

# User manual for the psych package Version 1.0.27

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## Description

The psych package has been developed at Northwestern University to include functions most useful for personality and psychological research. Some of the functions (e.g., `describe` and `pairs.panels`) are useful for basic descriptive analyses.

Psychometric applications include routines for Very Simple Structure (`VSS`), Item Cluster Analysis (`ICLUST`) and principal axes factor analysis (`factor.pa`), as well as functions to do Schmid Leiman transformations (`schmid`) to transform a hierarchical factor structure into a bifactor solution and to graph both structures (`omega.graph`) and to calculate reliability coefficients alpha (`score.items`), beta (`ICLUST`) and McDonald's omega (`omega` and `omega.graph`).

Additional functions make for more convenient descriptions of item characteristics. Functions under development include 1 and 2 parameter Item Response measures.

A number of procedures have been developed as part of the Synthetic Aperture Personality Assessment (SAPA) project. These routines facilitate forming and analyzing composite scales equivalent to using the raw data but doing so by adding within and between cluster/scale item correlations. These functions include extracting clusters from factor loading matrices (`factor2cluster`), synthetically forming clusters from correlation matrices (`cluster.cor`), and finding multiple correlation from correlation matrices (`mat.regress`).

The most recent development version of the package is always available for download as a source file from the repository at <http://personality-project.org/r/src/contrib/>

## Details

The psych package is a combination of multiple source files maintained at the <http://personality-project.org/r> repository: “useful.r”, `VSS.r`, `ICLUST.r`, `omega.r`, etc. “useful.r” is a set of routines for easy data entry (`read.clipboard`), simple descriptive statistics (`describe`), and splom plots combined with correlations (`pairs.panels`, adapted from the help files of `pairs`).

The `VSS` routines allow for testing the number of factors (`VSS`), showing plots (`VSS.plot`) of goodness of fit, and basic routines for estimating the number of factors/components to extract by examining the scree plot (`VSS.scree`) or comparing with the scree of an equivalent matrix of random numbers (`VSS.parallel`).

In addition, there are routines for hierarchical factor analysis using Schmid Leiman transformations (`omega`, `omega.graph`) as well as Item Cluster analysis (`ICLUST`, `ICLUST.graph`).

The more important functions in the package are for the analysis of multivariate data, with an emphasis upon those functions useful in scale construction of item composites.

When given a set of items from a personality inventory, one goal is to combine these into higher level item composites. This leads to several questions:

1) What are the basic properties of the data? `describe` reports basic summary statistics (mean, sd, median, mad, range, minimum, maximum, skew, kurtosis, standard error) for

vectors, columns of matrices, or data.frames. [describe.by](#) provides descriptive statistics, organized by a grouping variable. [pairs.panels](#) shows scatter plot matrices (SPLOMs) as well as histograms and the Pearson correlation for scales or items.

2) What is the most appropriate number of item composites to form? After finding either standard Pearson correlations, or finding tetrachoric or polychoric correlations using a wrapper ([poly.mat](#)) for John Fox's `hetcor` function, the dimensionality of the correlation matrix may be examined. The number of factors/components problem is a standard question of factor analysis, cluster analysis, or principal components analysis. Unfortunately, there is no agreed upon answer. The Very Simple Structure ([VSS](#)) set of procedures has been proposed as an answer to the question of the optimal number of factors. Other procedures ([VSS.scree](#), [VSS.parallel](#), and [fa.parallel](#)) also address this question.

3) What are the best composites to form? Although this may be answered using principal components ([principal](#)) or factor analysis ([factor.pa](#)) and to show the results graphically ([fa.graph](#)), it is sometimes more useful to address this question using cluster analytic techniques. (Better yet is to use maximum likelihood factor analysis using [factanal](#) from the stats package.) Previous versions of [ICLUST](#) (e.g., Revelle, 1979) have been shown to be particularly successful at doing this. Graphical output from [ICLUST.graph](#) uses the Graphviz dot language and allows one to write files suitable for Graphviz. (As of May, 2007, Rgraphviz now works on the Intel-Mac and two graphic functions take advantage of this. Graphviz still produces cleaner output.)

4) How well does a particular item composite reflect a single construct? This is a question of reliability and general factor saturation. Multiple solutions for this problem result in (Cronbach's) alpha ([score.items](#)), (Revelle's) Beta ([ICLUST](#)), and (McDonald's) [omega](#). Functions to estimate all three of these are included in `psych`.

5) For some applications, data matrices are synthetically combined from sampling different items for different people. So called Synthetic Aperture Personality Assessment (SAPA) techniques allow the formation of large correlation or covariance matrices even though no one person has taken all of the items. To analyze such data sets, it is easy to form item composites based upon the covariance matrix of the items, rather than original data set. These matrices may then be analyzed using a number of functions (e.g., [cluster.cor](#), [factor.pa](#), [ICLUST](#), [principal](#), [mat.regress](#), and [factor2cluster](#)).

6) More typically, one has a raw data set to analyze. [score.items](#) will score data sets on multiple scales, reporting the scale scores, item-scale and scale-scale correlations, as well as coefficient alpha and alpha-1. Using a 'keys' matrix, scales can have overlapping or independent items.

Package:	psych
Type:	Package
Version:	1.0-27
Date:	2007-7-31
License:	GPL version 2 or newer

Index:

[psych](#) A package for personality, psychometric, and psychological research.

Useful data entry and descriptive statistics

[describe](#) Basic descriptive statistics useful for psychometrics  
[describe.by](#) Find summary statistics by groups  
[read.clipboard](#) shortcut for reading from the clipboard  
[read.clipboard.csv](#) shortcut for reading comma delimited files from clipboard

[pairs.panels](#) SPLOM and correlations for a data matrix  
[multi.hist](#) Histograms of multiple variables arranged in matrix form  
[skew](#) Calculate skew for a vector, each column of a matrix, or data.frame  
[kurtosi](#) Calculate kurtosis for a vector, each column of a matrix or dataframe

[error.crosses](#) Two way error bars  
[geometric.mean](#) Find the geometric mean of a vector or columns of a data.frame  
[harmonic.mean](#) Find the harmonic mean of a vector or columns of a data.frame

Data reduction through cluster and factor analysis

[factor.pa](#) Do a principal Axis factor analysis  
[fa.graph](#) Show the results of a factor analysis or principal components analysis graphically  
[principal](#) Do an eigen value decomposition to find the principal components of a matrix  
[fa.parallel](#) Scree test and Parallel analysis  
[ICLUST](#) Apply the ICLUST algorithm  
[ICLUST.graph](#) Graph the output from ICLUST using the dot language  
[ICLUST.rgraph](#) Graph the output from ICLUST using rgraphviz  
[poly.mat](#) Find the polychoric correlations for items (uses J. Fox's hetcor)  
[omega](#) Calculate the omega estimate of factor saturation (requires the GPArotation package)  
[omega.graph](#) Draw a hierarchical or SL orthogonalized solution  
[schmid](#) Apply the Schmid Leiman transformation to a correlation matrix

[score.items](#) Combine items into multiple scales and find alpha  
[VSS](#) Apply the Very Simple Structure criterion to determine the appropriate number of factors.  
[VSS.parallel](#) Do a parallel analysis to determine the number of factors for a random matrix  
[VSS.plot](#) Plot VSS output  
[VSS.scree](#) Show the scree plot of the factor/principal components  
[VSS.simulate](#) Generate simulated data for the factor model  
[make.hierarchical](#) Generate simulated correlation matrices with hierarchical structure

Procedures particularly useful for Synthetic Aperture Personality Assessment

[alpha.scale](#) Find coefficient alpha for a scale (see also [score.items](#))  
[correct.cor](#) Correct a correlation matrix for unreliability  
[count.pairwise](#) Count the number of complete cases when doing pair wise correlations  
[cluster.cor](#) find correlations of composite variables from larger matrix  
[cluster.loadings](#) find correlations of items with composite variables from a larger matrix  
[eigen.loadings](#) Find the loadings when doing an eigen value decomposition  
[factor.pa](#) Do a principal Axis factor analysis  
[factor2cluster](#) extract cluster definitions from factor loadings

[factor.congruence](#) Factor congruence coefficient  
[factor.fit](#) How well does a factor model fit a correlation matrix  
[factor.model](#) Reproduce a correlation matrix based upon the factor model  
[factor.residuals](#) Fit = data - model  
[factor.rotate](#) “hand rotate” factors  
[mat.regress](#) multiple regression from matrix input  
[principal](#) Do an eigen value decomposition to find the principal components of a matrix

Functions for generating simulated data sets

[circ.sim](#) Generate a two dimensional circumplex item structure  
[item.sim](#) Generate a two dimensional simple structure with particular item characteristics  
[congeneric.sim](#) Generate a one factor congeneric reliability structure  
[psycho.demo](#) Create artificial data matrices for teaching purposes

Miscellaneous functions

[fisherz](#) Apply the Fisher r to z transform  
[paired.r](#) Test for the difference of two paired correlations  
[phi2poly](#) Given a phi coefficient, what is the polychoric correlation  
[poly.mat](#) Use John Fox’s hetcor to create a matrix of correlations from a data.frame or matrix of integer values  
[polychor.matrix](#) Use John Fox’s polycor to create a matrix of correlations (not yet very useful)

Functions that are under development and not recommended for casual use

[irt.item.diff.rasch](#) IRT estimate of item difficulty with assumption that  $\theta = 0$   
[irt.person.rasch](#) Item Response Theory estimates of  $\theta$  (ability) using a Rasch like model

[test.psych](#) Run a test of the major functions on 5 different data sets. Primarily for development purposes. Although the output can be used as a demo of the various functions.

## Note

Development versions of this package are maintained at the local repository <http://personality-project.org/r> along with further documentation. Specify that you are downloading a source package.

Some functions require other packages. Specifically, `omega` and `schmid` require the `GPArotation` package, and `poly.mat`, `phi2poly` and `polychor.matrix` requires John Fox’s `polychor` package. `ICLUST.rgraph` and `fa.graph` require `Rgraphviz`. i.e.:

function	requires
<a href="#">omega</a>	<code>GPArotation</code>
<a href="#">schmid</a>	<code>GPArotation</code>
<a href="#">poly.mat</a>	<code>polychor</code>
<a href="#">phi2poly</a>	<code>polychor</code>
<a href="#">polychor.matrix</a>	<code>polychor</code>
<a href="#">ICLUST.rgraph</a>	<code>Rgraphviz</code>

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**References**

A general guide to personality theory and research may be found at the personality-project <http://personality-project.org>. See also the short guide to R at <http://personality-project.org/r>. In addition, see An Introduction to Psychometric Theory with applications in R (Revelle, in preparation) at <http://personality-project.org/r/book/>

**Examples**

```
#See the separate man pages
```

---

<code>alpha.scale</code>	<i>Cronbach alpha for a scale</i>
--------------------------	-----------------------------------

---

**Description**

Find Cronbach's coefficient alpha given a scale and a data.frame of the items in the scale. For X, a total score composed of items in the data.frame Y, find Cronbach's alpha.

**Usage**

```
alpha.scale(x, y)
```

**Arguments**

x	A scale made by summing items
y	A data frame of items

## Details

Alpha is one of several estimates of the internal consistency reliability of a test. Perhaps because it is so easy to calculate, it is without doubt the most frequently reported measure of internal consistency reliability. Alpha is the mean of all possible split half reliabilities (corrected for test length). For a unifactorial test, it is a reasonable estimate of the first factor saturation, although if the test has any microstructure (i.e., if it is “lumpy”) coefficients beta and omega are more appropriate estimates of the general factor saturation.

Alpha is a positive function of the number of items in a test as well as the average inter-correlation of the items in the test. When calculated from the item variances and total test variance, as is done here, alpha is sensitive to differences in the item variances. Alternative functions `score.items` and `cluster.cor` will also score multiple scales and report more useful statistics. “Standardized” alpha is calculated from the inter-item correlations and will differ from raw alpha. Standardized alpha can be found by using `cluster.cor`.

## Value

Coefficient alpha

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## References

<http://personality-project.org/revelle/syllabi/405.syllabus.html>

## See Also

`score.items`, `cluster.cor`, `ICLUST`, `omega`, `describe.pairs.panels`, alpha in psychometrics

## Examples

```
y <- attitude      #from the datasets package
x <- rowSums(y)     #find the sum of the seven attitudes
alpha.scale(x,y)
#[1] 0.8431428
#compare with standardized alpha:
st.alpha <- cluster.cor(rep(1,7),cor(attitude),digits=4)
st.alpha
#compare with score.items
si <- score.items(rep(1,7), attitude,digits=3)
si$alpha
```



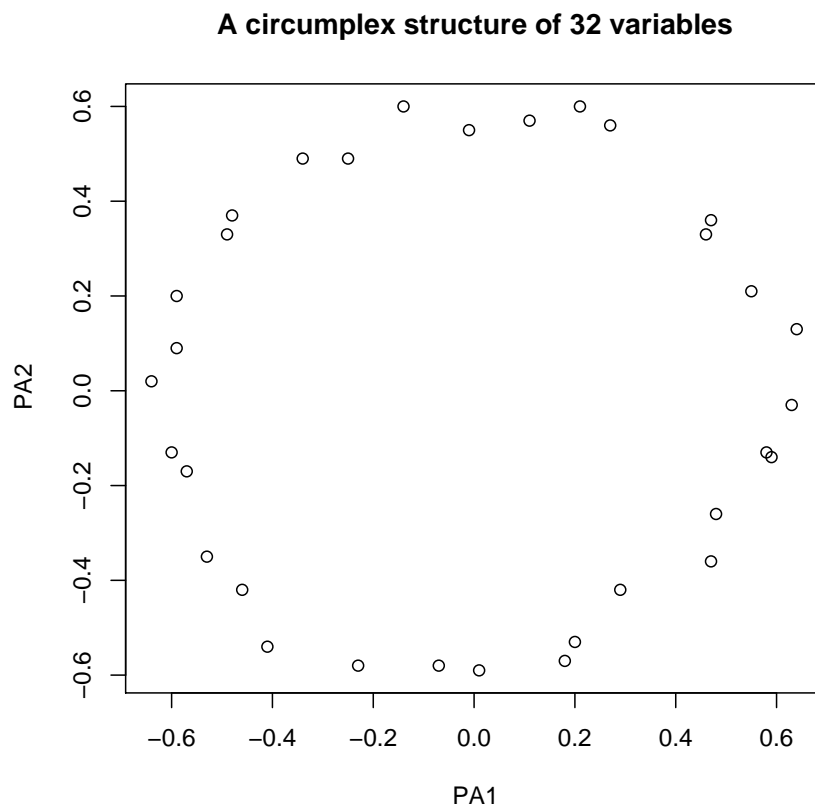


Figure 1: A circumplex structure is typically found for emotional words or for interpersonal behaviors. Tests of circumplex structure can be validated by analyzing simulated structures. Compare this structure to a simple structure [Figure 7](#)

## Description

Rotations of factor analysis and principal components analysis solutions typically try to represent correlation matrices as simple structured. An alternative structure, appealing to some, is a circumplex structure where the variables are uniformly spaced on the perimeter of a circle in a two dimensional space. Generating these data is straightforward, and is useful for exploring alternative solutions to affect and personality structure.

## Usage

```
circ.simulation(samplesize=c(100,200,400,800), numberofvariables=c(16,32,48,72))
```

## Arguments

`samplesize`        a vector of sample sizes to simulate  
`numberofvariables`  
                      vector of the number of variables to simulate

## Details

“A common model for representing psychological data is simple structure (Thurstone, 1947). According to one common interpretation, data are simple structured when items or scales have non-zero factor loadings on one and only one factor (Revelle & Rocklin, 1979). Despite the commonplace application of simple structure, some psychological models are defined by a lack of simple structure. Circumplexes (Guttman, 1954) are one kind of model in which simple structure is lacking.

“A number of elementary requirements can be teased out of the idea of circumplex structure. First, circumplex structure implies minimally that variables are interrelated; random noise does not a circumplex make. Second, circumplex structure implies that the domain in question is optimally represented by two and only two dimensions. Third, circumplex structure implies that variables do not group or clump along the two axes, as in simple structure, but rather that there are always interstitial variables between any orthogonal pair of axes (Saucier, 1992). In the ideal case, this quality will be reflected in equal spacing of variables along the circumference of the circle (Gurtman, 1994; Wiggins, Steiger, & Gaelick, 1981). Fourth, circumplex structure implies that variables have a constant radius from the center of the circle, which implies that all variables have equal communality on the two circumplex dimensions (Fisher, 1997; Gurtman, 1994). Fifth, circumplex structure implies that all rotations are equally good representations of the domain (Conte & Plutchik, 1981; Larsen & Diener, 1992).” (Acton and Revelle, 2004)

Acton and Revelle reviewed the effectiveness of 10 tests of circumplex structure and found that four did a particularly good job of discriminating circumplex structure from simple structure, or circumplexes from ellipsoidal structures. Unfortunately, their work was done in Pascal and is not easily available. Here we release R code to do the four most useful tests:

The Gap test of equal spacing

Fisher's test of equality of axes

A test of indifference to Rotation

A test of equal Variance of squared factor loadings across arbitrary rotations.

Included in this set of functions are simple procedure to generate circumplex structured or simple structured data, the four test statistics, and a simple simulation showing the effectiveness of the four procedures.

### Value

A data.frame with simulation results for circumplex, ellipsoid, and simple structure data sets for each of the four tests.

### Note

The simulations default values are for sample sizes of 100,200, 400, and 800 cases, with 16, 32, 48 and 72 items.

### Author(s)

William Revelle

### References

Acton, G. S. and Revelle, W. (2004) Evaluation of Ten Psychometric Criteria for Circumplex Structure. Methods of Psychological Research Online, Vol. 9, No. 1 [http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110\\_10.pdf](http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110_10.pdf)

### See Also

See Also as [circ.tests](#), [circ.sim](#)

### Examples

```
demo <- circ.simulation()
boxplot(demo[3:14])
title("4 tests of Circumplex Structure",sub="Circumplex, Ellipsoid, Simple Structure")
```

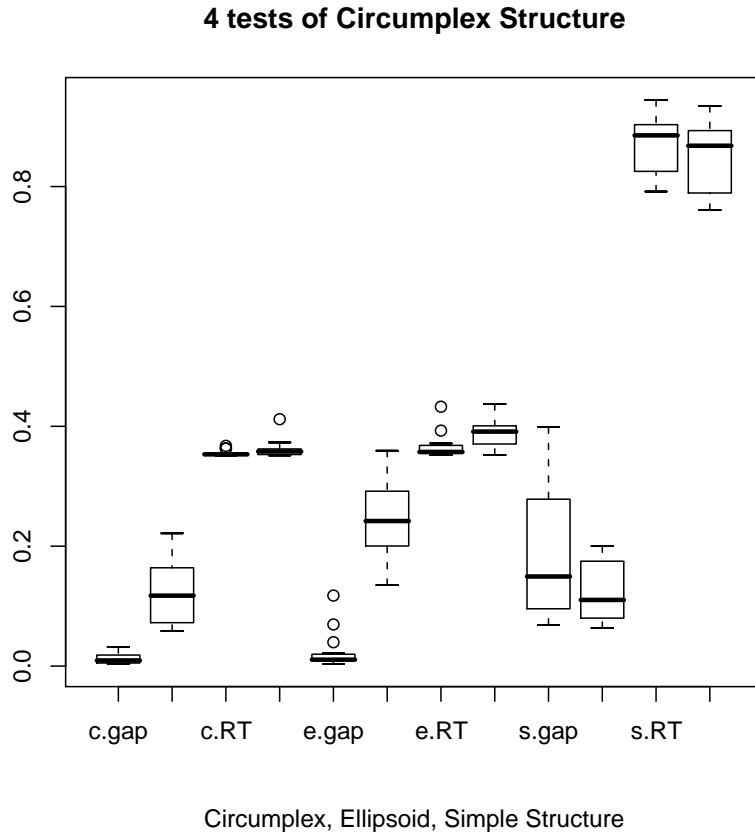


Figure 2: Tests for circumplex structure can be shown to discriminate circumplex and simple structure. Here we compare four tests of circumplex structure, the Gap Test, the Fisher Test, the Rotation Test, and the Variance Test of Circumplex structure. The data sets vary on whether or not they have a circumplex or simple structure. See Acton and Revelle (2004) for details.

## Description

Rotations of factor analysis and principal components analysis solutions typically try to represent correlation matrices as simple structured. An alternative structure, appealing to some, is a circumplex structure where the variables are uniformly spaced on the perimeter of a circle in a two dimensional space. Generating these data is straightforward, and is useful for exploring alternative solutions to affect and personality structure.

## Usage

```
circ.tests(loads, loading = TRUE, sorting = TRUE)
```

## Arguments

<code>loads</code>	A matrix of loadings <code>loads</code> here
<code>loading</code>	Are these loadings or a correlation matrix <code>loading</code>
<code>sorting</code>	Should the variables be sorted <code>sorting</code>

## Details

“A common model for representing psychological data is simple structure (Thurstone, 1947). According to one common interpretation, data are simple structured when items or scales have non-zero factor loadings on one and only one factor (Revelle & Rocklin, 1979). Despite the commonplace application of simple structure, some psychological models are defined by a lack of simple structure. Circumplexes (Guttman, 1954) are one kind of model in which simple structure is lacking.

“A number of elementary requirements can be teased out of the idea of circumplex structure. First, circumplex structure implies minimally that variables are interrelated; random noise does not a circumplex make. Second, circumplex structure implies that the domain in question is optimally represented by two and only two dimensions. Third, circumplex structure implies that variables do not group or clump along the two axes, as in simple structure, but rather that there are always interstitial variables between any orthogonal pair of axes (Saucier, 1992). In the ideal case, this quality will be reflected in equal spacing of variables along the circumference of the circle (Gurtman, 1994; Wiggins, Steiger, & Gaelick, 1981). Fourth, circumplex structure implies that variables have a constant radius from the center of the circle, which implies that all variables have equal communality on the two circumplex dimensions (Fisher, 1997; Gurtman, 1994). Fifth, circumplex structure implies that all rotations are equally good representations of the domain (Conte & Plutchik, 1981; Larsen & Diener, 1992). (Acton and Revelle, 2004)

Acton and Revelle reviewed the effectiveness of 10 tests of circumplex structure and found that four did a particularly good job of discriminating circumplex structure from simple structure, or circumplexes from ellipsoidal structures. Unfortunately, their work was done

in Pascal and is not easily available. Here we release R code to do the four most useful tests:

- 1 The Gap test of equal spacing
- 2 Fisher's test of equality of axes
- 3 A test of indifference to Rotation
- 4 A test of equal Variance of squared factor loadings across arbitrary rotations.

## Value

A list of four items is returned. These are the gap, fisher, rotation and variance test results.

<b>gaps</b>	gap.test
<b>fisher</b>	fisher.test
<b>RT</b>	rotation.test
<b>VT</b>	variance.test

## Note

Of the 10 criterion discussed in Acton and Revelle (2004), these tests operationalize the four most useful.

## Author(s)

William Revelle

## References

Acton, G. S. and Revelle, W. (2004) Evaluation of Ten Psychometric Criteria for Circumplex Structure. Methods of Psychological Research Online, Vol. 9, No. 1 [http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110\\_10.pdf](http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110_10.pdf)

## See Also

[circ.simulation](#), [circ.sim](#)

## Examples

```
circ.data <- circ.sim(24,500)
circ.fa <- factor.pa(circ.data,2)
#plot(circ.fa$loadings)
ct <- circ.tests(circ.fa)
#compare with non-circumplex data
simp.data <- item.sim(24,500)
simp.fa <- factor.pa(simp.data,2)
#plot(simp.fa$loadings)
st <- circ.tests(simp.fa)
print(rbind(ct,st),digits=2)
```

---

`cluster.cor`*Find correlations of composite variables from a larger matrix*

---

## Description

Given a  $n \times c$  cluster definition matrix of -1s, 0s, and 1s (the keys) , and a  $n \times n$  correlation matrix, find the correlations of the composite clusters. The keys matrix can be entered by hand, copied from the clipboard ([read.clipboard](#)), or taken as output from the [factor2cluster](#) function.

## Usage

```
cluster.cor(keys, r.mat, correct = TRUE,digits=2)
```

## Arguments

<code>keys</code>	A matrix of cluster keys
<code>r.mat</code>	A correlation matrix
<code>correct</code>	TRUE shows both raw and corrected for attenuation correlations
<code>digits</code>	round off answer to digits

## Details

This is one of the functions used in the SAPA procedures to form synthetic correlation matrices. Given any correlation matrix of items, it is easy to find the correlation matrix of scales made up of those items. This can also be done from the original data matrix using [score.items](#).

A typical use in the SAPA project is to form item composites by clustering or factoring (see [ICLUST](#), [principal](#)), extract the clusters from these results ([factor2cluster](#)), and then form the composite correlation matrix using [cluster.cor](#). The variables in this reduced matrix may then be used in multiple R procedures using [mat.regress](#).

The original correlation is pre and post multiplied by the (transpose) of the keys matrix.

If some correlations are missing from the original matrix this will lead to missing values (NA) for scale intercorrelations based upon those lower level correlations.

## Value

<code>cor</code>	the (raw) correlation matrix of the clusters
<code>sd</code>	standard deviation of the cluster scores
<code>corrected</code>	raw correlations below the diagonal, alphas on diagonal, disattenuated above diagonal
<code>size</code>	How many items are in each cluster?

## Note

See SAPA e.g., Revelle, W. (2006) Synthetic Aperture Personality Assessment. Invited paper at the Midwestern Psychological Association Annual Meeting, Chicago, May, 2006. pdf available at <http://personality-project.org/revelle/publications/sapa.mpa.key.pdf>

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## See Also

[factor2cluster](#), [mat.regress](#), [alpha.scale](#), [score.items](#)

## Examples

```
## Not run:
data(attitude)
keys <- matrix(c(1,1,1,0,0,0,0,
                 0,0,0,1,1,1,1),ncol=2)
colnames(keys) <- c("first","second")
r.mat <- cor(attitude)
cluster.cor(keys,r.mat)
## End(Not run)
##$cor
#      first second
#first   1.0    0.6
#second  0.6    1.0
#
##$sd
# first second
#  2.57    3.01
#
##$corrected
#      first second
#first  0.82    0.77
#second 0.60    0.74
#
##$size
# first second
#      3      4
```

---

`cluster.fit`

*cluster Fit: fit of the cluster model to a correlation matrix*

---



## Description

How well does the cluster model found by [ICLUST](#) fit the original correlation matrix? A similar algorithm [factor.fit](#) is found in [VSS](#). This function is internal to ICLUST but has more general use as well.

In general, the cluster model is a Very Simple Structure model of complexity one. That is, every item is assumed to represent only one factor/cluster. Cluster fit is an analysis of how well this model reproduces a correlation matrix. Two measures of fit are given: cluster fit and factor fit. Cluster fit assumes that variables that define different clusters are orthogonal. Factor fit takes the loadings generated by a cluster model, finds the cluster loadings on all clusters, and measures the degree of fit of this somewhat more complicated model. Because the cluster loadings are similar to, but not identical to factor loadings, the factor fits found here and by [factor.fit](#) will be similar.

## Usage

```
cluster.fit(original, load, clusters, diagonal = FALSE, digits = 2)
```

## Arguments

<code>original</code>	The original correlation matrix being fit
<code>load</code>	Cluster loadings – that is, the correlation of individual items with the clusters, corrected for item overlap
<code>clusters</code>	The cluster structure
<code>diagonal</code>	Should we fit the diagonal as well?
<code>digits</code>	Number of digits to be used in the output

## Details

The cluster model is similar to the factor model:  $R$  is fitted by  $C'C$ . Where  $C$  <- Cluster definition matrix x the loading matrix. How well does this model approximate the original correlation matrix and how does this compare to a factor model?

The fit statistic is a comparison of the original (squared) correlations to the residual correlations.  $\text{Fit} = 1 - r^2/r^2$  where  $r^*$  is the residual correlation of data - model and model =  $C'C$ .

## Value

<code>clusterfit</code>	The cluster model is a reduced form of the factor loading matrix. That is, it is the product of the elements of the cluster matrix * the loading matrix.
<code>factorfit</code>	How well does the complete loading matrix reproduce the correlation matrix?

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## References

<http://personality-project.org/r/r.ICLUST.html>

## See Also

[VSS](#), [ICLUST](#), [factor2cluster](#), [cluster.cor](#), [factor.fit](#)

## Examples

```
r.mat<- Harman74.cor$cov
iq.clus <- ICLUST(r.mat,nclusters =2)
fit <- cluster.fit(r.mat,iq.clus$loadings,iq.clus$clusters)
fit
```

---

<code>cluster.loadings</code>	<i>Find item by cluster correlations, corrected for overlap and reliability</i>
-------------------------------	---

---

## Description

Given a  $n \times n$  correlation matrix and a  $n \times c$  matrix of -1,0,1 cluster weights for those  $n$  items on  $c$  clusters, find the correlation of each item with each cluster. If the item is part of the cluster, correct for item overlap. Part of the [ICLUST](#) set of functions, but useful for many item analysis problems.

## Usage

```
cluster.loadings(keys, r.mat, correct = TRUE, digits = 2)
```

## Arguments

<code>keys</code>	Cluster keys: a matrix of -1,0,1 cluster weights
<code>r.mat</code>	A correlation matrix
<code>correct</code>	Correct for reliability
<code>digits</code>	Number of digits output

## Details

Given a set of items to be scored as (perhaps overlapping) clusters and the intercorrelation matrix of the items, find the clusters and then the correlations of each item with each cluster. Correct for item overlap by replacing the item variance with its average within cluster inter-item correlation.

Although part of ICLUST, this may be used in any SAPA application where we are interested in item- whole correlations of items and composite scales.

These loadings are particularly interpretable when sorted by absolute magnitude for each cluster (see [ICLUST.sort](#)).

## Value

<code>loadings</code>	A matrix of item-cluster correlations (loadings)
<code>cor</code>	Correlation matrix of the clusters
<code>corrected</code>	Correlation matrix of the clusters, raw correlations below the diagonal, alpha on diagonal, corrected for reliability above the diagonal
<code>sd</code>	Cluster standard deviations
<code>alpha</code>	alpha reliabilities of the clusters
<code>count</code>	Number of items in the cluster

## Note

Although part of ICLUST, this may be used in any SAPA application where we are interested in item- whole correlations of items and composite scales.

## Author(s)

Maintainer: William Revelle ([revelle@northwestern.edu](mailto:revelle@northwestern.edu))

## References

ICLUST: <http://personality-project.org/r/r.iclust.html>

## See Also

[ICLUST](#), [factor2cluster](#), [cluster.cor](#)

## Examples

```
## Not run:
r.mat<- Harman74.cor$cov
clusters <- matrix(c(1,1,1,rep(0,24),1,1,1,1,rep(0,17)),ncol=2)
cluster.loadings(clusters,r.mat)
## End(Not run)
```

---

<code>congeneric.sim</code>	<i>Simulate a congeneric data set</i>
-----------------------------	---------------------------------------

---

## Description

Classical Test Theory (CTT) considers four or more tests to be congenerically equivalent if all tests may be expressed in terms of one factor and a residual error. Parallel tests are the special case where (usually two) tests have equal factor loadings. Tau equivalent tests have equal factor loadings but may have unequal errors. Congeneric tests may differ in both factor loading and error variances.

## Usage

```
congeneric.sim(N = 1000, loads = c(0.8, 0.7, 0.6, 0.5), err=NULL, short = TRUE)
```

## Arguments

<b>N</b>	How many subjects to simulate
<b>loads</b>	A vector of factor loadings for the tests
<b>err</b>	A vector of error variances – if NULL then error = 1 - loading 2
<b>short</b>	short=TRUE: Just give the test scores, short=FALSE, report observed test scores as well as the implied pattern matrix

## Details

When constructing examples for reliability analysis, it is convenient to simulate congeneric data structures. These are the most simple of item structures, having just one factor.

The implied covariance matrix is just  $\text{pattern \%*\% t(pattern)}$ .

## Value

<b>observed</b>	a matrix of test scores for n tests
<b>pattern</b>	The pattern matrix implied by the loadings and error variances

## Author(s)

William Revelle

## References

## See Also

[item.sim](#)

## Examples

```
#test <- congeneric.sim()
```

---

`correct.cor`*Find dis-attenuated correlations and give alpha reliabilities*

---

## Description

Given a raw correlation matrix and a vector of reliabilities, report the disattenuated correlations above the diagonal.

## Usage

```
correct.cor(x, y)
```

## Arguments

<code>x</code>	A raw correlation matrix
<code>y</code>	Vector of reliabilities

## Details

Disattenuated correlations may be thought of as correlations between the latent variables measured by a set of observed variables. That is, what would the correlation be between two (unreliable) variables be if both variables were measured perfectly reliably.

This function is mainly used if importing correlations and reliabilities from somewhere else. If the raw data are available, use `score.items`, or `cluster.loadings` or `cluster.cor`.

Examples of the output of this function are seen in `cluster.loadings` and `cluster.cor`

## Value

Raw correlations below the diagonal, reliabilities on the diagonal, disattenuated above the diagonal.

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## References

<http://personality-project.org/revelle/syllabi/405.syllabus.html>

## See Also

`cluster.loadings` and `cluster.cor`

## Examples

```
# attitude from the datasets package
#example 1 is a rather clunky way of doing things
## Not run:
a1 <- attitude[,c(1:3)]
a2 <- attitude[,c(4:7)]
x1 <- rowSums(a1) #find the sum of the first 3 attitudes
x2 <- rowSums(a2) #find the sum of the last 4 attitudes
alpha1 <- alpha.scale(x1,a1)
alpha2 <- alpha.scale(x2,a2)
x <- matrix(c(x1,x2),ncol=2)
x.cor <- cor(x)
alpha <- c(alpha1,alpha2)
round(correct.cor(x.cor,alpha),2)
#
#much better - although uses standardized alpha
clusters <- matrix(c(rep(1,3),rep(0,7),rep(1,4)),ncol=2)
cluster.loadings(clusters,cor(attitude))
# or
clusters <- matrix(c(rep(1,3),rep(0,7),rep(1,4)),ncol=2)
cluster.cor(clusters,cor(attitude))
#
#best
scores <- score.items(matrix(c(rep(1,3),rep(0,7),rep(1,4)),ncol=2),attitude)
scores$corrected
## End(Not run)
```

---

count.pairwise	<i>Count number of pairwise cases for a data set with missing (NA) data.</i>
----------------	--

---

## Description

When doing `cor(x, use= "pairwise")`, it is nice to know the number of cases for each pairwise correlation. This is particularly useful when doing SAPA type analyses.

## Usage

```
count.pairwise(x, y = NULL)
```

## Arguments

x	An input matrix, typically a data matrix ready to be correlated.
y	An optional second input matrix

## Value

result = matrix of counts of pairwise observations

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## Examples

```
## Not run:
x <- matrix(rnorm(1000),ncol=6)
y <- matrix(rnorm(500),ncol=3)
x[x < 0] <- NA
y[y> 1] <- NA

count.pairwise(x)
count.pairwise(y)
count.pairwise(x,y)
## End(Not run)
```

---

describe.by

*Basic summary statistics by group*

---

## Description

Report basic summary statistics by a grouping variable. Useful if the grouping variable is some experimental variable and data are to be aggregated for plotting. Just a wrapper for `by` and [describe](#).

## Usage

```
describe.by(x, group,...)
```

## Arguments

x	a data.frame or matrix
group	a grouping variable
...	parameters to be passed to describe

## Details

## Value

A data.frame of the relevant statistics broken down by group:

item name  
item number  
number of valid cases  
mean  
standard deviation  
median  
mad: median absolute deviation (from the median)  
minimum  
maximum  
skew  
standard error

## Author(s)

William Revelle

## See Also

[describe](#)

## Examples

```
set.seed(42) #to get the same values each time
x.df <- data.frame(group=sample(2,20,replace=TRUE), matrix(rnorm(100),ncol=5))
x <- describe.by(x.df,x.df$group)
x #shows all the results
x[1] #shows just the first group
x <- matrix(sample(4,200,replace=TRUE),ncol=5)
y <- describe.by(x,x[,1])
y
#
#group: 1
#   var n mean  sd median mad   min max range  se
#group  1 7  1.00 0.00   1.00 0.00  1.00 1.00  0.00 0.00
#X1     2 7 -0.36 1.66  -0.31 2.16 -2.66 1.90  4.55 0.63
#X2     3 7 -0.10 1.04  -0.43 0.95 -1.37 1.44  2.81 0.39
#...
#-----
#group: 2
#   var n mean  sd median mad   min max range  se
#group  1 13  2.00 0.00   2.00 0.00  2.00 2.00  0.00 0.00
#X1     2 13 -0.07 1.26  -0.26 0.57 -2.44 2.29  4.73 0.35
#...
#> x[1] #shows just the first group
#`$`1`
#   var n mean  sd median mad   min max range  se
#group  1 7  1.00 0.00   1.00 0.00  1.00 1.00  0.00 0.00
#X1     2 7 -0.36 1.66  -0.31 2.16 -2.66 1.90  4.55 0.63
```



#X2	3	7	-0.10	1.04	-0.43	0.95	-1.37	1.44	2.81	0.39
#X3	4	7	0.56	0.90	0.92	0.97	-0.73	1.58	2.30	0.34
#X4	5	7	0.29	0.88	0.46	0.87	-1.19	1.51	2.71	0.33
#X5	6	7	0.23	1.12	0.19	1.27	-1.46	1.85	3.31	0.42

---

describe

*Basic descriptive statistics useful for psychometrics*

---

## Description

There are many summary statistics available in R; this function provides the ones most useful for scale construction and item analysis in classic psychometrics. Range is most useful for the first pass in a data set, to check for coding errors.

## Usage

```
describe(x, digits = 2, na.rm = TRUE, skew = TRUE, ranges = TRUE)
```

## Arguments

<b>x</b>	A data frame or matrix
<b>digits</b>	How many significant digits to report
<b>na.rm</b>	The default is to delete missing data
<b>skew</b>	Should the skew and kurtosis be calculated?
<b>ranges</b>	Should the range be calculated?

## Details

In basic data analysis it is vital to get basic descriptive statistics. Procedures such as [summary](#) and `hmisc::describe` do so. The describe function in the [psych](#) package is meant to produce the most frequently requested stats in psychometric and psychology studies, and to produce them in an easy to read data.frame. The results from describe can be used in graphics functions (e.g., [error.crosses](#)).

The range statistics (min, max, range) are most useful for data checking to detect coding errors, and should be found in early analyses of the data.

Although describe will work on data frames as well as matrices, it is important to realize that for data frames, descriptive statistics will be reported only for those variables where this makes sense (i.e., not for factors or for alphanumeric data).

In a typical study, one might read the data in from the clipboard ([read.clipboard](#)), show the splom plot of the correlations ([pairs.panels](#)), and then describe the data.

## Value

A data.frame of the relevant statistics:  
item name  
item number  
number of valid cases  
mean  
standard deviation  
median  
mad: median absolute deviation (from the median)  
minimum  
maximum  
skew  
kurtosis  
standard error

## Note

Describe uses either the mean or colMeans functions depending upon whether the data are a data.frame or a matrix. The mean function supplies means for the columns of a data.frame, but the overall mean for a matrix. Mean will throw a warning for non-numeric data, but colMeans stops with non-numeric data. Thus, the describe function uses either mean (for data frames) or colMeans (for matrices). This is true for skew and kurtosi as well.

## Author(s)

<http://personality-project.org/revelle.html>

Maintainer: William Revelle <revelle@northwestern.edu>

## See Also

[describe.by](#), [skew](#), [kurtosi](#), [pairs.panels](#), [read.clipboard](#), [error.crosses](#)

## Examples

```
describe(attitude)
#           var  n mean    sd median   mad min max range  skew kurtosis  se
#rating      1 30 64.63 12.17   65.5 10.38  40  85   45 -0.36   -0.77 2.22
#complaints   2 30 66.60 13.31   65.0 14.83  37  90   53 -0.22   -0.68 2.43
#privileges   3 30 53.13 12.24   51.5 10.38  30  83   53  0.38   -0.41 2.23
#learning     4 30 56.37 11.74   56.5 14.83  34  75   41 -0.05   -1.22 2.14
#raises       5 30 64.63 10.40   63.5 11.12  43  88   45  0.20   -0.60 1.90
#critical     6 30 74.77  9.89   77.5  7.41  49  92   43 -0.87    0.17 1.81
#advance      7 30 42.93 10.29   41.0  8.90  25  72   47  0.85    0.47 1.88

describe(attitude,skew=FALSE)  #attitude is taken from R data sets
```

#	var	n	mean	sd	median	mad	min	max	range	se
#rating	1	30	64.63	12.17	65.5	10.38	40	85	45	2.22
#complaints	2	30	66.60	13.31	65.0	14.83	37	90	53	2.43
#privileges	3	30	53.13	12.24	51.5	10.38	30	83	53	2.23
#learning	4	30	56.37	11.74	56.5	14.83	34	75	41	2.14
#raises	5	30	64.63	10.40	63.5	11.12	43	88	45	1.90
#critical	6	30	74.77	9.89	77.5	7.41	49	92	43	1.81
#advance	7	30	42.93	10.29	41.0	8.90	25	72	47	1.88

---

<code>eigen.loadings</code>	<i>Convert eigen vectors and eigen values to the more normal (for psychologists) component loadings</i>
-----------------------------	---

---

## Description

The default procedures for principal component returns values not immediately equivalent to the loadings from a factor analysis. `eigen.loadings` translates them into the more typical metric of eigen vectors multiplied by the squareroot of the eigenvalues. This lets us find pseudo factor loadings if we have used `princomp` or `princ`. If we use `principal` to do our principal components analysis, then we do not need this routine.

## Usage

```
eigen.loadings(x)
```

## Arguments

`x` a matrix of loadings from `princ` or `princomp`

## Value

A matrix of PA loadings more typical for what is expected in psychometrics.

## Note

Useful for SAPA analyses

## Author(s)

( `revelle@northwestern.edu` )  
<http://personality-project.org/revelle.html>

## Examples

---

<code>error.crosses</code>	<i>Plot x and y error bars</i>
----------------------------	--------------------------------

---

## Description

Given two vectors of data, plot the means and show standard errors in both X and Y directions.

## Usage

```
error.crosses(x, y, labels = NULL, pos = NULL, arrow.len = 0.2, ...)
```

## Arguments

<code>x</code>	A vector of summary statistics (from Describe)
<code>y</code>	A second vector of summary statistics (also from Describe)
<code>labels</code>	name the pair
<code>pos</code>	Labels are located where with respect to the mean?
<code>arrow.len</code>	Arrow length
<code>...</code>	Other parameters for plot

## Details

For an example of two way error bars describing the effects of mood manipulations upon positive and negative affect, see <http://personality-project.org/revelle/publications/happy-sad-appendix/FIG.A-6.pdf>)

## Author(s)

William Revelle  
<revelle@northwestern.edu>

## Examples

```
#desc <- describe(attitude)
#x <- desc[1,]
#y <- desc[2,]
#plot(x$mean,y$mean)      #in graphics window
#error.crosses(x,y)       #in graphics window
```

A graphNEL graph with directed edges  
Number of Nodes = 5  
Number of Edges = 4

**4 congeneric measures**

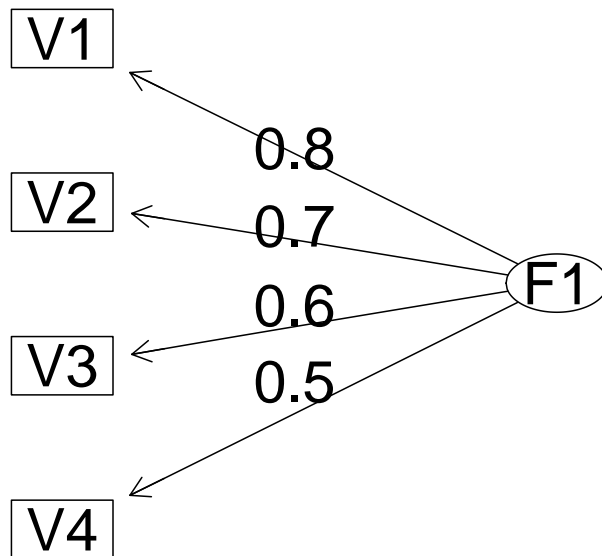


Figure 3: Congeneric tests have one common factor and can differ in their error variances. A demonstration of functions `congeneric.sim`, `factor.pa`, and `fa.graph`.

---

`fa.graph`

---

*Graph factor loading matrices*

## Description

Factor analysis or principal components analysis results are typically interpreted in terms of the major loadings on each factor. These structures may be represented as a table of loadings or graphically, where all loadings with an absolute value  $>$  some cut point are represented as an edge (path).

## Usage

```
fa.graph(fa.results, out.file = NULL, labels = NULL, cut = 0.3, simple = TRUE, size = c(8, 6),
```

## Arguments

<code>fa.results</code>	The output of factor analysis or principal components analysis
<code>out.file</code>	If it exists, a dot representation of the graph will be stored here
<code>labels</code>	Variable labels
<code>cut</code>	Loadings with $\text{abs}(\text{loading}) > \text{cut}$ will be shown
<code>simple</code>	Only the biggest loading per item is shown
<code>size</code>	
<code>node.font</code>	
<code>edge.font</code>	
<code>rank.direction</code>	
<code>digits</code>	Number of digits to show as an edgelable
<code>title</code>	Graphic title
<code>...</code>	other parameters

## Details

Path diagram representations have become standard in confirmatory factor analysis, but are not yet common in exploratory factor analysis. Representing factor structures graphically helps some people understand the structure.

## Value

A graph is drawn using `rgraphviz`. If an output file is specified, the graph instructions are also saved in the dot language.

## Note

Requires `Rgraphviz`.

As of June 1, there is an occasionally strange result when using the `simple=FALSE` option in Sweave.

### Parallel Analysis Scree Plots

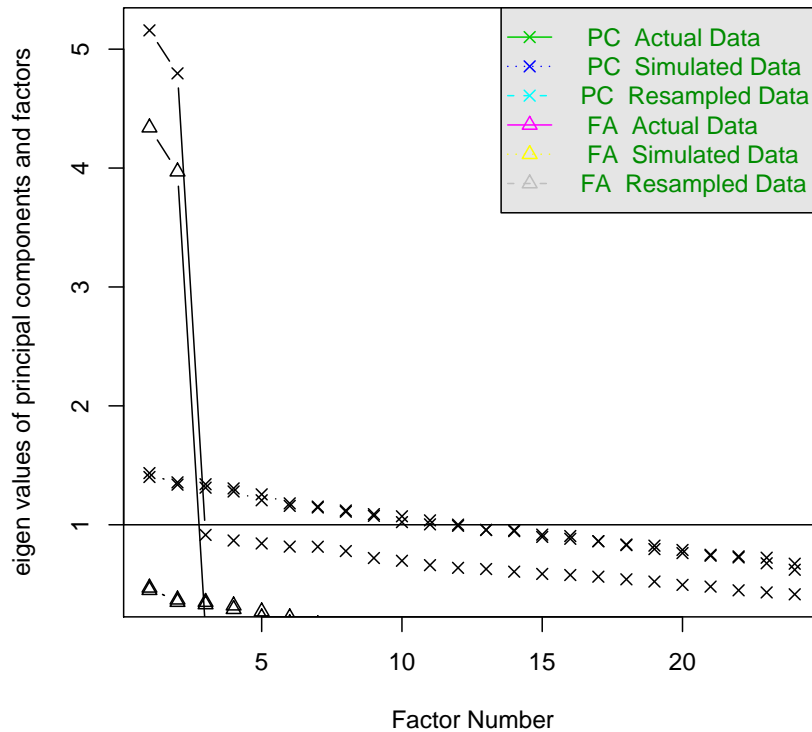


Figure 4: Plots of eigen values for real and simulated “parallel” data give two ways to determine the optimal number of factors: Examine the scree for large breaks or find when the eigen values of random data are greater than for the real data. For this data set, both methods suggest two factors are optimal.

### Author(s)

William Revelle

### See Also

[omega.graph](#), [ICLUST.graph](#)

### Examples

```
test.simple <- factor.pa(item.sim(16),2)
if(require(Rgraphviz)) {fa.graph(test.simple) }
```

---

<code>fa.parallel</code>	<i>Scree plots of data or correlation matrix compared to random “parallel” matrices</i>
--------------------------	---

---

## Description

One way to determine the number of factors or components in a data matrix or a correlation matrix is to examine the “scree” plot of the successive eigenvalues. Sharp breaks in the plot suggest the appropriate number of components or factors to extract. “Parallel” analysis is an alternative technique that compares the scree of the observed data with that of a random data matrix of the same size as the original.

## Usage

```
fa.parallel(x, n.obs = NULL, fa="both", main = "Parallel Analysis Scree Plots")
```

## Arguments

<code>x</code>	A data.frame or data matrix of scores. If the matrix is square, it is assumed to be a correlation matrix. Otherwise, correlations (with pairwise deletion) will be found
<code>n.obs</code>	n.obs=0 implies a data matrix/data.frame. Otherwise, how many cases were used to find the correlations.
<code>fa</code>	show the eigen values for a principal components (fa="pc") or a principal axis factor analysis (fa="fa") or both principal components and principal factors (fa="both")
<code>main</code>	a title for the analysis

## Details

Cattell’s “scree” test is one of most simple tests for the number of factors problem. Humphreys and Montanelli’s “parallel” analysis is an equally compelling procedure. Other procedures for determining the most optimal number of factors include finding the Very Simple Structure (VSS) criterion ([VSS](#)). `fa.parallel` plots the eigen values for a principal components and/or principal factor solution and does the same for random matrices of the same size as the original data matrix. For raw data, the random matrices are 1) a matrix of univariate normal data and 2) random samples (randomized across rows) of the original data.

## Value

A plot of the eigen values for the original data, a resampling of the original data, and of a equivalent size matrix of random normal deviates. If the data are a correlation matrix, specify the number of observations.

## Author(s)

William Revelle



## References

## See Also

[VSS](#), [VSS.plot](#), [VSS.parallel](#)

## Examples

```
#not run
#test.data <- Harman74.cor$cov
#fa.parallel(test.data,n.obs=200)
#
#fa.parallel(attitude)
#
```

---

<code>factor.congruence</code>	<i>Coefficient of factor congruence</i>
--------------------------------	---

---

## Description

Given two sets of factor loadings, report their degree of congruence.

## Usage

```
factor.congruence(x, y, loading=TRUE)
```

## Arguments

<code>x</code>	A matrix of factor loadings
<code>y</code>	A second matrix of factor loadings
<code>loading</code>	Is the input a loading matrix from a factor analysis or principal components, or is it just a matrix

## Details

Find the coefficient of factor congruence between two sets of factor loadings.

It is an interesting exercise to compare factor congruences with the correlations of factor loadings. Factor congruences are based upon the raw cross products, while correlations are based upon centered cross products.

## Value

A matrix of factor congruences.

## Author(s)

<revelle@northwestern.edu>  
<http://personality-project.org/revelle.html>

## References

Gorsuch, Richard, (1983) Factor Analysis. Lawrence Erlbaum Associates.  
Revelle, W. (In preparation) An Introduction to Psychometric Theory with applications in R (<http://personality-project.org/r/book/>)

## See Also

[principal](#), [factor.pa](#)

## Examples

```
#fa <- factanal(x,4,covmat=Harman74.cor$cov)
#pc <- principal(Harman74.cor$cov,4)
#pcv <- varimax(pc$loading)
#round(factor.congruence(fa,pcv),2)
#
#round(factor.congruence(pcv,fa),2)
#   Factor1 Factor2 Factor3 Factor4
#PC1    1.00    0.60    0.45    0.55
#PC2    0.44    0.49    1.00    0.56
#PC3    0.54    0.99    0.44    0.55
#PC4    0.47    0.52    0.48    0.99

#compare with
#round(cor(fa$loading,pcv$loading),2)
```

---

<code>factor.fit</code>	<i>How well does the factor model fit a correlation matrix. Part of the VSS package</i>
-------------------------	---

---

## Description

The basic factor or principal components model is that a correlation or covariance matrix may be reproduced by the product of a factor loading matrix times its transpose:  $F'F$  or  $P'P$ . One simple index of fit is the 1 - sum squared residuals/sum squared original correlations. This fit index is used by [VSS](#), [ICLUST](#), etc.

## Usage

```
factor.fit(r, f)
```

## Arguments

**r** a correlation matrix  
**f** A factor matrix of loadings.

## Details

There are probably as many fit indices as there are psychometricians. This fit is a plausible estimate of the amount of reduction in a correlation matrix given a factor model. Note that it is sensitive to the size of the original correlations. That is, if the residuals are small but the original correlations are small, that is a bad fit.

## Value

fit

## Author(s)

William Revelle

## See Also

[VSS](#), [ICLUST](#)

## Examples

```
## Not run:
#compare the fit of 4 to 3 factors for the Harman 24 variables
fa4 <- factanal(x,4,covmat=Harman74.cor$cov)
round(factor.fit(Harman74.cor$cov,fa4$loading),2)
#[1] 0.9
fa3 <- factanal(x,3,covmat=Harman74.cor$cov)
round(factor.fit(Harman74.cor$cov,fa3$loading),2)
#[1] 0.88

## End(Not run)
```

---

`factor.model`

*Find  $R = F F' + U2$  is the basic factor model*

---

## Description

The basic factor or principal components model is that a correlation or covariance matrix may be reproduced by the product of a factor loading matrix times its transpose. Find this reproduced matrix. Used by [factor.fit](#), [VSS](#), [ICLUST](#), etc.

## Usage

```
factor.model(f)
```

## Arguments

**f**                      A matrix of loadings.

## Details

## Value

A correlation or covariance matrix.

## Author(s)

`<revelle@northwestern.edu >`  
<http://personality-project.org/revelle.html>

## References

Gorsuch, Richard, (1983) Factor Analysis. Lawrence Erlbaum Associates.  
Revelle, W. In preparation) An Introduction to Psychometric Theory with applications in R (<http://personality-project.org/r/book/>)

## See Also

[ICLUST.graph](#), [ICLUST.cluster](#), [cluster.fit](#) , [VSS](#), [omega](#)

## Examples

```
f2 <- matrix(c(.9,.8,.7,rep(0,6),.6,.7,.8),ncol=2)
mod <- factor.model(f2)
round(mod,2)
```

---

<code>factor.pa</code>	<i>Principal Axis Factor Analysis</i>
------------------------	---------------------------------------

---

## Description

Among the many ways to do factor analysis, one of the most conventional is principal axes. An eigen value decomposition of a correlation matrix is done and then the communalities for each variable are estimated by the first n factors. These communalities are entered onto the diagonal and the procedure is repeated until the `sum(diag(r))` does not vary. For well behaved matrices, maximum likelihood factor analysis (`factanal`) is probably preferred.

## Usage

```
factor.pa(r, nfactors=1, residuals = FALSE, rotate = "varimax", n.obs = NULL,
min.err = 0.001, digits = 2, max.iter = 50)
```

A graphNEL graph with directed edges  
 Number of Nodes = 28  
 Number of Edges = 35

### Harman's 24 tests of mental ability

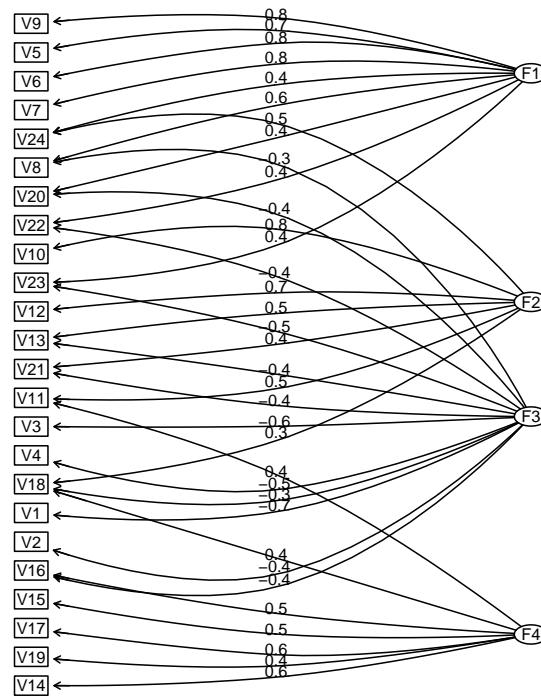


Figure 5: Four factors of the Harman 24 mental tests suggest that a simple structure is difficult to obtain. Note the cross loadings for several items. This figure was created with factor.pa and omega.graph. Compare with a hierarchical solution using the omega function (figure 8) or a cluster solution with ICLUST (figure 6).

## Arguments

<b>r</b>	A correlation matrix or a raw data matrix. If raw data, the correlation matrix will be found using pairwise deletion.
<b>nfactors</b>	Number of factors to extract, default is 1
<b>residuals</b>	Should the residual matrix be shown
<b>rotate</b>	
<b>n.obs</b>	Number of observations used to find the correlation matrix if using a correlation matrix. Used for finding the goodness of fit statistics.
<b>min.err</b>	Iterate until the change in communalities is less than min.err
<b>digits</b>	How many digits of output should be returned
<b>max.iter</b>	Maximum number of iterations for convergence

## Details

Factor analysis is an attempt to approximate a correlation or covariance matrix with one of lesser rank. The basic model is that  ${}_nR_n \approx {}_nF_{kk}F'_n + U^2$  where  $k$  is much less than  $n$ . There are many ways to do factor analysis, and maximum likelihood procedures are probably the most preferred (see [factanal](#)). The existence of uniquenesses is what distinguishes factor analysis from principal components analysis (e.g., [principal](#)).

Principal axes factor analysis has a long history in exploratory analysis and is a straight-forward procedure. Successive eigen value decompositions are done on a correlation matrix with the diagonal replaced with  $\text{diag}(FF')$  until  $\text{sum}(\text{diag}(FF'))$  does not change (very much). The current limit of `max.iter` = 50 seems to work for most problems, but the Holzinger-Harmon 24 variable problem needs about 203 iterations to converge for a 5 factor solution.

Principal axes may be used in cases when maximum likelihood solutions fail to converge.

The algorithm does not attempt to find the best (as defined by a maximum likelihood criterion) solution, but rather one that converges rapidly using successive eigen value decompositions. The maximum likelihood criterion of fit and the associated chi square value are reported, and will be worse than that found using maximum likelihood procedures.

## Value

<b>values</b>	Eigen values of the final solution
<b>communality</b>	Communality estimates for each item. These are merely the sum of squared factor loadings for that item.
<b>rotation</b>	which rotation was requested?
<b>n.obs</b>	number of observations specified or found
<b>loadings</b>	An item by factor loading matrix. Suitable for use in other programs (e.g., GPA rotation or <code>factor2cluster</code> ).
<b>fit</b>	How well does the factor model reproduce the correlation matrix. (See <a href="#">VSS</a> , <a href="#">ICLUST</a> , and <a href="#">principal</a> for this fit statistic.
<b>fit.off</b>	how well are the off diagonal elements reproduced?

<b>dof</b>	Degrees of Freedom for this model. This is the number of observed correlations minus the number of independent parameters. Let $n$ =Number of items, $nf$ = number of factors then $dof = n * (n - 1)/2 - n * nf + nf * (nf - 1)/2$
<b>objective</b>	value of the function that is minimized by maximum likelihood procedures. This is reported for comparison purposes and as a way to estimate chi square goodness of fit. The objective function is $f = \log(\text{trace}((FF' + U2)^{-1}R) - \log( (FF' + U2)^{-1}R ) - n.items.$
<b>STATISTIC</b>	If the number of observations is specified or found, this is a chi square based upon the objective function, $f$ . Using the formula from <a href="#">factanal</a> (which seems to be Bartlett's test) : $\chi^2 = (n.obs - 1 - (2 * p + 5)/6 - (2 * factors)/3)) * f$
<b>PVAL</b>	If $n.obs > 0$ , then what is the probability of observing a chisquare this large or larger?
<b>phi</b>	If oblique rotations (using oblimin from the GPArotation package) are requested, what is the interfactor correlation.
<b>communality.iterations</b>	The history of the communality estimates. Probably only useful for teaching what happens in the process of iterative fitting.
<b>residual</b>	If residuals are requested, this is the matrix of residual correlations after the factor model is applied.

## Author(s)

William Revelle

## References

Gorsuch, Richard, (1983) Factor Analysis. Lawrence Erlbaum Associates.

## See Also

[principal](#), [VSS](#), [ICLUST](#)

## Examples

```
#using the Harmon 24 mental tests, compare a principal factor with a principal components solution
pc <- principal(Harman74.cor$cov,4,rotate=TRUE)
pa <- factor.pa(Harman74.cor$cov,4,rotate="varimax")
round(factor.congruence(pc,pa),2)

#then compare with a maximum likelihood solution using factanal
mle <- factanal(x,4,covmat=Harman74.cor$cov)
round(factor.congruence(mle,pa),2)
#note that the order of factors and the sign of some of factors differ
```

---

<code>factor.residuals</code>	$R^* = R - F F'$
-------------------------------	------------------

---

## Description

The basic factor or principal components model is that a correlation or covariance matrix may be reproduced by the product of a factor loading matrix times its transpose. Find the residuals of the original minus the reproduced matrix. Used by [factor.fit](#), [VSS](#), [ICLUST](#), etc.

## Usage

```
factor.residuals(r, f)
```

## Arguments

<code>r</code>	A correlation matrix
<code>f</code>	A factor model matrix

## Details

The basic factor equation is  ${}_nR_n \approx {}_nF_{kk}F'_n + U^2$ . Residuals are just  $R^* = R - F'F$ .

## Value

`rstar` is the residual correlation matrix.

## Author(s)

Maintainer: William Revelle <revelle@northwestern.edu>

## See Also

[factor.pa](#), [principal](#), [VSS](#), [ICLUST](#)

## Examples



---

<code>factor.rotate</code>	<i>“Hand” rotate a factor loading matrix</i>
----------------------------	--

---

### Description

Given a factor or components matrix, it is sometimes useful to do arbitrary rotations of particular pairs of variables. This supplements the much more powerful rotation package `GRProtation` and is meant for specific requirements to do unusual rotations.

### Usage

```
factor.rotate(f, angle, col1, col2)
```

### Arguments

<code>f</code>	original loading matrix
<code>angle</code>	angle (in degrees!) to rotate
<code>col1</code>	column in factor matrix defining the first variable
<code>col2</code>	column in factor matrix defining the second variable

### Details

Partly meant as a demonstration of how rotation works, `factor.rotate` is useful for those cases that require specific rotations that are not available in more advanced packages such as `GPArotation`.

### Value

the resulting rotated matrix of loadings.

### Note

For a complete rotation package, see `GPArotation`

### Author(s)

Maintainer: William Revelle ([revelle@northwestern.edu](mailto:revelle@northwestern.edu))

### References

<http://personality-project.org/revelle/syllabi/405.syllabus.html>

### Examples

---

**factor2cluster***Extract cluster definitions from factor loadings*

---

## Description

Given a factor or principal components loading matrix, assign each item to a cluster corresponding to the largest (signed) factor loading for that item. Essentially, this is a Very Simple Structure approach to cluster definition that corresponds to what most people actually do: highlight the largest loading for each item and ignore the rest.

## Usage

```
factor2cluster(loads, loading = TRUE, cut = 0)
```

## Arguments

<b>loads</b>	either a matrix of loadings, or the result of a factor analysis/principal components analysis with a loading component
<b>loading</b>	TRUE if taking the output of a FA/PC or ICLUST, FALSE if just a matrix
<b>cut</b>	Extract items with absolute loadings > cut

## Details

A factor/principal components analysis loading matrix is converted to a cluster (-1,0,1) definition matrix where each item is assigned to one and only one cluster. This is a fast way to extract items that will be unit weighted to form cluster composites. Use this function in combination with `cluster.cor` to find the correlations of these composite scores.

A typical use in the SAPA project is to form item composites by clustering or factoring (see [ICLUST](#), [principal](#)), extract the clusters from these results ([factor2cluster](#)), and then form the composite correlation matrix using [cluster.cor](#). The variables in this reduced matrix may then be used in multiple R procedures using `mat.regress`.

## Value

a matrix of -1,0,1 cluster definitions for each item.

## Author(s)

<http://personality-project.org/revelle.html>

Maintainer: William Revelle < [revelle@northwestern.edu](mailto:revelle@northwestern.edu) >

## References

<http://personality-project.org/r/r.vss.html>

See Also

[cluster.cor](#), [factor2cluster](#), [principal](#), [ICLUST](#)

## Examples

```
## Not run:
f <- factanal(x,4,covmat=Harman74.cor$cov)
factor2cluster(f)
## End(Not run)
#
```

	Factor1	Factor2	Factor3	Factor4
#VisualPerception	0	1	0	0
#Cubes	0	1	0	0
#PaperFormBoard	0	1	0	0
#Flags	0	1	0	0
#GeneralInformation	1	0	0	0
#ParagraphComprehension	1	0	0	0
#SentenceCompletion	1	0	0	0
#WordClassification	1	0	0	0
#WordMeaning	1	0	0	0
#Addition	0	0	1	0
#Code	0	0	1	0
#CountingDots	0	0	1	0
#StraightCurvedCapitals	0	0	1	0
#WordRecognition	0	0	0	1
#NumberRecognition	0	0	0	1
#FigureRecognition	0	0	0	1
#ObjectNumber	0	0	0	1
#NumberFigure	0	0	0	1
#FigureWord	0	0	0	1
#Deduction	0	1	0	0
#NumericalPuzzles	0	0	1	0
#ProblemReasoning	0	1	0	0
#SeriesCompletion	0	1	0	0
#ArithmeticProblems	0	0	1	0

---

fisherz

*Fisher z transform of r*

---

## Description

convert a correlation to a z score using the Fisher transformation.

## Usage

```
fisherz(rho)
```

## Arguments

rho                      a Pearson r

## Value

z value corresponding to r

## Author(s)

Maintainer: William Revelle (revelle@northwestern.edu )

## Examples

```
## Not run:
cors <- seq(-.9,.9,.1)
round(fisherz(cors),2)
## End(Not run)
```

---

<code>geometric.mean</code>	<i>Find the geometric mean of a vector or columns of a data.frame.</i>
-----------------------------	--

---

## Description

The geometric mean is the nth root of n products or e to the mean log of x. Useful for describing non-normal, i.e., geometric distributions.

## Usage

```
geometric.mean(x)
```

## Arguments

x                      a vector or data.frame

## Details

Useful for teaching how to write functions, also useful for showing the different ways of estimating central tendency.

## Value

geometric mean(s) of x or x.df.

## Note

Not particularly useful if there are elements that are  $\leq 0$ .

**Author(s)**

William Revelle

**See Also**

[harmonic.mean](#), [mean](#)

**Examples**

```
x <- seq(1,5)
x2 <- x^2
geometric.mean(x)
geometric.mean(x2)
```

---

<code>harmonic.mean</code>	<i>Find the harmonic mean of a vector, matrix, or columns of a data.frame</i>
----------------------------	---

---

**Description**

The harmonic mean is merely the reciprocal of the arithmetic mean of the reciprocals.

**Usage**

```
harmonic.mean(x)
```

**Arguments**

`x` a vector, matrix, or data.frame

**Details**

Included as an example for teaching about functions. As well as for a discussion of how to estimate central tendencies.

**Value**

The harmonic mean(s)

**Note**

Included as a simple demonstration of how to write a function

## Examples

```
x <- seq(1,5)
x2 <- x^2
harmonic.mean(x)
harmonic.mean(x2)
```

---

ICLUST.cluster

*Function to form hierarchical cluster analysis of items*

---

## Description

The guts of the [ICLUST](#) algorithm. Called by [ICLUST](#).

## Usage

```
ICLUST.cluster(r.mat, ICLUST.options)
```

## Arguments

<code>r.mat</code>	A correlation matrix
<code>ICLUST.options</code>	A list of options (see <a href="#">ICLUST</a> )

## Details

See [ICLUST](#)

## Value

A list of cluster statistics, described more fully in [ICLUST](#)

<code>comp1</code>	Description of 'comp1'
<code>comp2</code>	Description of 'comp2'
...	

## Note

Although the main code for ICLUST is here in ICLUST.cluster, the more extensive documentation is for [ICLUST](#).

## Author(s)

William Revelle  
Department of Psychology  
Northwestern University  
Evanston, Illinois  
( [revelle@northwestern.edu](mailto:revelle@northwestern.edu) )  
<http://personality-project.org/revelle.html>

## References

Revelle, W. 1979, Hierarchical Cluster Analysis and the Internal Structure of Tests. Multivariate Behavioral Research, 14, 57-74. <http://personality-project.org/revelle/publications/iclust.pdf>  
See also more extensive documentation at <http://personality-project.org/r/r.ICLUST.html>

## See Also

[ICLUST.graph](#), [ICLUST](#), [cluster.fit](#) , [VSS](#), [omega](#)

## Examples

---

<code>ICLUST.graph</code>	<i>create control code for ICLUST graphical output</i>
---------------------------	--

---

## Description

Given a cluster structure determined by [ICLUST](#), create dot code to describe the [ICLUST](#) output. To use the dot code, use either <http://www.graphviz.org/> Graphviz or a commercial viewer (e.g., OmniGraffle).

## Usage

```
ICLUST.graph(ic.results, out.file,min.size=1, short = FALSE,labels=NULL,
size = c(8, 6), node.font = c("Helvetica", 14), edge.font = c("Helvetica", 12),
rank.direction = "RL", digits = 2, title = "ICLUST", ...)
```

## Arguments

<code>ic.results</code>	output list from ICLUST
<code>out.file</code>	name of output file (defaults to console)
<code>min.size</code>	draw a smaller node (without all the information) for clusters < min.size – useful for large problems
<code>short</code>	if short==TRUE, don't use variable names
<code>labels</code>	vector of text labels (contents) for the variables
<code>size</code>	size of output
<code>node.font</code>	Font to use for nodes in the graph
<code>edge.font</code>	Font to use for the labels of the arrows (edges)
<code>rank.direction</code>	LR or RL
<code>digits</code>	number of digits to show
<code>title</code>	any title
<code>...</code>	other options to pass

## Details

Will create (or overwrite) an output file and print out the dot code to show a cluster structure. This dot file may be imported directly into a dot viewer (e.g., <http://www.graphviz.org/>). The "dot" language is a powerful graphic description language that is particularly appropriate for viewing cluster output. Commercial graphics programs (e.g., OmniGraffle) can also read (and clean up) dot files.

ICLUST.graph takes the output from [ICLUST](#) results and processes it to provide a pretty picture of the results. Original variables shown as rectangles and ordered on the left hand side (if rank direction is RL) of the graph. Clusters are drawn as ellipses and include the alpha, beta, and size of the cluster. Edges show the cluster intercorrelations.

It is possible to trim the output to not show all cluster information. Clusters < min.size are shown as small ovals without alpha, beta, and size information.

## Value

Output is a set of dot commands written either to console or to the output file. These commands may then be used as input to any "dot" viewer, e.g., Graphviz.

## Author(s)

`<revelle@northwestern.edu >`  
<http://personality-project.org/revelle.html>

## References

ICLUST: <http://personality-project.org/r/r.iclust.html>

## See Also

[VSS.plot](#), [ICLUST](#)

## Examples

```
## Not run:
test.data <- Harman74.cor$cov
ic.out <- ICLUST(test.data)
out.file <- file.choose(new=TRUE) #create a new file to write the plot commands to
ICLUST.graph(ic.out,out.file)
now go to graphviz (outside of R) and open the out.file you created
print(ic.out,digits=2)
## End(Not run)

#test.data <- Harman74.cor$cov
#my.iclust <- ICLUST(test.data)
#ICLUST.graph(my.iclust)
#
#
#digraph ICLUST {
#  rankdir=RL;
```



```

# size="8,8";
# node [fontname="Helvetica" fontsize=14 shape=box, width=2];
# edge [fontname="Helvetica" fontsize=12];
# label = "ICLUST";
#     fontsize=20;
#V1  [label = VisualPerception];
#V2  [label = Cubes];
#V3  [label = PaperFormBoard];
#V4  [label = Flags];
#V5  [label = GeneralInformation];
#V6  [label = ParagraphComprehension];
#V7  [label = SentenceCompletion];
#V8  [label = WordClassification];
#V9  [label = WordMeaning];
#V10 [label = Addition];
#V11 [label = Code];
#V12 [label = CountingDots];
#V13 [label = StraightCurvedCapitals];
#V14 [label = WordRecognition];
#V15 [label = NumberRecognition];
#V16 [label = FigureRecognition];
#V17 [label = ObjectNumber];
#V18 [label = NumberFigure];
#V19 [label = FigureWord];
#V20 [label = Deduction];
#V21 [label = NumericalPuzzles];
#V22 [label = ProblemReasoning];
#V23 [label = SeriesCompletion];
#V24 [label = ArithmeticProblems];
#node [shape=ellipse, width ="1"];
#C1-> V9 [ label = 0.78 ];
#C1-> V5 [ label = 0.78 ];
#C2-> V12 [ label = 0.66 ];
#C2-> V10 [ label = 0.66 ];
#C3-> V18 [ label = 0.53 ];
#C3-> V17 [ label = 0.53 ];
#C4-> V23 [ label = 0.59 ];
#C4-> V20 [ label = 0.59 ];
#C5-> V13 [ label = 0.61 ];
#C5-> V11 [ label = 0.61 ];
#C6-> V7 [ label = 0.78 ];
#C6-> V6 [ label = 0.78 ];
#C7-> V4 [ label = 0.55 ];
#C7-> V1 [ label = 0.55 ];
#C8-> V16 [ label = 0.5 ];
#C8-> V14 [ label = 0.49 ];
#C9-> C1 [ label = 0.86 ];
#C9-> C6 [ label = 0.86 ];
#C10-> C4 [ label = 0.71 ];
#C10-> V22 [ label = 0.62 ];
#C11-> V21 [ label = 0.56 ];
#C11-> V24 [ label = 0.58 ];
#C12-> C10 [ label = 0.76 ];

```

```

#C12-> C11 [ label = 0.67 ];
#C13-> C8 [ label = 0.61 ];
#C13-> V15 [ label = 0.49 ];
#C14-> C2 [ label = 0.74 ];
#C14-> C5 [ label = 0.72 ];
#C15-> V3 [ label = 0.48 ];
#C15-> C7 [ label = 0.65 ];
#C16-> V19 [ label = 0.48 ];
#C16-> C3 [ label = 0.64 ];
#C17-> V8 [ label = 0.62 ];
#C17-> C12 [ label = 0.8 ];
#C18-> C17 [ label = 0.82 ];
#C18-> C15 [ label = 0.68 ];
#C19-> C16 [ label = 0.66 ];
#C19-> C13 [ label = 0.65 ];
#C20-> C19 [ label = 0.72 ];
#C20-> C18 [ label = 0.83 ];
#C21-> C20 [ label = 0.87 ];
#C21-> C9 [ label = 0.76 ];
#C22-> 0 [ label = 0 ];
#C22-> 0 [ label = 0 ];
#C23-> 0 [ label = 0 ];
#C23-> 0 [ label = 0 ];
#C1 [label = "C1\n alpha= 0.84\n beta= 0.84\nN= 2"] ;
#C2 [label = "C2\n alpha= 0.74\n beta= 0.74\nN= 2"] ;
#C3 [label = "C3\n alpha= 0.62\n beta= 0.62\nN= 2"] ;
#C4 [label = "C4\n alpha= 0.67\n beta= 0.67\nN= 2"] ;
#C5 [label = "C5\n alpha= 0.7\n beta= 0.7\nN= 2"] ;
#C6 [label = "C6\n alpha= 0.84\n beta= 0.84\nN= 2"] ;
#C7 [label = "C7\n alpha= 0.64\n beta= 0.64\nN= 2"] ;
#C8 [label = "C8\n alpha= 0.58\n beta= 0.58\nN= 2"] ;
#C9 [label = "C9\n alpha= 0.9\n beta= 0.87\nN= 4"] ;
#C10 [label = "C10\n alpha= 0.74\n beta= 0.71\nN= 3"] ;
#C11 [label = "C11\n alpha= 0.62\n beta= 0.62\nN= 2"] ;
#C12 [label = "C12\n alpha= 0.79\n beta= 0.74\nN= 5"] ;
#C13 [label = "C13\n alpha= 0.64\n beta= 0.59\nN= 3"] ;
#C14 [label = "C14\n alpha= 0.79\n beta= 0.74\nN= 4"] ;
#C15 [label = "C15\n alpha= 0.66\n beta= 0.58\nN= 3"] ;
#C16 [label = "C16\n alpha= 0.65\n beta= 0.57\nN= 3"] ;
#C17 [label = "C17\n alpha= 0.81\n beta= 0.71\nN= 6"] ;
#C18 [label = "C18\n alpha= 0.84\n beta= 0.75\nN= 9"] ;
#C19 [label = "C19\n alpha= 0.74\n beta= 0.65\nN= 6"] ;
#C20 [label = "C20\n alpha= 0.87\n beta= 0.74\nN= 15"] ;
#C21 [label = "C21\n alpha= 0.9\n beta= 0.77\nN= 19"] ;
#C22 [label = "C22\n alpha= 0\n beta= 0\nN= 0"] ;
#C23 [label = "C23\n alpha= 0\n beta= 0\nN= 0"] ;
#{ rank=same;
#V1;V2;V3;V4;V5;V6;V7;V8;V9;V10;V11;V12;V13;V14;V15;V16;V17;V18;V19;V20;V21;V22;V23;V24;}}
#
#copy the above output to Graphviz and draw it
#see \url{http://personality-project.org/r/ICLUST.html} for an example.

```

## Description

Given a cluster structure determined by [ICLUST](#), create a rgraphic directly using Rgraphviz. To create dot code to describe the [ICLUST](#) output with more precision, use [ICLUST.graph](#). As an option, dot code is also generated and saved in a file. To use the dot code, use either <http://www.graphviz.org/> Graphviz or a commercial viewer (e.g., OmniGraffle).

## Usage

```
ICLUST.rgraph(ic.results, out.file = NULL, min.size = 1, short = FALSE, labels = NULL, size = c
```

## Arguments

<code>ic.results</code>	output list from ICLUST
<code>out.file</code>	File name to save optional dot code.
<code>min.size</code>	draw a smaller node (without all the information) for clusters < min.size – useful for large problems
<code>short</code>	if short==TRUE, don't use variable names
<code>labels</code>	vector of text labels (contents) for the variables
<code>size</code>	size of output
<code>node.font</code>	Font to use for nodes in the graph
<code>edge.font</code>	Font to use for the labels of the arrows (edges)
<code>rank.direction</code>	LR or RL
<code>digits</code>	number of digits to show
<code>title</code>	any title
<code>...</code>	other options to pass

## Details

Will create (or overwrite) an output file and print out the dot code to show a cluster structure. This dot file may be imported directly into a dot viewer (e.g., <http://www.graphviz.org/>). The "dot" language is a powerful graphic description language that is particularly appropriate for viewing cluster output. Commercial graphics programs (e.g., OmniGraffle) can also read (and clean up) dot files.

ICLUST.rgraph takes the output from [ICLUST](#) results and processes it to provide a pretty picture of the results. Original variables shown as rectangles and ordered on the left hand side (if rank direction is RL) of the graph. Clusters are drawn as ellipses and include the alpha, beta, and size of the cluster. Edges show the cluster intercorrelations.

It is possible to trim the output to not show all cluster information. Clusters < min.size are shown as small ovals without alpha, beta, and size information.

## Value

Output is a set of dot commands written either to console or to the output file. These commands may then be used as input to any "dot" viewer, e.g., Graphviz.

ICLUST.rgraph is a version of [ICLUST.graph](#) that uses Rgraphviz to draw on the screen as well.

Additional output is drawn to main graphics screen.

## Note

Requires Rgraphviz

## Author(s)

`<revelle@northwestern.edu >`

<http://personality-project.org/revelle.html>

## References

ICLUST: <http://personality-project.org/r/r.iclust.html>

## See Also

[VSS.plot](#), [ICLUST](#)

## Examples

```
test.data <- Harman74.cor$cov
ic.out <- ICLUST(test.data)
if(require(Rgraphviz) ) {ICLUST.rgraph(ic.out) }
```

---

ICLUST.sort

*sort items by absolute size of cluster loadings*

---

## Description

Given a cluster analysis or factor analysis loadings matrix, sort the items by the (absolute) size of each column of loadings. Used as part of ICLUST and SAPA analyses.

## Usage

```
ICLUST.sort(ic.load, cut = 0, labels = NULL, loading=TRUE)
```

## Arguments

<code>ic.load</code>	A loading matrix from a factor or principal components analysis, or from ICLUST.
<code>cut</code>	Do not include items in clusters with absolute loadings less than cut
<code>labels</code>	labels for each item.
<code>loading</code>	if coming from a factor analysis or ICLUST output, use <code>loading=TRUE</code> , but, if just sorting a matrix use <code>loading=FALSE</code>

## Details

When interpreting cluster or factor analysis outputs, it is useful to group the items in terms of which items have their biggest loading on each factor/cluster and then to sort the items by size of the absolute factor loading.

A stable cluster solution will be one in which the output of these cluster definitions does not vary when clusters are formed from the clusters so defined.

## Value

<code>cluster</code>	A matrix of -1, 0, 1s defining each item by the factor/cluster with the row wise largest absolute loading.
<code>load</code>	A data.frame of item numbers, item contents, and item x factor loadings.
...	

## Note

Although part of the ICLUST set of programs, this is generally more useful for factor or principal components analysis.

## Author(s)

William Revelle

## References

<http://personality-project.org/r/r.ICLUST.html>

## See Also

[ICLUST.graph](#), [ICLUST.cluster](#), [cluster.fit](#) , [VSS](#), [factor2cluster](#)

## Examples

## ICLUST of 24 mental ability tests

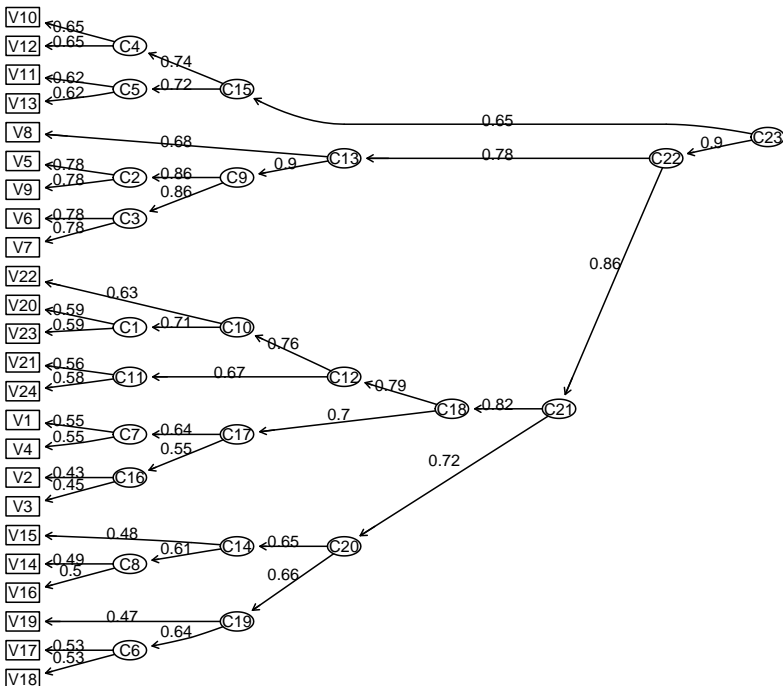


Figure 6: The Harman 24 variables may also be cluster analyzed to highlight the hierarchical structure of abilities. Compare to the hierarchical factor solution using omega (Figure 8) or the orthogonal factor solution (Figure 5)

## Description

A common data reduction technique is to cluster cases (subjects). Less common, but particularly useful in psychological research, is to cluster items (variables). This may be thought of as an alternative to factor analysis, based upon a much simpler model. The cluster model is that the correlations between variables reflect that each item loads on at most one cluster, and that items that load on those clusters correlate as a function of their respective loadings on that cluster and items that define different clusters correlate as a function of their respective cluster loadings and the intercluster correlations. Essentially, the cluster model is a Very Simple Structure factor model of complexity one (see [VSS](#)).

This function applies the ICLUST algorithm to hierarchically cluster items to form composite scales. Clusters are combined if coefficients alpha and beta will increase in the new cluster.

Alpha, the mean split half correlation, and beta, the worst split half correlation, are estimates of the reliability and general factor saturation of the test. (See also the [omega](#) function to estimate McDonald's coefficient omega.)

## Usage

```
ICLUST(r.mat, nclusters=1, alpha=3, beta=1, beta.size=4, alpha.size=3,
correct=TRUE, reverse=TRUE, beta.min=.5, output=1, digits=2, labels=NULL, cut=0,
n.iterations = 0, title="ICLUST")

#ICLUST(r.mat)      #use all defaults
#ICLUST(r.mat, nclusters = 3)    #use all defaults and if possible stop at 3 clusters
#ICLUST(r.mat, output = 3)      #long output shows clustering history
#ICLUST(r.mat, n.iterations = 3) #clean up solution by item reassignment
```

## Arguments

<b>r.mat</b>	A correlation matrix or data matrix/data.frame. (If r.mat is not square i.e, a correlation matrix, the data are correlated using pairwise deletion.)
<b>nclusters</b>	Extract clusters until nclusters remain (default = 1)
<b>alpha</b>	Apply the increase in alpha criterion (0) never or for (1) the smaller, 2) the average, or 3) the greater of the separate alphas. (default = 3)
<b>beta</b>	Apply the increase in beta criterion (0) never or for (1) the smaller, 2) the average, or 3) the greater of the separate betas. (default = 1)
<b>beta.size</b>	Apply the beta criterion after clusters are of beta.size (default = 4)
<b>alpha.size</b>	Apply the alpha criterion after clusters are of size alpha.size (default = 3)
<b>correct</b>	Correct correlations for reliability (default = TRUE)

<code>reverse</code>	Reverse negative keyed items (default = TRUE)
<code>beta.min</code>	Stop clustering if the beta is not greater than beta.min (default = .5)
<code>output</code>	1) short, 2) medium, 3 ) long output (default =1)
<code>labels</code>	vector of item content or labels
<code>cut</code>	sort cluster loadings > absolute(cut) (default = 0)
<code>n.iterations</code>	
<code>digits</code>	Precision of digits of output (default = 2)
<code>title</code>	Title for this run

## Details

Extensive documentation and justification of the algorithm is available in the original MBR 1979 <http://personality-project.org/revelle/publications/iclust.pdf> paper. Further discussion of the algorithm and sample output is available on the personality-project.org web page: <http://personality-project.org/r/r.ICLUST.html>

The results are best visualized using [ICLUST.graph](#), the results of which can be saved as a dot file for the Graphviz program. <http://www.graphviz.org/>

A common problem in the social sciences is to construct scales or composites of items to measure constructs of theoretical interest and practical importance. This process frequently involves administering a battery of items from which those that meet certain criteria are selected. These criteria might be rational, empirical, or factorial. A similar problem is to analyze the adequacy of scales that already have been formed and to decide whether the putative constructs are measured properly. Both of these problems have been discussed in numerous texts, as well as in myriad articles. Proponents of various methods have argued for the importance of face validity, discriminant validity, construct validity, factorial homogeneity, and theoretical importance.

Revelle (1979) proposed that hierarchical cluster analysis could be used to estimate a new coefficient (beta) that was an estimate of the general factor saturation of a test. More recently, Zinbarg, Revelle, Yovel and Li (2005) compared McDonald's Omega to Chronbach's alpha and Revelle's beta. They conclude that omega is the best estimate. An algorithm for estimating [omega](#) is available as part of this package.

This R version is a completely new version of ICLUST. Although early testing suggests it is stable, let me know if you have problems. Please email me if you want help with this version of ICLUST or if you desire more features.

The program currently has three primary functions: cluster, loadings, and graphics.

Clustering 24 tests of mental ability

A sample output using the 24 variable problem by Harman can be represented both graphically and in terms of the cluster order. Note that the graphic is created using GraphViz in the dot language. [ICLUST.graph](#) produces the dot code for Graphviz. Somewhat lower resolution graphs with fewer options are available in the [ICLUST.rgraph](#) function which requires Rgraphviz. Dot code can be viewed directly in Graphviz or can be tweaked using commercial software packages (e.g., OmniGraffle)

Note that for this problem, with these parameters, the data formed one large cluster. (This is consistent with the Very Simple Structure ([VSS](#)) output as well, which shows a clear one



factor solution for complexity 1 data.) See below for an example with this same data set, but with more stringent parameter settings.

To see the graphic output go to <http://personality-project.org/r/r.ICLUST.html> or use [ICLUST.rgraph](#) (requires Rgraphviz).

## Value

<code>title</code>	Name of this run
<code>results</code>	A list containing
<code>clusters</code>	a matrix of -1,0, and 1 values to define cluster membership.
<code>corrected</code>	The raw and corrected for alpha reliability cluster intercorrelations.
<code>purified</code>	A list of the cluster definitions and cluster loadings of the purified solution

## Author(s)

William Revelle  
Department of Psychology  
Northwestern University  
Evanston, Illinois  
< [revelle@northwestern.edu](mailto:revelle@northwestern.edu) >  
<http://personality-project.org/revelle.html>

## References

Revelle, W. Hierarchical Cluster Analysis and the Internal Structure of Tests. *Multivariate Behavioral Research*, 1979, 14, 57-74. <http://personality-project.org/revelle/publications/iclust.pdf>  
See also more extensive documentation at <http://personality-project.org/r/r.ICLUST.html>

## See Also

[ICLUST.graph](#), [ICLUST.cluster](#), [cluster.fit](#) , [VSS](#), [omega](#)

## Examples

```
## Not run:
test.data <- Harman74.cor$cov
ic.out <- ICLUST(test.data)          #use all defaults
out.file <- file.choose(new=TRUE)    #create a new file to write the plot commands to
ICLUST.graph(ic.out,out.file,title = "ICLUST of Harman's 24 mental variables" )

ic.out <- ICLUST(test.data,nclusters =3) #use all defaults and if possible stop at 3 clusters
ICLUST.graph(ic.out,out.file,title = "ICLUST of 24 mental variables with forced 3 cluster solution")

ICLUST(test.data, output =3)         #long output shows clustering history

ic.out <- ICLUST(test.data,,nclusters=4, n.iterations =3) #clean up solution by item reassignment
ICLUST.graph(ic.out,out.file,title = "ICLUST of 24 mental variables with forced 4 cluster solution")
ic.out      #shows the output on the console
```

```

## End(Not run)

#produces this output
#ICLUST(Harman74.cor$cov)
#$title
#[1] "ICLUST"
#
#clusters
#      VisualPerception          Cubes      PaperFormBoard          Flags      GeneralI
#              1              1              1              1
# ParagraphComprehension SentenceCompletion WordClassification WordMeaning
#              1              1              1              1
#              Code      CountingDots StraightCurvedCapitals WordRecognition      NumberR
#              1              1              1              1
#      FigureRecognition      ObjectNumber      NumberFigure      FigureWord
#              1              1              1              1
#      NumericalPuzzles      ProblemReasoning      SeriesCompletion      ArithmeticProblems
#              1              1              1              1
#
#corrected
#      [,1]
#[1,]      1
#
#loadings
#              [,1]
#VisualPerception      0.57
#Cubes                  0.36
#PaperFormBoard        0.40
#Flags                  0.46
#GeneralInformation    0.62
#ParagraphComprehension 0.62
#SentenceCompletion    0.60
#WordClassification    0.63
#WordMeaning           0.62
#Addition              0.43
#Code                  0.54
#CountingDots          0.44
#StraightCurvedCapitals 0.57
#WordRecognition       0.41
#NumberRecognition     0.38
#FigureRecognition     0.50
#ObjectNumber          0.45
#NumberFigure          0.51
#FigureWord            0.44
#Deduction             0.59
#NumericalPuzzles      0.58
#ProblemReasoning      0.58
#SeriesCompletion      0.66
#ArithmeticProblems    0.62
#
#fit
#$fit
#$fit$clusterfit
#[1] 0.78

```

```

#
#fit$factorfit
#[1] 0.78
#
#
#$results
#   Item/Cluster Item/Cluster similarity correlation alpha1 alpha2 beta1 beta2 size1 size2 rbar1 rbar2
#C1          V23          V20      1.00      0.51  0.51  0.51  0.51  0.51      1      1  0.51  0.51 0
#C2           V9           V5      1.00      0.72  0.72  0.72  0.72  0.72      1      1  0.72  0.72 0
#C3           V7           V6      1.00      0.72  0.72  0.72  0.72  0.72      1      1  0.72  0.72 0
#C4          V12          V10      1.00      0.58  0.58  0.58  0.58  0.58      1      1  0.58  0.58 0
#C5          V13          V11      1.00      0.54  0.54  0.54  0.54  0.54      1      1  0.54  0.54 0
#C6          V18          V17      1.00      0.45  0.45  0.45  0.45  0.45      1      1  0.45  0.45 0
#C7           V4           V1      0.99      0.47  0.47  0.48  0.47  0.48      1      1  0.47  0.48 0
#C8          V16          V14      0.98      0.41  0.43  0.41  0.43  0.41      1      1  0.43  0.41 0
#C9           C2           C3      0.93      0.78  0.84  0.84  0.84  0.84      2      2  0.72  0.72 0
#C10          C1          V22      0.91      0.56  0.67  0.56  0.68  0.56      2      1  0.51  0.56 0
#C11          V21          V24      0.87      0.45  0.51  0.53  0.51  0.53      1      1  0.51  0.53 0
#C12          C10          C11      0.86      0.58  0.74  0.62  0.72  0.62      3      2  0.49  0.45 0
#C13           C9           V8      0.84      0.64  0.90  0.64  0.88  0.64      4      1  0.69  0.64 0
#C14           C8          V15      0.84      0.41  0.58  0.41  0.58  0.41      2      1  0.41  0.41 0
#C15           C5           C4      0.82      0.59  0.70  0.74  0.70  0.73      2      2  0.54  0.58 0
#C16           V3           V2      0.81      0.32  0.41  0.38  0.41  0.38      1      1  0.41  0.38 0
#C17          C16           C7      0.81      0.45  0.48  0.64  0.48  0.64      2      2  0.32  0.47 0
#C18          C12          C17      0.81      0.59  0.79  0.67  0.73  0.62      5      4  0.43  0.34 0
#C19          V19           C6      0.80      0.40  0.40  0.62  0.40  0.62      1      2  0.40  0.45 0
#C20          C19          C14      0.77      0.49  0.64  0.64  0.57  0.58      3      3  0.38  0.37 0
#C21          C18          C20      0.74      0.58  0.83  0.74  0.74  0.66      9      6  0.35  0.32 0
#C22          C21          C13      0.70      0.62  0.86  0.90  0.73  0.78     15      5  0.29  0.64 0
#C23          C22          C15      0.65      0.55  0.90  0.79  0.77  0.74     20      4  0.31  0.49 0
#   beta rbar size
#C1  0.68 0.51    2
#C2  0.84 0.72    2
#C3  0.84 0.72    2
#C4  0.73 0.58    2
#C5  0.70 0.54    2
#C6  0.62 0.45    2
#C7  0.64 0.47    2
#C8  0.58 0.41    2
#C9  0.88 0.69    4
#C10 0.72 0.49    3
#C11 0.62 0.45    2
#C12 0.73 0.43    5
#C13 0.78 0.64    5
#C14 0.58 0.37    3
#C15 0.74 0.49    4
#C16 0.48 0.32    2
#C17 0.62 0.34    4
#C18 0.74 0.35    9
#C19 0.57 0.38    3
#C20 0.66 0.32    6
#C21 0.73 0.29   15
#C22 0.77 0.31   20

```

```

#C23 0.71 0.30 24
#
# $cor
#      [,1]
# [1,]    1
#
# $alpha
# [1] 0.91
#
# $size
# [1] 24
#
# $sorted
# $sorted$sorted
#
# item content cluster loadings
#SeriesCompletion 23 SeriesCompletion 1 0.66
#WordClassification 8 WordClassification 1 0.63
#GeneralInformation 5 GeneralInformation 1 0.62
#ParagraphComprehension 6 ParagraphComprehension 1 0.62
#WordMeaning 9 WordMeaning 1 0.62
#ArithmeticProblems 24 ArithmeticProblems 1 0.62
#SentenceCompletion 7 SentenceCompletion 1 0.60
#Deduction 20 Deduction 1 0.59
#NumericalPuzzles 21 NumericalPuzzles 1 0.58
#ProblemReasoning 22 ProblemReasoning 1 0.58
#VisualPerception 1 VisualPerception 1 0.57
#StraightCurvedCapitals 13 StraightCurvedCapitals 1 0.57
#Code 11 Code 1 0.54
#NumberFigure 18 NumberFigure 1 0.51
#FigureRecognition 16 FigureRecognition 1 0.50
#Flags 4 Flags 1 0.46
#ObjectNumber 17 ObjectNumber 1 0.45
#CountingDots 12 CountingDots 1 0.44
#FigureWord 19 FigureWord 1 0.44
#Addition 10 Addition 1 0.43
#WordRecognition 14 WordRecognition 1 0.41
#PaperFormBoard 3 PaperFormBoard 1 0.40
#NumberRecognition 15 NumberRecognition 1 0.38
#Cubes 2 Cubes 1 0.36
#
#
# $p.fit
# $p.fit$clusterfit
# [1] 0.78
#
# $p.fit$factorfit
# [1] 0.78
#
#
# $p.sorted
# $p.sorted$sorted
#
# item content cluster loadings
#SeriesCompletion 23 SeriesCompletion 1 0.66

```

#WordClassification	8	WordClassification	1	0.63
#GeneralInformation	5	GeneralInformation	1	0.62
#ParagraphComprehension	6	ParagraphComprehension	1	0.62
#WordMeaning	9	WordMeaning	1	0.62
#ArithmeticProblems	24	ArithmeticProblems	1	0.62
#SentenceCompletion	7	SentenceCompletion	1	0.60
#Deduction	20	Deduction	1	0.59
#NumericalPuzzles	21	NumericalPuzzles	1	0.58
#ProblemReasoning	22	ProblemReasoning	1	0.58
#VisualPerception	1	VisualPerception	1	0.57
#StraightCurvedCapitals	13	StraightCurvedCapitals	1	0.57
#Code	11	Code	1	0.54
#NumberFigure	18	NumberFigure	1	0.51
#FigureRecognition	16	FigureRecognition	1	0.50
#Flags	4	Flags	1	0.46
#ObjectNumber	17	ObjectNumber	1	0.45
#CountingDots	12	CountingDots	1	0.44
#FigureWord	19	FigureWord	1	0.44
#Addition	10	Addition	1	0.43
#WordRecognition	14	WordRecognition	1	0.41
#PaperFormBoard	3	PaperFormBoard	1	0.40
#NumberRecognition	15	NumberRecognition	1	0.38
#Cubes	2	Cubes	1	0.36
#				
#				
##\$purified				
##\$purified\$cor				
# [ ,1]				
#[1,] 1				
#				
##\$purified\$sd				
#[1] 13.79				
#				
##\$purified\$corrected				
# [ ,1]				
#[1,] 0.91				
#				
##\$purified\$size				
#[1] 24				
#				
#				

---

`irt.item.diff.rasch` *Simple function to estimate item difficulties using IRT concepts*

---

## Description

Steps toward a very crude and preliminary IRT program. These two functions estimate item difficulty and discrimination parameters.

## Usage

```
irt.item.diff.rasch(items)
irt.discrim(item.diff,theta,items)
```

## Arguments

<code>items</code>	a matrix of items
<code>item.diff</code>	a vector of item difficulties (found by <code>irt.item.diff</code> )
<code>theta</code>	ability estimate from <code>irt.person.theta</code>

## Details

Item Response Theory (aka "The new psychometrics") models individual responses to items with a logistic function and an individual ( $\theta$ ) and item difficulty ( $\text{diff}$ ) parameter.

`irt.item.diff.rasch` finds item difficulties with the assumption of  $\theta=0$  for all subjects and that all items are equally discriminating.

`irt.discrim` takes those difficulties and  $\theta$  estimates from [irt.person.rasch](#) to find item discrimination ( $\beta$ ) parameters.

A far better package with these features is the `ltm` package. The IRT functions in the `psych`-package are for pedagogical rather than production purposes. They are believed to be accurate, but are not guaranteed. They do seem to be slightly more robust to missing data structures associated with SAPA data sets than the `ltm` package.

## Value

a vector of item difficulties or item discriminations.

## Note

Under development. Not recommended for public consumption.

## Author(s)

William Revelle

## References

## See Also

[irt.person.rasch](#)

## Examples

---

<code>irt.1p</code>	<i>Item Response Theory estimate of theta (ability) using a Rasch (like) model</i>
---------------------	--

---

## Description

Item Response Theory models individual responses to items by estimating individual ability (theta) and item difficulty (diff) parameters. This is an early and crude attempt to capture this modeling procedure.

## Usage

```
irt.person.rasch(diff, items)
irt.0p(items,possible=20)
irt.1p(delta,items)
irt.2p(delta,beta,items)
```

## Arguments

<b>diff</b>	A vector of item difficulties –probably taken from <code>irt.item.diff.rasch</code>
<b>items</b>	A matrix of 0,1 items nrow = number of subjects, ncol = number of items
<b>possible</b>	Number of items in the scale – used to determine values of all wrong or all right
<b>delta</b>	delta is the same as diff and is the item difficulty parameter
<b>beta</b>	beta is the item discrimination parameter found in <a href="#">irt.discrim</a>

## Details

A very preliminary IRT estimation procedure. Given scores  $x_{ij}$  for  $i$ th individual on  $j$ th item

Classical Test Theory ignores item difficulty and defines ability as expected score :  $ability_i = \theta(i) = x(i)$ . A zero parameter model rescales these mean scores from 0 to 1 to a quasi logistic scale ranging from - 4 to 4 This is merely a non-linear transform of the raw data to reflect a logistic mapping.

Basic 1 parameter (Rasch) model considers item difficulties ( $\delta_j$ ):  $p(\text{correct on item } j \text{ for the } i\text{th subject} | \theta_i, \delta_j) = 1/(1+\exp(\delta_j - \theta_i))$  If we have estimates of item difficulty ( $\delta$ ), then we can find  $\theta_i$  by optimization

Two parameter model adds item sensitivity ( $\beta_j$ ):  $p(\text{correct on item } j \text{ for subject } i | \theta_i, \delta_j, \beta_j) = 1/(1+\exp(\beta_j * (\delta_j - \theta_i)))$  Estimate  $\delta$ ,  $\beta$ , and  $\theta$  to maximize fit of model to data.

The procedure used here is to first find the item difficulties assuming  $\theta = 0$  Then find  $\theta$  given those  $\delta$ s Then find  $\beta$  given  $\delta$  and  $\theta$ .

This is not an "official" way to do IRT, but is useful for basic item development.

**Value**

a data.frame with estimated ability (theta) and quality of fit. (for irt.person.rasch)  
a data.frame with the raw means, theta0, and the number of items completed

**Note**

Not recommended for serious use. This code is under development.

**Author(s)**

William Revelle

**References****See Also**

[irt.item.diff.rasch](#)

**Examples**

---

<code>item.sim</code>	<i>Generate simulated data structures for circumplex or simple structure</i>
-----------------------	--

---

**Description**

Rotations of factor analysis and principal components analysis solutions typically try to represent correlation matrices as simple structured. An alternative structure, appealing to some, is a circumplex structure where the variables are uniformly spaced on the perimeter of a circle in a two dimensional space. Generating these data is straightforward, and is useful for exploring alternative solutions to affect and personality structure.

**Usage**

```
item.sim(nvar = 72, nsub = 500, circum = FALSE, xloading = 0.6, yloading = 0.6, gloading = 0, x  
circ.sim(nvar = 72, nsub = 500, circum = TRUE, xloading = 0.6, yloading = 0.6, gloading = 0, x
```



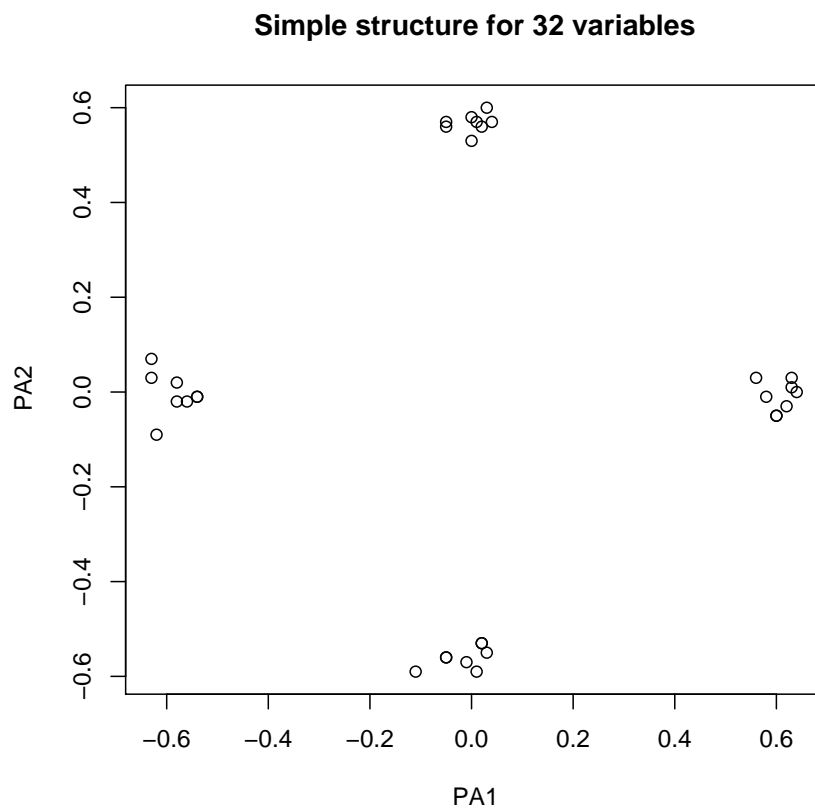


Figure 7: Simple structure is a goal for many factor rotation and extraction procedures. Data may be generated with simple or circumplex structure with varying degrees of skew and correlation. (Compare with Figure 1)

## Arguments

<code>nvar</code>	Number of variables to simulate
<code>nsub</code>	Number of subjects to simulate
<code>circum</code>	<code>circum=TRUE</code> is circumplex structure, <code>FALSE</code> is simple structure
<code>xloading</code>	the average loading on the first dimension
<code>yloading</code>	Average loading on the second dimension
<code>gloading</code>	Average loading on a general factor (default=0)
<code>xbias</code>	To introduce skew, how far off center is the first dimension
<code>ybias</code>	To introduce skew on the second dimension
<code>categorical</code>	continuous or categorical variables.
<code>low</code>	values less than low are forced to low
<code>high</code>	values greater than high are forced to high
<code>truncate</code>	Change all values less than cutpoint to cutpoint.
<code>cutpoint</code>	What is the cutpoint

## Details

This simulation was originally developed to compare the effect of skew on the measurement of affect (see Rafaeli and Revelle, 2005). It has been extended to allow for a general simulation of affect or personality items with either a simple structure or a circumplex structure. Items can be continuous normally distributed, or broken down into  $n$  categories (e.g., -2, -1, 0, 1, 2). Items can be distorted by limiting them to these ranges, even though the items have a mean of (e.g., 1).

## Value

A data matrix of (`nsub`) subjects by (`nvar`) variables.

## Author(s)

William Revelle

## References

Variations of a routine used in Rafaeli and Revelle, 2006; Rafaeli, E. & Revelle, W. (2006). A premature consensus: Are happiness and sadness truly opposite affects? Motivation and Emotion.

Acton, G. S. and Revelle, W. (2004) Evaluation of Ten Psychometric Criteria for Circumplex Structure. Methods of Psychological Research Online, Vol. 9, No. 1 [http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110\\_10.pdf](http://www.dgps.de/fachgruppen/methoden/mpr-online/issue22/mpr110_10.pdf)

## See Also

See Also the implementation in this to generate numerous simulations. [circ.simulation](#), [circ.tests](#)

## Examples

```
round(cor(circ.sim(nvar=8,nsup=200)),2)
plot(factor.pa(circ.sim(16,500),2)$loadings) #circumplex structure
#
#
plot(factor.pa(item.sim(16,500),2)$loadings) #simple structure
#
```

---

kurtosi	<i>Kurtosis of a vector, matrix, or data frame</i>
---------	--

---

## Description

Find the kurtosis of a vector, matrix, or dataframe.

## Usage

```
kurtosi(x, na.rm = TRUE)
```

## Arguments

x	vector, matrix, or data frame
na.rm	na.rm =TRUE removes missing data from the column

## Details

Kurtosis in the E1071 package finds the kurtosis for a single vector. This does it for matrices and dataframes. Used in the describe function.

## Value

kurtosi	a vector of the kurtosis for each column of the matrix
---------	--

## Note

The mean function supplies means for the columns of a data.frame, but the overall mean for a matrix. Mean will throw a warning for non-numeric data, but colMeans stops with non-numeric data. Thus, the function uses either mean (for data frames) or colMeans (for matrices). This is true for skew and kurtosi as well.

## Author(s)

William Revelle

## See Also

[skew](#), [describe](#)

## Examples

```
round(kurtosi(attitude),2)
```

---

<code>make.hierarchical</code>	<i>Create a population or sample correlation matrix with hierachical structure.</i>
--------------------------------	---

---

## Description

Create a population hierarchical correlation matrix from a set of factor loadings and factor intercorrelations. Samples of size n may be then be drawn from this population. Return either the sample data, sample correlations, or population correlations. This is used to create sample data sets for instruction and demonstration.

## Usage

```
make.hierarchical(gload=NULL, fload=NULL, n = 0, raw = FALSE)
```

## Arguments

<code>gload</code>	Loadings of group factors on a general factor
<code>fload</code>	Loadings of items on the group factor
<code>n</code>	Number of subjects to generate: N=0 => population values
<code>raw</code>	raw=TRUE, report the raw data, raw=FALSE, report the sample correlation matrix.

## Details

Many personality and cognitive tests have a hierarchical factor structure. For demonstration purposes, it is useful to be able to create such matrices, either with population values, or sample values.

Given a matrix of item factor loadings (`fload`) and of loadings of these factors on a general factor (`gload`), we create a population correlation matrix by using the general factor law ( $R = F' \theta F$  where  $\theta = g'g$ ).

To create sample values, we use the [mvrnorm](#) function from MASS.

The default is to return population correlation matrices. Sample correlation matrices are generated if `n > 0`. Raw data are returned if `raw = TRUE`.

The default values for `gload` and `fload` create a data matrix discussed by Jensen and Weng, 1994.

## Value

a matrix of correlations or a data matrix

## Author(s)

William Revelle

## References

<http://personality-project.org/r/r.omega.html>

Jensen, A.R., Weng, L.J. (1994) What is a Good g? *Intelligence*, 18, 231-258.

## See Also

[omega](#), [schmid](#), [ICLUST](#), [VSS](#), [mvrnorm](#)

## Examples

```
## Not run:
gload <- gload<-matrix(c(.9,.8,.7),nrow=3)    # a higher order factor matrix
fload <-matrix(c(                                #a lower order (oblique) factor matrix
  .8,0,0,
  .7,0,.0,
  .6,0,.0,
  0,.7,.0,
  0,.6,.0,
  0,.5,0,
  0,0,.6,
  0,0,.5,
  0,0,.4),    ncol=3,byrow=TRUE)

jensen <- make.hierarchical(gload,fload)    #the test set used by omega
round(jensen,2)
## End(Not run)
```

---

`mat.regress`

*Multiple Regression from matrix input*

---

## Description

This function extracts subsets of variables (x and y) from a correlation matrix (m) and then find the multiple correlation and beta weights of the (x) set predicting each member of the (y) set.

## Usage

```
mat.regress(m, x, y,digits=2)
```

## Arguments

<code>m</code>	a matrix of correlations
<code>x</code>	the column numbers of the x set (e.g., <code>c(1,3,5)</code> )
<code>y</code>	the column numbers of the y set (e.g., <code>c(2,4,6)</code> )
<code>digits</code>	round the answer to digits

## Details

Although it is more common to calculate multiple regression from raw data, it is, of course, possible to do so from a set of correlations. The input to the function is a square covariance or correlation matrix, as well as the column numbers of the x (predictor) and y (criterion) variables.

The output is a set of multiple correlations, one for each dependent variable in the y set.

A typical use in the SAPA project is to form item composites by clustering or factoring (see [ICLUST](#), [principal](#)), extract the clusters from these results ([factor2cluster](#)), and then form the composite correlation matrix using [cluster.cor](#). The variables in this reduced matrix may then be used in multiple R procedures using `mat.regress`.

Although the overall matrix can have missing correlations, the correlations in the subset of the matrix used for prediction must exist.

## Value

<code>beta</code>	the beta weights for each variable in X for each variable in Y
<code>R</code>	The multiple R for each equation (the amount of change a unit in the predictor set leads to in the criterion set).
<code>R2</code>	The multiple R2 (% variance accounted for) for each equation

## Author(s)

William Revelle

Maintainer: William Revelle <revell@northwestern.edu>

## See Also

[cluster.cor](#), [factor2cluster](#), [principal](#), [ICLUST](#)

## Examples

```
## Not run:
test.data <- Harman74.cor$cov      #24 mental variables
#choose 3 of them to regress against another 4 -- arbitrary choice of variables
print(mat.regress(test.data,c(1,2,3),c(4,5,10,12)),digits=2)
## End(Not run)
#gives this output
#print(mat.regress(test.data,c(1,2,3),c(4,5,10,12)),digits=2)
```

```

##$beta
#
#           Flags GeneralInformation Addition CountingDots
#VisualPerception 0.40           0.22    0.16    0.30
#Cubes           0.06           0.18    0.06    0.05
#PaperFormBoard  0.12           0.10   -0.16    0.00
#
##$R
#
#           Flags GeneralInformation           Addition           CountingDots
#           0.49           0.38           0.18           0.32
#
##$R2
#
#           Flags GeneralInformation           Addition           CountingDots
#           0.24           0.15           0.03           0.10
#
#
#

```

---

matrix.addition	<i>A function to add two vectors or matrices</i>
-----------------	--

---

**Description**

It is sometimes convenient to add two vectors or matrices in an operation analogous to matrix multiplication. For matrices nXm and mYp, the matrix sum of the i,jth element of nSp = sum(over m) of iXm + mYj.

**Usage**

x %+% y

**Arguments**

x                    a n by m matrix (or vector if m= 1)  
y                    a m by p matrix (or vector if m = 1)

**Details**

Used in such problems as Thurstonian scaling. Although not technically matrix addition, as pointed out by Krus, there are many applications where the sum or difference of two vectors or matrices is a useful operation. This can be done, of course, through

**Value**

a n by p matix of sums

**Author(s)**

William Revelle

## References

Krus, D. J. (2001) Matrix addition. Journal of Visual Statistics, 1, (February, 2001).

## Examples

```
x <- seq(1,4)
z <- x %+% -t(x)
x
z
x <- matrix(seq(1,6),ncol=2)
y <- matrix(seq(1,10),nrow=2)
z <- x %+% y
x
y
z
```

---

`multi.hist`

*Multiple histograms on one screen*

---

## Description

Given a matrix or data.frame, produce histograms for each variable in a "matrix" form.

## Usage

```
multi.hist(x)
```

## Arguments

`x`                      matrix or data.frame

## Author(s)

William Revelle  
Northwestern University  
Evanston, Illinois  
< [revelle@northwestern.edu](mailto:revelle@northwestern.edu) >  
<http://personality-project.org/revelle.html>

## Examples

```
multi.hist(attitude) #see graphics window
```



### Harman's 24 tests of mental ability

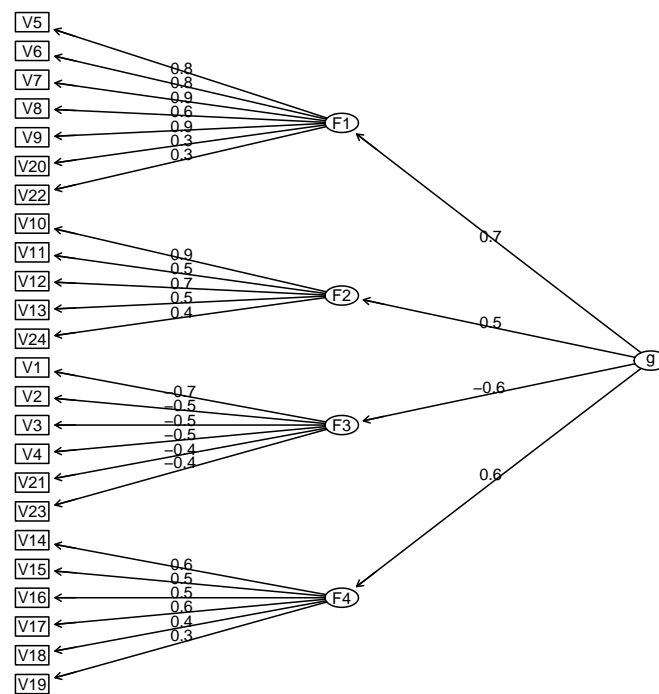


Figure 8: Hierarchical factor solutions are typical in the ability domain where a g factor is thought to reflect the correlations among lower level factors. An alternative transformation is to orthogonalize the g factor from the residual group factors using the Schmid-Leiman transformation (Figure 9)

### Harman's 24 tests of mental ability

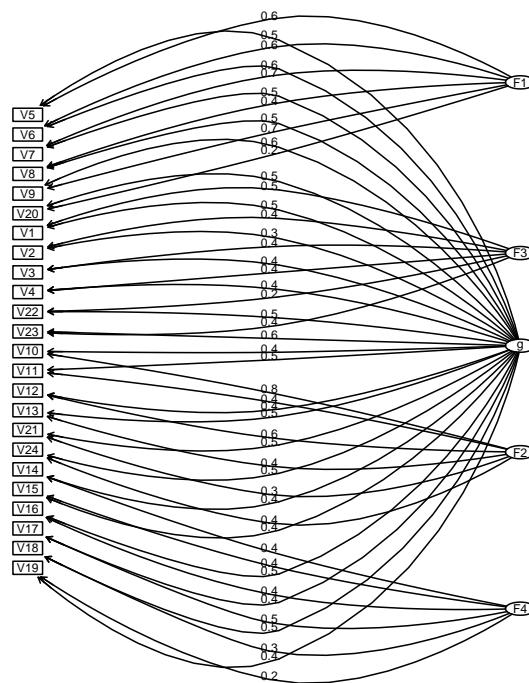


Figure 9: An alternative to the standard hierarchical factor solutions which are typical in the ability domain is to orthogonalize the g factor from the residual group factors using the Schmid-Leiman transformation. For the hierarchical solution, see Figure 8

## Description

Hierarchical factor structures represent the correlations between variables in terms of a smaller set of correlated factors which themselves can be represented by a higher order factor.

Two alternative solutions to such structures are found by the omega function. The correlated factors solutions represents the effect of the higher level, general factor, through its effect on the correlated factors. The other representation makes use of the Schmid Leiman transformation to find the direct effect of the general factor upon the original variables as well as the effect of orthogonal residual group factors upon the items.

Graphic presentations of these two alternatives are helpful in understanding the structure. omega.graph draws both such structures. Graphs are drawn directly onto the graphics window or expressed in “dot” commands for conversion to graphics using implementations of Graphviz.

## Usage

```
omega.graph(om.results, out.file = NULL, sl = TRUE, labels = NULL, size = c(8, 6), node.font =
```

## Arguments

<code>om.results</code>	The output from the omega function
<code>out.file</code>	Optional output file for off line analysis using Graphviz
<code>sl</code>	Orthogonal clusters using the Schmid-Leiman transform (sl=TRUE) or oblique clusters
<code>labels</code>	variable labels
<code>size</code>	size of graphics window
<code>node.font</code>	What font to use for the items
<code>edge.font</code>	What font to use for the edge labels
<code>rank.direction</code>	Defaults to left to right
<code>digits</code>	Precision of labels
<code>title</code>	Figure title
<code>...</code>	Other options to pass into the graphics packages

## Details

Requires the Rgraphviz package. omega requires the GPArotation package.

## Value

`clust.graph` A graph object

## Note

Requires rgraphviz. – omega requires GPARotation

## Author(s)

<http://personality-project.org/revelle.html>

Maintainer: William Revelle `<revelle@northwestern.edu>`

## References

<http://personality-project.org/r/r.omega.html>

Revelle, W. (1979). Hierarchical cluster analysis and the internal structure of tests. Multivariate Behavioral Research, 14, 57-74. (<http://personality-project.org/revelle/publications/iclust.pdf>)

Zinbarg, R.E., Revelle, W., Yovel, I., & Li, W. (2005). Cronbach's Alpha, Revelle's Beta, McDonald's Omega: Their relations with each and two alternative conceptualizations of reliability. Psychometrika, 70, 123-133. <http://personality-project.org/revelle/publications/zinbarg.revelle.pmet.05.pdf>

Zinbarg, R., Yovel, I., Revelle, W. & McDonald, R. (2006). Estimating generalizability to a universe of indicators that all have one attribute in common: A comparison of estimators for omega. Applied Psychological Measurement, 30, 121-144. DOI: 10.1177/0146621605278814 <http://apm.sagepub.com/cgi/reprint/30/2/121>

## See Also

[omega](#), [make.hierarchical](#), [ICLUST.rgraph](#)

## Examples

```
#24 mental tests from Holzinger-Swineford-Harman
if(require(GPARotation) ) {om24 <- omega(Harman74.cor$cov,4) } #run omega
if(require(Rgraphviz) ){om24pn <- omega.graph(om24,sl=FALSE)} #show the structure
#
#example hierarchical structure from Jensen and Weng
if(require(GPARotation) ) {jen.omega <- omega(make.hierarchical())}
if(require(Rgraphviz) ) {om.jen <- omega.graph(jen.omega,sl=FALSE) }
```

---

omega

*Calculate the omega estimate of factor saturation*

---

## Description

McDonald has proposed coefficient omega as an estimate of the general factor saturation of a test. One way to find omega is to do a factor analysis of the original data set, rotate the factors obliquely, do a Schmid Leiman transformation, and then find omega. This function estimates omega as suggested by McDonald by using hierarchical factor analysis (following Jensen).

## Usage

```
omega(m, nfactors, pc = "mle", key = NULL, ...)
```

## Arguments

<b>m</b>	A correlation matrix or a data.frame/matrix of data
<b>nfactors</b>	Number of factors believed to be group factors
<b>pc</b>	pc="pa" for principal axes, pc="pc" for principal components, pc="mle" for maximum likelihood.
<b>key</b>	a vector of +/- 1s to specify the direction of scoring of items. The default is to assume all items are positively keyed, but if some items are reversed scored, then key should be specified.
<b>...</b>	Allows additional parameters to be passed through to the factor routines

## Details

"Many scales are assumed by their developers and users to be primarily a measure of one latent variable. When it is also assumed that the scale conforms to the effect indicator model of measurement (as is almost always the case in psychological assessment), it is important to support such an interpretation with evidence regarding the internal structure of that scale. In particular, it is important to examine two related properties pertaining to the internal structure of such a scale. The first property relates to whether all the indicators forming the scale measure a latent variable in common.

The second internal structural property pertains to the proportion of variance in the scale scores (derived from summing or averaging the indicators) accounted for by this latent variable that is common to all the indicators (Cronbach, 1951; McDonald, 1999; Revelle, 1979). That is, if an effect indicator scale is primarily a measure of one latent variable common to all the indicators forming the scale, then that latent variable should account for the majority of the variance in the scale scores. Put differently, this variance ratio provides important information about the sampling fluctuations when estimating individuals' standing on a latent variable common to all the indicators arising from the sampling of indicators (i.e., when dealing with either Type 2 or Type 12 sampling, to use the terminology of Lord, 1956). That is, this variance proportion can be interpreted as the square of the correlation between the scale score and the latent variable common to all the indicators in the infinite universe of indicators of which the scale indicators are a subset. Put yet another way, this variance ratio is important both as reliability and a validity coefficient. This is a reliability issue as the larger this variance ratio is, the more accurately one can predict an individual's relative standing on the latent variable common to all the scale's indicators based on his or her observed scale score. At the same time, this variance ratio also bears on the construct validity of the scale given that construct validity encompasses the internal structure of a scale." (Zinbarg, Yovel, Revelle, and McDonald, 2006).

McDonald has proposed coefficient omega as an estimate of the general factor saturation of a test. Zinbarg, Revelle, Yovel and Li (2005) <http://personality-project.org/revelle/publications/zinbarg.revelle.pmet.05.pdf> compare McDonald's Omega to Cronbach's alpha and Revelle's beta. They conclude that omega is the best estimate. (See also Zinbarg et al., 2006)

One way to find omega is to do a factor analysis of the original data set, rotate the factors obliquely, do a Schmid-Leiman ([schmid](#)) transformation, and then find omega. Here we present code to do that.

Omega differs as a function of how the factors are estimated. Three options are available, `pc="pa"` does a principle axes factor analysis ([factor.pa](#)), `pc="mle"` uses the factanal function, and `pc="pc"` does a principal components analysis ([principal](#)).

Output from omega can be shown using the [omega.graph](#) function.

Beta, an alternative to omega, is defined as the worst split half reliability. It can be estimated by using [ICLUST](#) (a hierarchical clustering algorithm originally developed for main frames and written in Fortran and that is now available in R. (For a very complimentary review of why the ICLUST algorithm is useful in scale construction, see Cooksey and Soutar, 2005).

## Value

<code>alpha</code>	Cronbach's alpha
<code>schmid</code>	The Schmid Leiman transformed factor matrix and associated matrices
<code>schmid\$sl</code>	The g factor loadings as well as the residualized factors
<code>schmid\$orthog</code>	Varimax rotated solution of the original factors
<code>schmid\$oblique</code>	The oblimin transformed factors
<code>schmid\$fcor</code>	the correlation matrix of the oblique factors
<code>schmid\$gloading</code>	The loadings on the higher order, g, factor of the oblimin factors

## Note

Requires the GPArotation package

## Author(s)

<http://personality-project.org/revelle.html>

Maintainer: William Revelle < [revelle@northwestern.edu](mailto:revelle@northwestern.edu) >

## References

<http://personality-project.org/r/r.omega.html>

Revelle, W. (1979). Hierarchical cluster analysis and the internal structure of tests. Multivariate Behavioral Research, 14, 57-74. (<http://personality-project.org/revelle/publications/iclust.pdf>)

Zinbarg, R.E., Revelle, W., Yovel, I., & Li, W. (2005). Cronbach's Alpha, Revelle's Beta, McDonald's Omega: Their relations with each and two alternative conceptualizations of reliability. Psychometrika. 70, 123-133. <http://personality-project.org/revelle/publications/zinbarg.revelle.pmet.05.pdf>

Zinbarg, R., Yovel, I., Revelle, W. & McDonald, R. (2006). Estimating generalizability to a universe of indicators that all have one attribute in common: A comparison of estimators for omega. *Applied Psychological Measurement*, 30, 121-144. DOI: 10.1177/0146621605278814  
<http://apm.sagepub.com/cgi/reprint/30/2/121>

## See Also

[omega.graph](#), [ICLUST](#), [ICLUST.graph](#), [VSS](#), [schmid](#), [make.hierarchical](#)

## Examples

```
## Not run:
test.data <- Harman74.cor$cov
my.omega <- omega(test.data,3)
print(my.omega,digits=2)
#
jen.omega <- omega(make.hierarchical())
## End(Not run)
#produces this output

#$omega
#[1] 0.64
#
#$alpha
#[1] 0.91
#
#$schmid
#$schmid$sl
#
# g factor Factor1 Factor2 Factor3 h2 u2
#VisualPerception 0.53 0.018 0.4688 0.02089 0.494 0.51
#Cubes 0.34 0.022 0.3029 0.04544 0.209 0.79
#PaperFormBoard 0.38 0.033 0.3971 0.18505 0.398 0.60
#Flags 0.43 0.109 0.3233 0.06148 0.261 0.74
#GeneralInformation 0.57 0.564 0.0078 0.09900 0.606 0.39
#ParagraphComprehension 0.57 0.599 0.0244 0.02960 0.671 0.33
#SentenceCompletion 0.56 0.624 0.0352 0.02783 0.730 0.27
#WordClassification 0.56 0.394 0.1343 0.09255 0.341 0.66
#WordMeaning 0.58 0.637 0.0135 0.06034 0.762 0.24
#Addition 0.35 0.047 0.0939 0.81706 0.858 0.14
#Code 0.44 0.061 0.1442 0.44377 0.300 0.70
#CountingDots 0.37 0.100 0.1576 0.57732 0.491 0.51
#StraightCurvedCapitals 0.50 0.036 0.2677 0.33168 0.301 0.70
#WordRecognition 0.34 0.127 0.1506 0.10015 0.093 0.91
#NumberRecognition 0.32 0.060 0.2005 0.07972 0.105 0.90
#FigureRecognition 0.44 0.022 0.4080 0.00192 0.374 0.63
#ObjectNumber 0.37 0.064 0.1769 0.22163 0.139 0.86
#NumberFigure 0.43 0.076 0.3290 0.26170 0.339 0.66
#FigureWord 0.37 0.053 0.2431 0.10447 0.151 0.85
#Deduction 0.53 0.231 0.2814 0.00299 0.277 0.72
#NumericalPuzzles 0.50 0.025 0.2877 0.30211 0.301 0.70
#ProblemReasoning 0.52 0.222 0.2840 0.00067 0.272 0.73
#SeriesCompletion 0.59 0.198 0.3304 0.07553 0.325 0.68
#ArithmeticProblems 0.52 0.211 0.1106 0.40982 0.320 0.68
```

```

#
##$schmid$orthog
#
#VisualPerception      Factor1 Factor2  Factor3
#Cubes                 0.025   0.702 -0.02336
#PaperFormBoard        0.030   0.454 -0.05080
#Flags                 0.045   0.595 -0.20689
#GeneralInformation     0.149   0.484 -0.06874
#ParagraphComprehension 0.771  -0.012  0.11068
#SentenceCompletion     0.817   0.037 -0.03310
#WordClassification    0.852  -0.053  0.03111
#WordMeaning           0.538   0.201  0.10348
#Addition              0.870   0.020 -0.06746
#Code                  0.065  -0.141  0.91350
#CountingDots          0.083   0.216  0.49615
#StraightCurvedCapitals -0.136  0.236  0.64546
#WordRecognition       0.049   0.401  0.37082
#NumberRecognition     0.173   0.226  0.11197
#FigureRecognition     0.082   0.300  0.08912
#ObjectNumber          -0.029   0.611  0.00214
#NumberFigure          0.087   0.265  0.24779
#FigureWord            -0.104   0.493  0.29259
#Deduction             0.073   0.364  0.11681
#NumericalPuzzles      0.315   0.421 -0.00334
#ProblemReasoning      0.035   0.431  0.33777
#SeriesCompletion      0.303   0.425 -0.00074
#ArithmeticProblems    0.270   0.495  0.08445
#
##$schmid$fcor
#      [,1] [,2] [,3]
#[1,] 1.00 0.51 0.30
#[2,] 0.51 1.00 0.33
#[3,] 0.30 0.33 1.00
#
##$schmid$gloading
#
#Loadings:
#      Factor1
#[1,] 0.681
#[2,] 0.744
#[3,] 0.447
#
#
#      Factor1
#SS loadings      1.218
#Proportion Var   0.406
#

```

---

paired.r

*Test the difference between paired correlations*

---



## Description

Test the difference between paired correlations. Given 3 variables, x, y, z, is the correlation between xy different than that between xz? If y and z are independent, this is a simple t-test. But, if they are dependent, it is a bit more complicated.

## Usage

```
paired.r(xy, xz, yz, n)
```

## Arguments

xy	r(xy)
xz	r(xz)
yz	r(yz)
n	Number of subjects

## Details

Find a t2 statistic for the difference of two dependent correlations.

## Value

t2

## Author(s)

William Revelle  
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<http://personality-project.org/revelle.html>

## References

## Examples

```
paired.r(.5,.3, .4, 100)
```

---

`pairs.panels`*SPLOM, histograms and correlations for a data matrix*

---

## Description

Adapted from the help page for `pairs`, `pairs.panels` shows a scatter plot of matrices (SPLOM), with bivariate scatter plots below the diagonal, histograms on the diagonal, and the Pearson correlation above the diagonal. Useful for descriptive statistics of small data sets.

## Usage

```
pairs.panels(x, smooth = TRUE, scale = FALSE, digits = 2, pch=20, ...)
```

## Arguments

<code>x</code>	a data.frame or matrix
<code>smooth</code>	TRUE draws loess smooths
<code>scale</code>	TRUE scales the correlation font by the size of the absolute correlation.
<code>digits</code>	the number of digits to show
<code>pch</code>	The plot character (defaults to 20 which is a '.').
<code>...</code>	other options for <code>pairs</code>

## Details

Shamelessly adapted from the `pairs` help page. Uses `panel.cor`, `panel.cor.scale`, and `panel.hist`, all taken from the help pages for `pairs`.

## Value

a scatter plot matrix (SPLOM) is drawn in the graphic window. The lower off diagonal draws scatter plots, the diagonal histograms, the upper off diagonal reports the Pearson correlation (with pairwise deletion).

## See Also

[pairs](#)

## Examples

```
#pairs.panels(attitude)  #see the graphics window
```

---

<code>phi</code>	<i>Find the phi coefficient of correlation between two dichotomous variables</i>
------------------	--

---

## Description

Given a 1 x 4 vector or a 2 x 2 matrix of frequencies, find the phi coefficient of correlation. Typical use is in the case of predicting a dichotomous criterion from a dichotomous predictor.

## Usage

```
phi(t, digits = 2)
```

## Arguments

<code>t</code>	a 1 x 4 vector or a 2 x 2 matrix
<code>digits</code>	round the result to digits

## Details

In many prediction situations, a dichotomous predictor (accept/reject) is validated against a dichotomous criterion (success/failure). Although a polychoric correlation estimates the underlying Pearson correlation as if the predictor and criteria were continuous and bivariate normal variables, the phi coefficient is the Pearson applied to a matrix of 0's and 1s.

The calculation follows J. Wiggins discussion of personality assessment.

## Value

phi coefficient of correlation

## Author(s)

William Revelle with modifications by Leo Gurtler

## See Also

[phi2poly](#)

## Examples

```
phi(c(30,20,20,30))
phi(c(40,10,10,40))
x <- matrix(c(40,5,20,20),ncol=2)
phi(x)
```

---

**phi2poly***Convert a phi coefficient to a polychoric correlation*

---

**Description**

Given a phi coefficient (a Pearson r calculated on two dichotomous variables), and the marginal frequencies (in percentages), what is the corresponding estimate of the polychoric correlation?

**Usage**

```
phi2poly(ph, cp, cc)
```

**Arguments**

ph	phi
cp	probability of the predictor – the so called selection ratio
cc	probability of the criterion – the so called success rate.

**Details**

Uses John Fox's polycor package, which in turn requires the mvtnorm package

**Value**

a polychoric correlation

**Author(s)**

William Revelle

**See Also**

[polychor.matrix](#)

**Examples**

```
#phi2poly(.3,.5,.5)
#phi2poly(.3,.3,.7)
```

---

`poly.mat`

*Find polychoric correlations of item data*

---

## Description

Uses John Fox's `hetcor` function (from `polychor` package) to find a matrix of polychoric correlations for integer data. Essentially a wrapper for `hetcor` to convert integer item data into factor (categorical) data and then use `hetcor`. Just a useful shortcut for subsequent factor analysis.

## Usage

```
poly.mat(x, short = TRUE, std.err = FALSE, ML = FALSE)
```

## Arguments

<code>x</code>	A matrix or data frame of integer data
<code>short</code>	<code>short=TRUE</code> , just show the correlations, <code>short=FALSE</code> give the full <code>hetcor</code> output
<code>std.err</code>	<code>std.err=FALSE</code> does not report the standard errors (faster)
<code>ML</code>	<code>ML=FALSE</code> do a quick two step procedure, <code>ML=TRUE</code> , do longer maximum likelihood

## Details

Typical personality and item data are integer values (0,1 for ability; 1,2, 3, 4 for attitude scales). The normal correlation procedures will find Pearson correlations (`cor`). The `polycor` and `hetcor` functions from John Fox's `polychor` package will find polychoric correlations for categorical data. This wrapper function converts integer data to categorical data and then calls `hetcor`.

## Value

A matrix of polychoric correlations (if `short=TRUE`), otherwise a list of various estimates (see `hetcor`).

## Note

requires `polycor`

## Author(s)

William Revelle

## Examples

---

<code>polychor.matrix</code>	<i>Actually, what does this do? Convert a matrix of phi coefficients to a matrix of polycoric correlations</i>
------------------------------	--

---

## Description

Given a vector of a vector of frequencies, use John Fox's polycor function to convert these to polychoric correlations.

Not ready for public consumption.

## Usage

```
polychor.matrix(x, y = NULL)
```

## Arguments

<code>x</code>	a matrix of phi coefficients
<code>y</code>	perhaps

## Details

This is a stub of a function to convert a matrix of phi coefficients with an accompanying vector of frequencies into polychoric correlations. Clearly it is not there yet, but is included in the package as a stopgap.

Please do not use.

## Value

## Author(s)

William Revelle

## Examples

**Description**

Does an eigen value decomposition and returns eigen values, loadings, and degree of fit for a specified number of components. Basically it is just doing a principal components for n principal components. Can show the residual correlations as well. The quality of reduction in the squared correlations is reported by comparing residual correlations to original correlations. Unlike princomp, this returns a subset of just the best nfactors. The eigen vectors are rescaled by the sqrt of the eigen values to produce the component loadings more typical in factor analysis.

**Usage**

```
principal(r, nfactors = 1, residuals = FALSE, rotate="varimax", n.obs=NULL, digits=2)
```

**Arguments**

<b>r</b>	a correlation matrix. If a raw data matrix is used, the correlations will be found using pairwise deletions for missing values.
<b>nfactors</b>	Number of components to extract
<b>residuals</b>	FALSE, do not show residuals, TRUE, report residuals
<b>rotate</b>	
<b>n.obs</b>	Number of observations used to find the correlation matrix if using a correlation matrix. Used for finding the goodness of fit statistics.
<b>digits</b>	digits =2. Accuracy of answers as well as display

**Details**

Useful for those cases where the correlation matrix is improper (perhaps because of SAPA techniques).

There are a number of data reduction techniques including principal components and factor analysis. Both PC and FA attempt to approximate a given correlation or covariance matrix of rank n with matrix of lower rank (p).  ${}_nR_n \approx_n F_{kk}F'_n + U^2$  where k is much less than n. For principal components, the item uniqueness is assumed to be zero and all elements of the correlation matrix are fitted. That is,  ${}_nR_n \approx_n F_{kk}F'_n$  The primary empirical difference between a components versus a factor model is the treatment of the variances for each item. Philosophically, components are weighted composites of observed variables while in the factor model, variables are weighted composites of the factors.

For a n x n correlation matrix, the n principal components completely reproduce the correlation matrix. However, if just the first k principal components are extracted, this is the best k dimensional approximation of the matrix.

Some of the statistics reported are more appropriate for maximum likelihood factor analysis rather than principal components analysis, and are reported to allow comparisons with these other models.

## Value

<code>values</code>	Eigen Values of all components – useful for a scree plot
<code>rotation</code>	which rotation was requested?
<code>n.obs</code>	number of observations specified or found
<code>communality</code>	Communality estimates for each item. These are merely the sum of squared factor loadings for that item.
<code>loadings</code>	A standard loading matrix
<code>fit</code>	Fit of the model to the correlation matrix
<code>fit.off</code>	how well are the off diagonal elements reproduced?
<code>residual</code>	Residual matrix – if requested
<code>communality</code>	The history of the communality estimates. Probably only useful for teaching what happens in the process of iterative fitting.
<code>dof</code>	Degrees of Freedom for this model. This is the number of observed correlations minus the number of independent parameters (number of items * number of factors - $nf*(nf-1)/2$ ). That is, $dof = niI * (ni-1)/2 - ni * nf + nf*(nf-1)/2$ .
<code>objective</code>	value of the function that is minimized by maximum likelihood procedures. This is reported for comparison purposes and as a way to estimate chi square goodness of fit. The objective function is $f = \log(\text{trace}((FF' + U2)^{-1}R) - \log( (FF' + U2)^{-1}R ) - n.items$ .
<code>STATISTIC</code>	If the number of observations is specified or found, this is a chi square based upon the objective function, f. Using the formula from <a href="#">factanal</a> : $\chi^2 = (n.obs - 1 - (2 * p + 5)/6 - (2 * factors)/3)) * f$
<code>PVAL</code>	If <code>n.obs &gt; 0</code> , then what is the probability of observing a chisquare this large or larger?
<code>phi</code>	If oblique rotations (using oblimin from the <code>GPArotation</code> package) are requested, what is the interfactor correlation.

## Author(s)

William Revelle

## See Also

[VSS](#), [factor2cluster](#), [factor.pa](#), [factor.congruence](#)

## Examples

```
#Four principal components of the Harman 24 variable problem
#compare to a four factor principal axes solution using factor.congruence
pc <- principal(Harman74.cor$cov,4,rotate=TRUE)
pa <- factor.pa(Harman74.cor$cov,4,rotate=TRUE)
round(factor.congruence(pc,pa),2)
```



---

`psycho.demo`

*Create demo data for psychometrics*

---

## Description

A not very interesting demo of what happens if bivariate continuous data are dichotomized. Bascially a demo of  $r$ ,  $\phi$ , and polychor.

## Usage

```
psycho.demo()
```

## Details

Not one of my more interesting demonstrations. See <http://personality-project.org/r/simulating-personality.html> and <http://personality-project.org/r/r.datageneration.html> for better demonstrations of data generation.

## Value

a matrix of correlations)

## Author(s)

William Revelle

## References

<http://personality-project.org/r/simulating-personality.html> and <http://personality-project.org/r/r.datageneration.html> for better demonstrations of data generation.

## See Also

[VSS.simulate](#)

## Examples

---

<code>read.clipboard</code>	<i>shortcut for reading from the clipboard</i>
-----------------------------	--

---

## Description

input from the keyboard is easy but a bit obscure, particularly for Mac users. This is just an easier mnemonic to do so.

## Usage

```
read.clipboard(header = TRUE, ...)

#my.data <- read.clipboard()           #assumes headers and tab delimited
#my.data <- read.clipboard.csv()       #assumes heades and comma delimited
```

## Arguments

<code>header</code>	Does the first row have variable labels
<code>...</code>	Other parameters to pass to read

## Details

A typical session of R might involve data stored in text files, generated on line, etc. Although it is easy to just read from a file (particularly if using `file.locate()`, copying from the file to the clipboard and then reading from the clipboard is also very convenient (and somewhat more intuitive to the naive user.)

Based upon a suggestion by Ken Knoblauch to the R-help listserve.

## Value

the contents of the clipboard.

## Author(s)

William Revelle

## Examples

```
#my.data <- read.clipboard()
#my.data <- read.clipboard.csv()
#my.data <- read.clipboard(header=FALSE)
```

---

`schmid`

*Apply the Schmid Leiman transformation to a correlation matrix*

---

## Description

One way to find omega is to do a factor analysis of the original data set, rotate the factors obliquely, do a Schmid Leiman transformation, and then find omega. Here is the code for Schmid Leiman. The S-L transform takes a factor or PC solution, transforms it to an oblique solution, factors the oblique solution to find a higher order (g ) factor, and then residualizes g out of the the group factors.

## Usage

```
schmid(model, nfactors = 3, pc = "pa",...)
```

## Arguments

<code>model</code>	A correlation matrix
<code>nfactors</code>	Number of factors to extract
<code>pc</code>	pc="pa" for principal axes, pc="pc" for principal components, pc="mle" for maximum likelihood
<code>...</code>	Allows additional parameters to be passed to the factoring routines

## Details

Schmid Leiman orthogonalizations are typical in the ability domain, but are not seen as often in the non-cognitive personality domain. S-L is one way of finding the loadings of items on the general factor for estimating omega.

A typical example would be in the study of anxiety and depression. A general neuroticism factor (g) accounts for much of the variance, but smaller group factors of tense anxiety, panic disorder, depression, etc. also need to be considered.

An alternative model is to consider hierarchical cluster analysis techniques such as [ICLUST](#). Requires the GPArotation package.

## Value

<code>sl</code>	loadings on g + nfactors group factors, communalities, uniqueness
<code>orthog</code>	original orthogonal factor loadings
<code>oblique</code>	oblique factor loadings
<code>fcor</code>	correlations among the transformed factors
<code>gload</code>	loadings of the lower order factors on g
<code>...</code>	

## Author(s)

William Revelle

## References

<http://personality-project.org/r/r.omega.html> gives an example taken from Jensen and Weng, 1994 of a S-L transformation.

## See Also

[omega](#), [omega.graph](#), [fa.graph](#), [ICLUST](#), [VSS](#)

## Examples

---

<code>score.alpha</code>	<i>Score scales and find Cronbach's alpha as well as associated statistics</i>
--------------------------	--

---

## Description

Given a matrix or data.frame of k keys for m items (-1, 0, 1), and a matrix or data.frame of items scores for m items and n people, find the sum scores or average scores for each person and each scale. In addition, report Cronbach's alpha, the average r, the scale intercorrelations, and the item by scale correlations.

## Usage

```
score.alpha(keys, items, labels = NULL, totals=TRUE,digits = 2)
```

## Arguments

<code>keys</code>	A matrix or dataframe of -1, 0, or 1 weights for each item on each scale
<code>items</code>	Data frame or matrix of raw item scores
<code>labels</code>	column names for the resulting scales
<code>totals</code>	Find sum scores (default) or average score
<code>digits</code>	Number of digits for answer (default =2)

## Details

The process of finding sum or average scores for a set of scales given a larger set of items is a typical problem in psychometric research. Although the structure of scales can be determined from the item intercorrelations, to find scale means, variances, and do further analyses, it is typical to find the sum or the average scale score.

Various estimates of scale reliability include “Cronbach’s alpha”, and the average interitem correlation. For  $k$  = number of items in a scale, and  $av.r$  = average correlation between items in the scale,  $\alpha = k * av.r / (1 + (k-1)*av.r)$ . Thus, alpha is an increasing function of test length as well as the test homogeneity.

Alpha is a poor estimate of the general factor saturation of a test (see Zinbarg et al., 2005) for it can seriously overestimate the size of a general factor, and a better but not perfect estimate of total test reliability because it underestimates total reliability. None the less, it is a useful statistic to report.

## Value

<code>scores</code>	Sum or average scores for each subject on the $k$ scales
<code>alpha</code>	Cronbach’s coefficient alpha. A simple (but non-optimal) measure of the internal consistency of a test. See also <code>beta</code> and <code>omega</code> .
<code>av.r</code>	The average correlation within a scale, also known as $\alpha 1$ is a useful index of the internal consistency of a domain.
<code>n.items</code>	Number of items on each scale
<code>cor</code>	The intercorrelation of all the scales
<code>item.cor</code>	The correlation of each item with each scale. Because this is not corrected for item overlap, it will overestimate the amount that an item correlates with the other items in a scale.

## Author(s)

William Revelle

## References

An introduction to psychometric theory with applications in R (in preparation). <http://personality-project.org/r/book>

## See Also

[alpha.scale](#), [correct.cor](#), [alpha.scale](#), [cluster.loadings](#), [omega](#)

## Examples

```
y <- attitude      #from the datasets package
keys <- matrix(c(rep(1,7),rep(1,4),rep(0,7),rep(-1,3)),ncol=3)
labels <- c("first","second","third")
x <- score.alpha(keys,y,labels)
```

---

<code>score.items</code>	<i>Score item composite scales and find Cronbach's alpha as well as associated statistics</i>
--------------------------	---

---

### Description

Given a matrix or data.frame of k keys for m items (-1, 0, 1), and a matrix or data.frame of items scores for m items and n people, find the sum scores or average scores for each person and each scale. In addition, report Cronbach's alpha, the average r, the scale intercorrelations, and the item by scale correlations. Replace missing values with the item median or mean if desired. Will adjust scores for reverse scored items.

### Usage

```
score.items(keys, items, totals = FALSE, ilabels = NULL, missing = TRUE, impute="median", min =
```

### Arguments

<code>keys</code>	A matrix or dataframe of -1, 0, or 1 weights for each item on each scale
<code>items</code>	Matrix or dataframe of raw item scores
<code>totals</code>	if TRUE (default) find total scores, if FALSE, find average scores
<code>ilabels</code>	a vector of item labels
<code>missing</code>	TRUE: Replace missing values with the corresponding item median or mean. FALSE: do not score that subject
<code>impute</code>	impute="median" replaces missing values with the item median, impute = "mean" replaces values with the mean response.
<code>min</code>	May be specified as minimum item score allowed, else will be calculated from data
<code>max</code>	May be specified as maximum item score allowed, else will be calculated from data
<code>digits</code>	Number of digits to report

### Details

The process of finding sum or average scores for a set of scales given a larger set of items is a typical problem in psychometric research. Although the structure of scales can be determined from the item intercorrelations, to find scale means, variances, and do further analyses, it is typical to find scores based upon the sum or the average item score. For some strange reason, personality scale scores are typically given as totals, but attitude scores as averages. The default for `score.items` is the average.

Various estimates of scale reliability include "Cronbach's alpha", and the average interitem correlation. For  $k$  = number of items in a scale, and  $av.r$  = average correlation between items in the scale,  $\alpha = k * av.r / (1 + (k-1)*av.r)$ . Thus, alpha is an increasing function of test length as well as the test homogeneity.

Alpha is a poor estimate of the general factor saturation of a test (see Zinbarg et al., 2005) for it can seriously overestimate the size of a general factor, and a better but not perfect estimate of total test reliability because it underestimates total reliability. None the less, it is a useful statistic to report. To estimate the omega coefficient, use the [omega](#) function.

Correlations between scales are attenuated by a lack of reliability. Correcting correlations for reliability (by dividing by the square roots of the reliabilities of each scale) sometimes help show structure.

By default, missing values are replaced with the corresponding median value for that item. Means can be used instead (impute="mean"), or subjects with missing data can just be dropped (missing = FALSE).

## Value

<code>scores</code>	Sum or average scores for each subject on the k scales
<code>alpha</code>	Cronbach's coefficient alpha. A simple (but non-optimal) measure of the internal consistency of a test. See also beta and omega. Set to 1 for scales of length 1.
<code>av.r</code>	The average correlation within a scale, also known as alpha 1 is a useful index of the internal consistency of a domain. Set to 1 for scales with 1 item.
<code>n.items</code>	Number of items on each scale
<code>item.cor</code>	The correlation of each item with each scale. Because this is not corrected for item overlap, it will overestimate the amount that an item correlates with the other items in a scale.
<code>cor</code>	The intercorrelation of all the scales
<code>corrected</code>	The correlations of all scales (below the diagonal), alpha on the diagonal, and the unattenuated correlations (above the diagonal)

## Author(s)

William Revelle

## References

An introduction to psychometric theory with applications in R (in preparation). <http://personality-project.org/r/book>

## See Also

[alpha.scale](#), [correct.cor](#), [cluster.cor](#) , [cluster.loadings](#), [omega](#)

## Examples

```
y <- attitude      #from the datasets package
keys <- matrix(c(rep(1,7),rep(1,4),rep(0,7),rep(-1,3)),ncol=3)
colnames(keys) <- c("first","second","third")
x <- score.items(keys,y)
#x      #to see the output
```

---

**skew***Calculate skew for a vector, matrix, or data.frame*

---

### Description

Find the skew for each variable in a data.frame or matrix. Unlike skew in e1071, this calculates a different skew for each variable or column of a data.frame/matrix.

### Usage

```
skew(x, na.rm = TRUE)
```

### Arguments

<b>x</b>	A data.frame or matrix
<b>na.rm</b>	how to treat missing data

### Details

given a matrix or data.frame x, find the skew for each column.

### Value

if input is a matrix or data.frame, skew is a vector of skews

### Note

The mean function supplies means for the columns of a data.frame, but the overall mean for a matrix. Mean will throw a warning for non-numeric data, but colMeans stops with non-numeric data. Thus, the function uses either mean (for data frames) or colMeans (for matrices). This is true for skew and kurtosi as well.

### Author(s)

William Revelle

### See Also

[describe](#), [describe.by](#), [kurtosi](#)

### Examples

```
round(skew(attitude),2)
```



---

`test.psych`

---

*Testing of functions in the psych package*

---

## Description

Test to make sure the psych functions run on basic test data sets

## Usage

```
test.psych(first=1,last=5,short=TRUE)
```

## Arguments

<code>first</code>	first=1: start with dataset first
<code>last</code>	last=5: test for datasets until last
<code>short</code>	short=TRUE - don't return any analyses

## Details

When modifying the psych package, it is useful to make sure that adding some code does not break something else. This function tests the major functions on various standard data sets. This function also shows off a number of the capabilities of the psych package.

## Value

<code>out</code>	if short=FALSE, then list of the output from all functions tested
------------------	---

## Warning

Warning messages will be thrown by `fa.parallel`

## Note

Not for general consumption – used to make sure functions throw error messages or correct for weird conditions.

The datasets tested are part of the standard R data sets and represent some of the basic problems encountered.

## Author(s)

William Revelle

## Examples

```
#test <- psych.test()
```

---

**VSS.parallel***Compare real and random VSS solutions*

---

### Description

Another useful test for the number of factors is when the eigen values of a random matrix are greater than the eigen values of a real matrix. Here we show VSS solutions to random data.

### Usage

```
VSS.parallel(ncases, nvariables, scree=FALSE, rotate="none")
```

### Arguments

<b>ncases</b>	Number of simulated cases
<b>nvariables</b>	number of simulated variables
<b>scree</b>	Show a scree plot for random data – see <a href="#">omega</a>
<b>rotate</b>	rotate="none" or rotate="varimax"

### Value

VSS like output to be plotted by VSS.plot

### Author(s)

William Revelle

### References

Very Simple Structure (VSS)

### See Also

[fa.parallel](#), [VSS.plot](#), [ICLUST](#), [omega](#)

### Examples

```
#VSS.plot(VSS.parallel(200,24))
```

---

VSS.plot

*Plot VSS fits*

---

## Description

The Very Simple Structure criterion ( [VSS](#)) for estimating the optimal number of factors is plotted as a function of the increasing complexity and increasing number of factors.

## Usage

```
VSS.plot(x, plottitle = "Very Simple Structure", line = FALSE)
```

## Arguments

<code>x</code>	output from VSS
<code>plottitle</code>	any title
<code>line</code>	connect different complexities

## Details

Item-factor models differ in their "complexity". Complexity 1 means that all except the greatest (absolute) loading for an item are ignored. Basically a cluster model (e.g., [ICLUST](#)). Complexity 2 implies all except the greatest two, etc.

Different complexities can suggest different number of optimal number of factors to extract. For personality items, complexity 1 and 2 are probably the most meaningful.

The Very Simple Structure criterion will tend to peak at the number of factors that are most interpretable for a given level of complexity. Note that some problems, the most interpretable number of factors will differ as a function of complexity. For instance, when doing the Harman 24 psychological variable problems, an unrotated solution of complexity one suggests one factor (g), while a complexity two solution suggests that a four factor solution is most appropriate. This latter probably reflects a bi-factor structure.

For examples of VSS.plot output, see <http://personality-project.org/r/r.vss.html>

## Value

A plot window showing the VSS criterion varying as the number of factors and the complexity of the items.

## Author(s)

Maintainer: William Revelle ([revelle@northwestern.edu](mailto:revelle@northwestern.edu))

## References

<http://personality-project.org/r/r.vss.html>

## See Also

[VSS](#), [ICLUST](#), [omega](#)

## Examples

```
#test.data <- Harman74.cor$cov
#my.vss <- VSS(test.data)           #suggests that 4 factor complexity two solution is optimal
#VSS.plot(my.vss)                   #see the graphics window
```

---

VSS.scree

*Plot a scree test*

---

## Description

Cattell's scree test is one of most simple ways of testing the number of components in a correlation matrix. Here we plot the eigen values of a correlation matrix.

## Usage

```
VSS.scree(rx, main = "scree plot")
```

## Arguments

<b>rx</b>	a correlation matrix or a data matrix. If data, then correlations are found using pairwise deletions.
<b>main</b>	Title

## Author(s)

William Revelle  
Department of Psychology  
Northwestern University  
Evanston, Illinois

Maintainer: William Revelle <revelle@northwestern.edu>

## References

<http://personality-project.org/r/vss.html>

## See Also

[VSS.plot](#), [ICLUST](#), [omega](#)

## Examples

```
#VSS.scree(attitude)
#VSS.scree(cor(attitude))
```

---

<code>VSS.simulate</code>	<i>create VSS like data</i>
---------------------------	-----------------------------

---

## Description

Simulation is one of most useful techniques in statistics and psychometrics. Here we simulate a correlation matrix with a simple structure composed of a specified number of factors. Each item is assumed to have complexity one.

## Usage

```
VSS.simulate(ncases=1000, nvariables=16, nfactors=4, meanloading=.5)
```

## Arguments

<code>ncases</code>	number of simulated subjects
<code>nvariables</code>	Number of variables
<code>nfactors</code>	Number of factors to generate
<code>meanloading</code>	with a mean loading

## Value

a `ncases` x `nvariables` matrix

## Author(s)

William Revelle

## See Also

[VSS](#), [ICLUST](#)

## Examples

```
## Not run:
simulated <- VSS.simulate(1000,20,4,.6)
vss <- VSS(simulated,rotate="varimax")
VSS.plot(vss)
## End(Not run)
```

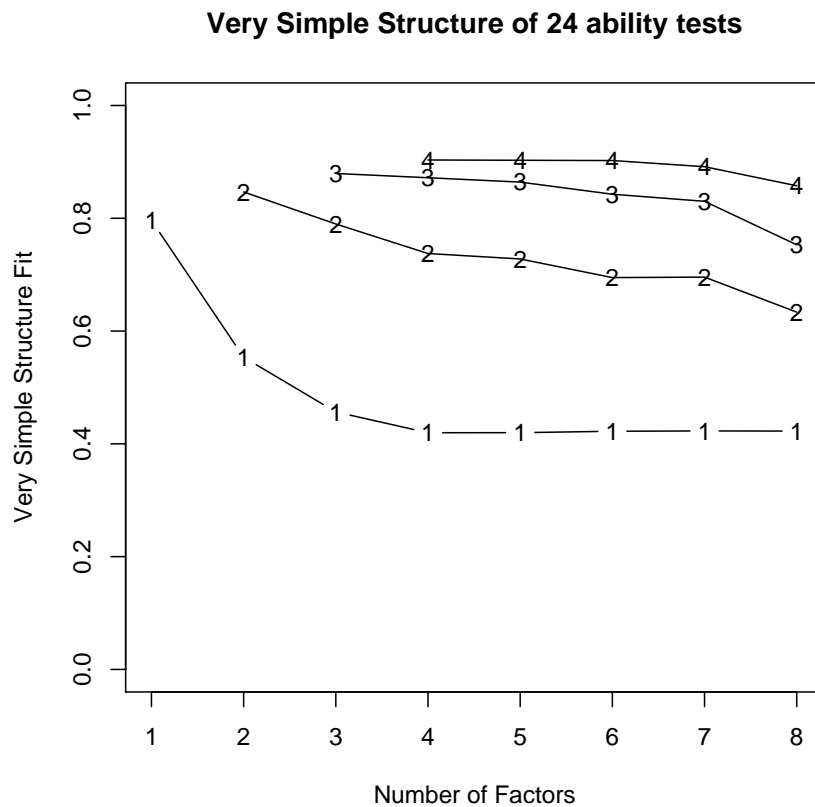


Figure 10: The Very Simple Structure criterion plots goodness of fit as a function of the number of factors extracted and the factorial complexity for each item/test. Note how a complexity one solution best fits the data if only one factor is extracted, but a complexity three solution is optimized at 4 factors.

---

VSS	<i>Apply the Very Simple Structure criterion to determine the appropriate number of factors.</i>
-----	--

---

## Description

There are multiple ways to determine the appropriate number of factors in exploratory factor analysis. Routines for the Very Simple Structure (VSS) criterion allow one to compare solutions of varying complexity and for different number of factors. Graphic output indicates the "optimal" number of factors for different levels of complexity.

## Usage

```
VSS(x, n = 8, rotate = "varimax", diagonal = FALSE, pc = "pa", n.obs=NULL,...)
```

## Arguments

<b>x</b>	a correlation matrix or a data matrix
<b>n</b>	Number of factors to extract – should be more than hypothesized!
<b>rotate</b>	what rotation to use c("none", "varimax", "promax")
<b>diagonal</b>	Should we fit the diagonal as well
<b>pc</b>	pc="pa" Principal Axis Factor Analysis, pc="mle" Maximum Likelihood FA, pc="pc" Principal Components
<b>n.obs</b>	Number of observations if doing a factor analysis of correlation matrix. This value is ignored by VSS but is necessary for the ML factor analysis package.
<b>...</b>	parameters to pass to the factor analysis program The most important of these is if using a correlation matrix is covmat= xx

## Details

Determining the most interpretable number of factors from a factor analysis is perhaps one of the greatest challenges in factor analysis. There are many solutions to this problem, none of which is uniformly the best. "Solving the number of factors problem is easy, I do it everyday before breakfast. But knowing the right solution is harder" (Kaiser, 195x).

Techniques most commonly used include

- 1) Extracting factors until the chi square of the residual matrix is not significant.
- 2) Extracting factors until the change in chi square from factor n to factor n+1 is not significant.
- 3) Extracting factors until the eigen values of the real data are less than the corresponding eigen values of a random data set of the same size (parallel analysis).
- 4) Plotting the magnitude of the successive eigen values and applying the scree test (a sudden drop in eigen values analogous to the change in slope seen when scrambling up the talus slope of a mountain and approaching the rock face.

- 5) Extracting principal components until the eigen value  $<1$ .  $<$
- 6) Extracting factors as long as they are interpretable.
- 7) Using the Very Structure Criterion.

Each of the procedures has its advantages and disadvantages. Using either the chi square test or the change in square test is, of course, sensitive to the number of subjects and leads to the nonsensical condition that if one wants to find many factors, one simply runs more subjects. Parallel analysis is partially sensitive to sample size in that for large samples the eigen values of random factors will be very small. The scree test is quite appealing but can lead to differences of interpretation as to when the scree "breaks". The eigen value of 1 rule, although the default for many programs, seems to be a rough way of dividing the number of variables by 3. Extracting interpretable factors means that the number of factors reflects the investigators creativity more than the data. VSS, while very simple to understand, will not work very well if the data are very factorially complex. (Simulations suggests it will work fine if the complexities of some of the items are no more than 2).

Most users of factor analysis tend to interpret factor output by focusing their attention on the largest loadings for every variable and ignoring the smaller ones. Very Simple Structure operationalizes this tendency by comparing the original correlation matrix to that reproduced by a simplified version (S) of the original factor matrix (F).  $R = SS' + U2$ . S is composed of just the c greatest (in absolute value) loadings for each variable. C (or complexity) is a parameter of the model and may vary from 1 to the number of factors.

The VSS criterion compares the fit of the simplified model to the original correlations:  $VSS = 1 - \text{sumsquares}(r^*) / \text{sumsquares}(r)$  where  $R^*$  is the residual matrix  $R^* = R - SS'$  and  $r^*$  and  $r$  are the elements of  $R^*$  and  $R$  respectively.

VSS for a given complexity will tend to peak at the optimal (most interpretable) number of factors (Revelle and Rocklin, 1979).

Although originally written in Fortran for main frame computers, VSS has been adapted to micro computers (e.g., Macintosh OS 6-9) using Pascal. We now release R code for calculating VSS.

Note that if using a correlation matrix (e.g., `my.matrix`) and doing a factor analysis, the parameters `n.obs` should be specified for the factor analysis: the call is `VSS(my.matrix,n.obs=500)`. Otherwise it defaults to 1000.

## Value

A data.frame with entries: `dof`: degrees of freedom (if using FA)  
`chisq`: chi square (from the factor analysis output (if using FA)  
`prob`: probability of residual matrix  $> 0$  (if using FA)  
`squaresid`: squared residual correlations  
`fit`: factor fit of the complete model  
`cfit.1`: VSS fit of complexity 1  
`cfit.2`: VSS fit of complexity 2  
...  
`cfit.8`: VSS fit of complexity 8  
`cresidual.1`: sum squared residual correlations for complexity 1  
...: sum squared residual correlations for complexity 2 ..8



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## References

<http://personality-project.org/r/vss.html> see also Revelle, W. and Rocklin, T. 1979, Very Simple Structure: an Alternative Procedure for Estimating the Optimal Number of Interpretable Factors, Multivariate Behavioral Research, 14, 403-414. <http://personality-project.org/revelle/publications/vss.pdf>

## See Also

[VSS.plot](#), [ICLUST](#), [omega](#)

## Examples

```
## Not run:
test.data <- Harman74.cor$cov
my.vss <- VSS(test.data)           #suggests that 4 factor complexity two solution is optimal
print(my.vss[,1:12],digits =2)
VSS.plot(my.vss)                  #see graphic window for a plot

#produces this output
#  dof chisq      prob sqresid  fit cfit.1 cfit.2 cfit.3 cfit.4 cfit.5 cfit.6 cfit.7
#1 252 4583  0.0e+00   17.2 0.79  0.79  0.00  0.00  0.00  0.00  0.00  0.00
#2 229 3105  0.0e+00   12.9 0.84  0.75  0.84  0.00  0.00  0.00  0.00  0.00
#4 186 1689 2.3e-240    8.0 0.90  0.66  0.87  0.90  0.90  0.00  0.00  0.00
#5 166 1398 9.3e-194    7.3 0.91  0.68  0.86  0.90  0.91  0.91  0.00  0.00
#6 147 1183 2.9e-161    6.5 0.92  0.53  0.83  0.88  0.91  0.92  0.92  0.00
#7 129 1002 5.8e-135    5.7 0.93  0.47  0.78  0.88  0.91  0.92  0.93  0.93
#8 112  803 5.3e-105    5.3 0.94  0.49  0.76  0.86  0.90  0.92  0.93  0.93

#compare the above solution to a "varimax" rotated solution which suggests 1 factor (g)

my.vss <- VSS(test.data,rotate="varimax")           #suggests that 1 factor complexity one solution is optimal
print(my.vss[,1:14],digits =2)
VSS.plot(my.vss)                  #see graphic window for a plot

#  dof chisq      prob sqresid  fit cfit.1 cfit.2 cfit.3 cfit.4 cfit.5 cfit.6 cfit.7 cfit.8 cresidual.1
#1 252 4583  0.0e+00   17.2 0.79  0.79  0.00  0.00  0.0  0.00  0.00  0.00  0.00  17
#2 229 3105  0.0e+00   12.9 0.84  0.54  0.84  0.00  0.0  0.00  0.00  0.00  0.00  38
#3 207 2193  0.0e+00   10.1 0.88  0.46  0.79  0.88  0.0  0.00  0.00  0.00  0.00  45
#4 186 1689 2.3e-240    8.0 0.90  0.42  0.73  0.87  0.9  0.00  0.00  0.00  0.00  48
#5 166 1398 9.3e-194    7.3 0.91  0.40  0.70  0.86  0.9  0.91  0.00  0.00  0.00  50
#6 147 1183 2.9e-161    6.5 0.92  0.39  0.69  0.86  0.9  0.92  0.92  0.00  0.00  51
#7 129 1002 5.8e-135    5.7 0.93  0.39  0.70  0.84  0.9  0.92  0.93  0.93  0.00  50
```

```
#8 112 803 5.3e-105 5.3 0.94 0.39 0.69 0.83 0.9 0.92 0.93 0.93 0.94 50
## End(Not run)
```