

# Package ‘FactorAssumptions’

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**Type** Package

**Title** Set of Assumptions for Factor and Principal Component Analysis

**Version** 2.0.1

**Maintainer** Jose Storopoli <jstoropoli@protonmail.com>

**Description** Tests for Kaiser-Meyer-Olkin (KMO) and communalities in a dataset. It provides a final sample by removing variables in a iterable manner while keeping account of the variables that were removed in each step. It follows the best practices and assumptions according to Hair, Black, Babin & Anderson (2018, ISBN:9781473756540).

**License** GPL-3

**URL** <https://github.com/storopoli/FactorAssumptions>

**BugReports** <https://github.com/storopoli/FactorAssumptions/issues>

**Depends** R (>= 3.6.0), MASS, psych

**Suggests** knitr, rmarkdown, testthat (>= 2.1.0)

**VignetteBuilder** knitr

**Encoding** UTF-8

**RoxygenNote** 7.0.2

**NeedsCompilation** no

**Author** Jose Storopoli [aut, cre] (<<https://orcid.org/0000-0002-0559-5176>>)

**Repository** CRAN

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 communalities\_optimal\_solution

*Calculates the Optimal Solution for Communalities in a Dataframe*


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### Description

communalities\_optimal\_solution() call upon either the `principal` or `fa` functions from psych package to iterate over the variables of a dataframe.

### Usage

```
communalities_optimal_solution(
  df,
  nfactors,
  type,
  rotate = "varimax",
  fm = "minres",
  squared = TRUE
)
```

### Arguments

df	a dataframe with only int or num type of variables
nfactors	number of factors to extract in principal components or factor analysis
type	either principal for <i>Principal Components Analysis</i> or fa for <i>Factor Analysis</i>
rotate	rotation to be employed (default is <i>varimax</i> ). "none", "varimax", "quartimax", "bentlerT", "equamax", "varimin", "geominT" and "bifactor" are orthogonal rotations. "Promax", "promax", "oblimin", "simplimax", "bentlerQ", "geominQ" and "biquartimin" and "cluster" are possible oblique transformations of the solution. The default is to do a oblimin transformation, although versions prior to 2009 defaulted to varimax. SPSS seems to do a Kaiser normalization before doing Promax, this is done here by the call to "promax" which does the normalization before calling Promax in GPArotation.
fm	Factoring method fm="minres" (default) will do a minimum residual as will fm="uls". Both of these use a first derivative. fm="ols" differs very slightly from "minres" in that it minimizes the entire residual matrix using an OLS procedure but uses the empirical first derivative. This will be slower. fm="wls" will do a weighted least squares (WLS) solution, fm="gls" does a generalized weighted least squares (GLS), fm="pa" will do the principal factor solution, fm="ml" will do a maximum likelihood factor analysis. fm="minchi" will minimize the sample size weighted chi square when treating pairwise correlations with different number of subjects per pair. fm="minrank" will do a minimum rank factor analysis. "old.min" will do minimal residual the way it was done prior to April, 2017 (see discussion below). fm="alpha" will do alpha factor analysis as described in Kaiser and Coffey (1965)
squared	TRUE if matrix is squared (such as adjacency matrices), FALSE otherwise

## Details

If finds any individual communality below the optimal value of 0.5 then removes the lowest communality value variable until no more variable has not-optimal communality values.

## Value

A list with

1. `df` - A dataframe that has reached its optimal solution in terms of KMO values
2. `removed` - A list of removed variables ordered by the first to last removed during the procedure
3. `loadings` - A table with the communalities loadings from the variables final iteration
4. `results` - Results of the final iteration of either the `principal` or `fa` functions from psych package

## See Also

`principal` the PCA function from psych and `fa` the Factor Analysis function from psych

## Examples

```
set.seed(123)
df <- as.data.frame(matrix(rnorm(100*10, 1, .5), ncol=10))
communalities_optimal_solution(df, nfactors = 2, type = "principal", squared = FALSE)
```

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kmo

*Calculates the Kayser-Meyer-Olkin (KMO)*

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## Description

`kmo()` handles both positive definite and not-positive definite matrix by employing the *Moore-Penrose* inverse (pseudoinverse)

## Usage

```
kmo(x, squared = TRUE)
```

## Arguments

<code>x</code>	a matrix or dataframe
<code>squared</code>	TRUE if matrix is squared (such as adjacency matrices), FALSE otherwise

**Value**

A list with

1. overall - Overall KMO value
2. individual - Individual KMO's dataframe
3. AIS - Anti-image Covariance Matrix
4. AIR - Anti-image Correlation Matrix

**Examples**

```
set.seed(123)
df <- as.data.frame(matrix(rnorm(100*10), 1, .5), ncol=10))
kmo(df, squared = FALSE)
```

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kmo\_optimal\_solution *Calculates the Optimal Solution for Kayser-Meyer-Olkin (KMO) in a Dataframe*

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**Description**

kmo\_optimal\_solution() call upon the [kmo](#) function to iterate over the variables of a dataframe.

**Usage**

```
kmo_optimal_solution(df, squared = TRUE)
```

**Arguments**

df                    a dataframe with only int or num type of variables  
squared                TRUE if matrix is squared (such as adjacency matrices), FALSE otherwise

**Details**

If finds any individual KMO's below the optimal value of 0.5 then removes the lowest KMO value variable until no more variable has not-optimal KMO values.

**Value**

A list with

1. df - A dataframe that has reached its optimal solution in terms of KMO values
2. removed - A list of removed variables ordered by the first to last removed during the procedure
3. kmo\_results - Results of the final iteration of the [kmo](#) function

**See Also**

[kmo](#) for kmo computation function

**Examples**

```
set.seed(123)
df <- as.data.frame(matrix(rnorm(100*10, 1, .5), ncol=10))
kmo_optimal_solution(df, squared = FALSE)
```

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