

Maximum Likelihood Module

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See the `max.e` files in the directory `c:\gauss\examples` for examples of how to use MAXLIK.

Procedure For Computing Likelihood Function

The user must provide a procedure for computing the log-likelihood for either one observation, or for a matrix of observations. The procedure must have two input arguments, first, a vector of parameter values, and second, one or more rows of the data matrix. The output argument is the log-likelihood for the observation or observations in the second argument evaluated at the parameters values in the first argument. Suppose that the function procedure has been named `fct`, the following considerations apply:

```

FORMAT
    logl=fct(x,y)

INPUT
    x - vector of parameters of model

    y - one or more rows of the data set (if the data set has
        been transformed, or if vars /= 0, i.e., there is
        selection, then y is a transformed, selected observation)
        if __row == 1, one row of the data set
        if __row >= 2, if data set is stored in memory then
            all of the data set will be passed to FCT;
            if data set is stored in GAUSS data file
            then __row will be passed to passed to
            FCT.
        if __row <= 0, For data set is stored in memory same as __row>= 2,
            for GAUSS data file the maximum number of
            rows that will fit in memory will be
            computed by MAXLIK.

    if _max_Lag >= 1, a matrix of observations, the first is
        the i-_max_Lag row, and the final row is
        the i-th row.

OUTPUT
    logl - the log-likelihood
        if __row == 1 or _max_Lag >= 1, a scalar value for
        a given observation, otherwise a vector of
        log-likelihoods.
  
```

REMARKS

If you have written the procedure such that it must compute the log-likelihood of one observation at a time then you must set `__row = 1`. But if you are able to write the procedure so that a vector of log-likelihoods may be returned then set `__row=0`; If you are getting "insufficient memory" messages when the data are being read from a GAUSS data file then either set `__row ==1` or to some positive value. Also, if the data set is stored in a GAUSS data set and the selected data set will fit into memory, then MAXLIK will read it in and store it before beginning the iterations. In this case the setting of `__row` will follow the rules of a data set stored in memory. Significant reduction in computation time may be achieved when the data set can be stored in memory and procedure is written to compute vectors of log-probabilities.

Calling MAXLIK Recursively

The procedure that computes the log-likelihood may itself call MAXLIK. When calling MAXLIK recursively the following considerations apply:

If a data set is being analyzed and it is to be transformed or deleted for missing data or cases are to be selected, then this can be done only on the outermost version of MAXLIK, i.e., the version called in the original command file. Variable selection (as opposed to case selection) can be done on any level through the second argument in the call to each version of MAXLIK. Data sets can be opened by nested versions of MAXLIK. If a nested version of MAXLIK is going to use the data set opened by the outer version of MAXLIK then pass a null string (i.e., "") in the first argument in the call. If it is going to analyze a different data set from the outer version then pass it the data set name in a string. You may also load and store a data set in memory in the command file and pass it as the first argument in the nested call to MAXLIK.

Before the call to the nested version of MAXLIK, the global variables may be re-set by calling MAXCLR. You must not use MAXSET because that will clear information about the data sets opened and processed in the outer version. The only differences between MAXSET and MAXCLR are references to these globals.

You may also want to disable the keyboard control of the nested versions. This is done by setting the global `_max_key = 0` after the call to MAXCLR and before the call to the nested MAXLIK.

PROC MAXLIK

FORMAT

```
{ x, f, g, cov, retcode } = MAXLIK(dataset, vars, &fct, start)
```

INPUT

`dataset` - string containing name of GAUSS data set, or
name of data matrix stored in memory

`vars` - character vector of labels selected for analysis, or
numeric vector of column numbers in data set
of variables selected for analysis

`fct` - the name of a procedure that returns either
the log-likelihood for one observation or a vector of
log-likelihoods for a matrix of observations

`start` - a Kx1 vector of start values

OUTPUT

```

x - Kx1 vector, estimated parameters
f - scalar, function at minimum (mean log-likelihood)
g - Kx1 vector, gradient evaluated at x
cov - KxK matrix, covariance matrix of the parameters
retcode - scalar, return code:

0 normal convergence
1 forced exit
2 maximum number of iterations exceeded
3 function calculation failed
4 gradient calculation failed
5 Hessian calculation failed
6 step length calculation failed
7 function cannot be evaluated at initial parameter values
8 number of elements in the gradient vector inconsistent
  with number of starting values
9 gradient function returned a column vector rather than
  the required row vector
10 secant update failed
11 maximum time exceeded
12 weights could not be found
20 Hessian failed to invert
34 data set could not be opened
99 termination condition unknown

```

GLOBAL VARIABLES

```

----- Options -----
_max_Options - string array, specification of options,
              default is equivalent to:

              string _max_Options = { bfgs stepbt forward info screen }

----- Descent -----
_max_Algorithm - scalar, determines descent algorithm (2)
_max_Delta     - scalar, floor for Hessian Eigenvalues in Newton (.1)

----- Line Search -----
_max_LineSearch - scalar, determines line search method (2)
_max_MaxTry     - scalar, maximum # of tries in step length methods (10)
_max_Extrap     - scalar, extrapolation constant for BRENT (2.0)
_max_Interp     - scalar, interpolation constant for BRENT (.25)
_max_RandRadius - scalar, radius of random direction (0)
_max_UserSearch - scalar, enables user defined line search (0)

----- Covariance Matrix of Parameters -----
_max_CovPar     - scalar, determines type of covariance matrix of
              parameters (1)
_max_XprodCov  - KxK matrix, cross-product covariance matrix of
              parameters when _max_CovPar = 3
_max_HessCov   - KxK matrix, information matrix covariance matrix
              of parameters when _max_CovPar = 3
_max_FinalHess - KxK matrix, stores hessian used for covariance

----- Gradients -----
_max_GradMethod - determines type of numerical gradient (1)
_max_GradProc   - scalar, pointer to analytical gradient procedure (0)
_max_UserNumGrad - scalar, pointer to numerical gradient procedure (0)
_max_HessProc   - scalar, pointer to analytical hessian procedure (0)
_max_UserNumHess - scalar, pointer to numerical hessian procedure (0)
_max_GradStep   - scalar, increment size for computing gradient (0)
_max_GradCheck  - scalar, if nonzero, check analytical gradients (0)

----- Convergence Criteria -----
_max_GradTol    - scalar, convergence tolerance for gradient (1e-5)
_max_MaxIters  - scalar, maximum number of iterations (1e+5)

```

```

_max_MaxTime - scalar, maximum time in iterations in minutes (1e+5)

----- Data -----
_max_Active - vector, defines fixed/active coefficients (1)
__weight - vector, frequency of observations (1)
_max_Lag - scalar, number of lags in model (0)
_max_NumObs - scalar, rows of data matrix (output)
_max_ParNames - char. vector, parameter names (0)
__row - scalar, # of rows of data set passed to procedures (0)
__rowfac - scalar, proportion of rows of data set (1)

----- Miscellaneous -----
__title - string, title ("")
_max_IterData - 3x1 vector, elapsed time, # of iters, cov method
_max_Key - scalar, controls keyboard trapping (0)
_max_Diagnostic - scalar, records current information from iterations

```

Options

`_max_Options` - string array, specification of options. This global permits setting various MAXLIK options in a single global using string identifiers. For example,

```
string _max_Options = { brent newton central file };
```

sets the line search method to BRENT, the descent method to NEWTON, the numerical gradient method to central differences, and `__OUTPUT` = 1.

Algorithms: STEEP, BFGS, DFP, NEWTON, BHHH, PRCG

Line Search: ONE, STEPBT, HALF, BRENT, BHHHSTEP

Covariance Matrix : NOCOV, INFO, XPROD, HETCON

Gradient method: CENTRAL, FORWARD

Output method: NONE, FILE, SCREEN

Descent

```

_max_Algorithm - scalar, indicator for optimization method:
= 1, SD (steepest descent)
= 2, BFGS (Broyden, Fletcher, Goldfarb, Shanno)
= 3, DFP (Davidon, Fletcher, Powell)
= 4, NEWTON (Newton-Raphson)
= 5, BHHH
= 6, Polak-Ribiere Conjugate Gradient

```

`_max_Delta` - scalar, floor for eigenvalues of Hessian in the NEWTON algorithm. This will insure that the Hessian will be positive definite.

Line Search

`_max_LineSearch` - scalar, indicator determining the line search method.

```

= 1, steplength = 1
= 2, STEPBT (default)
= 3, HALF
= 4, BRENT
= 5, BHHHSTEP

```

Usually `_max_Step = 2` will be best. If the optimization bogs down try setting `_max_Step = 1` or `3`. `_max_Step = 3` will generate slow iterations but faster convergence and `_max_Step = 1` will generate fast iterations but slower convergence.

`_max_MaxTry` - scalar, maximum number of tries in BRENT and GOLDEN.

`_max_Extrap` - scalar, extrapolation constant in BRENT.

`_max_Interp` - scalar, interpolation constant in BRENT.

`_max_RandRadius` - scalar, if `_max_RandRadius` is set to a nonzero value (1e-2, say) and all other line search methods fail then OPTMUM will attempt a random direction with radius determined by `_max_RandRadius`.

`_max_UserSearch` - scalar, if nonzero and if all other line search methods fail MAXLIK will enter an interactive mode in which the user can select a line search parameter.

Covariance Matrix of Parameters

`_max_CovPar` - scalar, type of covariance matrix of parameters,
 = 0, the inverse of the final information matrix from the optimization is returned in `cov` (default).
 = 1, the inverse of the second derivatives is returned.
 = 2, the inverse of the cross-product of the first derivatives is returned.
 = 3, the hetereskedastic-consistent covariance matrix is returned.

`_max_XprodCov` - KxK matrix, when `_max_CovPar` is set to 3 the cross-product matrix covariance matrix of the parameters will be returned in `_max_XprodCov`.

`_max_HessCov` - KxK matrix, when `_max_CovPar` is set to 3 the information matrix covariance matrix of the parameters, i.e., the inverse of the matrix of second order partial derivatives of the log-likelihood, will be returned in `_max_HessCov`.

`_max_FinalHess` - KxK matrix, the Hessian used to compute the covariance matrix of the parameters will be stored in `_max_FinalHess`. This will be most useful if the inversion of the hessian fails, which is indicated when MAXLIK returns a missing value for the covariance matrix of the parameters. An analysis of the Hessian stored in `_max_FinalHess` can then reveal the source of the linear dependency responsible for the singularity.

Gradients

`_max_GradMethod` - scalar, method for computing numerical gradient.
 = 0, central difference
 = 1, forward difference (default)

`_max_GradProc` - scalar, pointer to a procedure that computes the gradient of the function with respect to the parameters. For example, the instruction:

```
_max_GradProc=&gradproc
```

tells MAXLIK that a gradient procedure exists as well where to find it. The user-provided procedure has two input arguments, a Kx1 vector of parameter values and an NxP matrix of data. The procedure returns a single output argument, an NxK matrix of gradients of the log-likelihood function with respect to the parameters evaluated at the vector of parameter values.

Default = 0, i.e., no gradient procedure has been provided.

`_max_UserNumGrad` - scalar, pointer to user provided numerical gradient procedure. The instruction

```
_max_GradProc=&gradproc
```

tells MAXLIK that a procedure for computing the numerical gradients exists. The user-provided procedure three input arguments, a pointer to a function that computes the log-likelihood function, a Kx1 vector of parameter values, and an NxP matrix of data. The procedure returns a single output argument, an NxK matrix of gradients of each row of the input data matrix with respect to each parameter.

`_max_HessProc` - scalar, pointer to a procedure that computes the hessian, i.e., the matrix of second order partial derivatives of the function with respect to the parameters. For example, the instruction:

```
_max_HessProc=&hessproc
```

tells OPTMUM that a procedure has been provided for the computation of the hessian and where to find it. The procedure that is provided by the user has two input arguments, a Kx1 vector of parameter values and an NxK data matrix. The procedure returns a single output argument, the KxK symmetric matrix of second order derivatives of the function evaluated at the parameter values.

`_max_UserNumHess` - scalar, pointer to user provided numerical Hessian procedure. The instruction

```
_max_GradProc=&hessproc
```

tells MAXLIK that a procedure for computing the numerical Hessian exists. The user-provided procedure has three input arguments, a pointer to a function that computes the log-likelihood function, a Kx1 vector of parameter values, and an NxK matrix of data. The procedure returns a single output argument, a KxK Hessian matrix of the function with respect to the parameters.

`_max_GradStep` - increment size for computing numerical gradient.

`_max_GradCheck` - scalar, if nonzero and if proc's exist for computing the gradient or Hessian, their calculations will be compared with numerical gradients and Hessians in order to determine their correctness.

Convergence Criteria

`_max_GradTol` - scalar, convergence tolerance for gradient of estimated coefficients. Default = 1e-5. When this criterion has been

satisfied OPTMUM will exit the iterations.

`_max_MaxIters` - scalar, maximum number of iterations.

`_max_MaxTime` - scalar, maximum time in iterations in minutes.
Default = 1e+5, about 10 weeks.

Data

`_max_Active` - vector, 0 = fixed coefficient, 1 = active coefficient.
By default all coefficients are active.

`__weight` - vector, frequency of observations. By default all observations have a frequency of 1. zero frequencies are allowed. It is assumed that the elements of `__weight` sum to the number of observations.

`_max_Lag` - scalar, if the function includes lagged values of the variables `_max_Lag` may be set to the number of lags. When `_max_Lag` is set to a nonzero value then `__row` is set to 1 (that is, the function must be evaluated one observation at a time), and MAXLIK will pass a matrix to the user-provided function and gradient procedures. The first row in this matrix will be (i - `_max_Lag`)-th observation and the last row will be the i-th observation. The read loop will begin with the (`_max_Lag`+1)-th observation. Default = 0.

`_max_NumObs` - scalar, number of cases in the data set that was analyzed.

`_max_ParNames` - Kx1 character vector, parameter labels.

`__row` - determines the number of rows in the data set to be passed to the user-provided procedures. Default = 0.

`__rowfac` - If MAXLIK fails due to insufficient memory while attempting to read a GAUSS data set, then `__rowfac` may be set to some value between 0 and 1 to read a proportion of the original number of rows of the GAUSS data set.

Miscellaneous

`__title` - title of run

`_max_Diagnostic` - scalar. If 1, current estimates ("coeffs"), gradient ("gradient"), direction ("direct"), function value ("function"), Hessian ("Hessian"), and step length computed in the line search ("step") are printed to the screen. If 2, they are stored in `_nlpmax_Diagnostic` using VPUT. Use VREAD to extract. If 3, both 1 and 2 occur.

`_max_IterData` - 3x1 vector, contains information about the iterations. The first element contains the elapsed time in minutes of the iterations, the second element contains the # of iterations, and the third element contains a character variable indicating the type of covariance matrix of the parameters.

`_max_Key` - scalar, controls keyboard capture. Useful for recursively nested version of MAXLIK. Setting `_max_key` = 0 for the nested versions will turn off their key board captures permitting the outside version to retain control of the keyboard.
