

# CMAT Newsletter: December 2009

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# 1 General Remarks

A number of excellent methods for analyzing time series data, like EKG or EEG data, are based on the excellent *Oxford MFE Toolbox* written for Matlab users by Kevin Sheppard. We have modified some of his methods and used some of his ideas, especially for the reparametrization of constraints with the GARCH methods. With his permission we also use Kevin's tables of critical values for the (augmented) Dickey-Fuller testing.

## 1.1 New Functions

**arhetero** Heterogeneous Auto-Regressive modeling ([arhetero](#));

**arma** Auto-Regressive Moving-Average modeling ([arma](#));

**armafore** AutoRegressive Moving-Average forecasting ([armafore](#));

**affrma** RMA (Robust Multichip Average) by Bolstad et.al (2003) for the normalization of microarray data

**berkow** implements Berkowitz test of time series and cross sectional data for many data distributions ([berkow](#));

**garch** implements (G)ARCH, TARCH, AVARCH, ZARCH, EGARCH, APARCH, AGARCH, NAGARCH, IGARCH, FIGARCH time series modeling (based on MFE Toolbox) ([garch](#));

**jarbera** implements the Jarque-Bera test for normality based on the skewness and kurtosis of the data ([jarbera](#));

**mucomp** comparison of means, confirmatory ANOVA similar to Kuiper, Klugkist, & Hoijtink (2007) ([mucomp](#));

**shapwilk** implements Shapiro-Wilk and Shapiro-Francia tests for normal distributed data ([shapwilk](#)).

## 1.2 Fixed Bugs

# 2 Modifications of Features

## 2.1 Modification of the kstest Function

The old version was able only for testing uniform and normal distributions of data. The new version permits also testing of a large variety of other distributions, optional delivers some printed output, and returns additional arguments. Here is the new document for this function:

---

```
prob = kstest(y<,optn<,"dist"<,...>>)
```

```
< prob,stat > = kstest(y<,optn<,"dist"<,...>>)
```

**Purpose:** Different forms of the Cramer-vonMises-Kolmogorov-Smirnov test for a large variety of data distributions, including uniform and normal, are implemented.

**Input: y** The first input object  $y$  must be either a  $n$  vector or a  $n \times nc$  matrix. If  $y$  is a vector the scalar value of the probability of the CDF is returned, if  $y$  is a matrix then a vector of  $nc$  probabilities is returned by using each column of  $y$  separately.

**optn** this is a vector of options which should be initialized with missing values for defaults:

1. (int) amount of printed output (default is no output);
2. (real) probability for test (default is 0.05), must be in  $[0, 1]$ ;
3. (int) version number of algorithm (default is 0):
  - 0** the default version, for all implemented distributions (see below), returning all three results, probability, statistics, and hypothesis value.
  - 1** Anderson-Darling approximative, only for uniform distribution (i.e. data in  $[0, 1]$ ) which returns the probability only.
  - 2** Marsaglia-Tsang-Wang exact (JSS,2003), only for uniform distribution (i.e. data in  $[0, 1]$ ) which returns the probability only.
  - 3** the old implementation of the KS-Lilliefors method, only for normal distribution with  $\mu = 0$  and  $\sigma = 1$  and returning the probability only.

**"dist"** The following string options are valid:

**("bern",x,prob)** Bernoulli distribution

- $x$  is the upper limit for cumulation  $0 \leq x \leq 1$ .
- $p$  is the probability for success of 1,  $0 \leq prob \leq 1$ .

**("beta",x,a,b,<1>,<u>>)** Beta distribution

- $x$  is the upper limit for integration  $0 \leq x \leq 1$ .
- $a$  is the first shape parameter  $a \geq 0$ .
- $b$  is the second shape parameter  $b \geq 0$ .

**("bin",s,prob,n)** Binomial distribution

- $s$  is the number of successes,  $0 \leq s \leq n$ .
- $prob$  is the probability for success,  $0 \leq prob \leq 1$ .
- $n$  is the number of independent trials,  $n > 0$ .

**("cau",x,<μ>,<σ>>)** Cauchy distribution

- $x$  is the upper limit for integration  $x > 0$ .
- $\mu$  location parameter  $mu \geq 0$ .
- $\sigma$  scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

- (**"chis"**,**x**,**df**<**nc**>) (noncentral) ChiSquare distribution
- **x** is the upper limit for integration  $x > 0$ .
  - **df** degrees of freedom  $df > 0$ .
  - **nc** optional noncentrality  $nc \geq 0$ .
- (**"expo"**,**x**< **$\sigma$** >) Exponential distribution
- **x** is the upper limit for integration  $0 \leq x \leq 1$ .
  - **$\sigma$**  scale parameter ( $\sigma > 0$ ).
- Default is  $\sigma = 1$ .
- (**"f"**,**x**,**ndf**,**ddf**<**nc**>) (noncentral) *F* distribution
- **x** is the upper limit for integration  $x > 0$ .
  - **ndf** numerator degrees of freedom  $ndf > 0$ .
  - **ddf** denominator degrees of freedom  $ddf > 0$ .
  - **nc** optional noncentrality  $nc \geq 0$ .
- (**"gam"**,**x**,**shape**<**scale**>) • **x** is the upper limit for integration  $x > 0$ .
- **shape** shape parameter  $shape > 0$ .
  - **scale** is the scale parameter  $scale > 0$ .
- Default is  $scale = 1$ .
- (**"gaus"**,**x**< **$\mu$** >**,< $\sigma$ >>**) Gauss (Normal) distribution
- **x** is the upper limit for integration  $x > 0$ .
  - **$\mu$**  shape parameter.
  - **$\sigma$**  scale parameter ( $\sigma > 0$ ).
- Default is  $\mu = 0$  and  $\sigma = 1$ .
- (**"geom"**,**m**,**p**) • **m** is the upper limit for cumulation  $m > 0$ .
- **p** is the probability for success of 1,  $0 \leq prob \leq 1$ .
- (**"hypg"**,**x**,**m**,**k**,**n**<**r**>) Hypergeometric distribution
- **x** is the upper limit for cumulation  $x > 0$ .
  - **m** is the population size ( $m \geq 1$ ).
  - **k** is the number of items ( $0 \leq k \leq m$ ).
  - **n** is the sample size ( $0 \leq n \leq m$ ).
  - **r** is an optional odds ratio ( $r > 0$ ).
- Default is  $r = 1$ .
- (**"igau"**,**x**,**d**) Inverse Gauss (Wald) distribution
- **x** is the upper limit for integration  $x > 0$ .
  - **d** shape parameter  $d > 0$ .
- (**"lapl"**,**x**< **$\mu$** >**,< $\sigma$ >>**) Laplace distribution
- **x** is the upper limit for integration  $x > 0$ .
  - **$\mu$**  shape parameter.
  - **$\sigma$**  scale parameter ( $\sigma > 0$ ).
- Default is  $\mu = 0$  and  $\sigma = 1$ .

(**"logi"**,**x**<,< $\mu$ >,< $\sigma$ >>) • **x** is the upper limit for integration  $x > 0$ .

- $\mu$  shape parameter.
- $\sigma$  scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

(**"logn"**,**x**<,< $\mu$ >,< $\sigma$ >>) LogNormal distribution

- **x** is the upper limit for integration  $x > 0$ .
- $\mu$  shape parameter.
- $\sigma$  scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

(**"negb"**,**x**,**prob**,**n**) • **x** is the number of failures,  $0 \leq f \leq n$ .

- **prob** is the probability for success,  $0 \leq prob \leq 1$ .
- **n** is the number of successes,  $n > 0$ .

(**"norm"**,**x**<,< $\mu$ >,< $\sigma$ >>) Normal (Gauss) distribution

- **x** is the upper limit for integration  $x > 0$ .
- $\mu$  shape parameter.
- $\sigma$  scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

Using `optn[3]` two different versions may be specified:

**0** : default algorithm allowing the specification of additional arguments for  $\mu$  and  $\sigma$  and which returns probability, statistics and hypothesis value.

**3** : the old implementation of the KS-Lilliefors method which assumes  $\mu = 0$  and  $\sigma = 1$  and returns the probability only.

(**"pare"**,**x**,**a**<,**k**>) Pareto distribution

- **x** is the upper limit for integration  $x > 0$ .
- **a** shape parameter ( $a > 0$ ).
- **k** scale parameter ( $k > 0$ ).

Default is  $k = 1$ .

(**"pois"**,**n**, $\lambda$ ) Poisson distribution

- **x** is the upper limit for cumulation  $x > 0$ .
- $\lambda$  shape parameter.

(**"t"**,**t**,**df**<,**nc**>) (noncentral)  $t$  distribution

- **t** is chosen from  $t$  distribution.
- **df** degrees of freedom  $df > 0$ .
- **nc** optional noncentrality  $nc \geq 0$ .

(**"unif"**,**x**<,<**l**>,<**u**>>) • **x** is the upper limit for integration  $x > 0$ .

- **l** left location.
- **u** right location ( $u > l$ ).

Default is  $l = 0$  and  $u = 1$ .

Using `optn[3]` three different versions may be specified:

**0** : default algorithm: which returns probability, statistics, and hypothesis value.

**1** : Anderson-Darling approximative, which returns the probability only.

**2** : Marsaglia-Tsang-Wang exact (JSS,2003), which returns the probability only.

("wald",x,d) • **x** is the upper limit for integration  $x > 0$ .

• **d** shape parameter  $d > 0$ .

("weib",x,shape<,scale>) Weibull distribution

• **t** is the upper limit for integration  $x > 0$ .

• **shape** shape parameter  $shape > 0$ .

• **scale** scale parameter  $scale > 0$ .

Default is  $scale = 1$ .

... There may be a variable number of additional input arguments specifying non-default distributional parameters.

**Output: prob** The result is either a scalar or an object  $p = Prob(d < D_n)$ , where

$$D_n = \max(x_1 - \frac{0}{n}, x_2 - \frac{1}{n}, \dots, x_n - \frac{n-1}{n}, \frac{1}{n} - x_1, \frac{2}{n} - x_2, \dots, \frac{n}{n} - x_n)$$

The Anderson-Darling version of the Cramer-von Mises-Kolmogorov-Smirnov test for univariate distribution uses

$$A_n = -n - \frac{1}{n} [\ln(x_1 z_1) + 3 \ln(x_2 z_2) + 5 \ln(x_3 z_3) + \dots + (2n-1) \ln(x_n z_n)]$$

with  $z_i = 1 - x_{n+1-i}$ .

**stat** The **stat** result has the same size as **prob** and returns the test statistics (only for **optn[3]=0**).

**Restrictions:** 1. Missing value are returned if the input  $y$  contains missing data.

2.

**Relationships:** `ksprob()`, `berkow()`, `jarbera()`, `shapwilk()`

**Examples:** 1. Test different Distributions:

The following are  $N = 1000$  uniform random generated data:

```
options NOECHO;
#include "..\tdata\matlb.dat"
options ECHO;
```

```

optn = [ 2 , /* print */
        .05 , /* alpha */
        0 ]; /* version=cdf */
< p31,s31 > = kstest(uf1000,optn,"unif");
print "KSTEST P: UNIF",p31;
print "KSTEST S: UNIF",s31;

```

Hypothesis H0: Data are uniformly distributed versus  
Hypothesis H1: Data are not uniformly distributed  
H0 accepted: Probability=0.212528 Statistics=0.0333409

```

optn = [ 2 , /* print */
        .05 , /* alpha */
        0 ]; /* version=cdf */
< p32,s32 > = kstest(uf1000,optn,"norm");
print "KSTEST P: NORM",p32;
print "KSTEST S: NORM",s32;

```

Hypothesis H0: Data are normally distributed versus  
Hypothesis H1: Data are not normally distributed  
H0 rejected: Probability=1.42564e-218 Statistics=0.4992

```

optn = [ 2 , /* print */
        .05 , /* alpha */
        0 ]; /* version=cdf */
< p32,s32 > = kstest(uf1000,optn,"expo");
print "KSTEST P: NORM",p32;
print "KSTEST S: NORM",s32;

```

Hypothesis H0: Data are Exponential distributed versus  
Hypothesis H1: Data are not Exponential distributed  
H0 rejected: Probability=2.26383e-118 Statistics=0.367064

```

optn = [ 2 , /* print */
        .05 , /* alpha */
        0 ]; /* version=cdf */
< p32,s32 > = kstest(uf1000,optn,"f");
print "KSTEST P: NORM",p32;
print "KSTEST S: NORM",s32;

```

Hypothesis H0: Data are F distributed versus



Hypothesis H1: Data are not F distributed  
H0 rejected: Probability=0 Statistics=0.999001

```
optn = [ 2 , /* print */  
        .05 , /* alpha */  
        0 ]; /* version=cdf */  
< p32,s32 > = kstest(uf1000,optn,"logi");  
print "KSTEST P: NORM",p32;  
print "KSTEST S: NORM",s32;
```

Hypothesis H0: Data are Logistic distributed versus  
Hypothesis H1: Data are not Logistic distributed  
H0 rejected: Probability=1.65613e-218 Statistics=0.499126

2. Compare with other versions for uniform and normal distribution:  
Note, that there is only one return argument for versions  $\neq 0$ . The  
default version=0 computed for uniform: Probability=0.212528:

```
optn = [ 2 , /* print */  
        .05 , /* alpha */  
        1 ]; /* version=and */  
p31 = kstest(uf1000,optn,"unif");  
print "KSTEST: UNIF",p31;
```

KSTEST: UNIF 0.8380

```
optn = [ 2 , /* print */  
        .05 , /* alpha */  
        2 ]; /* version=mtw */  
p32 = kstest(uf1000,optn,"unif");  
print "KSTEST: UNIF",p32;
```

KSTEST: UNIF 0.8117

The default version=0 computed for normal: Probability=2.26383e-118  
:

```
optn = [ 2 , /* print */  
        .05 , /* alpha */  
        3 ]; /* version=ori */  
p33 = kstest(uf1000,optn,"norm");  
print "KSTEST: NORM",p33;
```

```

Number of Data=1000 Mean=0.488832 Standard Deviation=0.283241
Hypothesis H0: Data are normally distributed versus
Hypothesis H1: Data are not normally distributed
KS-Lilliefors-Test on Normal Distribution: D= 0.154522
Critical Values for d (two-sided)
      10%      5%      2%      1%
0.0261 0.0284 0.0313 0.0332
H0 rejected: Probability=0

```

3. Artificial Examples: for normal distribution:

```

opt3 = [ ., ., 3];
options nolapack;
s1 = [ .02 1.15 2.55 0.77 .33 .20 ];
pr1 = kstest(s1,opt3,"nor");
s2 = [ 5 7 9 8 5 0 ];
pr2 = kstest(s2,opt3,"nor");
print " 1: Prob for NOR:", pr1, pr2;

```

Prob for NOR: 0.10000 0.10000

```

opt3 = [ ., ., 3];
options lapack;
s1 = [ .02 1.15 2.55 0.77 .33 .20 ];
pr1 = kstest(s1,opt3,"nor");
s2 = [ 5 7 9 8 5 0 ];
pr2 = kstest(s2,opt3,"nor");
print " 2: Prob for NOR:", pr1, pr2;

```

Prob for NOR: 0.6400 0.9900

```

nr = 5000; nc = 1;
mu = rand(nr,nc,'g',"duni");
mn = rand(nr,nc,'g',"nor");
mn2 = mn .** 2;

```

```

opt3 = [ ., ., 3];
options nolapack;
p1 = kstest(mu,opt3,"nor");
print "Prob for UNIF:",p1;
p2 = kstest(mn,opt3,"nor");
print " 1: Prob for NORM:",p2;

```

```
Prob UNIF for UNIF: 0.0000
Prob NORM for NORM: 0.0000
```

```
opt3 = [ .,.,3];
options lapack;
p1 = kstest(mu,opt3,"nor");
print "Prob for UNIF:",p1;
p2 = kstest(mn,opt3,"nor");
print " 2: Prob for NORM:",p2;
```

```
Prob UNIF for NORM: 0.0000
Prob NORM for NORM: 0.0000
```

4. Comparison of Uniform Random Generators: Here the sample size N=1000 is selected.

```
nr = 1000; nc = 5;
opt1 = [ .,.,1];
opt2 = [ .,.,2];
options RANDUNI;
mu1 = rand(nr,nc,'g',"duni");
p11 = kstest(mu1,opt1,"uni");
print "AND: Prob UNIF for RANDUNI:",p11;
p12 = kstest(mu1,opt2,"uni");
print "MTW: Prob for RANDUNI:",p12;
print "RANDUNI:", p11[+]/nc, p12[+]/nc, ssq(p11 - p12);
```

AND: Prob UNIF for RANDUNI:

```
  |          1
-----
1 |  0.92352
2 |  0.68155
3 |  0.44995
4 |  0.75229
5 |  0.82055
```

MTW: Prob for RANDUNI:

```
  |          1
-----
1 |  0.82039
2 |  0.61299
3 |  0.56562
4 |  0.82924
```

5 | 0.71396

RANDUNI: 0.7256 0.7084 0.04600

```
opt1 = [ .,.,1];
opt2 = [ .,.,2];
options RANDKIS;
mu2 = rand(nr,nc,'g',"duni");
p21 = kstest(mu2,opt1,"uni");
print "AND: Prob UNIF for RANDKIS:",p21;
p22 = kstest(mu2,opt2,"uni");
print "MTW: Prob for RANDKIS:",p22;
print "RANDKIS:", p21[+]/nc, p22[+]/nc, ssq(p21 - p22);
```

AND: Prob UNIF for RANDKIS:

	1
1	0.63250
2	0.31721
3	0.51274
4	0.56430
5	0.52699

MTW: Prob for RANDKIS:

	1
1	0.38966
2	0.69173
3	0.29469
4	0.55523
5	0.62822

RANDKIS: 0.5107 0.5119 0.2571

```
opt1 = [ .,.,1];
opt2 = [ .,.,2];
options RANLECU;
mu3 = rand(nr,nc,'g',"duni");
p31 = kstest(mu3,opt1,"uni");
print "AND: Prob UNIF for RANLECU:",p31;
p32 = kstest(mu3,opt2,"uni");
print "MTW: Prob for RANLECU:",p32;
```

```
print "RANDLECU:", p31[+]/nc, p32[+]/nc, ssq(p31 - p32);
```

```
AND: Prob UNIF for RANDLECU:
```

```
  |          1
-----
1 |    0.96694
2 |    0.36281
3 |    0.42487
4 |    0.07408
5 |    0.38351
```

```
MTW: Prob for RANDLECU:
```

```
  |          1
-----
1 |    0.98915
2 |    0.21339
3 |    0.32869
4 |    0.05316
5 |    0.65719
```

```
RANDLECU: 0.4424 0.4483 0.1074
```

```
opt1 = [ .,.,1];
opt2 = [ .,.,2];
options RANDXORWOW;
mu4 = rand(nr,nc,'g',"duni");
p41 = kstest(mu4,opt1,"uni");
print "AND: Prob UNIF for RANDXORWOW:",p41;
p42 = kstest(mu4,opt2,"uni");
print "MTW: Prob for RANXORWOW:",p42;
res = ssq(p41 - p42);
print "RANDXORWOW:", p41[+]/nc, p42[+]/nc, ssq(p41 - p42);
options RANDUNI;
```

```
AND: Prob UNIF for RANDXORWOW:
```

```
  |          1
-----
1 |    0.05877
2 |    0.85525
3 |    0.18640
4 |    0.41601
5 |    0.27656
```

```
MTW: Prob for RANXORWOW:
```

```
  |      1
-----
1 |  0.10718
2 |  0.90353
3 |  0.04704
4 |  0.20441
5 |  0.61463
```

```
RANXORWOW: 0.3586 0.3754 0.1832
```

## 2.2 Modifications of the `mvelps` Function

The `mvelps` function now has the form `v = mvelps(Cov,mu,U,c,optn)` which means that `Cov,mu,U,c` now are specified as single input arguments and no longer as entries of the option vector.

The example shows that the results are unchanged:

```
sigm = [ 1.,
         .5 1.,
         .5 .5 1. ];
mu = cons(1,3,0.);

u = [ .25 ,
      0. .2,
      0. 0. .1];
m = [ 1. .5 .25 ];

optn = [ "ndim"      3,
         "bin"       ,
         "seed"     357,
         "ndf"      20,
         "randir"   1000,
         "nprel"    10,
         "e_rhs"    1. ];
ovec = mvelps(sigm,mu,u,m,optn);
```

```
Number Dimensions           3
T Distribution DF           20
Preliminary Estimates       10
Number Random Directions    10000
Value of Integral           0.6792134
Estimated Standard Error    5.07e-004
```

Initial Seed Value	357
Run Time (in seconds)	2.000000000

### 3 Extensions to the Language

## 4 Extensions to Various Functions

### 4.1 Extension to the `tsmeas` Function

We have added a number of more functions:

”**pacor**” computes partial autocorrelations w.r.t. a scalar or vector of integer delays (lags)  $\tau$ ; in addition optional robust asymptotic standard errors are computed;

”**ljung**” Ljung-Box test for serial correlation for  $\tau = 1, \dots, nlag$  lags;

”**lmtst**” LM test for serial correlation for  $\tau = 1, \dots, nlag$  lags.

”**nwcov**” computing the Newey-West (Bartlett) covariance matrix for  $\tau = 1, \dots, nlag$  lags. This matrix is normally applied to the scores or an ARMA application.

”**nwreg**” ordinary least squares regression  $\min \| \mathbf{X}\beta - y \|$  with  $t$  values based on the Newey-West covariance matrix for  $\tau = 1, \dots, nlag$  lags.

”**diful**” (augmented) Dickey-Fuller testing with specified lags or automatic lag selection (using critical values by K. Sheppard).

”**grang**” Granger causality testing: for a specified vector  $\tau$  of lags; for homoskedastic and heteroskedastic with correlated or uncorrelated errors

”**arvec**” Vector AR estimation: for a specified vector  $\tau$  of lags; for homoskedastic and heteroskedastic with correlated or uncorrelated errors

”**iresp**” Impulse Response modeling: for a specified vector  $\tau$  of lags; for homoskedastic and heteroskedastic with correlated or uncorrelated errors.

#### 4.1.1 Partial Autocorrelations

```
< pacor,ase > = tsmeas(xdat,"pacor",tau<,optn>)
```

The following specific options could be specified:

**3** indicates whether robust (=1) or nonrobust (=0) asymptotic standard errors are being computed. The default is 1.

Option `optn[3]` must be specified to zero for returning nonrobust asymptotic standard errors (which are constant). An additional vector `tau` specifies a set of lags.

If `xdat` has more than one column, each column is treated as separate time series.

Some examples follow:



```

xd = [ 0.8147 0.9058 0.1270 0.9134 0.6324
       0.0975 0.2785 0.5469 0.9575 0.9649
       0.1576 0.9706 0.9572 0.4854 0.8003
       0.1419 0.4218 0.9157 0.7922 0.9595 ];

```

```

optn = [ 2 ]; tau = [ 1 2 3 ];
pcor = tsmeas(xd,"pacor",tau,optn);
print "PCor=",pcor;

```

PCor=

	1
1	-0.07778
2	-0.27478
3	-0.00080

```

print "MFE Implementation with robust ASE";
rob = 1; optn = [ 2 . rob ];
tau = [ 1 2 3 ];
< pcor,rase > = tsmeas(xd,"pacor",tau,optn);
print "PCor=",pcor;
print "RASE=",rase;

```

PCor=

	1	2	3
1	-0.07778	-0.27478	-0.00080

RASE=

	1	2	3
1	0.21846	0.20710	0.24579

```

print "MFE implementation: NonRobust ASE";
tau = 3; rob = 0;
optn = [ 2 tau rob ];
< parcor,nase > = tsmeas(xd,"paco2",optn);
print "ParCor=",parcor;
print "Nonrobust ASE=",nase;

```

ParCor=

	1	2	3
1	-0.07778	-0.27478	-0.00080

Nonrobust ASE=

	1	2	3
1	0.22361	0.22361	0.22361

#### 4.1.2 Ljung-Box Test for Serial Correlation

```
< stat,prob > = tsmeas(xdat,"ljung"<,optn>)
```

The following specific options could be specified.

**2** integer `ntau`  $\geq 1$  specifying how many lags  $\tau = 1, \dots, ntau$  are to be considered. The default is 1.

If `xdat` has more than one column, each column is treated as separate time series.

The function returns two arguments, the test statistics and the probability for each of the `ntau` lags.

The following example illustrates the use of this function:

```
xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xc = xd2 - univar(xd2,"ari");

print "Ljung-Box test for serial correlation:";
tau = 5;
optn = [ 2 tau ];
< stat,prob > = tsmeas(xc,"ljung",optn);
print "LJUNG: stat=",stat;
print "LJUNG: prob=",prob;
```

Ljung-Box test for serial correlation:

```
LJUNG: stat=
|          1          2          3          4          5
-----
1 |  0.03534  0.07382  0.57096  2.0989  3.9281
```

```
LJUNG: prob=
|          1          2          3          4          5
-----
1 |  0.85089  0.96376  0.90305  0.71757  0.55982
```

### 4.1.3 (Robust) LM Test for Serial Correlation

```
< stat,prob > = tsmeas(xdat,"lmtst"<,optn>)
```

The following specific options could be specified.

- 2** an integer `ntau`  $\geq 1$  specifying how many lags  $\tau = 1, \dots, ntau$  are to be considered. The default is 1.
- 3** indicates whether a robust (=1) or a nonrobust (=0) statistic is being computed. The default is 1.

If `xdat` has more than one column, each column is treated as separate time series.

The function returns two arguments, the test statistics and the probability for each of the `ntau` lags.

The following example illustrates the use of this function:

```
xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xc = xd2 - univar(xd2,"ari");

print "LM test for serial correlation: Nonrobust";
tau = 5; rob = 0;
optn = [ 2 tau rob ];
< stat,prob > = tsmeas(xc,"lmtst",optn);
print "LMTST: stat=",stat;
print "LMTST: prob=",prob;
```

LM test for serial correlation: Nonrobust

```
LMTST: stat=
  |          1          2          3          4          5
-----
1 |  0.03455  0.14361  0.49659  2.3261  5.1064
```

```
LMTST: prob=
  |          1          2          3          4          5
-----
1 |  0.85255  0.93071  0.91964  0.67603  0.40303
```

```
print "LM test for serial correlation: Robust";
tau = 5; rob = 1;
optn = [ 2 tau rob ];
< stat,prob > = tsmeas(xc,"lmtst",optn);
print "LMTST: stat=",stat;
print "LMTST: prob=",prob;
```

LM test for serial correlation: Robust

```
LMTST: stat=
  |          1          2          3          4          5
-----
1 |  0.03603  0.14081  0.54035  2.2770  4.5774
```

```
LMTST: prob=
  |          1          2          3          4          5
-----
1 |  0.84945  0.93201  0.90994  0.68496  0.46960
```

#### 4.1.4 Newey-West (Bartlett) Covariance Matrix

```
nwcov = tsmeas(xdat,"nwcov"<,optn>)
```

This function is normally applied to the scores of an ARMA application. The following specific options could be specified.

- 2 an integer  $n\tau \geq 1$  specifying how many lags  $\tau = 1, \dots, n\tau$  are to be considered. The default is 1.
- 3 indicates whether the data are centered around the mean (=1) or the data are used as input (=0). The default is 0.

The function returns only one argument, the Newey-West covariance matrix. The following example shows not a typical use of the function:

```
xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xc = xd2 - univar(xd2,"ari");
```

```
print "Newey-West Covariance Matrix: Uncentered";
tau = 5; cent = 0;
optn = [ 2 tau cent ];
nwc = tsmeas(xd2,"nwcov",optn);
print "Newey-West COV: nwc=",nwc;
```

Newey-West COV: nwc= 1.9405

```
print "Newey-West Covariance Matrix: Centered";
tau = 5; cent = 1;
optn = [ 2 tau cent ];
nwc = tsmeas(xd2,"nwcov",optn);
print "Newey-West COV: nwc=",nwc;
```

Newey-West COV: nwc= 0.1060

#### 4.1.5 OLS Regression with Newey-West HAC Standard Errors

```
< gof,est,cov,yhat > = tsmeas(ydat,"nwreg",xdat<,optn>)
```

This function performs ordinary least squares (OLS) regression  $\min \| \mathbf{X}\beta - y \|$  and computes  $t$  values based on the Newey-West covariance matrix for  $\tau = 1, \dots, nlag$  lags. The first input argument `ydat` specifies usually a  $N$  vector of values of a dependent variable (response), whereas the third input argument `xdat` specifies a  $N \times n$  matrix of values of  $n$  independent variables (predictors). The number  $p$  of parameter estimates is then either  $p = n + 1$  if an intercept variable is added or not, depending on the setting of `optn[3]`.

If `ydat` has more than one column, each column is treated as separate time series.

The following specific options could be specified:

- 2** an integer `ntau`  $\geq 1$  specifying how many lags  $\tau = 1, \dots, ntau$  are to be considered. The default is 1.
- 3** indicates whether a constant is added (`=1`: default) or not (`=0`) to the set of parameter estimates.

Some examples follow:

1. OLS solution of  $40 \times 3$  linear system, without intercept:

```

print "Matlab 200 by 3 RandomData: ran203[200,3]";
options NOECHO;
%inc "..\tdata\matlb.dat";
options ECHO;

ran40 = ran203[1:40,];

xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xd2 = flip(xd2,"ud");

print "Newey-West Regression: w/o Intercept";
pri = 2; tau = 5; intc = 0;
optn = [ pri tau intc ];
< gof,est,nwcov,yhat > = tsmeas(xd2,"nwreg",ran40,optn);
print "Newey-West REG: gof=",gof;
print "Newey-West REG: est=",est;
print "Newey-West REG: nwcov=",nwcov;
print "Newey-West REG: yhat=",yhat;

```

Newey-West REG: gof=		Newey-West REG: yhat=				
	1		1			
-----						
1	0.09458	1	0.62073			
2	0.75065	2	0.49591			
3	0.73718	3	0.36245			
Newey-West REG: est=		4	0.70181			
	1	2	5	0.69115		
-----						
1	0.50690	4.8994	6	0.39841		
2	-0.04048	-0.19482	7	0.45876		
3	0.55293	4.6315	8	0.27353		
Newey-West REG: nwcov=		9	0.49820			
S	1	2	3	10	0.64206	
-----						
1	0.01070			11	0.36503	
2	-0.01403	0.04318			12	0.84164
3	-0.00069	-0.01440	0.01425	13	0.69152	
-----						
Newey-West REG: yhat=						
-----						
14   0.69012						
15   0.76872						
16   0.59961						
17   0.49844						
18   0.63702						
19   0.45074						
20   0.80658						
21   0.75040						
22   0.21488						
23   0.46320						
24   0.61332						
25   0.39239						
26   0.49979						
27   0.60226						
28   0.48576						
29   0.57474						
30   0.55427						
31   0.62026						
32   0.52725						
33   0.46856						
34   0.52417						
35   0.17334						
36   0.78652						
37   0.50004						
38   0.51929						
39   0.84883						
40   0.03448						

2. OLS solution of  $40 \times 3$  linear system, including intercept:

```
print "Newey-West Regression: With Intercept";
pri = 2; tau = 5; intc = 1;
optn = [ pri tau intc ];
< gof,est,nwcof,yhat > = tsmeas(xd2,"nwreg",ran40,optn);
print "Newey-West REG: gof=",gof;
print "Newey-West REG: est=",est;
print "Newey-West REG: nwcof=",nwcof;
print "Newey-West REG: yhat=",yhat;
```



Newey-West REG: gof=		Newey-West REG: yhat=		
	1		1	
-----				
1	0.07753	1	0.55774	
2	0.13355	2	0.59005	
3	0.06134	3	0.37505	
Newey-West REG: est=				
	1	2		
-----				
1	0.46282	2.5038	4	0.62702
2	0.24662	1.9244	5	0.66331
3	-0.30860	-1.5547	6	0.35072
4	0.21950	1.1649	7	0.40135
Newey-West REG: nwcov=				
S	1	2	3	4
-----				
1	0.03417			
2	-0.01549	0.01642		
3	-0.01478	-0.00474	0.03940	
4	-0.02892	0.01066	0.00371	0.03551
14				0.69139
15				0.55733
16				0.65031
17				0.61375
18				0.70733
19				0.61110
20				0.69911
21				0.69947
22				0.27962
23				0.55940
24				0.69463
25				0.38469
26				0.40903
27				0.60724
28				0.64096
29				0.64523
30				0.57107
31				0.56707
32				0.59687
33				0.48506
34				0.46495
35				0.47119
36				0.77809
37				0.60607
38				0.59010
39				0.71883
40				0.32949

#### 4.1.6 (Augmented) Dickey-Fuller Testing

```
< stat,pval,cval,lags,resi > = tsmeas(xdat,"diful"<,optn>)
```

This function performs (augmented) Dickey-Fuller testing with or without automatic lag selection. This function is based on Kevin Sheppard's implementation in the Oxford MFE Toolbox and is using the tables of critical values obtained by simulation with a Matlab script posted on that website. If automatic lag selection based on AIC or BIC is specified, the estimates are returned for the optimal number of lags.

If **xdat** has more than one column, each column is treated as separate time series.

It returns:

**stat** the test statistics (if **xdat** has more than one column this is a vector)

**pval** the probability (if **xdat** has more than one column this is a vector)

**cval** a vector of six critical values corresponding to the following probabilities: [.01.05.1.9.95.99] (if **xdat** has more than one column this is a matrix)

**lags** returns the number of lags to which the the other estimates correspond, i.e. for automatic lag selection the optimal number of lags (must be integer in  $[0, \text{opt}[2]]$ )

**resi** a vector or matrix of residuals.

Note, that the optimal solution may correspond to 0 lags. The following specific options could be specified:

**2** an integer **ntau**  $\geq 1$  specifying how many lags  $\tau = 1, \dots, \text{ntau}$  are to be considered. For automatic lags selection this is the maximum number of lags tried. The default is 1.

**3** specifies the degree of the polynomial included in the ADF regression, i.e. the order of augmentation:

**0** no deterministic terms included

**1** constant is included

**2** time trend is added to constant

**3** Constant, DGP assumed to have a time trend

**4** specifies which criterion is used for the automatic lag determination:

**0** no automatic lag determination (default);

**1** automatic lag determination using AIC;

**2** automatic lag determination using BIC.

All of the following examples use this data set:

```
xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
```

1.  $ndeg = 0$  with specified number of 5 lags:

```
lags = 5; ndeg = 0;
optn = [ 2 lags ndeg ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller: stat=",stat;
print "Dickey-Fuller: prob=",prob;
print "Dickey-Fuller: cval=",cval;
```

Dickey-Fuller Statistics -0.689827, Probability 0.4199 for 5 Lags

Dickey-Fuller: stat=-0.6898

Dickey-Fuller: prob= 0.4199

Dickey-Fuller: cval=

	1	2	3	4	5	6
1	-2.7328	-2.0186	-1.6648	0.94878	1.3721	2.1844

2.  $ndeg = 1$  with specified number of 5 lags:

```
lags = 5; ndeg = 1;
optn = [ 2 lags ndeg ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller: stat=",stat;
print "Dickey-Fuller: prob=",prob;
print "Dickey-Fuller: cval=",cval;
```

Dickey-Fuller Statistics -1.51482, Probability 0.5435 for 5 Lags

Dickey-Fuller: stat=-1.5148

Dickey-Fuller: prob= 0.5435

Dickey-Fuller: cval=

	1	2	3	4	5	6
1	-4.1091	-3.2310	-2.8221	-0.39517	-0.01339	0.71737

3.  $ndeg = 2$  with specified number of 5 lags:

```
lags = 5; ndeg = 2;
optn = [ 2 lags ndeg ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller: stat=",stat;
print "Dickey-Fuller: prob=",prob;
print "Dickey-Fuller: cval=",cval;
```

Dickey-Fuller Statistics -2.24125, Probability 0.5143 for 5 Lags

Dickey-Fuller: stat=-2.2413

Dickey-Fuller: prob= 0.5143

Dickey-Fuller: cval=

	1	2	3	4	5	6
1	-4.7563	-3.8449	-3.4375	-1.2658	-0.93043	-0.22962

4.  $ndeg = 3$  with specified number of 5 lags:

```
lags = 5; ndeg = 3;
optn = [ 2 lags ndeg ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller: stat=",stat;
print "Dickey-Fuller: prob=",prob;
print "Dickey-Fuller: cval=",cval;
```

Dickey-Fuller Statistics -1.78214, Probability 0.0374 for 5 Lags

Dickey-Fuller: stat=-1.7821

Dickey-Fuller: prob= 0.03736

Dickey-Fuller: cval=

	1	2	3	4	5	6
1	-2.3263	-1.6449	-1.2816	1.2816	1.6449	2.3263

5.  $ndeg = 0$  with BIC selection of a maximum of 8 lags:

```
lagmax = 8; ndeg = 0; bic = 2;
optn = [ 2 lagmax ndeg bic ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller BIC: stat=",stat;
print "Dickey-Fuller BIC: prob=",prob;
print "Dickey-Fuller BIC: cval=",cval;
print "Dickey-Fuller BIC: lags=",lags;
```

```
*****
Automatic Lag Search (BIC) for Dickey-Fuller Test
*****
```

Lags	SSE	AIC	BIC
0	0.15601843	-1.80778113	-1.76555914
1	0.14110212	-1.85827139	-1.77382741
2	0.12912306	-1.89698941	-1.77032345
3	0.12888016	-1.84887229	-1.67998434
4	0.10631328	-1.99136509	-1.78025516
5	0.10241622	-1.97871020	-1.72537828
6	0.09598876	-1.99352419	-1.69797029
7	0.08584639	-2.05519570	-1.71741980
8	0.08292584	-2.03980850	-1.65981063

The Smallest BIC Value of -1.78026 Determines Lag Size 4  
Dickey-Fuller Statistics -0.73469, Probability 0.4007 for 4 Lags

Dickey-Fuller BIC: stat=-0.7347

Dickey-Fuller BIC: prob= 0.4007

```

Dickey-Fuller BIC: cval=
|          1          2          3          4          5          6
-----
1 |  -2.7289  -2.0169  -1.6638   0.94707   1.3697   2.1797

Dickey-Fuller BIC: lags= 4.0000

```

6.  $ndeg = 1$  with BIC selection of a maximum of 8 lags:

```

lagmax = 8; ndeg = 1; bic = 2;
optn = [ 2 lagmax ndeg bic ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller BIC: stat=",stat;
print "Dickey-Fuller BIC: prob=",prob;
print "Dickey-Fuller BIC: cval=",cval;
print "Dickey-Fuller BIC: lags=",lags;

```

```

*****
Automatic Lag Search (BIC) for Dickey-Fuller Test
*****

```

Lags	SSE	AIC	BIC
0	0.11179578	-2.09108143	-2.00663745
1	0.11143637	-2.04430150	-1.91763554
2	0.11009370	-2.00642348	-1.83753553
3	0.10451435	-2.00843088	-1.79732094
4	0.09616274	-2.04171333	-1.78838142
5	0.09499131	-2.00396983	-1.70841592
6	0.08321125	-2.08637271	-1.74859682
7	0.07928279	-2.08473426	-1.70473638
8	0.07215951	-2.12887620	-1.70665634

The Smallest BIC Value of -2.00664 Determines Lag Size 0  
Dickey-Fuller Statistics -5.94067, Probability 0.0000 for 0 Lags

```

Dickey-Fuller BIC: stat=-5.9407

```

```

Dickey-Fuller BIC: prob= 0.0000

```

```

Dickey-Fuller BIC: cval=
|          1          2          3          4          5          6

```

```
-----
1 | -4.0210 -3.1847 -2.7907 -0.40155 -0.02262 0.70209

Dickey-Fuller BIC: lags= 0.0000
```

7.  $ndeg = 2$  with BIC selection of a maximum of 8 lags:

```
lagmax = 8; ndeg = 2; bic = 2;
optn = [ 2 lagmax ndeg bic ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller BIC: stat=",stat;
print "Dickey-Fuller BIC: prob=",prob;
print "Dickey-Fuller BIC: cval=",cval;
print "Dickey-Fuller BIC: lags=",lags;
```

```
*****
Automatic Lag Search (BIC) for Dickey-Fuller Test
*****
```

Lags	SSE	AIC	BIC
0	0.09891179	-2.16352685	-2.03686090
1	0.09831659	-2.11956245	-1.95067451
2	0.09828716	-2.06986186	-1.85875193
3	0.08904510	-2.11861232	-1.86528040
4	0.08625527	-2.10044407	-1.80489016
5	0.08613447	-2.05184556	-1.71406967
6	0.07481467	-2.14274130	-1.76274342
7	0.07231972	-2.12665843	-1.70443857
8	0.06426427	-2.19475143	-1.73030958

The Smallest BIC Value of -2.03686 Determines Lag Size 0  
Dickey-Fuller Statistics -6.26798, Probability 0.0000 for 0 Lags

```
Dickey-Fuller BIC: stat=-6.2680
```

```
Dickey-Fuller BIC: prob= 0.0000
```

```
Dickey-Fuller BIC: cval=
```

```

|          1          2          3          4          5          6
-----
1 | -4.6275 -3.7780 -3.3917 -1.2638 -0.93208 -0.24390
```

Dickey-Fuller BIC: lags= 0.0000

8. *ndeg* = 3 with BIC selection of a maximum of 8 lags:

```

lagmax = 8; ndeg = 3; bic = 2;
optn = [ 2 lagmax ndeg bic ];
< stat,prob,cval,lags,resi > = tsmeas(xd2,"diful",optn);
print "Dickey-Fuller BIC: stat=",stat;
print "Dickey-Fuller BIC: prob=",prob;
print "Dickey-Fuller BIC: cval=",cval;
print "Dickey-Fuller BIC: lags=",lags;

```

```

*****
Automatic Lag Search (BIC) for Dickey-Fuller Test
*****

```

	Lags	SSE	AIC	BIC
	0	0.11179578	-2.09108143	-2.00663745
	1	0.11143637	-2.04430150	-1.91763554
	2	0.11009370	-2.00642348	-1.83753553
	3	0.10451435	-2.00843088	-1.79732094
	4	0.09616274	-2.04171333	-1.78838142
	5	0.09499131	-2.00396983	-1.70841592
	6	0.08321125	-2.08637271	-1.74859682
	7	0.07928279	-2.08473426	-1.70473638
	8	0.07215951	-2.12887620	-1.70665634

The Smallest BIC Value of -2.00664 Determines Lag Size 0  
Dickey-Fuller Statistics -6.09299, Probability 0.0000 for 0 Lags

Dickey-Fuller BIC: stat=-6.0930

Dickey-Fuller BIC: prob= 5.541e-010

Dickey-Fuller BIC: cval=

	1	2	3	4	5	6
1	-2.3263	-1.6449	-1.2816	1.2816	1.6449	2.3263

Dickey-Fuller BIC: lags= 0.0000



#### 4.1.7 Granger Causality Testing

```
< stat,pval > = tsmeas(xdat,"grang",tau<,optn>)
```

The function implements the Granger causality test for an  $N$  *times*  $n$  data matrix and for a specified integer vector of  $K$  positive lags  $\tau = \tau_1, \dots, \tau_K$  where  $tmax = \max_k = 1^K(\tau_1, \dots, \tau_K)$ .

It returns:

**stat** the test statistics (if **xdat** has more than one column this is a vector)

**pval** the probability (if **xdat** has more than one column this is a vector)

The following specific options could be specified:

**2** an integer specifying the inference method:

**1** likelihood ratio test (default)

**2** LM test

**3** Wald test

**3** binary integer *intc* specifying whether (=1) or not (=0, is default) a constant intercept variable is used in the model;

**4** binary integer specifying Homoskedastic (=0) or Heteroskedastic (=1, is default);

**5** binary integer specifying uncorrelated (=0, is default) or correlated (=1) error;

All examples work with the same  $200 \times 3$  random generated data set:

```
print "Matlab 200 by 3 RandomData: ran203[200,3]";
options NOECHO;
%inc "..\tdata\matlb.dat";
options ECHO;
```

```
print "Ex 1: infer=1 ihet=1 uncor=0";
tau = [ 1 3 4 ];
pri = 2; infer = 1; intc = 1; ihet = 1; uncor = 0;
optn = [ pri infer intc ihet uncor ];
< stat,prob > = tsmeas(ran203,"grang",tau,optn);
print "STAT=",stat;
print "PROB=",prob;
```

```

STAT=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      3.2738      2.5597      0.85104
Var_2 |      0.28437     0.18616      3.6460
Var_3 |      2.6118      2.1632      1.9429

```

```

PROB=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      0.35130     0.46460     0.83722
Var_2 |      0.96294     0.97979     0.30231
Var_3 |      0.45542     0.53924     0.58435

```

```

print "Ex 2: infer=1 ihet=0 uncor=0";
pri = 2; infer = 1; intc = 1; ihet = 0; uncor = 0;
optn = [ pri infer intc ihet uncor ];
< stat,prob > = tsmeas(ran203,"grang",tau,optn);
print "STAT=",stat;
print "PROB=",prob;

```

```

STAT=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      2.9727      1.9965     0.67241
Var_2 |      0.26558     0.15035     3.0786
Var_3 |      2.3809      1.9809     1.4207

```

```

PROB=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      0.39586     0.57313     0.87967
Var_2 |      0.96637     0.98518     0.37966
Var_3 |      0.49720     0.57639     0.70069

```

```

print "Ex 3: infer=2 ihet=0 uncor=1";
pri = 2; infer = 2; intc = 1; ihet = 0; uncor = 1;
optn = [ pri infer intc ihet uncor ];
< stat,prob > = tsmeas(ran203,"grang",tau,optn);
print "STAT=",stat;
print "PROB=",prob;

```

```

STAT=
      |      Var_1      Var_2      Var_3
-----
Var_1 |  0.07282  0.01647  0.00751
Var_2 |  0.00216  0.00826  0.01171
Var_3 |  0.08615  0.03545  0.02027

```

```

PROB=
      |      Var_1      Var_2      Var_3
-----
Var_1 |  0.99489  0.99944  0.99983
Var_2 |  0.99997  0.99980  0.99966
Var_3 |  0.99345  0.99824  0.99924

```

```

print "Ex 4: infer=3 ihet=1 uncor=0";
pri = 2; infer = 3; intc = 1; ihet = 1; uncor = 0;
optn = [ pri infer intc ihet uncor ];
< stat,prob > = tsmeas(ran203,"grang",tau,optn);
print "STAT=",stat;
print "PROB=",prob;

```

```

STAT=
      |      Var_1      Var_2      Var_3
-----
Var_1 |  30.036  55.127  0.18754
Var_2 |   1.7373  4.0050  2.0567
Var_3 |   1.8878  1.7107  2.0502

```

```

PROB=
      |      Var_1      Var_2      Var_3
-----
Var_1 |  0.00000  6e-012  0.97958
Var_2 |  0.62866  0.26093  0.56071
Var_3 |  0.59603  0.63456  0.56205

```

```

print "Ex 5: infer=3 ihet=1 uncor=1";
pri = 2; infer = 3; intc = 1; ihet = 1; uncor = 1;
optn = [ pri infer intc ihet uncor ];
< stat,prob > = tsmeas(ran203,"grang",tau,optn);
print "STAT=",stat;
print "PROB=",prob;

```

```

STAT=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      30.036      55.127      0.18754
Var_2 |       1.7373       4.0050       2.0567
Var_3 |       1.8878       1.7107       2.0502

```

```

PROB=
      |      Var_1      Var_2      Var_3
-----
Var_1 |      0.00000      6e-012      0.97958
Var_2 |      0.62866      0.26093      0.56071
Var_3 |      0.59603      0.63456      0.56205

```

#### 4.1.8 Vector AR Modeling

```

< est,ase,r2s,cov,resi > = tsmeas(xdat,"arvec",tau<,optn>)

```

The function implements (multivariate) vector AR modeling for an  $N$  times  $n$  data matrix and for a specified integer vector of  $K$  positive lags  $\tau = \tau_1, \dots, \tau_K$  where  $tmax = \max_k = 1^K(\tau_1, \dots, \tau_K)$ .

There are  $p = (intc + K * n) * n$  parameters to fit.

It returns:

**est**  $(intc + K * n) \times n$  matrix of parameter estimates;

**ase**  $(intc + K * n) \times 3 * n$  matrix of asymptotic standard errors,  $t$  values, and probabilities;

**r2s**  $n \times n + 1$  matrix containing the  $n$  R-squared values in its first column and the  $n \times n$  matrix of sums of squared errors in its last  $n$  columns;

**cov** the symmetric  $p \times p$  covariance matrix of estimates;

**resi** a  $N$  vector or  $N \times n$  matrix of residuals.

The following specific options could be specified:

**2** unused;

**3** binary integer *intc* specifying whether (=1) or not (=0, is default) a constant intercept variable is used in the model;

**4** binary integer specifying Homoskedastic (=0) or Heteroskedastic (=1, is default);

**5** binary integer specifying uncorrelated (=0, is default) or correlated (=1) error;

All examples work with the same  $200 \times 3$  random generated data set:

```

print "Matlab 200 by 3 RandomData: ran203[200,3]";
options NOECHO;
%inc "..\tdata\matlb.dat";
options ECHO;

print "Ex 1.1: intc=1: ihet=1 uncor=0: tau=[ 1 4 ]";
tau = [ 1 4 ];
pri = 2; intc = 1; ihet = 1; uncor = 0;
optn = [ pri . intc ihet uncor ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"arvec",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

	Var_1	Var_2	Var_3
Constant	0.57015	0.45458	0.57867
Var_1_Lag_1	-0.02170	-0.02963	-0.05872
Var_1_Lag_2	-0.02145	-0.01881	-0.04751
Var_2_Lag_1	-0.03458	0.02628	0.05220
Var_2_Lag_2	0.01759	0.01559	-0.11212
Var_3_Lag_1	-0.11315	-0.02130	0.05411
Var_3_Lag_2	0.02703	0.10050	-0.05293

ASE=

	ASE_1	ASE_2	ASE_3	T_1	T_2
Constant	0.10354	0.09385	0.09510	5.5064	-0.20038
Var_1_Lag_1	0.07624	0.07069	0.06236	5.9625	-0.67202
Var_1_Lag_2	0.07035	0.07366	0.06548	8.2255	-0.46939
Var_2_Lag_1	0.07735	0.07770	0.07132	-0.28052	0.33827
Var_2_Lag_2	0.07406	0.06606	0.06734	-0.40005	0.79026
Var_3_Lag_1	0.07531	0.06917	0.06825	-0.77972	0.25424
Var_3_Lag_2	0.08528	0.07872	0.07435	-0.25154	0.19803

	T_3	P_1	P_2	P_3
Constant	-1.1790	4e-008	0.84118	0.23839

Var_1_Lag_1		-1.8146	2e-009	0.50157	0.06958
Var_1_Lag_2		-0.32534	2e-016	0.63879	0.74492
Var_2_Lag_1		0.75864	0.77908	0.73516	0.44807
Var_2_Lag_2		0.40134	0.68912	0.42938	0.68817
Var_3_Lag_1		1.4725	0.43556	0.79931	0.14088
Var_3_Lag_2		-0.71191	0.80140	0.84302	0.47652

R2S=

		R2	SE2_1	SE2_2	SE2_3
Constant		0.01380	0.08493	-0.00081	-0.01204
Var_1_Lag_1		0.01073	-0.00081	0.07861	0.00077
Var_1_Lag_2		0.02752	-0.01204	0.00077	0.06652

```

print "Ex 1.3: intc=1: ihet=1 uncor=1: tau=[ 1 4 ]";
tau = [ 1 4 ];
pri = 2; intc = 1; ihet = 1; uncor = 1;
optn = [ pri . intc ihet uncor ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"arvec",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

		Var_1	Var_2	Var_3
Constant		0.57015	0.45458	0.57867
Var_1_Lag_1		-0.02170	-0.02963	-0.05872
Var_1_Lag_2		-0.02145	-0.01881	-0.04751
Var_2_Lag_1		-0.03458	0.02628	0.05220
Var_2_Lag_2		0.01759	0.01559	-0.11212
Var_3_Lag_1		-0.11315	-0.02130	0.05411
Var_3_Lag_2		0.02703	0.10050	-0.05293

ASE=

		ASE_1	ASE_2	ASE_3	T_1	T_2
Constant		0.10354	0.09385	0.09510	5.5064	-0.20038
Var_1_Lag_1		0.07624	0.07069	0.06236	5.9625	-0.67202
Var_1_Lag_2		0.07035	0.07366	0.06548	8.2255	-0.46939
Var_2_Lag_1		0.07735	0.07770	0.07132	-0.28052	0.33827

Var_2_Lag_2		0.07406	0.06606	0.06734	-0.40005	0.79026
Var_3_Lag_1		0.07531	0.06917	0.06825	-0.77972	0.25424
Var_3_Lag_2		0.08528	0.07872	0.07435	-0.25154	0.19803
		T_3	P_1	P_2	P_3	
-----						
Constant		-1.1790	4e-008	0.84118	0.23839	
Var_1_Lag_1		-1.8146	2e-009	0.50157	0.06958	
Var_1_Lag_2		-0.32534	2e-016	0.63879	0.74492	
Var_2_Lag_1		0.75864	0.77908	0.73516	0.44807	
Var_2_Lag_2		0.40134	0.68912	0.42938	0.68817	
Var_3_Lag_1		1.4725	0.43556	0.79931	0.14088	
Var_3_Lag_2		-0.71191	0.80140	0.84302	0.47652	

```

print "Ex 2.1: intc=0: ihet=1 uncor=0: tau=[ 1 4 ]";
tau = [ 1 4 ];
pri = 2; intc = 0; ihet = 1; uncor = 0;
optn = [ pri . intc ihet uncor ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"arvec",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

		Var_1	Var_2	Var_3
-----				
Var_1_Lag_1		0.13654	0.09653	0.10188
Var_1_Lag_2		0.13901	0.10913	0.11536
Var_2_Lag_1		0.19367	0.20827	0.28386
Var_2_Lag_2		0.19434	0.15651	0.06727
Var_3_Lag_1		0.04117	0.10174	0.21073
Var_3_Lag_2		0.25120	0.27924	0.17460

ASE=

		ASE_1	ASE_2	ASE_3	T_1	T_2
-----						
Var_1_Lag_1		0.07368	0.06868	0.06076	1.8532	2.8201
Var_1_Lag_2		0.06605	0.06990	0.06476	1.4616	2.9796
Var_2_Lag_1		0.07002	0.06924	0.06170	1.4551	4.1000
Var_2_Lag_2		0.06978	0.06523	0.06153	1.9922	2.9791
Var_3_Lag_1		0.07550	0.06921	0.06660	1.4455	2.2614

Var_3_Lag_2		0.07819	0.07063	0.07340	1.4753	0.95240
		T_3	P_1	P_2	P_3	
-----						
Var_1_Lag_1		0.67757	0.06386	0.00480	0.49805	
Var_1_Lag_2		1.5710	0.14386	0.00289	0.11618	
Var_2_Lag_1		3.4152	0.14564	0.00004	0.00064	
Var_2_Lag_2		4.0824	0.04635	0.00289	0.00004	
Var_3_Lag_1		4.1930	0.14831	0.02373	0.00003	
Var_3_Lag_2		2.3786	0.14014	0.34089	0.01738	

```

print "Ex 2.3: intc=0: ihet=1 uncor=1: tau=[ 1 4 ]";
tau = [ 1 4 ];
pri = 2; intc = 0; ihet = 1; uncor = 1;
optn = [ pri . intc ihet uncor ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"arvec",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=		Var_1	Var_2	Var_3		
-----						
Var_1_Lag_1		0.13654	0.09653	0.10188		
Var_1_Lag_2		0.13901	0.10913	0.11536		
Var_2_Lag_1		0.19367	0.20827	0.28386		
Var_2_Lag_2		0.19434	0.15651	0.06727		
Var_3_Lag_1		0.04117	0.10174	0.21073		
Var_3_Lag_2		0.25120	0.27924	0.17460		
-----						
ASE=		ASE_1	ASE_2	ASE_3	T_1	T_2
-----						
Var_1_Lag_1		0.07368	0.06868	0.06076	1.8532	2.8201
Var_1_Lag_2		0.06605	0.06990	0.06476	1.4616	2.9796
Var_2_Lag_1		0.07002	0.06924	0.06170	1.4551	4.1000
Var_2_Lag_2		0.06978	0.06523	0.06153	1.9922	2.9791
Var_3_Lag_1		0.07550	0.06921	0.06660	1.4455	2.2614
Var_3_Lag_2		0.07819	0.07063	0.07340	1.4753	0.95240



	T_3	P_1	P_2	P_3
Var_1_Lag_1	0.67757	0.06386	0.00480	0.49805
Var_1_Lag_2	1.5710	0.14386	0.00289	0.11618
Var_2_Lag_1	3.4152	0.14564	0.00004	0.00064
Var_2_Lag_2	4.0824	0.04635	0.00289	0.00004
Var_3_Lag_1	4.1930	0.14831	0.02373	0.00003
Var_3_Lag_2	2.3786	0.14014	0.34089	0.01738

#### 4.1.9 Impulse Response Modeling

```
< est,ase,r2s,cov,resi > = tsmeas(xdat,"iresp",tau<,optn>)
```

The function implements (multivariate) impulse response modeling for an  $N \times n$  data matrix, for a specified value of lead, and for a specified integer vector of  $K$  positive lags  $\tau = \tau_1, \dots, \tau_K$  where  $tmax = \max_k = 1^K(\tau_1, \dots, \tau_K)$ .

There are  $p = (intc + tmax * n) * n$  parameters to fit.

It returns:

**est**  $(intc + tmax * n) \times n$  matrix of parameter estimates;

**ase**  $(intc + tmax * n) \times n$  matrix of asymptotic standard errors;

**r2s**  $n \times n + 1$  matrix containing the  $n$  R-squared values in its first column and the  $n \times n$  matrix used for **optn**[6] type scaling in its last  $n$  columns;

**cov** the symmetric  $p \times p$  covariance matrix of estimates;

**resi** a  $N$  vector or  $N \times n$  matrix of residuals.

The following specific options could be specified:

**2** positive integer specifying the number of leads (default is 1);

**3** binary integer *intc* specifying whether (=1) or not (=0, is default) a constant intercept variable is used in the model;

**4** binary integer specifying Homoskedastic (=0) or Heteroskedastic (=1, is default);

**5** binary integer specifying uncorrelated (=0, is default) or correlated (=1) error;

**6** positive integer specifying the type of shocks:

**0** unit (unscaled) shocks: use  $n \times n$  identity matrix;

**1** scaled but uncorrelated shocks: use square root of diagonal values of  $n \times n$  SSE matrix of vector AR model, this is the default;

- 2 scaled and correlated shocks: use Cholesky factor of  $n \times n$  SSE matrix of vector AR model,
- 3 scaled and correlated shocks (spectral decomposition): use square roots of all entries of  $n \times n$  SSE matrix of vector AR model.

All examples work with the same  $200 \times 3$  random generated data set:

```
print "Matlab 200 by 3 RandomData: ran203[200,3]";
options NOECHO;
%inc "..\tdata\matlb.dat";
options ECHO;

print "Ex 0.2: intc=1: ihet=1 uncor=0 sqtyp=0: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 0; sqtyp = 0;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;
```

EST=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	1.00000	0.00000	0.00000	0.00000	1.00000
Lead_1	-0.02467	-0.03246	-0.04451	-0.01073	-0.02738
Lead_2	0.00242	0.00030	0.00010	0.00195	-0.00023
Lead_3	-0.15225	0.02240	-0.00580	0.07529	-0.01507
Lead_4	0.00160	0.01018	0.01235	-0.00221	0.00093
Lead_5	-0.00089	0.00009	0.00010	-0.00026	0.00012

  

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	0.00000	0.00000	0.00000	1.00000
Lead_1	-0.04249	-0.03279	0.03131	0.05340
Lead_2	-0.00063	-0.00128	0.00188	0.00298
Lead_3	-0.09679	0.07659	-0.13703	-0.07969
Lead_4	-0.00178	0.01363	-0.00976	-0.00894
Lead_5	0.00003	0.00063	-0.00092	-0.00108

ASE=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	0.00000	0.00000	0.00000	0.00000	0.00000
Lead_1	0.07385	0.06916	0.07486	0.06948	0.07358
Lead_2	0.00529	0.00509	0.00540	0.00593	0.00619
Lead_3	0.07101	0.07005	0.07817	0.06932	0.06783
Lead_4	0.02534	0.01869	0.02245	0.01672	0.01554
Lead_5	0.00206	0.00175	0.00237	0.00188	0.00155

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	0.00000	0.00000	0.00000	0.00000
Lead_1	0.07843	0.06274	0.06483	0.07103
Lead_2	0.00597	0.00534	0.00546	0.00905
Lead_3	0.07831	0.06380	0.06626	0.06801
Lead_4	0.01606	0.01870	0.01507	0.01901
Lead_5	0.00219	0.00229	0.00205	0.00274

R2S=

	R2	SE2_1	SE2_2	SE2_3
Var_1	0.03679	1.00000	0.00000	0.00000
Var_2	0.01942	0.00000	1.00000	0.00000
Var_3	0.02485	0.00000	0.00000	1.00000

```

print "Ex 0.3: intc=1: ihet=1 uncor=1 sqtyp=0: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 1; sqtyp = 0;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	1.00000	0.00000	0.00000	0.00000	1.00000
Lead_1	-0.02467	-0.03246	-0.04451	-0.01073	-0.02738
Lead_2	0.00242	0.00030	0.00010	0.00195	-0.00023
Lead_3	-0.15225	0.02240	-0.00580	0.07529	-0.01507

Lead_4		0.00160	0.01018	0.01235	-0.00221	0.00093
Lead_5		-0.00089	0.00009	0.00010	-0.00026	0.00012
		Var_2_3	Var_3_1	Var_3_2	Var_3_3	
-----						
Lead_0		0.00000	0.00000	0.00000	1.00000	
Lead_1		-0.04249	-0.03279	0.03131	0.05340	
Lead_2		-0.00063	-0.00128	0.00188	0.00298	
Lead_3		-0.09679	0.07659	-0.13703	-0.07969	
Lead_4		-0.00178	0.01363	-0.00976	-0.00894	
Lead_5		0.00003	0.00063	-0.00092	-0.00108	

ASE=

		Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
-----						
Lead_0		0.00000	0.00000	0.00000	0.00000	0.00000
Lead_1		0.07385	0.06916	0.07486	0.06948	0.07358
Lead_2		0.00527	0.00518	0.00524	0.00591	0.00607
Lead_3		0.07101	0.07005	0.07816	0.06930	0.06781
Lead_4		0.02512	0.01799	0.02129	0.01645	0.01527
Lead_5		0.00203	0.00172	0.00228	0.00183	0.00141
		Var_2_3	Var_3_1	Var_3_2	Var_3_3	
-----						
Lead_0		0.00000	0.00000	0.00000	0.00000	
Lead_1		0.07843	0.06274	0.06483	0.07103	
Lead_2		0.00589	0.00520	0.00548	0.00895	
Lead_3		0.07833	0.06384	0.06627	0.06800	
Lead_4		0.01549	0.01780	0.01481	0.01812	
Lead_5		0.00202	0.00221	0.00193	0.00256	

```

print "Ex 1.2: intc=1: ihet=1 uncor=0 sqtyp=1: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 0; sqtyp = 1;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

```

EST=
      |  Var_1_1  Var_1_2  Var_1_3  Var_2_1  Var_2_2
-----
Lead_0 |  0.28801  0.00000  0.00000  0.00000  0.27914
Lead_1 | -0.00711 -0.00935 -0.01282 -0.00299 -0.00764
Lead_2 |  0.00070  0.00009  0.00003  0.00054 -0.00006
Lead_3 | -0.04385  0.00645 -0.00167  0.02102 -0.00421
Lead_4 |  0.00046  0.00293  0.00356 -0.00062  0.00026
Lead_5 | -0.00026  0.00003  0.00003 -0.00007  0.00003

      |  Var_2_3  Var_3_1  Var_3_2  Var_3_3
-----
Lead_0 |  0.00000  0.00000  0.00000  0.25827
Lead_1 | -0.01186 -0.00847  0.00809  0.01379
Lead_2 | -0.00018 -0.00033  0.00049  0.00077
Lead_3 | -0.02702  0.01978 -0.03539 -0.02058
Lead_4 | -0.00050  0.00352 -0.00252 -0.00231
Lead_5 |  0.00001  0.00016 -0.00024 -0.00028

```

```

ASE=
      |  Var_1_1  Var_1_2  Var_1_3  Var_2_1  Var_2_2
-----
Lead_0 |  0.00000  0.00000  0.00000  0.00000  0.00000
Lead_1 |  0.02127  0.01992  0.02156  0.01939  0.02054
Lead_2 |  0.00152  0.00147  0.00155  0.00166  0.00173
Lead_3 |  0.02045  0.02018  0.02251  0.01935  0.01894
Lead_4 |  0.00730  0.00538  0.00647  0.00467  0.00434
Lead_5 |  0.00059  0.00050  0.00068  0.00052  0.00043

      |  Var_2_3  Var_3_1  Var_3_2  Var_3_3
-----
Lead_0 |  0.00000  0.00000  0.00000  0.00000
Lead_1 |  0.02189  0.01620  0.01674  0.01834
Lead_2 |  0.00167  0.00138  0.00141  0.00234
Lead_3 |  0.02186  0.01648  0.01711  0.01756
Lead_4 |  0.00448  0.00483  0.00389  0.00491
Lead_5 |  0.00061  0.00059  0.00053  0.00071

```

```

print "Ex 1.3: intc=1: ihet=1 uncor=1 sqtyp=1: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 1; sqtyp = 1;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);

```

```

print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	0.28801	0.00000	0.00000	0.00000	0.27914
Lead_1	-0.00711	-0.00935	-0.01282	-0.00299	-0.00764
Lead_2	0.00070	0.00009	0.00003	0.00054	-0.00006
Lead_3	-0.04385	0.00645	-0.00167	0.02102	-0.00421
Lead_4	0.00046	0.00293	0.00356	-0.00062	0.00026
Lead_5	-0.00026	0.00003	0.00003	-0.00007	0.00003

  

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	0.00000	0.00000	0.00000	0.25827
Lead_1	-0.01186	-0.00847	0.00809	0.01379
Lead_2	-0.00018	-0.00033	0.00049	0.00077
Lead_3	-0.02702	0.01978	-0.03539	-0.02058
Lead_4	-0.00050	0.00352	-0.00252	-0.00231
Lead_5	0.00001	0.00016	-0.00024	-0.00028

ASE=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	0.00000	0.00000	0.00000	0.00000	0.00000
Lead_1	0.02127	0.01992	0.02156	0.01939	0.02054
Lead_2	0.00152	0.00149	0.00151	0.00165	0.00169
Lead_3	0.02045	0.02018	0.02251	0.01934	0.01893
Lead_4	0.00724	0.00518	0.00613	0.00459	0.00426
Lead_5	0.00059	0.00050	0.00066	0.00051	0.00039

  

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	0.00000	0.00000	0.00000	0.00000
Lead_1	0.02189	0.01620	0.01674	0.01834
Lead_2	0.00165	0.00134	0.00142	0.00231
Lead_3	0.02186	0.01649	0.01711	0.01756
Lead_4	0.00432	0.00460	0.00382	0.00468
Lead_5	0.00056	0.00057	0.00050	0.00066

```

print "Ex 2.2: intc=1: ihet=1 uncor=0 sqtyp=2: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 0; sqtyp = 2;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;

```

EST=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	0.28801	0.00311	-0.04061	0.00311	0.27913
Lead_1	-0.00581	-0.01071	-0.01512	-0.00302	-0.00779
Lead_2	0.00075	0.00001	-0.00009	0.00055	-0.00007
Lead_3	-0.04672	0.01197	0.00126	0.02043	-0.00394
Lead_4	-0.00010	0.00333	0.00391	-0.00063	0.00031
Lead_5	-0.00028	0.00006	0.00007	-0.00008	0.00004

  

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	-0.00143	-0.04061	-0.00143	0.25505
Lead_1	-0.01207	-0.00735	0.00934	0.01549
Lead_2	-0.00018	-0.00043	0.00047	0.00076
Lead_3	-0.02692	0.02561	-0.03584	-0.01995
Lead_4	-0.00045	0.00342	-0.00290	-0.00278
Lead_5	0.00001	0.00020	-0.00024	-0.00028

ASE=

	Var_1_1	Var_1_2	Var_1_3	Var_2_1	Var_2_2
Lead_0	0.00000	0.00000	0.00000	0.00000	0.00000
Lead_1	0.02153	0.02042	0.02216	0.01940	0.02053
Lead_2	0.00144	0.00142	0.00174	0.00166	0.00173
Lead_3	0.02087	0.02063	0.02300	0.01935	0.01894
Lead_4	0.00749	0.00563	0.00683	0.00462	0.00431
Lead_5	0.00061	0.00054	0.00076	0.00052	0.00043

  

	Var_2_3	Var_3_1	Var_3_2	Var_3_3
Lead_0	-0.00143	-0.04061	-0.00143	0.25505
Lead_1	-0.01207	-0.00735	0.00934	0.01549
Lead_2	-0.00018	-0.00043	0.00047	0.00076
Lead_3	-0.02692	0.02561	-0.03584	-0.01995
Lead_4	-0.00045	0.00342	-0.00290	-0.00278
Lead_5	0.00001	0.00020	-0.00024	-0.00028

```
-----
Lead_0 | 0.00000 0.00000 0.00000 0.00000
Lead_1 | 0.02188 0.01640 0.01712 0.01878
Lead_2 | 0.00166 0.00127 0.00134 0.00241
Lead_3 | 0.02187 0.01682 0.01742 0.01799
Lead_4 | 0.00442 0.00509 0.00418 0.00531
Lead_5 | 0.00060 0.00059 0.00056 0.00076
```

```
print "Ex 2.3: intc=1: ihet=1 uncor=1 sqtyp=2: tau=[ 1 3 ]";
tau = [ 1 3 ];
pri = 2; lead = 5; intc = 1; ihet = 1; uncor = 1; sqtyp = 2;
optn = [ pri lead intc ihet uncor sqtyp ];
< est,ase,r2s,cov,resi > = tsmeas(ran203,"iresp",tau,optn);
print "EST=",est;
print "ASE=",ase;
print "R2S=",r2s;
print "COV=",cov;
print "Resi=",resi;
```

EST=

```
-----
          |  Var_1_1  Var_1_2  Var_1_3  Var_2_1  Var_2_2
-----
Lead_0 |  0.28801  0.00311 -0.04061  0.00311  0.27913
Lead_1 | -0.00581 -0.01071 -0.01512 -0.00302 -0.00779
Lead_2 |  0.00075  0.00001 -0.00009  0.00055 -0.00007
Lead_3 | -0.04672  0.01197  0.00126  0.02043 -0.00394
Lead_4 | -0.00010  0.00333  0.00391 -0.00063  0.00031
Lead_5 | -0.00028  0.00006  0.00007 -0.00008  0.00004

          |  Var_2_3  Var_3_1  Var_3_2  Var_3_3
-----
Lead_0 | -0.00143 -0.04061 -0.00143  0.25505
Lead_1 | -0.01207 -0.00735  0.00934  0.01549
Lead_2 | -0.00018 -0.00043  0.00047  0.00076
Lead_3 | -0.02692  0.02561 -0.03584 -0.01995
Lead_4 | -0.00045  0.00342 -0.00290 -0.00278
Lead_5 |  0.00001  0.00020 -0.00024 -0.00028
```

ASE=

```
-----
          |  Var_1_1  Var_1_2  Var_1_3  Var_2_1  Var_2_2
-----
Lead_0 |  0.00000  0.00000  0.00000  0.00000  0.00000
Lead_1 |  0.02142  0.02009  0.02176  0.01940  0.02054
```



Lead_2		0.00143	0.00144	0.00167	0.00165	0.00169
Lead_3		0.02062	0.02036	0.02268	0.01934	0.01893
Lead_4		0.00737	0.00533	0.00638	0.00455	0.00424
Lead_5		0.00059	0.00053	0.00072	0.00051	0.00039
		Var_2_3	Var_3_1	Var_3_2	Var_3_3	
-----						
Lead_0		0.00000	0.00000	0.00000	0.00000	
Lead_1		0.02189	0.01628	0.01677	0.01837	
Lead_2		0.00164	0.00123	0.00134	0.00237	
Lead_3		0.02186	0.01654	0.01714	0.01763	
Lead_4		0.00428	0.00480	0.00399	0.00495	
Lead_5		0.00056	0.00057	0.00052	0.00071	

## 4.2 Extension to the `tstrans` Function

We have added a number of more functions:

”**boxc**” Box-Cox Transform

”**lagd**” Lag Difference Transform

”**logd**” Log Difference Transform

”**baki**” Baxter-King Filtering

”**hopr**” Hodrick-Prescott Filtering

### 4.2.1 Box-Cox, Lag- and Log-Difference Transforms

The old form of the `tstrans()` function

```
ydat = tstrans(xdat,"sopt"<,>optn>)
```

was extended with an optional input argument `add` to

```
ydat = tstrans(xdat,"sopt"<,>optn<,>add>>)
```

An additional argument was necessary to compute Box-Cox, Log and Lag Difference transformations. Only these three new transforms use the additional input argument. The following are the new "sopt" transformations:

”**boxc**” Box-Cox transform: the input argument specifies a real scalar or a vector with  $\lambda$  values:

$$y = \begin{cases} x^{\lambda-1}/\lambda & \text{for } \lambda \neq 0 \\ \log(x) & \text{for } \lambda = 0 \text{ and } x_{min} > 0 \\ \log(x + x_{min} + 1) & \text{for } \lambda = 0 \text{ and } x_{min} \leq 0 \end{cases}$$

”**lagd**” Lag difference transform: the last input argument specifies an integer scalar or a vector with lag values:

$$y_i = x_{i+lag} - x_i \quad , \quad i = 1, \dots, n - lag$$

”**logd**” Log difference transform: the last input argument specifies an integer scalar or a vector with lag values:

$$y_i = \log(x_{i+lag}) - \log(x_i) \quad , \quad i = 1, \dots, n - lag$$

”**boxc**” Example for Box-Cox transform:

```
xd1 = [ 0.9501  0.2311  0.6068  0.4860  0.8913
        0.7621  0.4565  0.0185  0.8214  0.4447
        0.6154  0.7919  0.9218  0.7382  0.1763
        0.4057  0.9355  0.9169  0.4103  0.8936 ];
```

```
optn = [ 2 ]; lamb = [ 1. 0. 2. ];
trans51 = tstrans(xd1,"boxc",optn,lamb);
print "Trans51=",trans51;
```

```
Trans51=
|          1          2          3
-----
1 | -0.04990 -0.05119 -0.04865
2 | -0.76890 -1.4649  -0.47330
3 | -0.39320 -0.49956 -0.31590
4 | -0.51400 -0.72155 -0.38190
5 | -0.10870 -0.11507 -0.10279
6 | -0.23790 -0.27168 -0.20960
7 | -0.54350 -0.78417 -0.39580
8 | -0.98150 -3.9900  -0.49983
9 | -0.17860 -0.19675 -0.16265
10 | -0.55530 -0.81036 -0.40112
11 | -0.38460 -0.48548 -0.31064
12 | -0.20810 -0.23332 -0.18645
13 | -0.07820 -0.08143 -0.07514
14 | -0.26180 -0.30354 -0.22753
15 | -0.82370 -1.7356  -0.48446
16 | -0.59430 -0.90214 -0.41770
17 | -0.06450 -0.06667 -0.06242
18 | -0.08310 -0.08676 -0.07965
19 | -0.58970 -0.89087 -0.41583
20 | -0.10640 -0.11250 -0.10074
```

**"lagd"** Example for Lag difference transform:

```
optn = [ 2 ]; lags = [ 0 1 2 ];  
trans61 = tstrans(xd1,"lagd",optn,lags);  
print "Trans61=",trans61;
```

```
Trans61=  
  |      1      2      3  
-----  
1 |  0.00000  -0.71900  -0.34330  
2 |  0.00000   0.37570   0.25490  
3 |  0.00000  -0.12080   0.28450  
4 |  0.00000   0.40530   0.27610  
5 |  0.00000  -0.12920  -0.43480  
6 |  0.00000  -0.30560  -0.74360  
7 |  0.00000  -0.43800   0.36490  
8 |  0.00000   0.80290   0.42620  
9 |  0.00000  -0.37670  -0.20600  
10 | 0.00000   0.17070   0.34720  
11 | 0.00000   0.17650   0.30640  
12 | 0.00000   0.12990  -0.05370  
13 | 0.00000  -0.18360  -0.74550  
14 | 0.00000  -0.56190  -0.33250  
15 | 0.00000   0.22940   0.75920  
16 | 0.00000   0.52980   0.51120  
17 | 0.00000  -0.01860  -0.52520  
18 | 0.00000  -0.50660  -0.02330  
19 | 0.00000   0.48330   .  
20 | 0.00000   .       .
```

**"logd"** Example for Log difference transform:

```
optn = [ 2 ]; lags = [ 0 1 2 ];  
trans71 = tstrans(xd1,"logd",optn,lags);  
print "Trans71=",trans71;
```

```
Trans71=  
  |      1      2      3  
-----  
1 | -1e-018  -1.4137  -0.44837  
2 | -5e-017   0.96535   0.74336
```

```

3 | 1e-017 -0.22199 0.38448
4 | 3e-017 0.60647 0.44987
5 | 6e-018 -0.15660 -0.66909
6 | 1e-017 -0.51249 -3.7183
7 | 2e-017 -3.2058 0.58742
8 | 1e-016 3.7932 3.1796
9 | 1e-017 -0.61361 -0.28874
10 | -3e-017 0.32487 0.57704
11 | -5e-018 0.25216 0.40406
12 | -6e-018 0.15189 -0.07022
13 | 6e-018 -0.22211 -1.6541
14 | -3e-017 -1.4320 -0.59860
15 | 6e-017 0.83343 1.6689
16 | -4e-017 0.83547 0.81538
17 | -2e-018 -0.02008 -0.82419
18 | 4e-019 -0.80411 -0.02574
19 | -2e-017 0.77837 .
20 | -5e-018 . .

```

#### 4.2.2 Baxter-King and Hodrick-Prescott Filtering

```

< ydat,cycl,noise > = tstrans(xdat,"baki"<,optn>)

```

The following specific options could be specified.

- 2 Number of periods to use in the higher frequency filter (e.g. 6 for quarterly data). Must be at least 2. The default is 6.
- 3 Number of periods to use in the lower frequency filter (e.g. 32 for quarterly data). There must be [3] >= [2]. The default is 32.
- 4 Number of points to use in the finite approximation bandpass filter. The filter throws away the first and last K points. The default is 12.

```

xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xc = xd2 - univar(xd2,"ari");

```

```

nphi = 6; nplo = 18; nkpt = 12;

```

```

optn = [ 2 nphi nplo nkpt ];
< trnd,cycl,nois > = tstrans(xc,"baki",optn);
print "Trend,Cycle,Noise=",trnd -> cycl -> nois;

```

Trend,Cycle,Noise=

L	1	2	3
1	0.24741	0	0
2	0.33851	0.00000	0
3	-0.44029	0.00000	0.00000
4	0.34611	0.00000	0.00000
5	0.06511	0.00000	0.00000
6	-0.46979	0.00000	0.00000
7	-0.28879	0.00000	0.00000
8	-0.02039	0.00000	0.00000
9	0.39021	0.00000	0.00000
10	0.39761	0.00000	0.00000
11	-0.40969	0.00000	0.00000
12	0.40331	0.00000	0.00000
13	0.06946	0.08202	0.23842
14	0.08122	-0.07580	-0.08732
15	0.09513	-0.19182	0.32970
16	0.09570	-0.17032	-0.35077
17	0.10864	-0.07158	-0.18256
18	0.12467	0.03503	0.18871
19	0.10820	0.12761	-0.01091
20	0.11169	0.08479	0.19573
21	0.11225	-0.01655	-0.00730
22	0.12558	-0.11698	-0.54019
23	0.13092	-0.11240	0.26328
24	0.08668	0.05017	0.22986
25	0.06173	0.15960	-0.10992
26	0.04178	0.16433	-0.01571
27	-0.01183	0.13531	0.05233
28	-0.17509	0.00000	0.00000
29	0.08821	0.00000	0.00000
30	-0.39609	0.00000	0.00000
31	0.13871	0.00000	0.00000
32	-0.53549	0.00000	0.00000
33	-0.29039	0.00000	0.00000
34	-0.52109	0.00000	0.00000
35	-0.47019	0.00000	0.00000
36	0.25621	0.00000	0.00000
37	0.12751	0.00000	0.00000

```

38 | -0.25019  0.00000  0.00000
39 |  0.38291  0.00000  0.00000
40 | -0.53289  0.00000  0.00000

```

```
< ydat,cycl > = tstrans(xdat,"hopr"<,optn>)
```

The following specific options could be specified.

**2** Positive, scalar integer containing the smoothing parameter of the HP filter. The default is 1600.

```

xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ]';
xc = xd2 - univar(xd2,"ari");

```

```

tlam = 100;
optn = [ 2 tlam ];
< trnd,cycl > = tstrans(xc,"hopr",optn);
print "Trend,Cycle=",trnd -> cycl;

```

Trend,Cycle=

	1	2
1	0.09799	0.14942
2	0.07174	0.26677
3	0.04698	-0.48727
4	0.02787	0.31824
5	0.01371	0.05140
6	0.00697	-0.47676
7	0.01063	-0.29942
8	0.02291	-0.04330
9	0.03903	0.35118
10	0.05379	0.34382
11	0.06548	-0.47518
12	0.07586	0.32745
13	0.08189	0.30801

14	0.08386	-0.16575
15	0.08510	0.14791
16	0.08730	-0.51269
17	0.09362	-0.23912
18	0.10211	0.24630
19	0.10841	0.11650
20	0.11062	0.28159
21	0.10803	-0.01962
22	0.10272	-0.63431
23	0.09659	0.18522
24	0.08519	0.28151
25	0.06593	0.04548
26	0.03902	0.15139
27	0.00514	0.17067
28	-0.03353	-0.14156
29	-0.07309	0.16130
30	-0.11107	-0.28503
31	-0.14336	0.28207
32	-0.16874	-0.36675
33	-0.18313	-0.10726
34	-0.18615	-0.33494
35	-0.17847	-0.29172
36	-0.16412	0.42033
37	-0.15004	0.27755
38	-0.13898	-0.11121
39	-0.13091	0.51381
40	-0.12689	-0.40600

### 4.3 Extension to the nlp Function

#### 4.3.1 New Techniques

**nondif** subgradient *bundle trust region* methods,

**simann** *Simulated Annealing* for global optimization (Goffe et.al, 1994).

#### 4.3.2 New Option

The **nlec** option was added as an alternative to the **nldb** option permitting a second form of nonlinear constraint specification: Assuming **nicon** is the name of a user specified function evaluating  $n_{nlc}$  nonlinear equality or inequality constraints  $c = (c_1, \dots, c_{n_{nlc}})^T$  for given variables (points)  $x = (x_1, \dots, x_n)^T$ . There are two different forms for specifying nonlinear constraints:

- The  $n_{nlc} = n_{nlec} + n_{nlic}$  nonlinear equality and inequality constraints can

always be specified with the function `nlcon` in the form:

$$\begin{aligned}c_j(x) &= 0 && \text{for } 1 \leq j \leq n_{nlec} \\c_j(x) &\geq 0 && \text{for } n_{nlec} < j \leq n_{nlc}\end{aligned}$$

If the function `nlcon` specifies the constraints in this form, the `nlec` option can be used for specifying the number  $n_{nlec}$  of constraints equal to zero. If there are no equality constraints,  $n_{nlec} = 0$ , the `nlec` option must not be specified since it defaults to zero.

- If for some reason you want to specify the constraints with the `nlcon` module in the form:

$$nllb_j \leq c_j(x) \leq nlub_j \quad \text{for } j = 1, \dots, n_{nlc}$$

then you can specify the lower  $nllb_j$  and upper  $nlub_j$  bounds using a two column matrix `nlcb`. This object must be specified using the `nlcb` option. The order of the  $n_{nlc}$  constraint functions specified in `nlcon` must correspond with the order of rows in the lower-upper-bound matrix specified with the `nlcb` option.

- If you specify both, the `nlec` and the `nlcb` options, the `nlcb` option will be used. However, using the `nlec` option and not `nlcb` is recommended.

The module `nlcon` must have the form `c = nlcon(x)`. During the execution of `nlp`, the argument `x` is set with trial values and the returned values of `y` are used to reduce the constraint violations when approximating the optimal location  $x^*$ . Note, that the definition of `nlcon` can have additional arguments `x2, ..., xm` defined with the `global` clause. No nonlinear constraints are imposed, if the `nlcon` function is not specified.



## 5 New Developments

### 5.1 Function arhetero

```
gof = arhetero(y,lagbnd<,optn>)
```

```
< gof,est,root,ase,nwcov,cov,resi > = arhet(y,lagbnd<,optn>)
```

**Purpose:** The `arhetero` function implements the modeling step for the Heterogeneous AutoRegressive method. There are two versions implemented (see MFE Toolbox):

**original** version using overlapping lag intervals,

**modified** version with non-overlapping lag intervals.

**Input:** `y` must be an  $N$  vector of time series data;

**lagbnd** can be a positive integer scalar or a  $p$  vector of positive upper bounds (assuming lower bounds of 1), or a  $p \times 2$  matrix with lower (in first column) and upper bounds (in second column);

**optn** The `optn` argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content.

Option	Second Column	Meaning
"alpha"	real	probability for confidence intervals, default=.05
"cent"		centers the $y$ data and as a consequence does not estimate the intercept
"meth"	string	method of estimation
	"ori"	original (overlapping) lag intervals (default)
	"mod"	modified (non-overlapping) lag intervals
"noint"		does not estimate the intercept
"print"	int	amount of printed output (def=0: no output)
"nopr"		do not print any output
"nwlag"	int	number of lags for Newey-West cov matrix (def=0: no lags)

**Output:** `gof` is a vector of scalar results currently containing:

1. indicator for problems in the algorithm (=0: no problems)
2. mean value  $\mu$  of input data  $y$
3. standard deviation  $\sigma$  of input data  $y$
4. variance  $\sigma^2$
5. computation time in seconds
6. error sum of squares
7. Akaike information criterion AIC

8. Bayesian (Schwartz) information criterion BIC

**est** is a vector of ML parameter estimates; the order of the estimates is:

1. intercept (optional)
2. nar parameters for the  $p$  lag intervals

**root** the magnitude (absolute values) of roots for the  $\max_j ar_j$  lags;

**ase** the asymptotic standard errors of the  $p$  estimates from the diagonal of the  $p \times p$  covariance matrix;

**nwcov** the  $p \times p$  (robust) Newey-West covariance matrix of the parameter estimates for a specified number of lags;

**cov** the  $p \times p$  asymptotic covariance matrix of the parameter estimates;

**resi** is an  $N$  vector of residuals.

**Restrictions:**

1. The time series data  $y$  cannot have any missing values or string data. If  $x$  is specified as nonmissing, then the same is true for the predictors  $x$ .
2. The lag numbers in `lagbnd` must be positive integers.

**Relationships:** `arma()`, `armafore()`, `arima()`, `burg()`, `tsmeas()`

**Examples:** 1. Lag Intervals with Vector Input, Original Version:

```
print "Robust NW Covariance Matrix without Lags";
plag = [ 1 5 22 ];
optn = [ "print"      2 ];
< gof,est,root,ase,nwcov,cov,resi > = arhetero(yy,plag,optn);
print "GOF=",gof;
print "PARM=",est;
print "Root=",root;
print "ASE=",ase;
print "NWCOV=",nwcov;
print "COV=",cov;
print "RESI[20:30]=",resi[20:30];
```

```
*****
Heterogeneous AR Modeling: Original Version (Uncentered)
*****
```

```
Number of Observations . . . . . 1000
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 20.7847
Standard Deviation Y . . . . . 1.1768
Mean Squared Error (MSE) . . . . . 1.3849
Number of Newey-West Lags . . . . . 0
```

Original (Overlapping) Lag Intervals

	Begin	End
AR_1	1	1
AR_2	1	5
AR_3	1	22

Standard Error Regression= 1.0037  
 BIC=0.00970143 AIC=-0.00992959

\*\*\*\*\*  
 AR Roots and Coefficients  
 \*\*\*\*\*

N	Coefficients	RealRoot	ImagRoot	Abs_Root
0	1.000000000			
1	-0.074546642	0.986396377	0.000000000	0.986396377
2	-0.086335537	-0.702542050	0.440720754	0.829337154
3	-0.086335537	0.852923558	-0.278322611	0.897185640
4	-0.086335537	-0.360801238	0.744403873	0.827233135
5	-0.086335537	0.319003199	-0.784422842	0.846807083
6	-0.027232902	0.527981456	0.664972164	0.849089158
7	-0.027232902	0.527981456	-0.664972164	0.849089158
8	-0.027232902	0.319003199	0.784422842	0.846807083
9	-0.027232902	0.709173134	-0.491738826	0.862979494
10	-0.027232902	0.094137249	-0.837687446	0.842960307
11	-0.027232902	-0.360801238	-0.744403873	0.827233135
12	-0.027232902	-0.137630194	-0.821719040	0.833165200
13	-0.027232902	-0.137630194	0.821719040	0.833165200
14	-0.027232902	0.709173134	0.491738826	0.862979494
15	-0.027232902	-0.791989818	-0.230843482	0.824946413
16	-0.027232902	-0.555251096	-0.614614010	0.828283865
17	-0.027232902	-0.791989818	0.230843482	0.824946413
18	-0.027232902	-0.702542050	-0.440720754	0.829337154
19	-0.027232902	0.852923558	0.278322611	0.897185640
20	-0.027232902	0.094137249	0.837687446	0.842960307
21	-0.027232902	-0.555251096	0.614614010	0.828283865
22	-0.027232902	-0.821858136	0.000000000	0.821858136

Parameter Estimates  
 \*\*\*\*\*

Dense Row Vector (ncol=4)

```

R | Intercept      AR_1      AR_2      AR_3
   2.4289563 -0.0117889  0.2955132  0.5991238

```

Robust (Newey-West) Asymptotic Standard Errors  
 \*\*\*\*\*

Dense Matrix (4 by 3)

	A_St_Err.	T_Value	P_Value
Intercept	0.9738120	2.4942763	0.0063913
AR_1	0.0360480	-0.3270335	0.6281444
AR_2	0.0820911	3.5998212	1.67e-004
AR_3	0.0859935	6.9670817	2.94e-012

Newey-West Covariance Matrix of Estimates (0 Lags)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept	0.9483098			
AR_1	-0.0019482	0.0012995		
AR_2	9.91e-005	-0.0013292	0.0067389	
AR_3	-0.0438759	1.24e-004	-0.0054079	0.0073949

Covariance Matrix of Parameter Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept	0.9181031			
AR_1	-2.56e-004	0.0013961		
AR_2	-0.0019961	-0.0014799	0.0074492	
AR_3	-0.0418698	9.62e-005	-0.0058720	0.0077889

## 2. Lag Intervals with Matrix Specification, Original Version:

```

print "Matrix Specification of LAG Intervals: > 1 Start";
plag = [ 1 1, 2 2, 10 22 ];

```

```

optn = [ "print"      2 ];
< gof,est,root,ase,nwcov,cov,resi > = arhetero(yy,plag,optn);
print "GOF=",gof;
print "PARM=",est;
print "Root=",root;
print "ASE=",ase;
print "NWCOV=",nwcov;
print "COV=",cov;
print "RESI[20:30]=",resi[20:30];

```

```

*****
Heterogeneous AR Modeling: Original Version (Uncentered)
*****

```

```

Number of Observations . . . . . 1000
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 20.7847
Standard Deviation Y . . . . . 1.1768
Mean Squared Error (MSE) . . . . . 1.3849
Number of Newey-West Lags . . . . . 0

```

Original (Overlapping) Lag Intervals

	Begin	End
AR_1	1	1
AR_2	2	2
AR_3	10	22

Standard Error Regression= 1.02833  
BIC=0.0581787 AIC=0.0385477

Parameter Estimates  
\*\*\*\*\*

Dense Row Vector (ncol=4)

R	Intercept	AR_1	AR_2	AR_3
	3.8495117	0.1228098	0.0785620	0.6130674

Robust (Newey-West) Asymptotic Standard Errors  
\*\*\*\*\*

Dense Matrix (4 by 3)

	A_St_Err.	T_Value	P_Value
Intercept	0.9693597	3.9711902	3.83e-005
AR_1	0.0309529	3.9676305	3.89e-005
AR_2	0.0319426	2.4594742	0.0070416
AR_3	0.0562102	10.906701	0.0000000

Newey-West Covariance Matrix of Estimates (0 Lags)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept	0.9396582			
AR_1	-0.0046478	9.58e-004		
AR_2	-0.0036824	-9.88e-005	0.0010203	
AR_3	-0.0369353	-6.33e-004	-7.46e-004	0.0031596

### 3. Newey West Covariance Matrix with Lags:

```
print "With Newey-West Covariance Matrix: Default Lags";
plag = [ 1 5 22 ];
tt = (real)ncol(yy); tt = ceil(pow(tt,1. / 3.));
optn = [ "print"      2 ,
        "nwlag"      tt ];
< gof,est,root,ase,nwcov,cov,resi > = arhetero(yy,plag,optn);
print "GOF=",gof;
print "PARM=",est;
print "Root=",root;
print "ASE=",ase;
print "NWCOV=",nwcov;
print "COV=",cov;
print "RESI[20:30]=",resi[20:30];
```

\*\*\*\*\*  
Heterogeneous AR Modeling: Original Version (Uncentered)  
\*\*\*\*\*

Number of Observations . . . . .	1000
Number of Parameter Estimates . . . . .	4
Mean Y . . . . .	20.7847
Standard Deviation Y . . . . .	1.1768

Mean Squared Error (MSE) . . . . . 1.3849  
 Number of Newey-West Lags . . . . . 10

Original (Overlapping) Lag Intervals

	Begin	End
AR_1	1	1
AR_2	1	5
AR_3	1	22

Standard Error Regression= 1.0037  
 BIC=0.00970143 AIC=-0.00992959

Parameter Estimates  
 \*\*\*\*\*

Dense Row Vector (ncol=4)

R	Intercept	AR_1	AR_2	AR_3
	2.4289563	-0.0117889	0.2955132	0.5991238

Robust (Newey-West) Asymptotic Standard Errors  
 \*\*\*\*\*

Dense Matrix (4 by 3)

		A_St_Err.	T_Value	P_Value
Intercept		0.9452535	2.5696347	0.0051625
AR_1		0.0319191	-0.3693364	0.6440222
AR_2		0.0807807	3.6582162	1.34e-004
AR_3		0.0898222	6.6701070	2.12e-011

Newey-West Covariance Matrix of Estimates (10 Lags)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept		0.8935043		
AR_1		-0.0032674	0.0010188	
AR_2		0.0064259	-9.20e-004	0.0065255

```
AR_3 | -0.0461011  5.98e-005 -0.0059129  0.0080680
```

Covariance Matrix of Parameter Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept	0.9181031			
AR_1	-2.56e-004	0.0013961		
AR_2	-0.0019961	-0.0014799	0.0074492	
AR_3	-0.0418698	9.62e-005	-0.0058720	0.0077889

4. Modified version: Non-overlapping Lag Intervals

```
print "Use Modified Version";
plag = [ 1 1, 1 5, 1 22 ];
optn = [ "print"    2 ,
         "meth"    "mod" ];
< gof,est,root,ase,nwcov,cov,resi > = arhetero(yy,plag,optn);
print "GOF=",gof;
print "PARM=",est;
print "Root=",root;
print "ASE=",ase;
print "NWCOV=",nwcov;
print "COV=",cov;
print "RESI[20:30]=",resi[20:30];
```

\*\*\*\*\*  
Heterogeneous AR Modeling: Modified Version (Uncentered)  
\*\*\*\*\*

Number of Observations . . . . .	1000
Number of Parameter Estimates . . . . .	4
Mean Y . . . . .	20.7847
Standard Deviation Y . . . . .	1.1768
Mean Squared Error (MSE) . . . . .	1.3849
Number of Newey-West Lags . . . . .	0

Original (Overlapping) Lag Intervals

	Begin	End
AR_1	1	1



AR_2	1	5
AR_3	1	22

Modified (Non-Overlapping) Lag Intervals

	Begin	End
AR_1	1	1
AR_2	2	5
AR_3	6	22

Standard Error Regression= 1.0037  
 BIC=0.00970143 AIC=-0.00992959

Parameter Estimates  
 \*\*\*\*\*

Dense Row Vector (ncol=4)

R	Intercept	AR_1	AR_2	AR_3
	2.4289563	0.0745466	0.3453421	0.4629593

Robust (Newey-West) Asymptotic Standard Errors  
 \*\*\*\*\*

Dense Matrix (4 by 3)

		A_St_Err.	T_Value	P_Value
Intercept		0.9738120	2.4942763	0.0063913
AR_1		0.0310735	2.3990420	0.0083105
AR_2		0.0546277	6.3217440	1.95e-010
AR_3		0.0664495	6.9670817	2.94e-012

Newey-West Covariance Matrix of Estimates (0 Lags)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept		0.9483098		
AR_1		-0.0039228	9.66e-004	
AR_2		-0.0078981	-2.95e-004	0.0029842

```
AR_3 | -0.0339041 -4.80e-004 -0.0023041 0.0044155
```

Covariance Matrix of Parameter Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	AR_1	AR_2	AR_3
Intercept	0.9181031			
AR_1	-0.0025589	0.0010202		
AR_2	-0.0092096	-3.37e-004	0.0033167	
AR_3	-0.0323540	-5.60e-004	-0.0025356	0.0046508

## 5.2 Function arma

```
gof = arma(y<,x<,ar<,ma<,optn<,p0> . >)
```

```
< gof,est,root,ase,cis,cov,sco > = arma(y<,x<,ar<,ma<,optn <,p0> . >)
```

**Purpose:** The `arma` function implements the modeling step for the AutoRegressive Moving-Average method. It computes ML estimates and the (robust) asymptotic covariance matrix of estimates. In a second step, the optimal estimates can be used for computing predicted values using the `armafore` function.

**Input:** `y` must be an  $N$  vector of time series data;

`x` [optional] could be a  $N \times n$  matrix of predictors; can be specified as missing;

`ar` can be a positive integer scalar or vector of positive integers with the *nar* AR lag numbers, default is 1;

`ma` can be a positive integer scalar or vector of positive integers with the *nma* MA lag numbers, default is 1;

`optn` The `optn` argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content.

`p0` specifies an optional  $p$  vector of starting values for the ML estimation. In addition to many of the optimization options, like termination criteria etc., the following options may be specified:

Option	Second Column	Meaning
"alpha"	real	probability for confidence intervals, default=.05
"cent"		centers the $y$ data and as a consequence does not estimate the intercept
"lstech"	string	optimization technique for least squares minimization, default is Levenberg-Marquardt;
"mltech"	string	optimization technique for maximum likelihood optimization, default is Newton Raphson;
"nacor"	int	specifies number of (partial) autocorrelations
"noint"		does not estimate the intercept
"print"	int	amount of printed output (def=0: no output)
"nopr"		do not print any output

**Output:** `gof` is a vector of scalar results currently containing:

1. indicator for problems in the algorithm (=0: no problems)
2. mean value  $\mu$  of input data  $y$
3. standard deviation  $\sigma$  of input data  $y$
4. variance  $\sigma^2$
5. computation time in seconds
6. Loglikelihood
7. number of iterations for optimization
8.  $L1$  norm of gradient at the optimal solution
9. error sum of squares
10. Akaike information criterion AIC
11. Bayesian (Schwartz) information criterion BIC

`est` is a vector of ML parameter estimates; the order of the estimates is:

1. intercept (optional)
2. `nar` parameters for the AR lags,
3. `nma` parameters for the MA lags.

`root` the magnitude (absolute values) of roots for the  $\max_j ar_j$  and  $\max_k ma_k$  lags;

`ase` the asymptotic standard errors of the  $p$  estimates from the diagonal of the  $p \times p$  covariance matrix;

`cis` confidence intervals based on the asymptotic standard errors of the  $p$  estimates

`cov` the  $p \times p$  asymptotic covariance matrix of the ML parameter estimates

`sco` is an  $N \times p$  matrix of scores.

**Restrictions:** 1. The time series data  $y$  cannot have any missing values or string data. If  $x$  is specified as nonmissing, then the same is true for the predictors  $x$ .

2. The lag numbers in `ar` and `ma` must be positive integers.

**Relationships:** arhetero(), armafore(), arima(), burg(), tsmeas()

**Examples:** The following three examples refer to the  $N = 40$  data of randomly generated values:

```
xd2 = [ 0.8147 0.9058 0.1270 0.9134 0.6324
        0.0975 0.2785 0.5469 0.9575 0.9649
        0.1576 0.9706 0.9572 0.4854 0.8003
        0.1419 0.4218 0.9157 0.7922 0.9595
        0.6557 0.0357 0.8491 0.9340 0.6787
        0.7577 0.7431 0.3922 0.6555 0.1712
        0.7060 0.0318 0.2769 0.0462 0.0971
        0.8235 0.6948 0.3171 0.9502 0.0344 ];
```

1. Use simple lags ar=1 and ma=1:

The algorithm is running two optimizations, the first is a least squares optimization computing excellent starting values for the ML optimization. That's why we able to specify two optimization methods:

```
print "Simple ARMA: LSTech=LEVMar, MLTech=TRUREG";
ar = 1; ma = 1;
optn = [ "lstech"  "lm" ,
         "mltech"  "tr" ,
         "print"   3 ];
< gof,parm,root,ase,cis,cov,scor > = arma(xd2,.,ar,ma,optn);
print "GOF=",gof;
print "PARM=",parm;
print "Root=",root;
print "ASE=",ase;
print "CIS=",cis;
print "COV=",cov;
print "SCOR=",scor;
```

```
*****
ARMA Modeling: N(AR)=1 N(MA)=1 (Uncentered)
*****
AR[1] = 1 MA[1] = 1
```

```
Number of Observations . . . . . 40
Number of Predictor Variables . . . . . 0
Number of Parameter Estimates . . . . . 3
Mean . . . . . 0.5673
Standard Deviation . . . . . 0.3323
Mean Squared Error (MSE) . . . . . 0.1104
LS Optimization Technique . . . . . LEVMAR
```

When specifying the "print" 3 option the optimization histories are printed:

\*\*\*\*\*  
 Optimization Start  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	0.34996599	0.1010283
2 AR_1	0.37292898	-0.0057028
3 MA_1	-0.41847023	-0.0743579

Value of Objective Function = 2.15852

Levenberg-Marquardt Optimization  
 Scaling Update of More (1978)  
 Gradient Computed by Finite Differences  
 CRP Jacobian Computed by Finite Differences

Iteration Start:

N. Variables	3	N. Equations	40
Criterion	2.158519520	Max Grad Entry	0.101028317
TR Radius	1.081565870		

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	rho
1	0	0	0	2.158312	2.1e-004	0.02017	0.00000	0.21832
2	0	0	0	2.157933	3.8e-004	0.01434	2e-003	0.61247
3	0	0	0	2.157910	2.2e-005	4e-004	0.00000	0.99724
4	0	0	0	2.157910	5.4e-008	7e-004	5e-004	0.10030
5	0	0	0	2.157910	1.7e-007	3e-004	2e-003	0.56175
6	0	0	0	2.157910	5.5e-009	8e-005	0.00000	0.40304
7	0	0	0	2.157910	4.7e-009	5e-005	2e-003	0.57227
8	0	0	0	2.157910	2.1e-010	1e-005	0.00000	0.63508

Successful Termination After 8 Iterations  
 GCONV convergence criterion satisfied.  
 Criterion 2.157910208 Max Grad Entry 1.0103e-005

Ridge (lambda)	0.00000000	TR Radius	0.000373462
Act.dF/Pred.dF	0.635080549		
N. Function Calls	0	N. Gradient Calls	3
N. Hessian Calls	10	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates

Parameter	Estimate	Gradient
1 Intercept	0.36010117	7.25e-007
2 AR_1	0.35313547	-3.04e-006
3 MA_1	-0.38641826	-1.01e-005

Value of Objective Function = 2.15791

The negative LogLikelihood is minimized:

Trust Region Optimization  
 Scaling Update of More (1978)  
 Gradient Computed by Finite Differences  
 Hessian Computed by Finite Differences (dense)  
 (Using Only Function Calls)

Iteration Start:

N. Variables	3		
Criterion	12.22570100	Max Grad Entry	9.3640e-005
TR Radius	1.000000000		

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1	0	2	0	12.22570	7.0e-010	8e-009	0.00000	1.00000

Successful Termination After 1 Iterations

GCONV convergence criterion satisfied.

Criterion	12.22570100	Max Grad Entry	7.6173e-009
Ridge (lambda)	0.000000000	TR Radius	1.000000000
Act.dF/Pred.dF	1.000119405		
N. Function Calls	3	N. Gradient Calls	7

N. Hessian Calls	3	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	0.36011365	-4.74e-009
2 AR_1	0.35311336	2.17e-009
3 MA_1	-0.38639573	-7.62e-009

Value of Objective Function = 12.2257

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3
1	933.34733	537.45465	10.248984
2	537.45465	354.55737	51.546290
3	10.248984	51.546290	49.056449

Log Likelihood= -12.2257  
 SBC=-1.89313 AIC=-2.0198

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (3 by 3)

S	Intercept	AR_1	MA_1
Intercept	0.1288180		
AR_1	-0.2258580	0.3993282	
MA_1	0.2104084	-0.3724091	0.3677363

Factor sigma=1 (nobs=40 df=3 vard=37)  
 Determinant = 8.74154e-006  
 Matrix has Only Positive Eigenvalues

Approximate Correlation Matrix of Parameter Estimates  
 \*\*\*\*\*

Dense Symmetric Matrix (3 by 3)

S	Intercept	AR_1	MA_1
Intercept	1.0000000		
AR_1	-0.9958231	1.0000000	
MA_1	0.9667328	-0.9718229	1.0000000

Determinant = 0.00046211  
 Matrix has Only Positive Eigenvalues

\*\*\*\*\*  
 Parameter Estimates and Asympt. Standard Errors  
 \*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.3601136	0.3589123	1.0033473	0.161105
AR_1	0.3531134	0.6319242	0.5587907	0.289836
MA_1	-0.3863957	0.6064126	-0.6371828	0.736035

\*\*\*\*\*  
 Wald Confidence Limits (Alpha=0.05)  
 \*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.3601136	-0.3433415	1.0635688
AR_1	0.3531134	-0.8854352	1.5916620
MA_1	-0.3863957	-1.5749427	0.8021512

Factor sigma=1 (nobs=40 df=3 vard=37)

GOF=

	g-of-fit
Error	0.00000
Y_Mean	0.56729
Y_Stdv	0.33231



MSE		0.11043
Time		0.00000
LogLik		12.226
N_Iter		9.0000
MaxGrad		8e-009
SEReg		0.35137
AIC		-2.0198
SBIC		-1.8931
unused		.

PARM=

		Estimates	Gradient
Intercept		0.36011	-5e-009
AR_1		0.35311	2e-009
MA_1		-0.38640	-8e-009

Root=

		1
AR_Root		0.35311
MA_Root		0.38642

ASE=

		A_St_Err.	T_Value	P_Value
Intercept		0.35891	1.0033	0.16111
AR_1		0.63192	0.55879	0.28984
MA_1		0.60641	-0.63718	0.73604

CI=

		Wald_LCI	Wald_UCI
Intercept		-0.34334	1.0636
AR_1		-0.88544	1.5917
MA_1		-1.5749	0.80215

COV=

	SYM		Intercept	AR_1	MA_1
Intercept		0.12882			
AR_1		-0.22586	0.39933		
MA_1		0.21041	-0.37241	0.36774	

SCOR=

	Intercept	AR_1	MA_1
1	1e-010	-4e-011	2e-010
2	2.3912	1.9481	9e-011
3	-5.8243	-5.1278	-1.0839
4	4.7439	1.8492	-1.0922
.....			
37	2.0787	1.1384	-0.06555
38	-3.5523	-2.2662	-0.25670
39	5.8491	2.5799	-0.68075
40	-7.7277	-5.8226	-1.4887

2. Use lags ar=1,2,3 and ma=1:

```
print "Simple ARMA: Tech=LEVMAR";
ar = [ 1:3 ]; ma = 1;
optn = [ "lstech"  "lm"  ,
         "mltech"  "tr"  ,
         "print"   3  ];
< gof,parm,root,ase,cis,cov > = arma(xd2,.,ar,ma,optn);
print "GOF=",gof;
print "PARM=",parm;
print "Root=",root;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;
```

Here we only report the final result:

Trust Region Optimization  
Scaling Update of More (1978)  
Gradient Computed by Finite Differences  
Hessian Computed by Finite Differences (dense)  
(Using Only Function Calls)

```
Iteration Start:
N. Variables      5
Criterion         9.861684378      Max Grad Entry  0.001189907
TR Radius         1.000000000
```

```

Iter rest nfun act  optcrit difcrit maxgrad lambda  radius
  1   0   2   0  9.861684 4.2e-008 3e-007 0.00000 1.00000

```

Successful Termination After 1 Iterations

GCONV convergence criterion satisfied.

```

Criterion          9.861684336          Max Grad Entry  2.8946e-007
Ridge (lambda)    0.000000000          TR Radius        1.000000000
Act.dF/Pred.dF    0.999854944
N. Function Calls      3          N. Gradient Calls      7
N. Hessian Calls      3          Preproces. Time        0
Time for Method       0          Effective Time         0

```

```

*****
Optimization Results
*****

```

Parameter Estimates

-----

Parameter	Estimate	Gradient
1 Intercept	0.72978139	5.02e-008
2 AR_1	-0.43982250	5.09e-008
3 AR_2	-0.03502382	-3.68e-008
4 AR_3	0.16555089	1.91e-007
5 MA_1	0.48770070	2.89e-007

Value of Objective Function = 9.86168

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4	5
1	178.94289	99.914095	103.26008	103.07296	4.0279008
2	99.914095	103.41070	37.594099	59.708318	44.177141
3	103.26008	37.594099	108.73105	41.247163	-13.278503
4	103.07296	59.708318	41.247163	106.80399	6.6848378
5	4.0279008	44.177141	-13.278503	6.6848378	55.149279

Log Likelihood= -9.86168

SBC=-1.74689 AIC=-1.958

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (5 by 5)

S	Intercept	AR_1	AR_2	AR_3	MA_1
Intercept	0.0677439				
AR_1	-0.0532225	0.0727391			
AR_2	-0.0323532	0.0162133	0.0295282		
AR_3	-0.0251912	0.0076546	0.0110593	0.0254186	
MA_1	0.0329496	-0.0514044	-0.0048556	-0.0047101	0.0563052

Factor sigma=1 (nobs=40 df=5 vard=35)  
 Determinant = 4.0632e-009  
 Matrix has Only Positive Eigenvalues

Approximate Correlation Matrix of Parameter Estimates  
 \*\*\*\*\*

Dense Symmetric Matrix (5 by 5)

S	Intercept	AR_1	AR_2	AR_3	MA_1
Intercept	1.0000000				
AR_1	-0.7581866	1.0000000			
AR_2	-0.7233759	0.3498402	1.0000000		
AR_3	-0.6070690	0.1780181	0.4036760	1.0000000	
MA_1	0.5335085	-0.8032331	-0.1190829	-0.1245039	1.0000000

Determinant = 0.0195116  
 Matrix has Only Positive Eigenvalues

\*\*\*\*\*  
 Parameter Estimates and Asympt. Standard Errors  
 \*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.7297814	0.2602766	2.8038686	0.004091
AR_1	-0.4398225	0.2697019	-1.6307726	0.944047
AR_2	-0.0350238	0.1718377	-0.2038192	0.580162
AR_3	0.1655509	0.1594322	1.0383780	0.153107
MA_1	0.4877007	0.2372872	2.0553185	0.023685

\*\*\*\*\*  
Wald Confidence Limits (Alpha=0.05)  
\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.7297814	0.2196486	1.2399142
AR_1	-0.4398225	-0.9684286	0.0887836
AR_2	-0.0350238	-0.3718195	0.3017719
AR_3	0.1655509	-0.1469305	0.4780323
MA_1	0.4877007	0.0226264	0.9527750

Factor sigma=1 (nobs=40 df=5 vard=35)

GOF=

	g-of-fit
Error	0.00000
Y_Mean	0.56729
Y_Stdv	0.33231
MSE	0.11043
Time	1.00000
LogLik	9.8617
N_Iter	12.000
MaxGrad	3e-007
SEReg	0.35443
AIC	-1.9580
SBIC	-1.7469
unused	.

PARM=

	Estimates	Gradient
Intercept	0.72978	5e-008
AR_1	-0.43982	5e-008
AR_2	-0.03502	-4e-008
AR_3	0.16555	2e-007
MA_1	0.48770	3e-007

Root=

	1	2	3
AR_Root	0.79557	0.79557	0.69489
MA_Root	0.48763	.	.

```

SCOR=
| Intercept AR_1 AR_2 AR_3 MA_1
-----
1 | -1e-009 -1e-009 9e-010 -5e-009 -7e-009
2 | -1e-009 -1e-009 9e-010 -5e-009 -7e-009
3 | -1e-009 -1e-009 9e-010 -5e-009 -7e-009
4 | 1.4220 0.18060 1.2881 1.1585 -6e-009
5 | 0.49355 0.82030 -0.30324 0.48986 0.13134
6 | -3.0377 -0.87932 -4.3205 0.48992 -0.10477
.....
36 | 1.7194 0.45974 -0.43340 1.1310 0.10133
37 | 1.4259 1.5610 0.38128 -0.35944 0.47922
38 | -1.3568 -0.67808 -1.4853 -0.36280 -0.18813
39 | 2.3998 0.54720 1.1993 2.6271 -0.85315
40 | -3.8445 -5.0071 -0.87660 -1.9213 -2.6241

```

3. Use lags ar=1 and ma=1,2,3:

```

ar = 1; ma = [ 1:3 ];
optn = [ "lstech" "lm" ,
         "mltech" "tr" ,
         "print" 3 ];
< gof,parm,root,ase,cis,cov > = arma(xd2,.,ar,ma,optn);
print "GOF=",gof;
print "PARM=",parm;
print "Root=",root;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

GOF=
| g-of-fit
-----
Error | 0.00000
Y_Mean | 0.56729
Y_Stdv | 0.33231
MSE | 0.11043
Time | 1.00000
LogLik | 11.734
N_Iter | 16.000
MaxGrad | 7e-008
SEReg | 0.34594
AIC | -2.0065
SBIC | -1.7954
unused | .

```

PARM=

	Estimates	Gradient
Intercept	0.32928	-2e-008
AR_1	0.40673	-6e-009
MA_1	-0.43325	1e-008
MA_2	-0.04240	-4e-008
MA_3	0.08842	7e-008

Root=

	1	2	3
AR_Root	0.40673	.	.
MA_Root	0.74631	0.76196	0.76196

ASE=

	A_St_Err.	T_Value	P_Value
Intercept	0.32817	1.0034	0.16128
AR_1	0.57655	0.70545	0.24260
MA_1	0.59611	-0.72679	0.76391
MA_2	0.24669	-0.17187	0.56774
MA_3	0.23102	0.38274	0.35211

CI=

	Wald_LCI	Wald_UCI
Intercept	-0.31392	0.97247
AR_1	-0.72329	1.5367
MA_1	-1.6016	0.73510
MA_2	-0.52590	0.44110
MA_3	-0.36438	0.54122

COV=

	Intercept	AR_1	MA_1
Intercept	0.10769		
AR_1	-0.18832	0.33241	
MA_1	0.18658	-0.32970	0.35535
MA_2	-0.00489	0.00882	-0.02982
MA_3	-0.01137	0.01971	-0.00963

	SYM	MA_2	MA_3		
	MA_2	0.06086			
	MA_3	-0.04140	0.05337		
SCOR=					
	Intercept	AR_1	MA_1	MA_2	MA_3
1	4e-010	1e-010	-3e-010	1e-009	-2e-009
2	2.3948	1.9511	-1e-010	4e-010	-7e-010
3	-6.5027	-5.7111	-1.1123	-1e-009	2e-009
4	5.5508	2.3590	-1.1955	0.81815	2e-010
5	0.63274	0.44886	0.07356	-0.13389	0.09163
6	-6.8298	-4.3627	-0.35289	-0.80142	1.4588
37	2.7454	1.4474	-0.09572	-1.2659	-1.2981
38	-2.7003	-1.7532	-0.30430	0.09414	1.2451
39	6.4223	3.1440	-0.10074	0.72373	-0.22391
40	-8.4361	-6.5401	-2.0917	0.13233	-0.95065

### 5.3 Function armafore

```

gof = armafore(np,nh,est,y<,x<,ar<,ma<,optn> . >)
< gof,yhat > = armafore(np,nh,est,y<,x<,ar<,ma<,optn> . >)

```

**Purpose:** The `armafore` function implements the forecasting step of the AutoRegressive Moving-Average method.

**Input:** `np` must be a positive integer with  $1 \leq np < N$  specifying the number of last observations of  $y$  should be forecast; that means, the first  $N - np$  observations are used for regression

`nh` must be a positive integer with  $1 \leq nh < N$  specifying the forecast "horizon";

`est` must be an  $p$  vector with the optimal ML parameters estimated from calling the `arma` function.

`y` must be an  $N$  vector of time series data.

`x` [optional] could be a  $N \times n$  matrix of predictors

`ar` can be an integer scalar or vector of integers with the  $nar$  AR lag numbers, default is 1;

`ma` can be an integer scalar or vector of integers with the  $nma$  MA lag numbers, default is 1;



**optn** The **optn** argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content.

Option	Second Column	Meaning
"center"		centers the $y$ data and as a consequence does not use an intercept parameter
"noint"		does not use an intercept parameter
"print"	int	amount of printed output (def=0: no output)
"nopr"		do not print any output
"sereg"	real	should be the standard error output from the <b>gof</b> result of <b>arma</b> .

**Output:** **gof** is a vector of scalar results currently containing:

1. indicator for problems in the algorithm (=0: no problems)
2. mean value  $\mu$  of input data  $y$
3. standard deviation  $\sigma$  of input data  $y$
4. variance  $\sigma^2$
5. computation time in seconds
6. theoretical standard deviation of the h-step ahead forecast (assuming homoskedasticity)

**yhat** is a  $N \times 3$  matrix containing in its columns:

1. the predicted  $\hat{y}$  values;
2. the shifted input data  $\bar{y}$ ;
3. the residuals  $\hat{y} - \bar{y}$ .

**Restrictions:** 1. The time series data  $y$  cannot have any missing values or string data. If  $x$  is specified as nonmissing, then the same is true for the predictors  $x$ .

2. There are methodological problems with ARMA forecasting of data including a regressor matrix  $\mathbf{X}$ .
3. The **ar** and **ma** specifications must agree with those used in the **arma** modeling step.
4. Some of the options should be compatible with those used in the **arma** modeling step, like "center" or "noint"!

**Relationships:** **arhetero()**, **arma()**, **arima()**, **burg()**, **garch()**, **tsmeas()**

**Examples:** Here we use the results from the three examples documented for the **arma** function:

1. Example 1 of **arma**: Use simple lags **ar**=1 and **ma**=1:

```

est = parm[,1]; sereg = gof[9];
np = 10; nh = 2;
optn = [ "print"      3 ,
        "sereg"    sereg ];
< gof,yhat > = armafore(np,nh,est,xd2,.,ar,ma,optn);
print "GOF=",gof;
print "YHAT=",yhat;

```

```

*****
ARMA Forecasting: N(AR)=1 N(MA)=1 (Uncentered)
*****
          AR[1] =      1 MA[1] =      1

```

```

Number of Observations . . . . . 40
Number of Predictor Variables . . . . . 0
Number of Parameter Estimates . . . . . 3
Mean . . . . . 0.5673
Standard Deviation . . . . . 0.3323
Mean Squared Error (MSE) . . . . . 0.1104
Standard Error Regression . . . . . 0.3514

```

GOF=

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	g-of-fit
Error	0.00000
Y_Mean	0.56729
Y_Stdv	0.33231
MSE	0.11043
Time	0.00000
YSTD	0.35156
unused	.

YHAT=

	YHAT	YTPH	FERR
1	.	.	.
2	.	.	.
3	.	.	.
.....			
27	.	.	.
28	.	.	.
29	.	.	.
30	0.56085	0.03180	-0.52905

```

31 | 0.55654 0.27690 -0.27964
32 | 0.56280 0.04620 -0.51660
33 | 0.56234 0.09710 -0.46524
34 | 0.56487 0.82350 0.25863
35 | 0.56525 0.69480 0.12955
36 | 0.55686 0.31710 -0.23976
37 | 0.55513 0.95020 0.39507
38 | 0.55890 0.03440 -0.52450
39 | 0.55292 . .
40 | 0.56137 . .

```

2. Example 2 of arma: Use lags ar=1,2,3 and ma=1:

```

est = parm[,1]; sereg = gof[9];
np = 10; nh = 2;
optn = [ "print"      3 ,
        "sereg"     sereg ];
< gof,yhat > = armafore(np,nh,est,xd2,.,ar,ma,optn);
print "GOF=",gof;
print "YHAT=",yhat;

```

```

GOF=
      | g-of-fit
-----
Error | 0.00000
Y_Mean | 0.56729
Y_Stdv | 0.33231
  MSE | 0.11043
  Time | 0.00000
  YSTD | 0.35484
unused | .
unused | .
unused | .
unused | .

```

```

YHAT=
      |      YHAT      YTPH      FERR
-----
  1 | .          .          .
  2 | .          .          .
  3 | .          .          .
.....
 27 | .          .          .
 28 | .          .          .

```

```

29 | . . .
30 | 0.61619 0.03180 -0.58439
31 | 0.45786 0.27690 -0.18096
32 | 0.65669 0.04620 -0.61049
33 | 0.43973 0.09710 -0.34263
34 | 0.59330 0.82350 0.23020
35 | 0.47969 0.69480 0.21511
36 | 0.50087 0.31710 -0.18377
37 | 0.61720 0.95020 0.33300
38 | 0.56632 0.03440 -0.53192
39 | 0.49271 . .
40 | 0.68072 . .

```

3. Example 3 of arma: Use lags ar=1 and ma=1,2,3:

```

est = parm[,1]; sereg = gof[9];
np = 10; nh = 2;
optn = [ "print"      3 ,
        "sereg"     sereg ];
< gof,yhat > = armafore(np,nh,est,xd2,.,ar,ma,optn);
print "GOF=",gof;
print "YHAT=",yhat;

```

```

GOF=
      | g-of-fit
-----|-----
Error | 0.00000
Y_Mean | 0.56729
Y_Stdv | 0.33231
  MSE | 0.11043
  Time | 0.00000
  YSTD | 0.34607
unused | .
unused | .
unused | .
unused | .

```

```

YHAT=
      |      YHAT      YTPH      FERR
-----|-----
1 | . . .
2 | . . .
3 | . . .
.....

```

27		.	.	.
28		.	.	.
29		.	.	.
30		0.58534	0.03180	-0.55354
31		0.54813	0.27690	-0.27123
32		0.58139	0.04620	-0.53519
33		0.52150	0.09710	-0.42440
34		0.58340	0.82350	0.24010
35		0.53209	0.69480	0.16271
36		0.55172	0.31710	-0.23462
37		0.52947	0.95020	0.42073
38		0.49106	0.03440	-0.45666
39		0.54282	.	.
40		0.57429	.	.

## 5.4 Function affrma

---

```
gof = affrma(data,optn<,ref>)
```

```
< gof,dnew > = affrma(data,optn<,ref>)
```

**Purpose:** The `affrma` function implements the Bioconductor RMA algorithm by Bolstad et.al (2003) for the normalization of microarray data. This algorithm is an alternative to the `affvsn` and `affarms` methods for obtaining column normalized microarray data. However, other than fitting the parameters of a Maximum Likelihood model, the RMA method applies robust Medianpolish (Tukey, 1977a, p. 179), see also the `mpolish` function in CMAT to the data.

**Input: data** this should be a  $N \times n$  matrix of microarray data where the rows correspond to features (genes) and the columns to samples which need to be normalized. The data may obtain a column for an ID variable specifying a strata. The column number of the ID variable must be specified by the `optn` argument.

**optn** The `optn` argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content.

**ref** not yet implemented.

Option	Second Column	Meaning
"bgcor"		perform background correction default is background correction
"idvar"	int	column number containing ID variable; default is no ID variable
"medp"	int	algorithm for mean or median polish default is =0: median polish, =1: mean polish
"mpit"	int	maximum number of median polish iterations; default is 10 for median and 1 for mean polish
"mpeps"	real	for terminating median polish; default is 1.e-2
"nobg"		do not perform background correction default is background correction
"nsamp"		size $n_r \leq N$ of a reduced data set, needed only when $N$ is too large for the computer resources. default is $n_r = 0$
"pdat"	int	print input and fitted data sets
"print"	int	amount of printed output =0 specifies no printed output, default is 1

**Output:** `gof` a vector of scalar results, see below for content;

`dnew` this is the normalized  $N \times n$  data set.

**Restrictions:** 1. The input data must be numeric and cannot contain any string or complex data.

2. Currently the input data must not have any missing values.

**Relationships:** `affarms()`, `affvsn()`

**Examples:** 1. :

## 5.5 Function berkow

---

```
prob = berkow(y<,optn<,"dist"<,>>)
```

```
< prob,stat > = berkow(y<,optn<,"dist"<,>>)
```

**Purpose:** Implements the Berkowitz test for a large variety of data distributions, including uniform and normal.

**Input: y** The first input object  $y$  must be either a  $n$  vector or a  $n \times nc$  matrix. If  $y$  is a vector the scalar value of the probability of the CDF is returned, if  $y$  is a matrix then a vector of  $nc$  probabilities is returned by using each column of  $y$  separately.

**optn** this is a vector of options which should be initialized with missing values for defaults:

1. (int) amount of printed output (default is no output);
2. (real) probability for test (default is 0.05), must be in  $[0, 1]$ ;
3. (int) version number of algorithm:  
0 for time series data (is default),  
1 for cross sectional data.

**"dist"** The following string options are valid:

**("bern",x,prob)** Bernoulli distribution

- $x$  is the upper limit for cumulation  $0 \leq x \leq 1$ .
- $p$  is the probability for success of 1,  $0 \leq prob \leq 1$ .

**("beta",x,a,b,<l>,<u>>)** Beta distribution

- $x$  is the upper limit for integration  $0 \leq x \leq 1$ .
- $a$  is the first shape parameter  $a \geq 0$ .
- $b$  is the second shape parameter  $b \geq 0$ .

**("bin",s,prob,n)** Binomial distribution

- $s$  is the number of successes,  $0 \leq s \leq n$ .
- $prob$  is the probability for success,  $0 \leq prob \leq 1$ .
- $n$  is the number of independent trials,  $n > 0$ .

**("cau",x,< $\mu$ >,< $\sigma$ >>)** Cauchy distribution

- $x$  is the upper limit for integration  $x > 0$ .
- $\mu$  location parameter  $mu \geq 0$ .
- $\sigma$  scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

**("chis",x,df,<nc>)** (noncentral) ChiSquare distribution

- $x$  is the upper limit for integration  $x > 0$ .
- $df$  degrees of freedom  $df > 0$ .
- $nc$  optional noncentrality  $nc \geq 0$ .

- ("expo",x,<, $\sigma$ >) Exponential distribution
- $x$  is the upper limit for integration  $0 \leq x \leq 1$ .
  - $\sigma$  scale parameter ( $\sigma > 0$ ).
- Default is  $\sigma = 1$ .
- ("f",x,ndf,ddf,<,nc>) (noncentral)  $F$  distribution
- $x$  is the upper limit for integration  $x > 0$ .
  - ndf numerator degrees of freedom  $ndf > 0$ .
  - ddf denominator degrees of freedom  $ddf > 0$ .
  - nc optional noncentrality  $nc \geq 0$ .
- ("gam",x,shape,<,scale>) •  $x$  is the upper limit for integration  $x > 0$ .
- shape shape parameter  $shape > 0$ .
  - scale is the scale parameter  $scale > 0$ .
- Default is  $scale = 1$ .
- ("gaus",x,<,< $\mu$ >,< $\sigma$ >>) Gauss (Normal) distribution
- $x$  is the upper limit for integration  $x > 0$ .
  - $\mu$  shape parameter.
  - $\sigma$  scale parameter ( $\sigma > 0$ ).
- Default is  $\mu = 0$  and  $\sigma = 1$ .
- ("geom",m,p) •  $m$  is the upper limit for cumulation  $m > 0$ .
- $p$  is the probability for success of 1,  $0 \leq prob \leq 1$ .
- ("hypg",x,m,k,n,<,r>) Hypergeometric distribution
- $x$  is the upper limit for cumulation  $x > 0$ .
  - $m$  is the population size ( $m \geq 1$ ).
  - $k$  is the number of items ( $0 \leq k \leq m$ ).
  - $n$  is the sample size ( $0 \leq n \leq m$ ).
  - $r$  is an optional odds ratio ( $r > 0$ ).
- Default is  $r = 1$ .
- ("igau",x,d) Inverse Gauss (Wald) distribution
- $x$  is the upper limit for integration  $x > 0$ .
  - $d$  shape parameter  $d > 0$ .
- ("lapl",x,<,< $\mu$ >,< $\sigma$ >>) Laplace distribution
- $x$  is the upper limit for integration  $x > 0$ .
  - $\mu$  shape parameter.
  - $\sigma$  scale parameter ( $\sigma > 0$ ).
- Default is  $\mu = 0$  and  $\sigma = 1$ .
- ("logi",x,<,< $\mu$ >,< $\sigma$ >>) •  $x$  is the upper limit for integration  $x > 0$ .
- $\mu$  shape parameter.
  - $\sigma$  scale parameter ( $\sigma > 0$ ).
- Default is  $\mu = 0$  and  $\sigma = 1$ .



("logn",x,<,<μ>,<σ>>) LogNormal distribution

- x is the upper limit for integration  $x > 0$ .
- μ shape parameter.
- σ scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

("negb",x,prob,n) • x is the number of failures,  $0 \leq f \leq n$ .

- prob is the probability for success,  $0 \leq prob \leq 1$ .
- n is the number of successes,  $n > 0$ .

("norm",x,<,<μ>,<σ>>) Normal (Gauss) distribution

- x is the upper limit for integration  $x > 0$ .
- μ shape parameter.
- σ scale parameter ( $\sigma > 0$ ).

Default is  $\mu = 0$  and  $\sigma = 1$ .

("pare",x,a,<k>) Pareto distribution

- x is the upper limit for integration  $x > 0$ .
- a shape parameter ( $a > 0$ ).
- k scale parameter ( $k > 0$ ).

Default is  $k = 1$ .

("pois",n,λ) Poisson distribution

- x is the upper limit for cumulation  $x > 0$ .
- λ shape parameter.

("t",t,df,<nc>) (noncentral) t distribution

- t is chosen from t distribution.
- df degrees of freedom  $df > 0$ .
- nc optional noncentrality  $nc \geq 0$ .

("unif",x,<l>,<u>>) • x is the upper limit for integration  $x > 0$ .

- l left location.
- u right location ( $u > l$ ).

Default is  $l = 0$  and  $u = 1$ .

("wald",x,d) • x is the upper limit for integration  $x > 0$ .

- d shape parameter  $d > 0$ .

("weib",x,shape,<scale>) Weibull distribution

- t is the upper limit for integration  $x > 0$ .
- shape shape parameter  $shape > 0$ .
- scale scale parameter  $scale > 0$ .

Default is  $scale = 1$ .

... There may be a variable number of additional input arguments specifying non-default distributional parameters.

**Output: prob** The first result is either a scalar or an object of the test probabilities.

**stat** The **stat** result has the same size as **prob** and returns the test statistics (only for `optn[3]=0`).

**Restrictions:** 1. Missing value are returned if the input  $y$  contains missing data.  
2.

**Relationships:** `jarbera()`, `ksprob()`, `kstest()`, `shapwilk()`

**Examples:** 1. Test  $N = 1000$  random uniform generated data:

```
options NOECHO;
#include "..\tdata\matlb.dat"
options ECHO;

optn = [ 2 , /* print */
        .05 , /* alpha */
        1 ]; /* version=cs */
< p21,s21 > = berkow(uf1000,optn,"unif");
print "BERKOW P: UNIF",p21;
print "BERKOW S: UNIF",s21;
```

```
Hypothesis H0: Data are uniformly distributed versus
Hypothesis H1: Data are not uniformly distributed
H0 accepted: Probability=0.0701486 Statistics=5.31428
```

```
optn = [ 2 , /* print */
        .05 , /* alpha */
        1 ]; /* version=cs */
< p22,s22 > = berkow(uf1000,optn,"norm");
print "BERKOW P: NORM",p22;
print "BERKOW S: NORM",s22;
```

```
Hypothesis H0: Data are normally distributed versus
Hypothesis H1: Data are not normally distributed
H0 rejected: Probability=0 Statistics=1843.02
```

## 5.6 Function garch

---

```
gof = garch(meth,y,x,o,q,p<,optn<x0>>)
```

```
< gof,est,ase,cis,cov,sco > = garch(meth,y,x,o,q,p<,optn<x0>>)
```

**Purpose:** The `garch` function implements ARCH, GARCH, TARCH, AVARCH, ZARCH, APARCH, EGARCH, AGARCH, NAGARCH, IGARCH, and FIGARCH time series modeling.

The methods 2, ..., 8 permit the specification of the error distribution using the "error" option:

"norm" Gaussian innovations (default); no additional parameter is estimated;

"stdt" Students  $T$  distributed error; one additional parameter  $\nu > 2$  is estimated;

"gend" generalized error distribution; one additional parameter  $\nu > 1$  is estimated;

"skew" skewed  $T$  error distribution; two additional parameters  $\nu > 2$  and  $-.99 < \lambda < .99$  are estimated.

**meth=1: (G)ARCH** implements the ARCH and GARCH: the `o` input argument specifies the AR lags; the predictor data `x` may be specified; the input `y` must not be centered; for  $p = 0$  the ARCH model is estimated; for  $o > 1$  the optimization is subject to nonlinear constraints; for  $o = 1$  optimization algorithm LMEMQN is default, also NRRIDG or TRUREG are among preferred; for  $o > 1$  optimization algorithm DQNSQP is default or COBYLA could be tried. There are

$$p = ni + nb + nar + nq + np$$

parameters to estimate, where the intercept  $ni = 0$  or  $1$  is optional,  $nb \geq 0$  is the number of columns of  $\mathbf{X}$ , and  $nar$  the number of entries of  $o$ ,  $nq$  the number of entries of  $q$ , and  $np$  the number of entries of  $p$ .

**meth=2: ARCH, GARCH, TARCH, AVGARCH, ZARCH** : The input `y` must be centered around the mean. Currently the argument `x` must be specified as missing.

- the `o` input argument specifies the  $no$  asymmetric innovations;
- the `q` input argument specifies the  $nq$  symmetric innovations;
- the `p` input argument specifies the  $np$  lagged variances.

For  $\delta = 2$  (default) and  $\delta = 1$ :

- for  $\delta = 2$  squared value estimation ARCH, GARCH, GJR-GARCH;
- for  $\delta = 1$  absolute value estimation TARCH, AVGARCH, ZARCH.

The conditional variance is given by:

$$\sigma_\delta^2 = \omega + \sum_{q=1}^Q \alpha_q |\epsilon_{t-q}|^\delta + \sum_{o=1}^O \gamma_o |\epsilon_{t-o}|^\delta I_{[\epsilon_{t-o} < 0]} + \sum_{p=1}^P \beta_p \sigma_{t-p}^\delta$$

There are

$$p = 1 + no + nq + np + nm$$

parameters to estimate, where  $no$  the number of entries of  $o$ ,  $nq$  the number of entries of  $q$ ,  $np$  the number of entries of  $p$ , and  $nm$  is either zero, or one or two depending on estimating additional  $\mu$  and  $\lambda$  parameters for nonnormal error distributions.

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\begin{aligned} \omega &\geq 0 \\ \alpha_q &\geq 0, \quad q = 1, \dots, Q \\ \beta_p &\geq 0, \quad p = 1, \dots, P \\ \alpha_j + \gamma_j &\geq 0, \quad j = 1, \dots, \max(Q, O) \\ \sum_q \alpha_q + .5 \sum_o \gamma_o + \sum_p \beta_p &< 1 \end{aligned}$$

**meth=3: APARCH** implements the APARCH method: which is similar to  $meth = 2$ , but there are two versions:

- a fixed power of  $\delta > .3$  can be specified; i.e. there are  $p = 1 + no + nq + np + nm$  parameters to estimate;
- the best  $\delta > .3$  is estimated as an additional parameter, i.e. there are  $p = 2 + no + nq + np + nm$  parameters to estimate;

The input  $\mathbf{y}$  must be centered around the mean. Currently the argument  $\mathbf{x}$  must be specified as missing.

- the  $o$  input argument specifies the  $no$  asymmetric innovations;
- the  $q$  input argument specifies the  $nq$  symmetric innovations;
- the  $p$  input argument specifies the  $np$  lagged variances.

The conditional variance is given by:

$$\sigma_\delta^2 = \omega + \sum_{j=1}^{\max(Q,O)} \alpha_j (|\epsilon_{t-j}| + \gamma_j \epsilon_{t-j})^\delta + \sum_{p=1}^P \beta_p \sigma_{t-p}^\delta$$

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\omega \geq 0$$

$$\begin{aligned}
\alpha_q &\geq 0, & q = 1, \dots, Q \\
\beta_p &\geq 0, & p = 1, \dots, P \\
-1 &\leq \gamma_o \leq 1, & o = 1, \dots, O \\
\delta &\geq .3 \\
\sum_q \alpha_q + \sum_p \beta_p &< 1
\end{aligned}$$

**meth=4: EGARCH** implements the EGARCH: The input  $y$  must be centered around the mean. Currently the argument  $x$  must be specified as missing.

- the  $o$  input argument specifies the  $no$  asymmetric innovations;
- the  $q$  input argument specifies the  $nq$  symmetric innovations;
- the  $p$  input argument specifies the  $np$  lagged variances.

The conditional variance is given by:

$$\ln(h_t) = \omega + \sum_{q=1}^Q \alpha_q (|\epsilon_{t-q}| - C) + \sum_{o=1}^O \gamma_o \epsilon_{t-o} + \sum_{p=1}^P \beta_p \ln(h_{t-p})$$

where  $e_t = r_t / \sqrt{h_t}$  and  $C = \sqrt{2} / \sqrt{\pi}$ .

There are

$$p = 1 + no + nq + np + nm$$

parameters to estimate, where  $no$  the number of entries of  $o$  etc.

For  $p > 0$  the optimization is subject to  $p$  nonlinear constraints on the absolute size of the roots of the characteristic polynomial of the  $\beta$  parameters. The optimization algorithm DQNSQP is default or COBYLA could be tried; there may be convergence problems if the model is not fitting the data well enough.

**meth=5: AGARCH** implements the AGARCH (Engle, 1990): The input  $y$  must be centered around the mean. Currently the argument  $x$  must be specified as missing. The input argument  $o$  must be specified as missing.

- the  $q$  input argument specifies the  $nq$  symmetric innovations;
- the  $p$  input argument specifies the  $np$  lagged variances.

An additional model parameter  $\gamma$  is estimated.

The volatility dynamics is given by:

$$h_t = \omega + \sum_{q=1}^Q (r_{t-q} - \gamma)^2 + \sum_{p=1}^P h_{t-p}$$

There are

$$p = 2 + nq + np + nm$$

parameters to estimate, where  $nq$  the number of entries of  $q$  etc. The input  $y$  must be centered around the mean.

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\begin{aligned}\omega &\geq 0 \\ \alpha_q &\geq 0, \quad q = 1, \dots, Q \\ \beta_p &\geq 0, \quad p = 1, \dots, P \\ q(.01, y) &\leq \gamma \leq q(.99, y) \\ \sum_q \alpha_q + \sum_p \beta_p &< 1\end{aligned}$$

**meth=6: NAGARCH** implements the NAGARCH (Engle & Ng, 1993): The NAGARCH model is very similar to AGARCH. The input  $y$  must be centered around the mean. Currently the argument  $x$  must be specified as missing.

- The input argument  $o$  must be specified as missing;
- the  $q$  input argument specifies the  $nq$  symmetric innovations;
- the  $p$  input argument specifies the  $np$  lagged variances.

An additional model parameter  $\gamma$  is estimated.

The volatility dynamics is given by:

$$h_t = \omega + \sum_{q=1}^Q (r_{t-q} - \gamma \sqrt{h_{t-q}})^2 + \sum_{p=1}^P h_{t-p}$$

There are

$$p = 2 + nq + np + nm$$

parameters to estimate, where  $nq$  the number of entries of  $q$  etc.

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\begin{aligned}\omega &\geq 0 \\ \alpha_q &\geq 0, \quad q = 1, \dots, Q \\ \beta_p &\geq 0, \quad p = 1, \dots, P \\ \sum_q \alpha_q + \sum_p \beta_p + (1 + \gamma^2) &< 1\end{aligned}$$

**meth=7: IGARCH** implements the IGARCH and IAVARCH methods: The input  $y$  must be centered around the mean. Currently the argument  $x$  must be specified as missing.

- The input argument `o` must be specified as missing;
- the `q` input argument specifies the  $nq$  symmetric innovations;
- the `p` input argument specifies the  $np$  lagged variances.

An intercept  $\omega$  is estimated optional. It is default and is suppressed by `NOINT` option. A  $\delta = 1$  or  $\delta = 2$  (default) can be specified using the options argument.

The conditional variance is given by:

$$g(h_t) = \omega + \sum_{q=1}^Q \alpha_q |r_{t-q}|^\delta + \sum_{p=1}^P \beta_p g(h_{t-p})$$

where

$$g(x) = \begin{cases} \sqrt{x} & \text{for } \delta = 1 \\ x & \text{for } \delta = 2 \end{cases}$$

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\begin{aligned} \omega &\geq 0 \\ \alpha_q &\geq 0, \quad q = 1, \dots, Q \\ \beta_p &\geq 0, \quad p = 1, \dots, P \\ \sum_q \alpha_q + \sum_p \beta_p &= 1 \end{aligned}$$

**meth=8: FIGARCH** implements the FIGARCH method: The input `y` must be centered around the mean. Currently the argument `x` must be specified as missing.

- The input argument `o` must be specified as missing;
- the `q` input argument specifies the order of the short memory autoregressive process, must be 0 or 1;
- the `p` input argument specifies the order of the moving average process, must be 0 or 1.

Additionally the integer truncation lag must be specified with the options argument.

The conditional variance is given by:

$$h_t = \omega + \sum_{i=1}^{\infty} \lambda_i \epsilon_{t-i}^2$$

where

$$\begin{aligned} \delta_1 &= d \\ \lambda_1 &= \Phi - \beta + d \\ \delta_i &= \frac{i-1-d}{i} \delta_{i-1}, \quad i = 2, \dots \\ \lambda_i &= \beta \lambda_{i-1} + \delta_i - \Phi \delta_{i-1}, \quad i = 2, \dots \end{aligned}$$

The following constraints are resolved by model reparametrization so that the optimization is unconstrained:

$$\omega \geq 0$$

$$0 \leq d \leq 1$$

$$0 \leq \Phi \leq (1 - d)/2$$

$$0 \leq \beta \leq d + \Phi$$

**Input: meth** should be an integer scalar  $meth \in \{1, \dots, 8\}$  (see above for method assignment).

**y** should be an  $N$  vector  $\mathbf{Y}$  of time series data.

**x** could be either a missing value or an  $N \times nb$  matrix  $\mathbf{X}$  of predictor data.

**o** should be either an integer scalar or vector of lag numbers depending on the estimation method.

**q** should be either an integer scalar or vector determining the lag numbers of the ARCH parameter.

**p** should be either an integer scalar or vector determining the lag numbers of the GARCH parameter.

**optn** The **optn** argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content.

**x0** a  $p$  vector specifying initial estimates for the optimization.

In addition to many options for the optimization algorithms the following can be specified:



Option	Second Column	Meaning
"alpha"	real	probability for confidence intervals, default=.05
"center"		centers the $y$ data
"delta"	real	power for methods=2,3,7, default is delta=2
"error"	string	error distribution
	"norm"	Gaussian innovations (default);
	"stdt"	$T$ distributed error;
	"gend"	generalized error distribution;
	"skew"	skewed $T$ error distribution;
"noint"		does not use an intercept parameter valid only for methods 1 and
"norob"		suppress robust (sandwich) covariance matrix
"nosco"		suppress computation of scores
"nopr"		do not print any output
"pcov"		print asymptotic covariance matrices
"phis"		print optimization history
"pscor"		print (large $N \times p$ ) score matrix
"print"	int	amount of printed output (def=0: no output)
"truncl"	int	size of truncation lag (default=1000) valid only for meth=8
"tech"	string	optimization technique see function <code>nlp()</code>

**Output:** `gof` a vector of scalar results, see below for content;

**est** a  $p \times 2$  matrix of estimates (in first column) and the gradient vector in its second column;

**ase** a  $p \times 3$  matrix with asymptotic standard error (based on the diagonal of the robust covariance matrix) in its first column, the  $t$  value in its second, and the  $p$  value in its third column;

**cis** a  $p \times 2$  matrix of (robust) lower and upper Wald confidence intervals of the estimates;

**cov** a  $p \times p$  (robust, sandwich, White) asymptotic covariance matrix;

**sco** is an  $N \times p$  matrix of scores.

The current content of the `gof` vector:

1. indicator for problems in the algorithm (=0: no problems)
2. mean value  $\mu$  of input data  $y$
3. standard deviation  $\sigma$  of input data  $y$
4. variance  $\sigma^2$
5. computation time in seconds
6. Loglikelihood
7. Akaike's information criterion AIC

8. Bayesian (Schwartz) information criterion BIC
9. number of iterations for optimization
10.  $L1$  norm of gradient at the optimal solution

**Restrictions:** 1. The input data  $y$  must be numeric and cannot contain any string or complex data.

2. Currently the input data  $y$  must not have any missing values.

**Relationships:** arma(), arima(), tsmeas()

**Examples:** 1. Data:

The following time series data has  $N = 300$  observations:

```
print "Read Data TWO by Minbo: from nlp213";
options NOECHO;
%inc "c:\cmat\tdata\minbo0.dat";
options ECHO;
```

2. Method 1:

```
m = q = p = 1;
print "Method 1: Unconstrained (GARAR)";
optn = [ "tech"      "lq" ,
        "cent"      ,
        "print"     3 ];
< gof,est,ase,cis,cov,scor > = garch(1,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;
```

```
*****
Model 1: (G)ARCH Modeling: AR=1 NQ=1 NP=1
*****
AR[1] =    1 Q[1] =    1 P[1] =    1

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Optimization Technique . . . . . LMEMQN
```

\*\*\*\*\*  
 Optimization Start  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient	Lower BC	Upper BC
1 AR_1	0.00000000	-24.019592	-0.9999000	0.9999000
2 ARCH_0	1.30926796	0.0016684	1.00e-008	.
3 ARCH_1	1.000e-006	-29.834384	1.00e-008	.
4 GARCH_1	1.000e-006	0.0015156	1.00e-008	.

Value of Objective Function = 466.102

Limited Memory Quasi Newton Method  
 Gradient Computed by Finite Differences

Iteration Start:

N. Variables	4		
N. Bound. Constr.	5	N. Mask Constr.	0
Criterion	466.1017562	Max Grad Entry	29.83438353
N. Active Constraints	0	N. Grad Storage	5

Iter	nfun	act	optcrit	difcrit	maxgrad	alpha	slope	gcrit
1	3	2	465.0617	1.040046	18.4056	7e-003	-914.108	1.30926
2	4	1	462.8073	2.254445	14.3460	1.00000	-2.78328	1.25336
3	5	1	460.9802	1.827046	9.98118	1.00000	-3.48729	20.0384
4	7	0	454.6460	6.334191	15.0617	5.00000	-1.30142	1.43248
5	10	0	454.1369	0.509175	11.4576	0.15347	-5.91310	1.48010
6	11	0	453.2002	0.936687	15.2952	1.00000	-1.84432	15.2952
7	12	0	452.7153	0.484894	7.43397	1.00000	-0.68535	7.43397
8	13	0	452.4463	0.268983	6.31992	1.00000	-0.51750	0.85465
9	14	0	452.1939	0.252427	2.23234	1.00000	-0.33597	1.82205
10	15	0	452.1174	0.076441	1.79658	1.00000	-0.21904	1.79658
11	16	0	452.1039	0.013526	0.29609	1.00000	-0.03298	0.29609
12	17	0	452.1032	6.6e-004	0.28550	1.00000	-2e-003	0.28550
13	19	0	452.1031	9.1e-005	0.06968	0.19375	-9e-004	0.06968
14	20	0	452.1031	2.1e-005	5e-003	1.00000	-4e-005	5e-003
15	21	0	452.1031	7.9e-008	3e-004	1.00000	-2e-007	3e-004
16	22	0	452.1031	5.5e-010	3e-004	1.00000	-7e-010	3e-004
17	23	0	452.1031	3.6e-010	9e-005	1.00000	-1e-009	9e-005
18	24	0	452.1031	6.3e-011	8e-006	1.00000	-1e-010	8e-006

Successful Termination After 18 Iterations  
 ABSGCONV convergence criterion satisfied.

Criterion	452.1031258	Max Grad Entry	8.0944e-006
N. Active Constraints	0	N. Grad Storage	5
Skipped Updates	0	Segments Explored	21
N. Function Calls	25	N. Gradient Calls	25
N. Line Searches	18	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient	Active BC
1 AR_1	0.26060427	8.09e-006	
2 ARCH_0	0.42786973	-6.27e-006	
3 ARCH_1	0.30722878	-7.83e-007	
4 GARCH_1	0.37519470	-4.16e-007	

Value of Objective Function = 452.103

Log Likelihood= -452.103  
 SBC=-881.391 AIC=-896.206

Covariance Matrix 2: H = (NOBS/d) inv(G)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	AR_1	ARCH_0	ARCH_1	GARCH_1
AR_1	0.0119455			
ARCH_0	0.0010376	0.0125162		
ARCH_1	-6.45e-004	0.0010024	0.0097181	
GARCH_1	-3.92e-004	-0.0095789	-0.0066432	0.0128294

Factor sigma=1 (nobs=300 df=4 vard=300)

Determinant = 2.72817e-009

Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	AR_1	ARCH_0	ARCH_1	GARCH_1
AR_1	0.0363805			
ARCH_0	-0.0023527	0.0077765		
ARCH_1	-0.0060095	-6.61e-004	0.0094990	
GARCH_1	0.0043184	-0.0043883	-0.0045065	0.0064840

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	AR_1	ARCH_0	ARCH_1	GARCH_1
AR_1	1.0000000			
ARCH_0	-0.1398734	1.0000000		
ARCH_1	-0.3232716	-0.0769177	1.0000000	
GARCH_1	0.2811663	-0.6179963	-0.5742187	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
AR_1	0.2606043	0.1907367	1.3663040	0.086440
ARCH_0	0.4278697	0.0881842	4.8520000	0.000001
ARCH_1	0.3072288	0.0974628	3.1522683	0.000893
GARCH_1	0.3751947	0.0805233	4.6594522	0.000002

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
AR_1	0.2606043	-0.1132327	0.6344413
ARCH_0	0.4278697	0.2550319	0.6007076
ARCH_1	0.3072288	0.1162053	0.4982523
GARCH_1	0.3751947	0.2173718	0.5330176

Factor sigma=1 (nobs=300 df=4 vard=300)

3. Method 2: GARCH: delta=2:

```

m = q = p = 1;
print "Method 2: Reparametrized Constraints: delta=2";
optn = [ "cent"          ,
         "delta"         2. ,
         "tech"          "tr" ,
         "error"         "norm" ,
         "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(2,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 2: (G)ARCH Modeling: NO=1 NQ=1 NP=1
*****
Modeling Squared Values (Delta=2)
O[1] = 1 Q[1] = 1 P[1] = 1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal
Optimization Technique . . . . . TRUREG

```

```

*****
Optimization Start
*****

```

Parameter Estimates  
-----

Parameter	Estimate	Gradient
1 Intercept	-1.33996974	-7.4183275
2 Alpha_1	-1.38629190	0.3059796
3 Gamma_1	-1.25276015	0.1473073

4 Beta\_1 0.91630058 -1.9482010

Value of Objective Function = 455.645

Hessian Matrix  
\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	47.491963	1.7283319	-4.3302148	26.269433
2	1.7283319	3.5187833	-2.9335752	1.6697438
3	-4.3302148	-2.9335752	7.3798025	-1.0955059
4	26.269433	1.6697438	-1.0955059	16.182565

Trust Region Optimization  
Without Parameter Scaling  
Gradient Computed by Finite Differences  
Hessian Computed by Finite Differences (dense)  
(Using Only Function Calls)

Iteration Start:

N. Variables	4		
Criterion	455.6448096	Max Grad Entry	7.418327504
TR Radius	1.000000000		

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1	0	2	0	454.4628	1.181992	4.81094	1.10038	1.00000
2	0	3	0	454.2628	0.200034	0.20377	0.00000	1.04053
3	0	4	0	454.2590	3.8e-003	7e-003	0.00000	0.18087
4	0	5	0	454.2590	1.3e-006	4e-006	0.00000	0.04976

Successful Termination After 4 Iterations

GCONV convergence criterion satisfied.

Criterion	454.2590151	Max Grad Entry	4.0547e-006
Ridge (lambda)	0.000000000	TR Radius	0.049761569
Act.dF/Pred.dF	0.999835589		
N. Function Calls	6	N. Gradient Calls	3
N. Hessian Calls	6	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*

Optimization Results

\*\*\*\*\*

Parameter Estimates

-----

Parameter	Estimate	Gradient
-----------	----------	----------

1 Intercept	-0.82432306	-4.05e-006
2 Alpha_1	-1.35239391	1.92e-007
3 Gamma_1	-1.08135794	4.42e-007
4 Beta_1	0.14335696	-3.94e-006

Value of Objective Function = 454.259

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	62.341578	3.1051598	-1.4674951	25.986019
2	3.1051598	2.7639938	-1.8008995	1.4816808
3	-1.4674951	-1.8008995	5.2601208	-0.0024428
4	25.986019	1.4816808	-0.0024428	13.691414

Log Likelihood= -454.259

SBC=-885.703 AIC=-900.518

Covariance Matrix 2: H = (NOBS/d) inv(G)

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0151475			
Alpha_1	5.65e-004	0.0133848		
Gamma_1	0.0029399	-0.0071469	0.0263580	
Beta_1	-0.0120534	-0.0071370	-0.0045574	0.0165277

Factor sigma=1 (nobs=300 df=4 vard=300)

Determinant = 8.94864e-009



Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0088788			
Alpha_1	2.27e-005	0.0124115		
Gamma_1	-7.70e-005	-0.0048557	0.0213510	
Beta_1	-0.0058128	-0.0068907	-0.0017048	0.0101131

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	1.0000000			
Alpha_1	0.0021639	1.0000000		
Gamma_1	-0.0055927	-0.2982873	1.0000000	
Beta_1	-0.6134338	-0.6150440	-0.1160142	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.4385318	0.0942275	4.6539678	0.000002
Alpha_1	0.2054382	0.1114069	1.8440345	0.033089
Gamma_1	0.2489317	0.1461199	1.7036131	0.044751
Beta_1	0.3589155	0.1005640	3.5690243	0.000209

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.4385318	0.2538492	0.6232143
Alpha_1	0.2054382	-0.0129153	0.4237917

```

Gamma_1  0.2489317 -0.0374580  0.5353214
Beta_1   0.3589155  0.1618136  0.5560174
Factor sigma=1 (nobs=300 df=4 vard=300)

```

4. Method 2: GARCH: delta=1:

```

m = q = p = 1;
print "Method 2: Reparametrized Constraints: delta=1";
optn = [ "cent"          ,
         "delta"         1. ,
         "tech"          "tr" ,
         "error"         "norm" ,
         "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(2,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 2: (G)ARCH Modeling: NO=1 NQ=1 NP=1
*****
Modeling Absolute Values (Delta=1)
O[1] = 1 Q[1] = 1 P[1] = 1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal
Optimization Technique . . . . . TRUREG

```

```

*****
Optimization Start
*****

```

Parameter Estimates

```

-----
Parameter      Estimate      Gradient

```

```

1 Intercept-1.53586564 -58.124474
2 Alpha_1 -1.38629190 3.6784135
3 Gamma_1 -2.02437929 -6.7925300
4 Beta_1 1.09025389 -34.811202

```

Value of Objective Function = 464.281

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

```

S |          1          2          3          4
-----
1 | 244.19334  4.7451953  20.045738  170.53481
2 |  4.7451953  12.384253 -6.9473193  7.9429617
3 |  20.045738 -6.9473193  15.343707  17.500619
4 |  170.53481  7.9429617  17.500619  115.50154

```

Trust Region Optimization

Without Parameter Scaling

Gradient Computed by Finite Differences

Hessian Computed by Finite Differences (dense)

(Using Only Function Calls)

Iteration Start:

```

N. Variables          4
Criterion             464.2813571      Max Grad Entry  58.12447438
TR Radius             1.000000000

```

```

Iter rest nfun act  optcrit  difcrit  maxgrad  lambda  radius
1*  0  2  0  456.3216  7.959783  20.3783  7.64481  1.00000
2  0  3  0  455.2247  1.096896  1.54743  0.00000  1.04789
3  0  4  0  455.1950  0.029695  0.14818  0.00000  0.30112
4  0  5  0  455.1947  2.8e-004  3e-003  0.00000  0.12380
5  0  6  0  455.1947  6.3e-008  1e-006  0.00000  0.01537

```

Successful Termination After 5 Iterations

GCONV convergence criterion satisfied.

```

Criterion             455.1947014      Max Grad Entry  1.2954e-006
Ridge (lambda)       0.000000000      TR Radius       0.015373816
Act.dF/Pred.dF      1.000341620
N. Function Calls          7      N. Gradient Calls          3

```

N. Hessian Calls	7	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	-0.84061128	1.30e-006
2 Alpha_1	-1.48322181	1.51e-008
3 Gamma_1	-1.57885226	2.15e-007
4 Beta_1	0.22287797	3.49e-007

Value of Objective Function = 455.195

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	287.78217	9.7874108	6.9416767	133.94190
2	9.7874108	5.3759445	-3.8963319	5.0234865
3	6.9416767	-3.8963319	11.843638	4.2857946
4	133.94190	5.0234865	4.2857946	65.293889

Log Likelihood= -455.195  
 SBC=-887.574 AIC=-902.389

Covariance Matrix 2:  $H = (NOBS/d) \text{inv}(G)$   
 \*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0151876			
Alpha_1	0.0018433	0.0068219		

```
Gamma_1 | 4.43e-004 -0.0035559 0.0077095
Beta_1 | -0.0148603 -0.0052871 -3.88e-004 0.0175127
```

```
Factor sigma=1 (nobs=300 df=4 vard=300)
Determinant = 3.03027e-010
Matrix has Only Positive Eigenvalues
```

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

```

S | Intercept Alpha_1 Gamma_1 Beta_1
-----
Intercept | 0.0097406
Alpha_1 | 4.07e-004 0.0062409
Gamma_1 | -0.0010235 -0.0027778 0.0062047
Beta_1 | -0.0084654 -0.0039304 0.0010388 0.0102681
```

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

```

S | Intercept Alpha_1 Gamma_1 Beta_1
-----
Intercept | 1.0000000
Alpha_1 | 0.0522012 1.0000000
Gamma_1 | -0.1316572 -0.4463881 1.0000000
Beta_1 | -0.8464653 -0.4909820 0.1301413 1.0000000
```

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

```
Parameter Estimate AStdErr T Value P Value
Intercept 0.4314467 0.0986943 4.3715476 0.000009
Alpha_1 0.1849043 0.0789995 2.3405741 0.009959
Gamma_1 0.1253326 0.0787700 1.5911220 0.056325
Beta_1 0.4178559 0.1013315 4.1236515 0.000024
```

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.4314467	0.2380095	0.6248839
Alpha_1	0.1849043	0.0300680	0.3397405
Gamma_1	0.1253326	-0.0290537	0.2797189
Beta_1	0.4178559	0.2192498	0.6164620

Factor sigma=1 (nobs=300 df=4 vard=300)

5. Method 3: APARCH: specify  $\delta = 1.5$ :

```

m = q = p = 1;
print "Method 3: APARCH: Reparametrized Constraints: Specify Delta";
optn = [ "cent"          ,
         "delta"         1.5 ,
         "tech"          "tr" ,
         "error"         "norm" ,
         "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(3,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

\*\*\*\*\*  
Model 3: APARCH Modeling: NO=1 NQ=1 NP=1  
\*\*\*\*\*

Modeling Constant Delta=1.5  
O[1] = 1 Q[1] = 1 P[1] = 1

Number of Observations . . . . .	300
Number of Parameter Estimates . . . . .	4
Mean Y . . . . .	0.0016
Standard Deviation Y . . . . .	1.1442
Mean Squared Error (MSE) . . . . .	1.3093
Number of Back Casts . . . . .	1
Error Distribution . . . . .	Normal
Optimization Technique . . . . .	TRUREG

\*\*\*\*\*  
Optimization Start  
\*\*\*\*\*

Parameter Estimates

Parameter	Estimate	Gradient
1 Intercept	-0.80807956	-67.266644
2 Alpha_1	-1.04456937	-24.750763
3 Gamma_1	0.00000000	1.9300417
4 Beta_1	0.14454518	-134.16255

Value of Objective Function = 484.494

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	104.67489	52.890431	-3.3000239	241.90207
2	52.890431	9.0992573	-0.2050936	61.237870
3	-3.3000239	-0.2050936	0.3314837	-4.9151948
4	241.90207	61.237870	-4.9151948	376.88691

Trust Region Optimization  
 Without Parameter Scaling  
 Gradient Computed by Finite Differences  
 Hessian Computed by Finite Differences (dense)  
 (Using Only Function Calls)

Iteration Start:

N. Variables 4  
 Criterion 484.4939750 Max Grad Entry 134.1625549  
 TR Radius 1.00000000

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1*	0	3	0	461.7268	22.76718	25.9680	66.6286	0.45258
2	0	5	0	458.1502	3.576632	9.88555	1.44608	1.42239
3	0	6	0	457.5656	0.584535	0.71054	0.08086	1.46799
4	0	7	0	457.4567	0.108951	0.16465	0.00000	1.51947
5	0	8	0	457.4223	0.034412	0.04680	0.00000	1.09544
6	0	9	0	457.4101	0.012209	0.01514	0.00000	1.01036
7	0	10	0	457.4055	4.5e-003	5e-003	0.00000	0.95878

8	0	11	0	457.4036	1.9e-003	3e-003	0.00000	0.90175
9	0	12	0	457.4029	7.6e-004	1e-003	0.00000	0.95438
10	0	13	0	457.4026	2.4e-004	2e-004	0.00000	1.02312
11	0	14	0	457.4026	5.2e-005	1e-004	0.00000	0.79537
12*	0	15	0	457.4025	9.5e-005	3e-004	2e-004	1.10559
13	0	16	0	457.4025	1.6e-005	2e-004	0.00000	1.11262
14	0	17	0	457.4025	3.8e-006	6e-005	0.00000	1.14367

Successful Termination After 14 Iterations

GCONV convergence criterion satisfied.

Criterion	457.4024572	Max Grad Entry	5.9941e-005
Ridge (lambda)	0.000000000	TR Radius	1.143670882
Act.dF/Pred.dF	1.871066805		
N. Function Calls	18	N. Gradient Calls	7
N. Hessian Calls	16	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates

-----

Parameter	Estimate	Gradient
1 Intercept	-1.13572475	-5.99e-005
2 Alpha_1	0.45005298	-3.25e-005
3 Gamma_1	-10.9289805	2.68e-005
4 Beta_1	0.34625681	-1.76e-005

Value of Objective Function = 457.402

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	31.066380	12.768076	-1.83e-004	110.73000
2	12.768076	8.3003477	2.24e-005	53.621172
3	-1.83e-004	2.24e-005	-7.53e-005	-2.41e-004
4	110.73000	53.621172	-2.41e-004	481.50746



Log Likelihood= -457.402  
SBC=-891.99 AIC=-906.805

Covariance Matrix 2: H = (NOBS/d) inv(G)  
\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0195339			
Alpha_1	-0.0019982	0.0249742		
Gamma_1	0.0060737	-0.0028987	0.0357288	
Beta_1	-0.0127947	-0.0100935	-0.0013812	0.0157775

Factor sigma=1 (nobs=300 df=4 vard=300)  
Determinant = 3.37945e-008  
Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates  
\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0175506			
Alpha_1	-0.0120204	0.0257178		
Gamma_1	0.0145278	-0.0151766	0.0377656	
Beta_1	-0.0055463	-0.0031582	-0.0029162	0.0061195

Robust Asymptotic Correlation Matrix of Estimates  
\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	1.0000000			
Alpha_1	-0.5657894	1.0000000		
Gamma_1	0.5642936	-0.4869768	1.0000000	
Beta_1	-0.5351819	-0.2517489	-0.1918264	1.0000000

```
*****
Parameter Estimates and Asympt. Standard Errors
*****
```

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.3211893	0.1324788	2.4244575	0.007965
Alpha_1	0.6105297	0.1603676	3.8070647	0.000085
Gamma_1	0.1047319	0.1943338	0.5389280	0.295171
Beta_1	0.3462568	0.0782271	4.4263051	0.000007

```
*****
Wald Confidence Limits (Alpha=0.05)
*****
```

Parameter	Estimate	Lower	Upper
Intercept	0.3211893	0.0615356	0.5808430
Alpha_1	0.6105297	0.2962150	0.9248444
Gamma_1	0.1047319	-0.2761553	0.4856192
Beta_1	0.3462568	0.1929346	0.4995790

Factor sigma=1 (nobs=300 df=4 vard=300)

6. Method 3: APARCH: estimate  $\delta$ :

```
m = q = p = 1;
print "Method 3: APARCH: Reparametrized Constraints: Estimate Delta";
optn = [ "cent"          ,
        "delta"         . ,
        "tech"          "tr" ,
        "error"         "norm" ,
        "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(3,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;
```

```
*****
Model 3: APARCH Modeling: NO=1 NQ=1 NP=1
*****
Modeling Variable Delta=2
O[1] = 1 Q[1] = 1 P[1] = 1
```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 5
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal
Optimization Technique . . . . . TRUREG

```

```

*****
Optimization Start
*****

```

Parameter Estimates  
-----

Parameter	Estimate	Gradient
1 Intercept	-0.80807956	-67.266644
2 Alpha_1	-1.04456937	-24.750763
3 Gamma_1	0.00000000	1.9300417
4 Beta_1	0.14454518	-134.16255
5 Delta	-1.45528723	-16.203738

Value of Objective Function = 484.494

Hessian Matrix  
\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4	5
1	104.67489	52.890431	-3.3000239	241.90207	68.519298
2	52.890431	9.0992573	-0.2050936	61.237870	21.760473
3	-3.3000239	-0.2050936	0.3314837	-4.9151948	-0.7538566
4	241.90207	61.237870	-4.9151948	376.88691	118.88276
5	68.519298	21.760473	-0.7538566	118.88276	12.505634

Trust Region Optimization  
Without Parameter Scaling  
Gradient Computed by Finite Differences  
Hessian Computed by Finite Differences (dense)  
(Using Only Function Calls)

Iteration Start:

N. Variables 5  
Criterion 484.4939750 Max Grad Entry 134.1625549  
TR Radius 1.000000000

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1*	0	3	0	462.6536	21.84039	35.7901	196.014	0.22429
2*	0	4	0	460.9077	1.745880	11.7648	17.8182	0.24322
3*	0	7	0	456.9692	3.938522	3.02933	0.84716	2.24672
4*	0	8	0	454.7810	2.188210	0.42557	0.44642	2.25214
5	0	9	0	454.4958	0.285173	0.53019	0.00000	2.32350
6	0	10	0	454.4208	0.074962	0.15293	0.00000	1.42516
7	0	11	0	454.3954	0.025435	0.06860	0.00000	1.29484
8	0	12	0	454.3861	9.3e-003	0.02121	0.00000	1.01182
9*	0	15	0	454.3843	1.9e-003	4e-003	-4e-003	2.14396
10*	0	17	0	454.3830	1.3e-003	2e-003	6e-003	1.69460
11*	0	18	0	454.3824	5.5e-004	2e-003	-2e-003	1.70251
12*	0	20	0	454.3819	5.4e-004	3e-003	3e-003	1.79430
13	0	21	0	454.3812	7.3e-004	2e-003	0.00000	1.79980
14*	0	24	0	454.3812	3.7e-006	3e-004	0.02389	0.03805

Successful Termination After 14 Iterations

GCONV convergence criterion satisfied.

Criterion 454.3811585 Max Grad Entry 0.000290528  
Ridge (lambda) 0.023888796 TR Radius 0.038049758  
Act.dF/Pred.dF 1.007922555  
N. Function Calls 25 N. Gradient Calls 7  
N. Hessian Calls 16 Preproces. Time 0  
Time for Method 0 Effective Time 0

\*\*\*\*\*  
Optimization Results  
\*\*\*\*\*

Parameter Estimates

-----  
Parameter Estimate Gradient  
1 Intercept -1.56352973 -8.09e-006  
2 Alpha\_1 9.76986195 -2.91e-004  
3 Gamma\_1 -2.95937948 1.38e-005

4 Beta\_1 0.23833660 1.44e-006  
 5 Delta 1.18702907 4.02e-005

Value of Objective Function = 454.381

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4	5
1	2.5715182	-1.40e-004	-1.91e-004	7.7087105	0.4908551
2	-1.40e-004	4.70e-004	0.0000000	8.43e-004	-2.63e-004
3	-1.91e-004	0.0000000	-2.39e-004	3.16e-004	5.82e-004
4	7.7087105	8.43e-004	3.16e-004	134.15224	4.4204392
5	0.4908551	-2.63e-004	5.82e-004	4.4204392	0.2892853

Log Likelihood= -454.381  
 SBC=-880.243 AIC=-898.762

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (5 by 5)

S	Intercept	Alpha_1	Gamma_1	Beta_1	Delta
Intercept	0.0336666				
Alpha_1	-0.0295690	0.2319315			
Gamma_1	0.0104679	-0.0055272	0.0208672		
Beta_1	0.0094309	-0.0646167	0.0047090	0.0335773	
Delta	-0.2785531	0.5637437	-0.1686795	-0.3327043	6.3247923

Factor sigma=1 (nobs=300 df=5 vard=300)  
 Determinant = 2.79972e-006  
 Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates  
 \*\*\*\*\*

Dense Symmetric Matrix (5 by 5)

S	Intercept	Alpha_1	Gamma_1	Beta_1	Delta
Intercept	0.0351807				
Alpha_1	-0.0283652	0.1171894			
Gamma_1	0.0222053	-0.0201169	0.0291104		
Beta_1	0.0133267	-0.0234699	0.0095243	0.0164274	
Delta	-0.3641835	0.2552282	-0.3449748	-0.2426870	6.6053951

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (5 by 5)

S	Intercept	Alpha_1	Gamma_1	Beta_1	Delta
Intercept	1.0000000				
Alpha_1	-0.4417626	1.0000000			
Gamma_1	0.6938757	-0.3444232	1.0000000		
Beta_1	0.5543539	-0.5349114	0.4355375	1.0000000	
Delta	-0.7554725	0.2900916	-0.7867091	-0.7367370	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.2093957	0.1875652	1.1163889	0.132582
Alpha_1	0.9997429	0.3423293	2.9204126	0.001883
Gamma_1	1.65e-005	0.1706176	9.67e-005	0.499961
Beta_1	0.2383366	0.1281695	1.8595424	0.031973
Delta	3.1349744	2.5700963	1.2197887	0.111760

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.2093957	-0.1582253	0.5770166
Alpha_1	0.9997429	0.3287897	1.6706960
Gamma_1	1.65e-005	-0.3343878	0.3344208
Beta_1	0.2383366	-0.0128710	0.4895442
Delta	3.1349744	-1.9023218	8.1722706

Factor sigma=1 (nobs=300 df=5 vard=300)

7. Method 4: EGARCH

```

m = q = p = 1;
print "Method 4: EGARCH: Nonlinear Constraints: DQNSQP";
optn = [ "cent"           ,
         "tech"          "sqp" ,
         "error"         "norm" ,
         "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(4,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIS=",cis;
print "COV=",cov;
print "SCOR=",scor;

```

```

*****
Model 4: EGARCH Modeling: NO=1 NQ=1 NP=1
*****
O[1] =    1 Q[1] =    1 P[1] =    1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal
Optimization Technique . . . . . DQNSQP

```

```

*****
Optimization Start
*****

```

Parameter Estimates  
-----

Parameter	Estimate	Gradient	Grad. LF	Lower BC	Upper BC
1 Intercept	0.11103067	5.20e-005	5.20e-005	.	.
2 Alpha_1	0.50000000	0.6583880	0.6583880	-3.0000000	3.0000000
3 Gamma_1	-0.20000000	-15.839860	-15.839860	-3.0000000	3.0000000
4 Beta_1	0.50000000	-7.3873198	-7.3873198	.	.

Value of Objective Function = 457.143

Value of Lagrange Function = 457.143

Values of Nonlinear Constraints

-----

Constraint	Residual
1 1	0.4999800

NLC Jacobian Matrix

\*\*\*\*\*

Row Vector: Dense Storage

R	1	2	3	4
	0.0000000	0.0000000	0.0000000	-1.0000000

DQNSQP (Modified VMCWD) Algorithm by Powell (1978, 1982)  
Lagrange Multiplier Update of Powell(1982)  
Gradient Computed by Finite Differences  
Jacobian Nonlinear Constraints (dense)  
(Computed by Finite Differences)

Iteration Start:

N. Variables	4		
N. Bound. Constr.	4	N. Mask Constr.	0
N. NonLin Constr.	1	NonLin EquConstr.	0
Criterion	457.1429772	Max Grad Entry	15.83985967
N. Active Constraints	0	Max Const Viol.	0.000000000

Iter	rest	nfun	objfunc	conmax	pred	alfa	lfgmax
1	0	4	456.134	0.00000	0.43037	0.06967	5.770163
2	0	5	455.898	0.00000	0.45711	1.00000	5.238383
3	0	6	455.683	0.00000	0.02384	1.00000	1.366673
4	0	7	455.670	0.00000	6e-003	1.00000	0.670068
5	0	8	455.666	0.00000	1e-003	1.00000	0.229266
6	0	9	455.665	0.00000	4e-005	1.00000	0.110952
7	0	10	455.665	0.00000	2e-006	1.00000	0.028218

Successful Termination After 7 Iterations

GCONV convergence criterion satisfied.

Criterion	455.6652474	Max Grad Entry	0.028217578
Lagrange Funct.	455.6652474	Max Grad LagF.	0.028217578



Max Const Viol.	0.00000000	N. Active Constraints	0
N. Function Calls	11	N. Gradient Calls	9
Preproces. Time	0	Time for Method	0
Effective Time	0		

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient	Grad. LF	Active BC
1 Intercept	0.06951210	-0.0282176	-0.0282176	
2 Alpha_1	0.45112468	-0.0015194	-0.0015194	
3 Gamma_1	-0.10095187	-0.0099995	-0.0099995	
4 Beta_1	0.64448549	-0.0046092	-0.0046092	

Value of Objective Function = 455.665  
 Value of Lagrange Function = 455.665

Values of Nonlinear Constraints  
 -----

Constraint	Residual
1 1	0.3554945

NLC Jacobian Matrix  
 \*\*\*\*\*

Row Vector: Dense Storage

R	1	2	3	4
	0.0000000	0.0000000	0.0000000	-1.0000000

Log Likelihood= -455.665  
 SBC=-888.515 AIC=-903.33

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0024646			
Alpha_1	0.0016399	0.0175085		
Gamma_1	-2.99e-005	2.80e-004	0.0050599	
Beta_1	-0.0026972	-0.0043345	-3.90e-004	0.0134336

Factor sigma=1 (nobs=300 df=4 vard=300)

Determinant = 2.05744e-009

Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	0.0020083			
Alpha_1	7.84e-004	0.0173408		
Gamma_1	5.04e-007	-1.67e-004	0.0038506	
Beta_1	-0.0019126	-8.94e-004	-0.0018338	0.0107178

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma_1	Beta_1
Intercept	1.0000000			
Alpha_1	0.1328165	1.0000000		
Gamma_1	1.81e-004	-0.0204093	1.0000000	
Beta_1	-0.4122337	-0.0656023	-0.2854585	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.0695121	0.0448144	1.5511109	0.060972
Alpha_1	0.4511247	0.1316844	3.4258023	0.000350

```
Gamma_1 -0.1009519  0.0620534 -1.6268539  0.947584
Beta_1  0.6444855  0.1035270  6.2252882  0.000000
```

```
*****
Wald Confidence Limits (Alpha=0.05)
*****
```

Parameter	Estimate	Lower	Upper
Intercept	0.0695121	-0.0183225	0.1573467
Alpha_1	0.4511247	0.1930280	0.7092213
Gamma_1	-0.1009519	-0.2225744	0.0206706
Beta_1	0.6444855	0.4415763	0.8473947

Factor sigma=1 (nobs=300 df=4 vard=300)

## 8. Method 5: AGARCH

```
m = .; q = p = 1;
print "Method 5: AGARCH: Reparametrized Constraints";
optn = [ "cent"      ,
         "tech"      "tr"  ,
         "error"     "norm" ,
         "print"     3  ];
< gof,est,ase,cis,cov,scor > = garch(5,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;
```

```
*****
Model 5: AGARCH Modeling: NQ=1 NP=1
*****
```

Q[1] = 1 P[1] = 1

Number of Observations . . . . .	300
Number of Parameter Estimates . . . . .	4
Mean Y . . . . .	0.0016
Standard Deviation Y . . . . .	1.1442
Mean Squared Error (MSE) . . . . .	1.3093
Number of Back Casts . . . . .	1
Error Distribution . . . . .	Normal
Optimization Technique . . . . .	TRUREG

```
*****
```

Optimization Start

\*\*\*\*\*

Parameter Estimates

-----

Parameter      Estimate      Gradient

1 Intercept -0.80893295 -0.0156817  
2 Alpha\_1 -0.65145797 -8.63e-004  
3 Gamma 0.12019977 -10.120248  
4 Beta\_1 0.08284350 -0.0055090

Value of Objective Function = 455.504

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	60.094827	2.1424727	9.0724793	24.498126
2	2.1424727	4.3238422	-2.4711464	1.3176409
3	9.0724793	-2.4711464	63.271550	4.0454594
4	24.498126	1.3176409	4.0454594	12.970194

Trust Region Optimization

Without Parameter Scaling

Gradient Computed by Finite Differences

Hessian Computed by Finite Differences (dense)

(Using Only Function Calls)

Iteration Start:

N. Variables 4  
Criterion 455.5035429      Max Grad Entry 10.12024794  
TR Radius 1.000000000

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1	0	2	0	454.7243	0.779193	0.90219	0.00000	1.00000
2	0	3	0	454.6731	0.051214	0.14599	0.00000	0.20814
3	0	4	0	454.6727	4.2e-004	1e-003	0.00000	0.15570
4	0	5	0	454.6727	4.0e-008	1e-006	0.00000	0.01078

Successful Termination After 4 Iterations  
 GCONV convergence criterion satisfied.

Criterion	454.6727129	Max Grad Entry	1.4851e-006
Ridge (lambda)	0.000000000	TR Radius	0.010779429
Act.dF/Pred.dF	0.997041601		
N. Function Calls	6	N. Gradient Calls	7
N. Hessian Calls	6	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	-0.82184321	7.21e-008
2 Alpha_1	-0.68460525	1.31e-006
3 Gamma	0.28751079	-2.37e-007
4 Beta_1	0.04295362	1.49e-006

Value of Objective Function = 454.673

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	59.097069	3.0961381	12.899948	24.465324
2	3.0961381	4.6931412	5.4377903	2.0330739
3	12.899948	5.4377903	61.523260	6.6195508
4	24.465324	2.0330739	6.6195508	13.145823

Log Likelihood= -454.673  
 SBC=-886.53 AIC=-901.345

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	0.0144828			
Alpha_1	0.0013327	0.0123990		
Gamma	-9.80e-004	-0.0054169	0.0302728	
Beta_1	-0.0109716	-0.0083462	0.0020905	0.0149192

Factor sigma=1 (nobs=300 df=4 vard=300)  
 Determinant = 1.0427e-008  
 Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	0.0087005			
Alpha_1	-0.0011078	0.0119599		
Gamma	-3.57e-004	-0.0048228	0.0235804	
Beta_1	-0.0048669	-0.0059886	0.0019479	0.0080261

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	1.0000000			
Alpha_1	-0.1085972	1.0000000		
Gamma	-0.0249095	-0.2871868	1.0000000	
Beta_1	-0.5824019	-0.6112349	0.1415926	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.4396206	0.0932766	4.7130849	0.000002

```

Alpha_1  0.3351672  0.1093611  3.0647750  0.001190
Gamma    0.2152178  0.1535592  1.4015295  0.081052
Beta_1   0.3394524  0.0895886  3.7890126  0.000092

```

```

*****
Wald Confidence Limits (Alpha=0.05)
*****

```

Parameter	Estimate	Lower	Upper
Intercept	0.4396206	0.2568018	0.6224394
Alpha_1	0.3351672	0.1208234	0.5495110
Gamma	0.2152178	-0.0857528	0.5161883
Beta_1	0.3394524	0.1638619	0.5150429

Factor sigma=1 (nobs=300 df=4 vard=300)

#### 9. Method 6: NAGARCH:

```

m = .; q = p = 1;
print "Method 6: NAGARCH: Reparametrized Constraints";
optn = [ "cent"
        "tech"      "tr"
        "error"     "norm"
        "print"     3 ];
< gof,est,ase,cis,cov,scor > = garch(6,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 6: NAGARCH Modeling: NQ=1 NP=1
*****
Q[1] = 1 P[1] = 1

```

Number of Observations . . . . .	300
Number of Parameter Estimates . . . . .	4
Mean Y . . . . .	0.0016
Standard Deviation Y . . . . .	1.1442
Mean Squared Error (MSE) . . . . .	1.3093
Number of Back Casts . . . . .	1
Error Distribution . . . . .	Normal
Optimization Technique . . . . .	TRUREG

\*\*\*\*\*  
 Optimization Start  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	-0.80893295	-0.0156817
2 Alpha_1	-0.65145797	-8.63e-004
3 Gamma	0.00000000	-6.3343419
4 Beta_1	0.08284350	-0.0055090

Value of Objective Function = 455.504

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	60.094788	2.1426024	3.5418457	24.498126
2	2.1426024	4.3231317	-1.1918934	1.3178576
3	3.5418457	-1.1918934	26.798373	2.0994381
4	24.498126	1.3178576	2.0994381	12.972177

Trust Region Optimization  
 Without Parameter Scaling  
 Gradient Computed by Finite Differences  
 Hessian Computed by Finite Differences (dense)  
 (Using Only Function Calls)

Iteration Start:

N. Variables	4		
Criterion	455.5035429	Max Grad Entry	6.334341932
TR Radius	1.000000000		

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1	0	2	0	454.7537	0.749798	0.42360	0.00000	1.00000
2	0	3	0	454.7329	0.020876	0.01323	0.00000	0.26611
3	0	4	0	454.7329	3.4e-006	8e-006	0.00000	0.09923



Successful Termination After 3 Iterations  
 GCONV convergence criterion satisfied.

Criterion	454.7328650	Max Grad Entry	8.2550e-006
Ridge (lambda)	0.000000000	TR Radius	0.099227568
Act.dF/Pred.dF	0.999981427		
N. Function Calls	5	N. Gradient Calls	7
N. Hessian Calls	5	Preproces. Time	0
Time for Method	0	Effective Time	0

\*\*\*\*\*  
 Optimization Results  
 \*\*\*\*\*

Parameter Estimates  
 -----

Parameter	Estimate	Gradient
1 Intercept	-0.79882912	3.31e-007
2 Alpha_1	-0.66456439	-3.95e-006
3 Gamma	0.25256726	-8.25e-006
4 Beta_1	0.04328314	1.30e-006

Value of Objective Function = 454.733

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
1	62.232247	2.7850828	-0.2077795	25.072832
2	2.7850828	4.6852822	1.7877625	1.8948846
3	-0.2077795	1.7877625	21.513094	0.2034453
4	25.072832	1.8948846	0.2034453	13.092182

Log Likelihood= -454.733  
 SBC=-886.651 AIC=-901.466

Covariance Matrix 2: H = (NOBS/d) inv(G)  
 \*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	0.0144757			
Alpha_1	0.0012657	0.0122713		
Gamma	3.14e-004	-0.0068194	0.0290811	
Beta_1	-0.0110178	-0.0081967	0.0017286	0.0149454

Factor sigma=1 (nobs=300 df=4 vard=300)

Determinant = 9.23947e-009

Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	0.0088774			
Alpha_1	-0.0010126	0.0120685		
Gamma	6.46e-004	-0.0073773	0.0258309	
Beta_1	-0.0051584	-0.0060528	0.0024665	0.0083331

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Alpha_1	Gamma	Beta_1
Intercept	1.0000000			
Alpha_1	-0.0978309	1.0000000		
Gamma	0.0426414	-0.4178324	1.0000000	
Beta_1	-0.5997496	-0.6035727	0.1681129	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.4498554	0.0942199	4.7745255	0.000001

```

Alpha_1  0.3267569  0.1098565  2.9743965  0.001589
Gamma    0.1986172  0.1607199  1.2357977  0.108757
Beta_1   0.3372187  0.0912858  3.6940963  0.000131

```

```

*****
Wald Confidence Limits (Alpha=0.05)
*****

```

```

Parameter  Estimate      Lower      Upper
Intercept  0.4498554  0.2651877  0.6345230
Alpha_1    0.3267569  0.1114420  0.5420718
Gamma      0.1986172 -0.1163879  0.5136224
Beta_1     0.3372187  0.1583017  0.5161357
Factor sigma=1 (nobs=300 df=4 vard=300)

```

10. Method 7: IGARCH use  $\delta = 2$ :

```

m = .; q = p = 1;
print "Method 7: IGARCH: Reparametrized Constraints: Delta=2";
optn = [ "cent"
         "delta"      2. ,
         "tech"       "tr" ,
         "error"      "norm" ,
         "print"      3 ];
< gof,est,ase,cis,cov,scor > = garch(7,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 7: IGARCH Modeling: NQ=1 NP=1
*****
Modeling Squared Values (Delta=2)
Q[1] = 1 P[1] = 1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 2
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal

```

Optimization Technique . . . . . TRUREG

\*\*\*\*\*  
Optimization Start  
\*\*\*\*\*

Parameter Estimates  
-----

Parameter	Estimate	Gradient
1 Intercept	-4.33236311	-0.1900006
2 Alpha_1	-2.19722238	-1.5994832

Value of Objective Function = 466.563

Hessian Matrix  
\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2
1	1.7309304	-3.4318452
2	-3.4318452	5.0825081

Trust Region Optimization  
Without Parameter Scaling  
Gradient Computed by Finite Differences  
Hessian Computed by Finite Differences (dense)  
(Using Only Function Calls)

Iteration Start:  
N. Variables 2  
Criterion 466.5625441 Max Grad Entry 1.599483213  
TR Radius 1.000000000

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1*	0	4	0	465.4522	1.110313	1.56396	0.65221	4.01012
2*	0	5	0	464.5735	0.878697	10.6363	1.09837	2.00506
3	0	6	0	460.5322	4.041335	5.78896	0.00000	2.00548
4	0	7	0	460.1095	0.422739	0.11949	0.00000	1.66177

```

5 0 8 0 460.1076 1.9e-003 6e-003 0.00000 0.14209
6 0 9 0 460.1076 6.1e-007 2e-006 0.00000 0.03490

```

```

Successful Termination After      6 Iterations
GCONV convergence criterion satisfied.
Criterion      460.1076084      Max Grad Entry  2.1990e-006
Ridge (lambda) 0.000000000      TR Radius      0.034904588
Act.dF/Pred.dF 0.999530876
N. Function Calls      10      N. Gradient Calls      3
N. Hessian Calls      8      Preproces. Time      0
Time for Method      0      Effective Time      0

```

```

*****
Optimization Results
*****

```

```

Parameter Estimates
-----

```

Parameter	Estimate	Gradient
1 Intercept	-1.13300176	7.31e-007
2 Alpha_1	0.74808044	-2.20e-006

```

Value of Objective Function =      460.108

```

```

Hessian Matrix
*****

```

```

Symmetric Matrix: Dense Storage

```

S	1	2
1	38.015479	-17.841882
2	-17.841882	11.906938

```

Log Likelihood= -460.108
SBC=-908.808 AIC=-916.215

```

```

Covariance Matrix 2: H = (NOBS/d) inv(G)
*****

```

```

Dense Symmetric Matrix (2 by 2)

```

	S	Intercept	Alpha_1
Intercept		0.0092191	
Alpha_1		0.0093608	0.0135003

Factor sigma=1 (nobs=300 df=2 vard=300)  
 Determinant = 3.68368e-005  
 Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates  
 \*\*\*\*\*

Dense Symmetric Matrix (2 by 2)

	S	Intercept	Alpha_1
Intercept		0.0074443	
Alpha_1		0.0070108	0.0101716

Robust Asymptotic Correlation Matrix of Estimates  
 \*\*\*\*\*

Dense Symmetric Matrix (2 by 2)

	S	Intercept	Alpha_1
Intercept		1.0000000	
Alpha_1		0.8056720	1.0000000

\*\*\*\*\*  
 Parameter Estimates and Asympt. Standard Errors  
 \*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.3220650	0.0862806	3.7327625	0.000113
Alpha_1	0.6786245	0.1008545	6.7287453	0.000000

\*\*\*\*\*  
 Wald Confidence Limits (Alpha=0.05)  
 \*\*\*\*\*

Parameter	Estimate	Lower	Upper
-----------	----------	-------	-------

```

Intercept  0.3220650  0.1529581  0.4911719
Alpha_1    0.6786245  0.4809533  0.8762958
Factor sigma=1 (nobs=300 df=2 vard=300)

```

11. Method 7: IGARCH use  $\delta = 1$ :

```

m = .; q = p = 1;
print "Method 7: IGARCH: Reparametrized Constraints: Delta=1";
optn = [ "cent"          ,
        "delta"         1. ,
        "tech"          "tr" ,
        "error"         "norm" ,
        "print"         3 ];
< gof,est,ase,cis,cov,scor > = garch(7,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 7: IGARCH Modeling: NQ=1 NP=1
*****
Modeling Absolute Values (Delta=1)
      Q[1] =      1 P[1] =      1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 2
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Back Casts . . . . . 1
Error Distribution . . . . . Normal
Optimization Technique . . . . . TRUREG

```

```

*****
Optimization Start
*****

```

Parameter Estimates  
-----

Parameter	Estimate	Gradient
-----------	----------	----------

1 Intercept-4.93814434 -8.5560129  
 2 Alpha\_1 -2.94443689 8.2633134

Value of Objective Function = 470.207

Hessian Matrix  
 \*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2
1	5.4931025	-8.2545947
2	-8.2545947	10.430961

Trust Region Optimization  
 Without Parameter Scaling  
 Gradient Computed by Finite Differences  
 Hessian Computed by Finite Differences (dense)  
 (Using Only Function Calls)

Iteration Start:

N. Variables	2		
Criterion	470.2067871	Max Grad Entry	8.556012903
TR Radius	1.000000000		

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1*	0	3	0	467.5957	2.611061	4.07006	9.42808	0.49607
2*	0	7	0	465.6413	1.954411	1.35533	0.32023	5.91678
3	0	9	0	464.8224	0.818955	0.33783	0.94923	0.85436
4	0	10	0	464.7775	0.044887	0.07254	0.00000	0.90206
5	0	11	0	464.7774	1.2e-004	4e-004	0.00000	0.25776
6	0	12	0	464.7774	1.7e-009	3e-008	0.00000	0.01028

Successful Termination After 6 Iterations

GCONV convergence criterion satisfied.

Criterion	464.7773539	Max Grad Entry	2.6987e-008
Ridge (lambda)	0.000000000	TR Radius	0.010284096
Act.dF/Pred.dF	0.999686729		
N. Function Calls	13	N. Gradient Calls	3
N. Hessian Calls	8	Preproces. Time	0
Time for Method	0	Effective Time	0



\*\*\*\*\*  
Optimization Results  
\*\*\*\*\*

Parameter Estimates  
-----

Parameter	Estimate	Gradient
1 Intercept	-2.69415957	-8.89e-009
2 Alpha_1	-1.26091968	-2.70e-008

Value of Objective Function = 464.777

Hessian Matrix  
\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2
1	48.759932	-46.781661
2	-46.781661	47.834734

Log Likelihood= -464.777  
SBC=-918.147 AIC=-925.555

Covariance Matrix 2: H = (NOBS/d) inv(G)  
\*\*\*\*\*

Dense Symmetric Matrix (2 by 2)

S	Intercept	Alpha_1
Intercept	0.0015155	
Alpha_1	0.0037710	0.0100023

Factor sigma=1 (nobs=300 df=2 vard=300)  
Determinant = 9.37374e-007  
Matrix has Only Positive Eigenvalues

Robust Asymptotic Covariance Matrix of Estimates  
\*\*\*\*\*

Dense Symmetric Matrix (2 by 2)

```
      S |   Intercept   Alpha_1
-----
Intercept |   0.0071639
Alpha_1   |   0.0190380   0.0512523
```

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (2 by 2)

```
      S |   Intercept   Alpha_1
-----
Intercept |   1.0000000
Alpha_1   |   0.9935531   1.0000000
```

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.0675992	0.0846398	0.7986692	0.212559
Alpha_1	0.2207715	0.2263898	0.9751828	0.165130

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.0675992	-0.0982917	0.2334901
Alpha_1	0.2207715	-0.2229444	0.6644873

Factor sigma=1 (nobs=300 df=2 vard=300)

## 12. Method 8: FIGARCH: Truncated Lags=100

```
m = .; q = p = 1;
print "Method 8: FIGARCH: Reparametrized Constraints: TRUNCL=100";
optn = [ "cent"
        "truncl"    100 ,
        "tech"      "tr" ,
        "error"     "norm" ,
        "print"     3 ];
```

```

< gof,est,ase,cis,cov,scor > = garch(8,two,.,m,q,p,optn);
print "GOF=",gof;
print "EST=",est;
print "ASE=",ase;
print "CIs=",cis;
print "COV=",cov;

```

```

*****
Model 8: FIGARCH Modeling: NQ=1 NP=1
*****
Q[1] = 1 P[1] = 1

```

```

Number of Observations . . . . . 300
Number of Parameter Estimates . . . . . 4
Mean Y . . . . . 0.0016
Standard Deviation Y . . . . . 1.1442
Mean Squared Error (MSE) . . . . . 1.3093
Number of Weights in ARCH(oo) . . . . . 100
Error Distribution . . . . . Normal
Optimization Technique . . . . . TRUREG

```

```

*****
Optimization Start
*****

```

Parameter Estimates  
-----

Parameter	Estimate	Gradient
1 Intercept	-0.88588323	0.3649173
2 Phi	1.38629436	-0.4245310
3 D	-1.38629436	0.2307636
4 Beta	0.00000000	1.1283417

Value of Objective Function = 457.638

```

Hessian Matrix
*****

```

Symmetric Matrix: Dense Storage

```

S | 1 2 3 4
-----

```

```

1 | 19.991728 0.3046813 10.991260 -0.7993952
2 | 0.3046813 0.2911064 0.5745450 -0.2276935
3 | 10.991260 0.5745450 7.0245531 -0.1297626
4 | -0.7993952 -0.2276935 -0.1297626 2.8163192

```

Trust Region Optimization  
Without Parameter Scaling  
Gradient Computed by Finite Differences  
Hessian Computed by Finite Differences (dense)  
(Using Only Function Calls)

Iteration Start:

```

N. Variables          4
Criterion             457.6376615      Max Grad Entry  1.128341659
TR Radius             1.000000000

```

Iter	rest	nfun	act	optcrit	difcrit	maxgrad	lambda	radius
1	0	2	0	457.1540	0.483708	0.21628	0.22278	1.00000
2	0	5	0	456.9815	0.172470	0.51927	0.01525	2.30785
3	0	6	0	456.8819	0.099625	0.15624	0.03489	1.00857
4	0	7	0	456.8402	0.041700	0.15303	8e-003	1.01691
5	0	8	0	456.8237	0.016415	9e-003	0.00000	1.01747
6	0	9	0	456.8183	5.5e-003	7e-003	0.00000	1.05950
7	0	10	0	456.8162	2.1e-003	4e-003	0.00000	0.98833
8	0	11	0	456.8155	7.6e-004	2e-003	2e-004	0.99528
9	0	12	0	456.8152	2.8e-004	4e-004	3e-004	0.99560
10	0	13	0	456.8151	4.6e-005	1e-004	0.00000	0.99560
11	0	14	0	456.8151	4.6e-005	1e-004	0.00000	1.09166
12	0	15	0	456.8151	3.6e-005	2e-004	0.00000	1.16833
13	0	16	0	456.8151	5.8e-006	4e-005	0.00000	0.66855

Successful Termination After 13 Iterations

GCONV convergence criterion satisfied.

```

Criterion             456.8150627      Max Grad Entry  3.8784e-005
Ridge (lambda)       0.000000000      TR Radius       0.668550915
Act.dF/Pred.dF      1.832806390
N. Function Calls    17      N. Gradient Calls  3
N. Hessian Calls     15      Preproces. Time    0
Time for Method      1      Effective Time     1

```

\*\*\*\*\*  
Optimization Results  
\*\*\*\*\*

Parameter Estimates

-----

Parameter	Estimate	Gradient
1 Intercept	-0.46297451	-3.88e-005
2 Phi	11.8851377	-3.17e-005
3 D	-2.60537231	-3.74e-007
4 Beta	-0.20489682	3.46e-005

Value of Objective Function = 456.815

Hessian Matrix

\*\*\*\*\*

Symmetric Matrix: Dense Storage

S	1	2	3	4
-----				
1	47.152502	6.17e-005	11.418389	-2.6275098
2	6.17e-005	1.26e-004	7.51e-005	7.49e-005
3	11.418389	7.51e-005	2.8779569	-0.5813992
4	-2.6275098	7.49e-005	-0.5813992	3.0395612

Log Likelihood= -456.815

SBC=-890.815 AIC=-905.63

Covariance Matrix 2:  $H = (NOBS/d) \text{inv}(G)$

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Phi	D	Beta
-----				
Intercept	-0.0345733			
Phi	-0.0806959	-0.0525482		
D	0.0294820	0.0394115	-0.0175934	
Beta	-0.0364651	-0.0143724	0.0169448	0.0052774

Factor sigma=1 (nobs=300 df=4 vard=300)

Determinant = -3.18551e-008

Matrix has 1 Negative Eigenvalue(s)

Robust Asymptotic Covariance Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Phi	D	Beta
Intercept	0.1283577			
Phi	0.0360338	0.0295502		
D	-0.0518958	-0.0182783	0.0226552	
Beta	0.0053435	0.0109712	-0.0034809	0.0075612

Robust Asymptotic Correlation Matrix of Estimates

\*\*\*\*\*

Dense Symmetric Matrix (4 by 4)

S	Intercept	Phi	D	Beta
Intercept	1.0000000			
Phi	0.5850856	1.0000000		
D	-0.9623608	-0.7064356	1.0000000	
Beta	0.1715219	0.7339742	-0.2659561	1.0000000

\*\*\*\*\*

Parameter Estimates and Asympt. Standard Errors

\*\*\*\*\*

Parameter	Estimate	AStdErr	T Value	P Value
Intercept	0.6294087	0.3582705	1.7567975	0.039993
Phi	0.4656001	0.1719017	2.7085257	0.003576
D	0.0687935	0.1505163	0.4570499	0.323985
Beta	0.2399182	0.0869549	2.7591106	0.003078

\*\*\*\*\*

Wald Confidence Limits (Alpha=0.05)

\*\*\*\*\*

Parameter	Estimate	Lower	Upper
Intercept	0.6294087	-0.0727886	1.3316059
Phi	0.4656001	0.1286790	0.8025211
D	0.0687935	-0.2262131	0.3638000
Beta	0.2399182	0.0694897	0.4103468

Factor sigma=1 (nobs=300 df=4 vard=300)

## 5.7 Function jarbera

---

```
prob = jarbera(y<,optn>)
```

```
< prob,stat > = jarbera(y<,optn>)
```

**Purpose:** Implements the Jarque-Bera test for normally distributed data.

**Input: y** The first input object  $y$  must be either a  $n$  vector or a  $n \times nc$  matrix. If  $y$  is a vector the scalar value of the probability of the CDF is returned, if  $y$  is a matrix then a vector of  $nc$  probabilities is returned by using each column of  $y$  separately.

**optn** this is a vector of options which should be initialized with missing values for defaults:

1. (int) amount of printed output (default is no output);
2. (real) probability for test (default is 0.05), must be in  $[0, 1]$ ;
3. (int) number of dependent variables (default=2).

**Output: prob** The first result is either a scalar or an object of the test probabilities.

**stat** The **stat** result has the same size as **prob** and returns the test statistics (only for `optn[3]=0`).

**Restrictions:** 1. Missing value are returned if the input  $y$  contains missing data.

2.

**Relationships:** `berkow()`, `ksprob()`, `kstest()`, `shapwilk()`

**Examples:** 1. Test  $N = 1000$  random uniform generated data:

```
options NOECHO;
#include "..\\tdata\\matlb.dat"
options ECHO;

optn = [ 2 , /* print */
        .05 , /* alpha */
        2 ]; /* nvar */
< p1,s1 > = jarbera(uf1000,optn);
print "JARBERA P:",p1;
print "JARBERA S:",s1;
```

```
Hypothesis H0: Data are uniformly distributed versus
Hypothesis H1: Data are not uniformly distributed
H0 rejected: Probability=7.38734e-014 Statistics=60.4728
```

2. Test  $N = 1000$  random  $(0, 1)$ -normal generated data:

```
nr = 500; nc = 5;
print "\n *** nr,nc=",nr,nc," ***\n";
options RANDUNI;
mu1 = rand(nr,nc,'g',"norm");
```

Hypothesis H0: Data are uniformly distributed versus  
Hypothesis H1: Data are not uniformly distributed

Col	Probability	Statistics	Rej/Acc
1	0.870686046	0.276947638	Accept
2	0.575339062	1.105591478	Accept
3	0.732156146	0.623522948	Accept
4	0.745872441	0.586401367	Accept
5	0.677028247	0.780084567	Accept



## 5.8 Function mucomp

---

```
< gof,BF > = mucomp(data,nres,modl<,optn>)
```

```
< gof,lik,est,fpbar,BF > = mucomp(data,modres,order,rest<,optn>)
```

**Purpose:** The `mucomp` function implements three algorithms for comparison of means (confirmatory ANOVA) as it is available in the Fortran program by Kuiper, Klugkist, & Hojtink (2009). The data set is usually a vector  $x$  of  $N$  observations grouped into  $g$  groups. The common 1-way ANOVA model,

$$y_{ij} = \mu_i d_{ij} + \epsilon_{ij}, \quad i = 1, \dots, g, j = 1, \dots, N_i$$

where  $d_{ij}$  is observation  $j$  belonging to group  $i$ , and the error is normal distributed  $\epsilon_{ij} \sim \mathcal{N}(0, \sigma^2)$ , tests the null hypothesis that all  $g$  means are equal  $\mu_1 = \dots = \mu_g$  against the alternative that the means are not equal:  $\mu_1, \dots, \mu_g$ . This method however, was extended to permit a wider variety of order relations among the means, e.g. testing models of the form:

$$\mu_5 = \mu_3 > (\mu_1, \mu_4) < \mu_2$$

As in the Fortran program by Kuiper, Klugkist, & Hojtink three different algorithms are implemented:

**Fbar** an algorithm for estimating the **Fbar** and **pFbar** criterion as shown in Silvapulle & Sen (2005).

**ORIC** an algorithm for computing the order restricted information criterion ORIC proposed by Anraku (1999).

**BMS** an algorithm for Bayesian model selection, see Klugkist, Laudy, & Hoijtink (2005).

For the **Fbar** and **ORIC** algorithms a set of optimal estimates (means satisfying the constraints) and log likelihood values for each model is computed. The **ORIC** and **BMS** algorithms permit to rank the models.

The method can be extended to deal with 2 or more way classification problems.

**Input: data** The data matrix consists of either one or two columns.

**2 column data** : One of the two columns (specified by an option) must contain an integer ID variable indicating the group membership. This ID variable must not have consecutive values. Groups are specified by identical ID values.

**1 column data** : Here we assume the total number of observations  $N$  is a multiple of the number  $g$  of groups. For the group membership the data must then be arranged by one of the orders  
:

**block** blockwise: [ 1 1 1 2 2 2 3 3 3 ],  
**split** splitwise: [ 1 2 3 1 2 3 1 2 3 ],  
**random** random: [ 3 1 2 2 1 3 3 2 1 ].

The user may specify in which order the data are arranged.

**nres** This must be a  $m$  vector  $nr$  of positive integers specifying how many restrictions are specified for each of the models  $k = 1, \dots, m$  with the third input argument. Let  $NR = \sum_k^m nr_k$  be the total number of restrictions for all models.

**modl** There are two different forms of the specification of restrictions on the means  $\mu_j$ , the *tensor* and the easier *string* specification:

**tensor** The tensor `tmod[2,NR,g]` consists of two matrices of integers and is the same as the matrix form developed and described in detail by Kuiper, Klugkist, & Hoijtink (2009).

1. The first  $NR \times g$  matrix of integers specifies the order of groups in the  $NR$  restrictions. The rows of this matrix have to ordered as specified by the `modres` input argument.
2. The second  $NR \times g$  matrix of integers specifies the  $NR$  model restrictions.

The meaning of the integers in the restriction matrix is the following:

- -1: defines the  $j$  relationship;
- -3: defines the  $i$  relationship;
- 0: defines a singleton (unused by the restriction);
- positive integers: specify the group number of equality constrained clusters of variables.

**string** This must be a vector of  $NR$  strings where each of the strings formulates restrictions in a clearly readable form.

Even though the string seems to be much nicer for applications, the matrix or tensor specification may have some use when the input arguments are created by some other software rather than writing them down by the user. See the example below for more details.

**optn** The `optn` argument is specified in form of a two column matrix where the first column defines the option as string value (in quotes) and the second column can be used for a numeric or string specification of the option. See table below for content. Many of the options have default values and must not be specified for many applications. However, using the string input of the model requires to specify the number  $g$  of groups, which must not necessarily be determined with the model input.

Option	Second Column	Meaning
"block"		block sorted 1-column input data
"bms"		perform BMS algorithm (not default)
"delta"	real	only for BMS: $\Delta \geq 0$ , default=0
"fbar"		perform Fbar algorithm (is default)
"group"		specifies number $g$ of groups (string model)
"idvar"	int	column number containing ID variable default is first column
"nopr"		suppress all printed output
"oric"		perform ORIC algorithm (not default)
"nsamp"	int	number of random samples for Fbar and ORIC default is $n_s = 100,000$
"pinp"		print the model specification in matrix form
"print"	int	amount of printed output =0 specifies no printed output, default is 1
"pv"	real	only for BMS: prior vagueness, $p_v > 0$ , default=2
"qpnuSP"		perform null space QP algorithm
"qpPogi"		perform Powell's Goldfarb and Idnani QP algorithm
"rand"		random sorted 1-column input data
"split"		split sorted 1-column input data
"seed"	int	seed for random generator, default=time of day
"sing"	real	singularity criterion, default=1.e-8

**Output:** **gof** a vector of scalar results, see below for content;

**lik** return only for the Fbar and ORIC methods:

- for Fbar method: the  $m \times 2$  matrix containing the log likelihood values and the residual sum of squares for all models in its two columns;
- for ORIC method:  $m \times 4$  matrix with an additional two matrices containing the penalty term and the ORIC value (which is the sum of log likelihood and penalty term). The value of the penalty term is in  $[2, g+1]$ , the lower value for the null hypothesis (all means are equal) and  $g + 1$  for the unconstrained solution.

**est** return only for Fbar and ORIC: this is the  $m \times g$  matrix of parameter estimates for the  $m$  models which should satisfy the specified restrictions.

**fpbar** only for Fbar: contains the Fbar and corresponding  $p$  values for the hypotheses.

**BF** only for BMS: this is the  $m \times 2$  matrix of Bayes factors  $BF_k$  in its first column and the *posterior model probabilities*  $pmp_k$ ,  $k = 1, \dots, m$ , which follow directly from the Bayes factors:

$$pmp_k = BF_k / \sum_k^M BF_k \quad , \quad k = 1, \dots, m$$

That means  $0 \leq pmp_k \leq 1$  and the model with the highest  $pmp_k$  value should be the most fitting w.r.t. to the data.

- Restrictions:**
1. The input data must be numeric and cannot contain any string or complex data.
  2. Currently the input data must not have any missing values.

**Relationships:** `anova()`, `hbanova()`

**Examples:** 1. Example by Lucas (2003): Tensor Input:

```
data = [ 1  3.58, 1 -0.15, 1  0.67, 1  2.22, 1  2.56, 1  1.7 , 1 -0.45, 1
         1  4.83, 1  2.44, 1  6.74, 1  1.4 , 1 -0.03, 1  0.71, 1  4.3 , 1
         1  2.44, 1  0.81, 1  3.65, 1  1.41, 1 -0.38, 1  3.11, 1  0.21, 1
         1  2.63, 1  4.74, 1  2.12, 1  4.36, 1  3.26, 1  0.94, 2  1.67, 2
         .....
         5  1.38, 5  4.58, 5  3.48, 5  0.39, 5  3.89, 5  4.18, 5  3.72, 5
         5  3.63, 5  1.55, 5 -0.46, 5  2.04, 5  2.26, 5  4.27, 5  5.67, 5
         5  2.17, 5  2.27, 5  4.1 , 5  3.74, 5  4.54, 5  1.71, 5  4.74, 5
         5  4.1 , 5  4.8 , 5  2.9 , 5  2.35, 5  4.09, 5  1.09 ];
```

The model specification contains one restriction row for all models (hypotheses) except for the second which takes two rows:

```
nres = [ 1  2  1  1 ];
nmod = dim(nres);
```

The following is the string model specification:

```
print "String Model Specification without Singletons";
modl = [ "1 = 2 = 3 = 4 = 5",
         "5 = 3 > 1 > 2 ",
         "3 > 4 > 2 ",
         "3 > 1 > 4 = 5 > 2",
         " " ];
```

The following string specification is equivalent but closer to the matrix specification used in Kuiper, Klugkist, and Hojtink (2009);

```
print "Model specification with Singletons";
modl = [ "1 = 2 = 3 = 4 = 5",
         "5 = 3 > 1 > 2,  4",
         "3 > 4 > 2,  1,  5",
         "3 > 1 > 4 = 5 > 2",
         "1 , 2 , 3 , 4 , 5" ];
```

The group number must be specified with the string model specification. The default sample size for the simulation with Fbar ( $p$  value) and ORIC is 100,000 but is here specified to only 1000. The "seed" specification ensures replicable results.

```
print "Fbar, ORIC, and BMS together";
optns = [ "bms"           ,
          "fbar"         ,
          "oric"         ,
          "group"        5 ,
          "samp"         1000 ,
          "seed"         123 ,
          "delt"         0.3 ,
          "pv"           2.0 ,
          "pinput"       ,
          "print"        2 ];
```

The function starts with the output of some data properties, where the means are the free (unconstrained) parameter estimates (as with specified model 4):

```
*****
Summary of Input Data
*****
```

Group	Nobs	Mean	StdDev
1	30	2.32933333	1.86012223
2	30	1.32800000	1.14947484
3	30	3.20000000	1.78986418
4	30	2.23066667	1.45044805
5	30	3.22900000	1.50012264

The following is the output generated for the pinp option:

```
*****
Ordering and Restrictions for 4 Models and 5 Groups
*****
```

Model NR Orders and Restrictions

1	1	1	2	3	4	5
		1	1	1	1	1
2	1	5	3	1	2	4
		1	1	-3	-3	0
	2	3	4	2	1	5

```

      1 -3 -3 0 0
3  1  3  1  4  5  2
      1 -3 -3  3 -3
4  1  1  2  3  4  5
      0  0  0  0  0

```

With Fbar and ORIC the optimal (constrained) estimates are computed:

Estimated Coefficients

\*\*\*\*\*

Dense Matrix (4 by 5)

	Group1	Group2	Group3	Group4	Group5
Model1	2.4634000	2.4634000	2.4634000	2.4634000	2.4634000
Model2	2.3293333	1.3280000	3.2145000	2.2306667	3.2145000
Model3	2.5963333	1.3280000	3.2000000	2.5963333	2.5963333
Model4	2.3293333	1.3280000	3.2000000	2.2306667	3.2290000

The program

Fbar Test for Null Hypothesis 1 and Unconstrained Hypothesis 4  
Hypothesis 1 versus 4: Fbar= 30.27 p-value= 0.0000

\*\*\*\*\*  
"Ordered Alternative" Fbar Test  
\*\*\*\*\*

Hypothesis	Fbar	p-value
H0 vs 2	30.2643855	0.00000000
H0 vs 3	22.9115843	0.00000000

\*\*\*\*\*  
"Ordered Null" Fbar Test  
\*\*\*\*\*

Hypothesis	Fbar	p-value
2 vs Ha	0.00511179	0.98600000
3 vs Ha	7.35791301	0.05600000

Fbar and Pvalues for Null and Alt Models

\*\*\*\*\*

Dense Matrix (4 by 5)

	Model1	Model2	Model3	Model4
H0_FbarProb	.	0.0051118	7.3579130	.
H0_Pval	.	0.9860000	0.0560000	.
Ha_Fbar	.	30.264385	22.911584	.
Ha_Pval	.	0.0000000	0.0000000	.

H0\_vs\_Hator

H0_FbarProb	.
H0_Pval	.
Ha_Fbar	30.269497
Ha_Pval	0.0000000

After Fbar the pecified ORIC is computed. The model with the smallest ORIC value is the most preferred:

\*\*\*\*\*  
 The value of the Order-Restricted Information Criterion  
 $ORIC = -2 * \log \text{likelihood} + 2 * \text{penalty}$   
 \*\*\*\*\*

Model	LogLik	Penalty	ORIC
1	-292.267801	2.00000000	588.535603
2	-278.051118	3.17300000	562.448236
3	-281.760877	3.15300000	569.827753
4	-278.048474	6.00000000	568.096948

The preferred hypothesis, according to the Order-Restricted Information Criterion of the hypotheses be compared is hypothesis number 2 with the following ordering(s) of means:

5	3	1	2	4
3	4	2	1	5
Corresponding Order of Restrictions:				
1	1	-3	-3	0
1	-3	-3	0	0

After ORIC the pecified BMS is computed. Here is it necessary to specify the `delta` and `pv` options which may have a huge effect on the computer time needed for the bayes estimation.

\*\*\*\*\*

The resulting Bayes factor (BF) values (of the order-restricted hypotheses versus the unconstrained hypothesis) and the posterior model probabilities (PMP) of the order-restricted hypotheses with respect to the whole set of hypotheses:

Hypothesis	BF	PMP
1	0.02643567	5.797e-004
2	39.5893012	0.86820503
3	4.98328638	0.10928494
4	1.00000000	0.02193029

\*\*\*\*\*

The preferred hypothesis, according to confirmatory Bayesian model selection of the hypotheses to be compared is hypothesis number 2 with the following ordering(s) of means:

3	4	2	1	5
3	1	4	5	2

Corresponding Order of Restrictions:

1	-3	-3	0	0
1	-3	-3	3	-3

\*\*\*\*\*

Specification of the Encompassing Prior:

For all means, the same normal prior with mean 2.28037 and variance 2.3046 is used.

For the residual variance, a scaled inverse chi-square, with degrees of freedom 1 and scale parameter 2.50256 is used.

The following is the printed output of the return arguments:

```
< gof,lik,est,fpbar,BF > = mucomp(data,nres,ord,rest,optns);
print "GOF=",gof;
print "LIK=",lik;
print "EST=",est;
print "FPBAR=",fpbar;
print "BF=",BF;
```



GOF=

	g-of-fit
Error	0.000000
Mu_Total	2.46340
SS_Total	357.835
Time_Tot	64.0000
Time_Opt	18.0000
Iter_Tot	17439.0
unused	.
unused	.
unused	.
unused	.

LIK=

	LogLik	ResiSS	PenTrm	ORIC
Model1	-292.27	432.53	2.0000	588.54
Model2	-278.05	357.85	3.1730	562.45
Model3	-281.76	375.99	3.1530	569.83
Model4	-278.05	357.83	6.0000	568.10

EST=

	Group1	Group2	Group3	Group4	Group5
Model1	2.4634	2.4634	2.4634	2.4634	2.4634
Model2	2.3293	1.3280	3.2145	2.2307	3.2145
Model3	2.5963	1.3280	3.2000	2.5963	2.5963
Model4	2.3293	1.3280	3.2000	2.2307	3.2290

FPBAR=

	Model1	Model2	Model3
HO_FbarProb	.	0.0051	7.3579
HO_Pval	.	0.98600	0.0560
Ha_Fbar	.	30.264	22.912
Ha_Pval	.	0.00000	0.00000

	Model4	HO_vs_Hator
HO_FbarProb	.	.
HO_Pval	.	.
Ha_Fbar	.	30.269

```
Ha_Pval | . 0.00000
```

```
BF=
      | BayesFactor PostModProb
-----|-----
Model1 | 0.0264 0.0006
Model2 | 39.589 0.86821
Model3 | 4.9833 0.10928
Model4 | 1.00000 0.0219
```

## 2. Example by Lucas (2003): Tensor Input:

The following specification is equivalent with that of the Kuiper, Klugkist, & Hojtink (2009) program: The first matrix specifies the order of variables in the restrictions (second matrix):

```
ord = [ 1 2 3 4 5 ,
        5 3 1 2 4 ,
        3 4 2 1 5 ,
        3 1 4 5 2 ,
        1 2 3 4 5 ];
print "Ordering of means in restriction",ord;
```

The second matrix specifies the restrictions of the four models:

```
rest = [ 1 1 1 1 1 ,
         1 1 -3 -3 0 ,
         1 -3 -3 0 0 ,
         1 -3 -3 3 -3 ,
         0 0 0 0 0 ];
print "(Order) Restrictions",rest;
```

The meaning of the integers in the restriction matrix is the following:

- -1: defines the  $i$  relationship;
- -3: defines the  $j$  relationship;
- 0: defines a singleton (unused by the restriction);
- positive integers: specify the group number of equality constrained clusters of variables.

The two matrices are merged in the tensor `tmodl`:

```
int tmodl[2,5,5];
tmodl[1,,] = ord; tmodl[2,,] = rest;
print "Tensor=",tmodl;
```

The results are identical to those obtained with the string specification of the model.

## 5.9 Function shapwilk

---

```
prob = shapwilk(y<,optn>)
```

```
< prob,stat > = shapwilk(y<,optn>)
```

**Purpose:** Implements the Shapiro-Wilks and the Shapiro-Francia tests for normally distributed data.

**Input:** **y** The first input object *y* must be either a *n* vector or a  $n \times nc$  matrix. If *y* is a vector the scalar value of the probability of the CDF is returned, if *y* is a matrix then a vector of *nc* probabilities is returned by using each column of *y* separately.

**optn** this is a vector of options which should be initialized with missing values for defaults:

1. (int) amount of printed output (default is no output);
2. (real) probability for test (default is 0.05), must be in [0, 1];
3. (int) version: =0: Shapiro-Wilks test (default); =1: Shapiro-Francia test.

**Output:** **prob** The first result is either a scalar or an object of the test probabilities.

**stat** The **stat** result has the same size as **prob** and returns the test statistics (only for `optn[3]=0`).

**Restrictions:** 1. Missing value are returned if the input *y* contains missing data.

2.

**Relationships:** `berkow()`, `jarbera()`, `ksprob()`, `kstest()`

**Examples:** 1. Shapiro-Wilks Test:  $N = 1000$

```
options NOECHO;
#include "..\\tdata\\matlb.dat"
options ECHO;

print "Shapiro-Wilks test for Normality";
optn = [ 2 , /* print */
        .05 , /* alpha */
        0 ]; /* vers */
< p1,s1 > = shapwilk(uf1000,optn);
print "SHAPIRO-Wilks P:",p1;
print "SHAPIRO-Wilks S:",s1;
```

Shapiro-Wilks Test for Normal Distribution  
Hypothesis H0: Data are normally distributed versus  
Hypothesis H1: Data are not normally distributed  
H0 rejected: Probability=5.41037e-017 Statistics=8.29541

SHAPIRO-Wilks P: 5.410e-017

SHAPIRO-Wilks S: 8.2954

```
print "Shapiro-Francia test for Normality";  
optn = [ 2 , /* print */  
        .05 , /* alpha */  
        1 ]; /* vers */  
< p1,s1 > = shapwilk(uf1000,optn);  
print "SHAPIRO-Francia P:",p1;  
print "SHAPIRO-Francia S:",s1;
```

Shapiro-Francia Test for Normal Distribution  
Hypothesis H0: Data are normally distributed versus  
Hypothesis H1: Data are not normally distributed  
H0 rejected: Probability=0.000253907 Statistics=3.4766

SHAPIRO-Francia P: 0.0002539

SHAPIRO-Francia S: 3.4766