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Maximum Likelihood (ML) (factor analysis algorithms)

Version 21.0.0 ▼

The maximum likelihood solutions of arLambda and $arphi^2$ are obtained by minimizing

$$F = tr[(\Lambda\Lambda' + \psi^2)^{-1}\mathbf{R}] - \log|(\Lambda\Lambda' + \psi^2)^{-1}\mathbf{R}| - p$$

with respect to Λ and ψ , where p is the number of variables, Λ is the factor loading matrix, and ψ^2 is the diagonal matrix of unique variances.

The minimization of F is performed by way of a two-step algorithm. First, the conditional minimum of F for a given y is found. This gives the function $f(\psi)$, which is minimized numerically using the Newton-Raphson procedure. Let $\mathbf{x}^{(s)}$ be the column vector containing the logarithm of the diagonal elements of y at the sth iteration; then

$$\mathbf{x}^{(s+1)} = \mathbf{x}^{(s)} - \mathbf{d}^{(s)}$$

where $\mathbf{d}^{\,(\,s\,)}$ is the solution to the system of linear equations

$$\mathbf{H}^{(s)} \mathbf{d}^{(s)} = \mathbf{h}^{(s)}$$

and where

$$\mathbf{H}^{(s)} = (\partial^2 f(\psi) / \partial x_i \partial x_j)$$

and $\mathbf{h}^{(s)}$ is the column vector containing $\partial f(\psi)/\partial x_i$. The starting point $\mathbf{x}^{(1)}$ is

$$\sqrt{1}$$
 $\log[(1-m/2p)/r^{ii}]$ for ML and GLS

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where m is the number of factors and r^{ii} is the ith diagonal element \square Email feedback to IBM Knowledge Center

The values of $f(\psi)$, $\partial f/\partial x_i$, and $\partial^2 f/\partial x_i \partial x_j$ can be expressed in coins or the eigenvalues

$$\gamma_1 \le \gamma_2 \le \dots \le \gamma_p$$

and corresponding eigenvectors

$$\omega_1$$
, ω_2 , ..., ω_p

of matrix $\psi \mathbf{R}^{-1} \psi$. That is,

$$f(\psi) = \sum_{k=m+1}^{p} (\log \gamma_k + \gamma_k^{-1} - 1)$$

$$\frac{\partial f}{\partial x_i} = \sum_{k=m+1}^{p} \left(1 - \gamma_k^{-1}\right) \omega_{ik}^2$$

$$\frac{\partial^{2} f}{\partial x_{i} \partial x_{j}} = -\delta_{ij} \frac{\partial f}{\partial x_{i}} + \sum_{k=m+1}^{p} \omega_{ik} \omega_{jk} \left(\sum_{n=1}^{m} \frac{\gamma_{k} + \gamma_{n} - 2}{\gamma_{k} - \gamma_{n}} \omega_{in} \omega_{jn} + \delta_{ij} \right)$$

where

$$\delta_{ij} = \begin{cases} 1 & \text{if } i = j \\ 0 & \text{if } i \neq j \end{cases}$$

The approximate second-order derivatives

$$\frac{\partial^2 f}{\partial x_i \partial x_j} \cong \left(\sum_{k=m+1}^p \omega_{ik} \omega_{jk}\right)^2$$

are used in the initial step and when the matrix of the exact second-vector ${\bf d}$ are greater than 0.1. If $\partial^2 f/\partial x_i^2 < 0.05$ (Heywood variables of that column and row are set to 0. If the value of $f(\psi)$ is not decrease of $f(\psi)$ decreases or 25 halvings fail to produce a decrease. (In this largest absolute value of the elements of ${\bf d}$ is less than the criterion

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the value
es until the

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largest absolute value of the elements of \mathbf{d} is less than the criterion value (default 0.001) or until the maximum number of iterations (default 25) is reached. Using the converged value of ψ (denoted by

ŵ

), the eigenanalysis is performed on the matrix



 \mathbf{R}^{-1}

 $\hat{\psi}$

. The factor loadings are computed as

? =

m =

ŵ

$$\Omega_m \left(\Gamma_m^{-1} - \mathbf{I}_m \right)^{1/2}$$

where

$$\Gamma_m = \text{diag} (\gamma_1, \gamma_2, \dots, \gamma_m)$$

 $\Omega_m = (\omega_1, \omega_2, \dots, \omega_m)$

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