender", "answer"), levels=universe, values=1:4)
t1
t2 <- cpt(~gender:answer, levels=universe, values=1:4)
t2
t3 <- cpt(~gender:answer, c(2, 2), values=1:4)
t3

```

cptable

*Create conditional probability tables (CPTs)***Description**

Creates conditional probability tables of the form  $p(v|pa(v))$ .

**Usage**

```
cptable(vpar, levels = NULL, values = NULL, normalize = TRUE, smooth = 0)
```

**Arguments**

vpar	Specifications of the names in $P(v pa1, \dots, pa_k)$ . See section 'details' for information about the form of the argument.
levels	See 'details' below.
values	Probabilities; recycled if necessary. Regarding the order, please see section 'details' and the examples.
normalize	See 'details' below.
smooth	See 'details' below.

**Details**

If `normalize=TRUE` then the probabilities are normalized to sum to one for each configuration of the parents.

If `smooth` is non-zero then zero entries of values are replaced with `smooth` **before** normalization takes place.

Regarding the form of the argument `vpar`: To specify  $P(a|b, c)$  one may write `~a|b:c`, `~a:b:c`, `~a|b+c`, `~a+b+c` or `c("a", "b", "c")`. Internally, the last form is used. Notice that the `+` and `:` operator is used as a separator only. The order of the variables IS important so the operators DO NOT commute.

If `a` has levels `a1, a2` and likewise for `b` and `c` then the order of values corresponds to the configurations `(a1, b1, c1)`, `(a2, b1, c1)`, `(a1, b2, c1)`, `(a2, b2, c1)` etc. That is, the first variable varies fastest. Hence the first two elements in values will be the conditional probabilities of a given `b=b1, c=c1`.

**Value**

A cptable object (a numeric vector with various attributes).

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[andtable](#), [ortable](#), [extractCPT](#), [compileCPT](#), [extractPOT](#), [compilePOT](#), [grain](#), [parray](#)

**Examples**

```
## See the wet grass example at
## https://en.wikipedia.org/wiki/Bayesian_network

yn <- c("yes", "no")
p.R   <- cptable(~R, values=c(.2, .8), levels=yn)
p.S_R <- cptable(~S:R, values=c(.01, .99, .4, .6), levels=yn)
p.G_SR <- cptable(~G:S:R, values=c(.99, .01, .8, .2, .9, .1, 0, 1), levels=yn)

# or
ssp <- list(R=yn, S=yn, G=yn) # state space
p.R   <- cptable(~R, values=c(.2, .8), levels=ssp)
p.S_R <- cptable(~S:R, values=c(.01, .99, .4, .6), levels=ssp)
p.G_SR <- cptable(~G:S:R, values=c(.99, .01, .8, .2, .9, .1, 0, 1), levels=ssp)

# components above are "intermediate representations" and are turned into arrays with
wet.cpt <- compileCPT(p.R, p.S_R, p.G_SR)
wet.cpt
wet.cpt$S # etc

# A Bayesian network is created with:
wet.bn <- grain(wet.cpt)

# Can also create arrays directly
## Not run:
ssp <- list(R=yn, S=yn, G=yn) # state space
p.R   <- c(.2, .8)
p.S_R <- c(.01, .99, .4, .6)
p.G_SR <- c(.99, .01, .8, .2, .9, .1, 0, 1)
dim(p.R) <- 2
dimnames(p.R) <- ssp["R"]
dim(p.S_R) <- c(2, 2)
dimnames(p.S_R) <- ssp[c("S", "R")]
```



```

dim(p.G_SR) <- c(2, 2, 2)
dimnames(p.G_SR) <- ssp[c("G", "S", "R")]

# Arrays can be created (easier?) with parray() from gRbase
p.R   <- parray("R", levels=ssp, values=c(.2, .8))
p.S_R <- parray(c("S", "R"), levels = ssp, values=c(.01, .99, .4, .6))
p.G_SR <- parray(~ G:S:R, levels = ssp, values=c(.99, .01, .8, .2, .9, .1, 0, 1))

## End(Not run)

```

---

evidence_object	<i>Evidence objects</i>
-----------------	-------------------------

---

## Description

Functions for defining and manipulating evidence.

## Usage

```

new_evi(evi_list = NULL, levels)

is.null_evi(object)

## S3 method for class 'grain_evidence'
as.data.frame(x, row.names = NULL, optional = FALSE, ...)

setdiff_evi(ev1, ev2)

union_evi(ev1, ev2)

```

## Arguments

evi_list	A named list with evidence; see 'examples' below.
levels	A named list with the levels of all variables.
object	Some R object.
x	An evidence object.
row.names	Not used.
optional	Not used.
...	Not used.
ev1, ev2	Evidence.

## Details

Evidence is specified as a list. Internally, evidence is represented as a grain evidence object which is a list with 4 elements.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**Examples**

```
## Define the universe
yn <- c("yes", "no")
uni <- list(asia = yn, tub = yn, smoke = yn, lung = yn,
           bronc = yn, either = yn, xray = yn, dysp = yn)

e1 <- list(dysp="no", xray="no")
eo1 <- new_evi(e1, levels=uni)
eo1 |> as.data.frame()

e2 <- list(dysp="no", xray=c(0, 1))
eo2 <- new_evi(e2, levels=uni)
eo2 |> as.data.frame()

# Above e1 and e2 specifies the same evidence but information about
# whether the state has been set definite or as a weight is
# maintained.

e3 <- list(dysp="yes", asia="yes")
eo3 <- new_evi(e3, uni)
eo3 |> as.data.frame()

# If evidence 'e1' is already set in the network and new evidence
# 'e3' emerges, then evidence in the network must be updated. But
# there is a conflict in that dysp="yes" in 'e1' and
# dysp="no" in 'e3'. The (arbitrary) convention is that
# existing evidence overrides new evidence so that the only new
# evidence in 'e3' is really asia="yes".

# To subtract existing evidence from new evidence we can do:
setdiff_evi(eo3, eo1) |> as.data.frame()

# Likewise the 'union' is
union_evi(eo3, eo1) |> as.data.frame()
```

---

finding

*Set, retrieve, and retract finding in Bayesian network.*

---

**Description**

Set, retrieve, and retract finding in Bayesian network. NOTICE: The functions described here are kept only for backward compatibility; please use the corresponding evidence-functions in the future.

**Usage**

```
setFinding(object, nodes = NULL, states = NULL, flist = NULL, propagate = TRUE)
```

**Arguments**

object	A "grain" object
nodes	A vector of nodes
states	A vector of states (of the nodes given by 'nodes')
flist	An alternative way of specifying findings, see examples below.
propagate	Should the network be propagated?

**Note**

NOTICE: The functions described here are kept only for backward compatibility; please use the corresponding evidence-functions in the future:

`setEvidence()` is an improvement of `setFinding()` (and as such `setFinding` is obsolete). Users are recommended to use `setEvidence()` in the future.

`setEvidence()` allows to specification of "hard evidence" (specific values for variables) and likelihood evidence (also known as virtual evidence) for variables.

The syntax of `setEvidence()` may change in the future.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[setEvidence](#), [getEvidence](#), [retractEvidence](#), [pEvidence](#), [querygrain](#)

**Examples**

```
## setFindings
yn <- c("yes", "no")
a <- cptable(~asia, values=c(1, 99), levels=yn)
t.a <- cptable(~tub+asia, values=c(5, 95, 1, 99), levels=yn)
s <- cptable(~smoke, values=c(5,5), levels=yn)
l.s <- cptable(~lung+smoke, values=c(1, 9, 1, 99), levels=yn)
b.s <- cptable(~bronc+smoke, values=c(6, 4, 3, 7), levels=yn)
e.lt <- cptable(~either+lung+tub, values=c(1, 0, 1, 0, 1, 0, 0, 1), levels=yn)
x.e <- cptable(~xray+either, values=c(98, 2, 5, 95), levels=yn)
d.be <- cptable(~dysp+bronc+either, values=c(9, 1, 7, 3, 8, 2, 1, 9), levels=yn)
chest.cpt <- compileCPT(a, t.a, s, l.s, b.s, e.lt, x.e, d.be)
```

```
chest.bn <- grain(chest.cpt)

## These two forms are equivalent
bn1 <- setFinding(chest.bn, nodes=c("chest", "xray"), states=c("yes", "yes"))
bn2 <- setFinding(chest.bn, flist=list(c("chest", "yes"), c("xray", "yes")))

getFinding(bn1)
getFinding(bn2)

pFinding(bn1)
pFinding(bn2)

bn1 <- retractFinding(bn1, nodes="asia")
bn2 <- retractFinding(bn2, nodes="asia")

getFinding(bn1)
getFinding(bn2)

pFinding(bn1)
pFinding(bn2)
```

---

generics

*gRain generics*

---

### **Description**

Generic functions etc for the gRain package

### **Usage**

```
nodeNames(object)

## S3 method for class 'grain'
nodeNames(object)

nodeStates(object, nodes = nodeNames(object))

## S3 method for class 'grain'
nodeStates(object, nodes = nodeNames(object))

universe(object, ...)

## S3 method for class 'grain'
universe(object, ...)

isCompiled(object)
```

```

isPropagated(object)

isCompiled(object) <- value

isPropagated(object) <- value

## S3 method for class 'cpt_spec'
vpar(object, ...)

## S3 method for class 'cpt_grain'
vpar(object, ...)

## S3 method for class 'grain'
rip(object, ...)

## S3 method for class 'grainEvidence_'
varNames(x)

```

### Arguments

nodes	Some nodes of the object.
...	Additional arguments; currently not used.
value	Value to be set for slot in object.
x, object	A relevant object.

---

get_superset_list	<i>Get superset for each element in a list</i>
-------------------	--

---

### Description

For each element (vector)  $x$  in  $x\_set$ , find the first element (vector)  $y$  in  $y\_set$  such that  $x$  is contained in  $y$

### Usage

```
get_superset_list(x_set, y_set, warn = FALSE)
```

### Arguments

x_set	Vector or list of vectors.
y_set	Vector or list of vectors.
warn	Should a warning be made if an element is found.

**Examples**

```

x_set <- list(c("a", "b"), "e", c("b", "a"))
y_set <- list(c("f","u", "e"), c("a", "b", "c", "a"), c("b", "c", "a"))
get_superset_list(x_set, y_set)
get_superset_list(letters[1:4], y_set)
get_superset_list(letters[1:4], letters[1:10])
get_superset_list(x_set, letters[1:10])
x_set <- list(c("a", "b"), "e", c("b", "a"), "o")
y_set <- list(c("f","u", "e"), c("a", "b", "c", "a"), c("b", "c", "a"))
get_superset_list(x_set, y_set, warn=TRUE)
get_superset_list(x_set, y_set, warn=FALSE)

```

grain-main

*Graphical Independence Network***Description**

Creating grain objects (graphical independence network).

**Usage**

```

grain(x, ...)

## S3 method for class 'cpt_spec'
grain(x, control = list(), smooth = 0, compile = TRUE, details = 0, ...)

## S3 method for class 'CPTspec'
grain(x, control = list(), smooth = 0, compile = TRUE, details = 0, ...)

## S3 method for class 'pot_spec'
grain(x, control = list(), smooth = 0, compile = TRUE, details = 0, ...)

## S3 method for class 'igraph'
grain(
  x,
  control = list(),
  smooth = 0,
  compile = TRUE,
  details = 0,
  data = NULL,
  ...
)

## S3 method for class 'dModel'
grain(
  x,

```

```

control = list(),
smooth = 0,
compile = TRUE,
details = 0,
data = NULL,
...
)

```

### Arguments

x	An argument to build an independence network from. Typically a list of conditional probability tables, a DAG or an undirected graph. In the two latter cases, data must also be provided.
...	Additional arguments, currently not used.
control	A list defining controls, see 'details' below.
smooth	A (usually small) number to add to the counts of a table if the grain is built from a graph plus a dataset.
compile	Should network be compiled.
details	Debugging information.
data	An optional data set (currently must be an array/table)

### Details

If 'smooth' is non-zero then entries of 'values' which a zero are replaced by the value of 'smooth' - BEFORE any normalization takes place.

### Value

An object of class "grain"

### Note

A change from earlier versions of this package is that grain objects are now compiled upon creation.

### Author(s)

Søren Højsgaard, <sorenh@math.aau.dk>

### References

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

### See Also

[cptable](#), [compile.grain](#), [propagate.grain](#), [setFinding](#), [setEvidence](#), [getFinding](#), [pFinding](#), [retractFinding](#), [extractCPT](#), [extractPOT](#), [compileCPT](#), [compilePOT](#)

**Examples**

```

## Create network from conditional probability tables CPTs:

yn <- c("yes", "no")
a <- cpt(~asia, values=c(1,99), levels=yn)
t.a <- cpt(~tub+asia, values=c(5,95,1,99), levels=yn)
s <- cpt(~smoke, values=c(5,5), levels=yn)
l.s <- cpt(~lung+smoke, values=c(1,9,1,99), levels=yn)
b.s <- cpt(~bronc+smoke, values=c(6,4,3,7), levels=yn)
e.lt <- cpt(~either+lung+tub, values=c(1,0,1,0,1,0,0,1), levels=yn)
x.e <- cpt(~xray+either, values=c(98,2,5,95), levels=yn)
d.be <- cpt(~dysp+bronc+either, values=c(9,1,7,3,8,2,1,9), levels=yn)
cpt_list <- list(a, t.a, s, l.s, b.s, e.lt, x.e, d.be)
chest_cpt <- compileCPT(cpt_list)
## Alternative: chest_cpt <- compileCPT(a, t.a, s, l.s, b.s, e.lt, x.e, d.be)

chest_bn <- grain(chest_cpt)

## Create network from data and graph specification.

data(lizard, package="gRbase")

## From a DAG: height <- species -> diam
daG <- dag(~species + height:species + diam:species, result="igraph")

## From an undirected graph UG : [height:species][diam:species]
uG <- ug(~height:species + diam:species, result="igraph")

liz.ug <- grain(uG, data=lizard)
liz.dag <- grain(daG, data=lizard)

```

---

grain-simulate

*Simulate from an independence network*


---

**Description**

Simulate data from an independence network.

**Usage**

```

## S3 method for class 'grain'
simulate(object, nsim = 1, seed = NULL, ...)

```

**Arguments**

object	An independence network.
nsim	Number of cases to simulate.
seed	An optional integer controlling the random number generation.
...	Not used.



**Value**

A data frame

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**Examples**

```
tf <- system.file("huginex", "chest_clinic.net", package = "gRain")

chest <- loadHuginNet(tf, details=1)
simulate(chest, n=10)

chest2 <- setFinding(chest, c("VisitToAsia", "Dyspnoea"),
                    c("yes", "yes"))
simulate(chest2, n=10)
```

---

grain\_compile

*Compile a graphical independence network (a Bayesian network)*

---

**Description**

Compiles a Bayesian network. This means creating a junction tree and establishing clique potentials.

**Usage**

```
## S3 method for class 'grain'
compile(
  object,
  propagate = FALSE,
  tug = NULL,
  root = NULL,
  control = object$control,
  details = 0,
  ...
)
```

**Arguments**

object	A grain object.
propagate	If TRUE the network is also propagated meaning that the cliques of the junction tree are calibrated to each other.
tug	A triangulated undirected graph.
root	A set of variables which must be in the root of the junction tree
control	Controlling the compilation process.
details	For debugging info. Do not use.
...	Currently not used.

**Value**

A compiled Bayesian network; an object of class `grain`.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[grain](#), [propagate](#), [propagate.grain](#), [triangulate](#), [rip](#), [junctionTree](#)

---

`grain_evidence`      *Set, update and remove evidence.*

---

**Description**

Set, update and remove evidence.

**Usage**

```
setEvidence(
  object,
  nodes = NULL,
  states = NULL,
  evidence = NULL,
  propagate = TRUE,
  details = 0
)
```

```

retractEvidence(object, nodes = NULL, propagate = TRUE)

absorbEvidence(object, propagate = TRUE)

pEvidence(object, evidence = NULL)

getEvidence(object, short = TRUE)

evidence(object, short = TRUE)

## S3 method for class 'grain'
evidence(object, short = TRUE)

evidence(object) <- value

## S3 replacement method for class 'grain'
evidence(object) <- value

```

### Arguments

object	A "grain" object
nodes	A vector of nodes; those nodes for which the (conditional) distribution is requested.
states	A vector of states (of the nodes given by 'nodes')
evidence	An alternative way of specifying findings (evidence), see examples below.
propagate	Should the network be propagated?
details	Debugging information
short	If TRUE a dataframe with a summary is returned; otherwise a list with all details.
value	The evidence in the form of a named list or an evidence-object.

### Value

A list of tables with potentials.

### Note

`setEvidence()` is an improvement of `setFinding()` (and as such `setFinding` is obsolete). Users are recommended to use `setEvidence()` in the future.

`setEvidence()` allows to specification of "hard evidence" (specific values for variables) and likelihood evidence (also known as virtual evidence) for variables.

The syntax of `setEvidence()` may change in the future.

### Author(s)

Søren Højsgaard, <sorenh@math.aau.dk>

## References

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

## See Also

[setFinding](#), [getFinding](#), [retractFinding](#), [pFinding](#)

## Examples

```
data(chest_cpt)
chest.bn <- grain(compileCPT(chest_cpt))
chest.bn <- compile(chest.bn)

## 1) These two forms are identical
setEvidence(chest.bn, c("asia", "xray"), c("yes", "yes"))
setFinding(chest.bn, c("asia", "xray"), c("yes", "yes"))

## 2) Suppose we do not know with certainty whether a patient has
## recently been to Asia. We can then introduce a new variable
## "guess.asia" with "asia" as its only parent. Suppose
##  $p(\text{guess.asia}=\text{yes}|\text{asia}=\text{yes})=0.8$  and  $p(\text{guess.asia}=\text{yes}|\text{asia}=\text{no})=0.1$ 
## If the patient is e.g. unusually tanned we may set
## guess.asia=yes and propagate.
##
## This corresponds to modifying the model by the likelihood (0.8,
## 0.1) as

setEvidence(chest.bn, c("asia", "xray"), list(c(0.8, 0.1), "yes"))

## 3) Hence, the same result as in 1) can be obtained with
setEvidence(chest.bn, c("asia", "xray"), list(c(1, 0), "yes"))

## 4) An alternative specification using evidence is
setEvidence(chest.bn, evidence=list(asia=c(1, 0), xray="yes"))
```

---

grain\_joint\_evidence *Set joint evidence in grain objects*

---

## Description

Setting and removing joint evidence in grain objects.

**Usage**

```

setJEvidence(object, evidence = NULL, propagate = TRUE, details = 0)

retractJEvidence(object, items = NULL, propagate = TRUE, details = 0)

new_jev(ev, levels)

```

**Arguments**

object	A "grain" object.
evidence	A list of evidence. Each element is a named array.
propagate	Should evidence be absorbed once entered; defaults to TRUE.
details	Amount of printing; for debugging.
items	Items in the evidence list to be removed. Here, NULL means remove everything, 0 means nothing is removed. Otherwise items is a numeric vector.
ev	A named list.
levels	A named list.

**Note**

All the joint evidence functionality should be used with great care.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**Examples**

```

data(chest_cpt)
chest.bn <- grain(compileCPT(chest_cpt))
chest.bn <- compile(chest.bn)

uni <- list(asia = c("yes", "no"), tub = c("yes", "no"),
           smoke = c("yes", "no"), lung = c("yes", "no"),
           bronc = c("yes", "no"), either = c("yes", "no"),
           xray = c("yes", "no"), dysp = c("yes", "no"))

ev <- list(tabNew("asia", levels=uni, values=c(1,0)),
          tabNew("dysp", levels=uni, values=c(1,0)),
          tabNew(c("dysp","bronc"), levels=uni, values=c(.1, .2, .9, .8)) )

chest.bn
chest.bn2 <- setJEvidence(chest.bn, evidence=ev)
chest.bn2
getEvidence(chest.bn2)

# Notice: The evidence is defined on (subsets of) cliques of the junction tree
# and therefore evidence can readily be absorbed:
getgrain(chest.bn, "rip")$cliques %>% str

```

```

# On the other hand, below is evidence which is not defined cliques
# of the junction tree and therefore evidence can not easily be
# absorbed. Hence this will fail:

## Not run:
ev.fail <- list(tabNew(c("dysp", "smoke"), levels=uni, values=c(.1, .2, .9, .8)) )
setJEvidence(chest.bn, evidence=ev.fail)

## End(Not run)

## Evidence can be removed with

retractJEvidence(chest.bn2)      ## All evidence removed.
retractJEvidence(chest.bn2, 0)   ## No evidence removed.
retractJEvidence(chest.bn2, 1:2) ## Evidence items 1 and 2 are removed.

# Setting additional joint evidence to an object where joint
# evidence already is set will cause an error. Hence this will fail:

## Not run:
ev2 <- list(smoke="yes")
setJEvidence(chest.bn2, evidence=ev2)

## End(Not run)

## Instead we can do
new.ev <- c(getEvidence(chest.bn2), list(smoke="yes"))
chest.bn
setJEvidence(chest.bn, evidence=new.ev)

## Create joint evidence object:
yn <- c("yes", "no")
db <- pararray(c("dysp", "bronc"), list(yn, yn), values=c(.1, .2, .9, .8))
db
ev <- list(asia=c(1, 0), dysp="yes", db)

jevi <- new_jev(ev, levels=uni)
jevi

chest.bn3 <- setJEvidence(chest.bn, evidence=jevi)
evidence(chest.bn3)

```

**Description**

Makes predictions (either as the most likely state or as the conditional distributions) of variables conditional on finding (evidence) on other variables in an independence network.

**Usage**

```
## S3 method for class 'grain'  
predict(  
  object,  
  response,  
  predictors = setdiff(names(newdata), response),  
  newdata,  
  type = "class",  
  ...  
)
```

**Arguments**

object	A grain object
response	A vector of response variables to make predictions on
predictors	A vector of predictor variables to make predictions from. Defaults to all variables that are not responses.
newdata	A data frame
type	If "class", the most probable class is returned; if "distribution" the conditional distribution is returned.
...	Not used

**Value**

A list with components

pred	A list with the predictions
pFinding	A vector with the probability of the finding (evidence) on which the prediction is based

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[grain](#)

**Examples**

```

data(chest_cpt)
data(chestSim500)

chest.bn <- grain(compileCPT(chest_cpt))
nd <- chestSim500[1:4]

predict(chest.bn, response="bronc", newdata=nd)
predict(chest.bn, response="bronc", newdata=nd, type="distribution")

```

---

grain_propagate	<i>Propagate a graphical independence network (a Bayesian network)</i>
-----------------	--

---

**Description**

Propagation refers to calibrating the cliques of the junction tree so that the clique potentials are consistent on their intersections; refer to the reference below for details.

**Usage**

```

## S3 method for class 'grain'
propagate(object, details = object$details, engine = "cpp", ...)

propagateLS(cq_pot_list, rip, initialize = TRUE, details = 0)

compute_p_evidence(object, details = object$details, engine = "cpp", ...)

```

**Arguments**

object	A grain object
details	For debugging info
engine	Either "R" or "cpp"; "cpp" is the default and the fastest.
...	Currently not used
cq_pot_list	List of clique potentials
rip	A rip ordering
initialize	Always true.

**Details**

The propagate method invokes propagateLS which is a pure R implementation of the Lauritzen-Spiegelhalter algorithm. The c++ based version is several times faster than the purely R based version.

**Value**

A compiled and propagated grain object.



**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[grain](#), [compile](#)

**Examples**

```
example("grain")

## Uncompiled and unpropagated network:
bn0 <- grain(chest_cpt, compile=FALSE)
bn0
## Compiled but unpropagated network:
bn1 <- compile(bn0, propagate=FALSE)
## Compiled and propagated network
bn2 <- propagate(bn1)
bn2
## Default is that networks are compiled but not propagated at creation time:
bn3 <- grain(chest_cpt)
bn3
```

---

grass

*Wet grass example*

---

**Description**

Conditional probability tables for the wet grass example.

**Usage**

```
data(grass_cpt)
```

**Format**

An object of class `list` of length 3.

## Examples

```
## 'data' generated with the following code fragment
yn <- c("yes", "no")
p.R <- cptable(~R, values=c(.2, .8), levels=yn)
p.S_R <- cptable(~S:R, values=c(.01, .99, .4, .6), levels=yn)
p.G_SR <- cptable(~G:S:R, values=c(.99, .01, .8, .2, .9, .1, 0, 1), levels=yn)

grain(compileCPT(p.R, p.S_R, p.G_SR))

# 'data' generated from
# grass_cpt <- list(p.R, p.S_R, p.G_SR)

data(grass_cpt)
```

---

load-save-hugin

*Load and save Hugin net files*

---

## Description

These functions can load a net file saved in the 'Hugin format' into R and save a network in R as a file in the 'Hugin format'.

## Usage

```
loadHuginNet(file, description = NULL, details = 0)
```

```
saveHuginNet(gin, file, details = 0)
```

## Arguments

file	Name of HUGIN net file. Convenient to give the file the extension '.net'
description	A text describing the network, defaults to file
details	Debugging information.
gin	An independence network

## Value

An object of class grain.

## Note

- In Hugin, it is possible to specify the potential of a node as a functional relation between other nodes. In a .net file, such a specification will appear as 'function' rather than as 'node'. Such a specification is not recognized by loadHuginNet.
- It is recommended to avoid the text node as part of the name of a node.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. *Journal of Statistical Software*, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[grain](#)

**Examples**

```
## Load HUGIN net file
tf <- system.file("huginex", "chest_clinic.net", package = "gRain")
chest <- loadHuginNet(tf, details=1)
chest

## Save a copy
td <- tempdir()
saveHuginNet(chest, paste(td, "/chest.net", sep=''))

## Load the copy
chest2 <- loadHuginNet(paste(td, "/chest.net", sep=''))

tf <- system.file("huginex", "golf.net", package = "gRain")
golf <- loadHuginNet(tf, details=1)

saveHuginNet(golf, paste(td, "/golf.net", sep=''))
golf2 <- loadHuginNet(paste(td, "/golf.net", sep=''))
```

---

logical

*Conditional probability tables based on logical dependencies*

---

**Description**

Generate conditional probability tables based on the logical expressions AND and OR.

**Usage**

```
booltab(vpa, levels = c(TRUE, FALSE), op = '&')
```

```
andtab(vpa, levels = c(TRUE, FALSE))
```

```
ortab(vpa, levels = c(TRUE, FALSE))
```

```
andtable(vpa, levels = c(TRUE, FALSE))
```

```
ortable(vpa, levels = c(TRUE, FALSE))
```

### Arguments

vpa	Node and two parents; as a formula or a character vector.
levels	The levels (or rather labels) of v, see 'examples' below.
op	A logical operator.

### Details

Regarding the form of the argument vpa: To specify  $P(a|b, c)$  one may write  $\sim a|b+c$  or  $\sim a+b+c$  or  $\sim a|b:c$  or  $\sim a:b:c$  or  $c("a", "b", "c")$ . Internally, the last form is used. Notice that the + and : operator are used as separators only. The order of the variables is important so + and : DO NOT commute.

### Value

An array.

### Note

andtable and ortable are aliases for andtab and ortab and are kept for backward compatibility.

### Author(s)

Søren Højsgaard, <sorenh@math.aau.dk>

### References

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

### See Also

[cptable](#)

### Examples

```
## Logical OR:

## A variable v is TRUE if either of its parents pa1 and pa2 are TRUE:
ortab( c("v", "pa1", "pa2") ) %>% ftable(row.vars="v")
## TRUE and FALSE can be recoded to e.g. yes and no:
ortab( c("v", "pa1", "pa2"), levels=c("yes", "no") ) %>% ftable(row.vars="v")

## Logical AND:
```

```
## Same story here:
andtab(c("v", "pa1", "pa2") ) %>% ftable(row.vars="v")
andtab(c("v", "pa1", "pa2"), levels=c("yes", "no") ) %>% ftable(row.vars="v")

## Combined approach

booltab(c("v", "pa1", "pa2"), op=`&`) %>% ftable(row.vars="v") ## AND
booltab(c("v", "pa1", "pa2"), op=`|`) %>% ftable(row.vars="v") ## OR

booltab(~v + pa1 + pa2, op=`&`) %>% ftable(row.vars="v") ## AND
booltab(~v + pa1 + pa2, op=`|`) %>% ftable(row.vars="v") ## OR
```

---

mendel

*Mendelian segregation*


---

## Description

Generate conditional probability table for mendelian segregation.

## Usage

```
mendel(allele, names = c("child", "father", "mother"))
```

## Arguments

allele	A character vector.
names	Names of columns in dataframe.

## Note

No error checking at all on the input.

## Examples

```
## Inheritance of the alleles "y" and "g"

men <- mendel(c("y", "g"), names=c("ch", "fa", "mo"))
men
```

---

 querygrain

*Query a network*


---

### Description

Query an independence network, i.e. obtain the conditional distribution of a set of variables - possibly (and typically) given finding (evidence) on other variables.

### Usage

```
querygrain(
  object,
  nodes = nodeNames(object),
  type = "marginal",
  evidence = NULL,
  exclude = TRUE,
  normalize = TRUE,
  simplify = FALSE,
  result = "array",
  details = 0
)
```

### Arguments

object	A grain object.
nodes	A vector of nodes; those nodes for which the (conditional) distribution is requested.
type	Valid choices are "marginal" which gives the marginal distribution for each node in nodes; "joint" which gives the joint distribution for nodes and "conditional" which gives the conditional distribution for the first variable in nodes given the other variables in nodes.
evidence	An alternative way of specifying findings (evidence), see examples below.
exclude	If TRUE then nodes on which evidence is given will be excluded from nodes (see above).
normalize	Should the results be normalized to sum to one.
simplify	Should the result be simplified (to a dataframe) if possible.
result	If "data.frame" the result is returned as a data frame (or possibly as a list of dataframes).
details	Debugging information

### Value

A list of tables with potentials.

**Note**

setEvidence() is an improvement of setFinding() (and as such setFinding is obsolete). Users are recommended to use setEvidence() in the future.

setEvidence() allows to specification of "hard evidence" (specific values for variables) and likelihood evidence (also known as virtual evidence) for variables.

The syntax of setEvidence() may change in the future.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[setEvidence](#), [getEvidence](#), [retractEvidence](#), [pEvidence](#)

**Examples**

```
testfile <- system.file("huginex", "chest_clinic.net", package = "gRain")
chest <- loadHuginNet(testfile, details=0)
qb <- querygrain(chest)
qb

lapply(qb, as.numeric) # Safe
sapply(qb, as.numeric) # Risky
```

---

repeatPattern

*Create repeated patterns in Bayesian networks*

---

**Description**

Repeated patterns is a useful model specification short cut for Bayesian networks

**Usage**

```
repeatPattern(plist, instances, unlist = TRUE, data = NULL)
```

**Arguments**

plist	A list of conditional probability tables. The variable names must have the form name[i] and the i will be substituted by the values given in instances below. See also the data argument.
instances	A vector of distinct integers
unlist	If FALSE the result is a list in which each element is a copy of plist in which name[i] are substituted. If TRUE the result is the result of applying unlist().
data	A two column matrix. The first column is the index / name of a node; the second column is the index / name of the node's parent.

**Author(s)**

Søren Højsgaard, <sorenh@math.aau.dk>

**References**

Søren Højsgaard (2012). Graphical Independence Networks with the gRain Package for R. Journal of Statistical Software, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

**See Also**

[grain](#), [compileCPT](#)

**Examples**

```
## Example: Markov chain
yn <- c("yes", "no")

x.0 <- cptable(~x0, values=c(1, 9), levels=yn)          ## p(x0)
x.x <- cptable(~x[i]|x[i-1], values=c(1, 99, 2, 98), levels=yn) ## p(x[i]|x[i-1])

x.0 <- cpt(~x0, values=c(1, 9), levels=yn)            ## p(x0)
x.x <- cpt(~x[i]|x[i-1], values=c(1, 99, 2, 98), levels=yn) ## p(x[i]|x[i-1])

pat <- list(x.x) ## Pattern to be repeated

n <- 5
rep.pat <- repeatPattern(pat, instances=1:n)
mc <- compileCPT(c(list(x.0), rep.pat)) |> grain()
if (interactive()) iplot(mc)

## Example: Hidden markov model: The x[i]'s are unobserved, the
## y[i]'s can be observed.

x.0 <- cptable(~x0, values=c(1, 9), levels=yn)          ## p(x0)
x.x <- cptable(~x[i]|x[i-1], values=c(1, 99, 2, 98), levels=yn) ## p(x[i]|x[i-1])
y.x <- cptable(~y[i]|x[i], values=c(10, 90, 20, 80), levels=yn) ## p(y[i]|x[i])

x.0 <- cpt(~x0, values=c(1, 9), levels=yn)            ## p(x0)
```



```

x.x <- cpt(~x[i]|x[i-1], values=c(1, 99, 2, 98), levels=yn) ## p(x[i]|x[i-1])
y.x <- cpt(~y[i]|x[i], values=c(10, 90, 20, 80), levels=yn) ## p(y[i]|x[i])

pat <- list(x.x, y.x) ## Pattern to be repeated

rep.pat <- repeatPattern(pat, instances=1:n)
hmm <- compileCPT(c(list(x.0), rep.pat)) |> grain()
hmm

if (interactive()) iplot(hmm)

## Data-driven variable names
dep <- data.frame(i=c(1, 2, 3, 4, 5, 6, 7, 8),
                 p=c(0, 1, 2, 2, 3, 3, 4, 4))

z0 <- cptable(~z0, values=c(0.5, 0.5), levels=yn)
za <- cptable(~z[i] | z[data[i, "p"]], values=c(0.5, 0.5), levels=yn)
zb <- repeatPattern(list(za), instances=1:nrow(dep), data=dep)
tree <- compileCPT(c(list(z0), zb)) |> grain()

x0 <- cpt(~x0, values=c(0.5, 0.5), levels=yn)
xa <- cpt(~x[i] | x[data[i, "p"]], values=c(0.5, 0.5), levels=yn)
xb <- repeatPattern(list(xa), instances=1:nrow(dep), data=dep)
tree <- compileCPT(c(list(x0), xb)) |> grain()
tree

if (interactive()) iplot(tree)

```

---

replace-cpt

*Replace CPTs in Bayesian network*


---

## Description

Replace CPTs of Bayesian network.

## Usage

```

replaceCPT(object, value)

## S3 method for class 'cpt_grain'
replaceCPT(object, value)

```

## Arguments

object            A grain object.  
value             A named list, see examples below.

## Details

When a Bayesian network (BN) is constructed from a list of conditional probability tables (CPTs) (e.g. using the function `grain()`), various actions are taken:

1. It is checked that the list of CPTs define a directed acyclic graph (DAG).
2. The DAG is moralized and triangulated.
3. A list of clique potentials (one for each clique in the triangulated graph) is created from the list of CPTs.
4. The clique potentials are, by default, calibrated to each other so that the potentials contain marginal distributions.

The function described here bypass the first two steps which can provide an important gain in speed compared to constructing a new BN with a new set of CPTs with the same DAG.

## Author(s)

Søren Højsgaard, <sorenh@math.aau.dk>

## References

Søren Højsgaard (2012). Graphical Independence Networks with the `gRain` Package for R. *Journal of Statistical Software*, 46(10), 1-26. <https://www.jstatsoft.org/v46/i10/>.

## See Also

[grain](#), [propagate](#), [triangulate](#), [rip](#), [junctionTree](#)

## Examples

```
## See the wet grass example at
## https://en.wikipedia.org/wiki/Bayesian_network

yn <- c("yes", "no")
p.R <- cptable(~R, values=c(.2, .8), levels=yn)
p.S_R <- cptable(~S:R, values=c(.01, .99, .4, .6), levels=yn)
p.G_SR <- cptable(~G:S:R, values=c(.99, .01, .8, .2, .9, .1, 0, 1), levels=yn)

wet.bn <- compileCPT(p.R, p.S_R, p.G_SR) |> grain()
getgrain(wet.bn, "cpt")[c("R", "S")]

# Update some CPTs
wet.bn <- replaceCPT(wet.bn, list(R=c(.3, .7), S=c(.1, .9, .7, .3)))
getgrain(wet.bn, "cpt")[c("R", "S")]
```

---

simplify_query	<i>Simplify output query to a Bayesian network</i>
----------------	--

---

**Description**

Simplify output query to a Bayesian network to a dataframe provided that each node has the same levels.

**Usage**

```
simplify_query(b)
```

**Arguments**

b                    Result from running querygrain.

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