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Description

`xtgls` fits panel-data linear models by using feasible generalized least squares. This command allows estimation in the presence of AR(1) autocorrelation within panels and cross-sectional correlation and heteroskedasticity across panels.

Quick start

GLS regression of `y` on `x1`, `x2`, and [indicators](#) for levels of categorical variable `a` using [xtset](#) data

```
xtgls y x1 x2 i.a
```

With heteroskedastic but uncorrelated errors across panels

```
xtgls y x1 x2 i.a, panels(heteroskedastic)
```

With heteroskedastic and correlated errors across panels

```
xtgls y x1 x2 i.a, panels(correlated)
```

Three-stage GLS with a common first-order autocorrelation within panels

```
xtgls y x1 x2 i.a, panels(correlated) corr(ar1)
```

Same as above, but let autocorrelation structure be panel-specific

```
xtgls y x1 x2 i.a, panels(correlated) corr(pсар1)
```

Same as above, but estimate by iterated GLS

```
xtgls y x1 x2 i.a, panels(correlated) corr(pсар1) igls
```

Menu

Statistics > Longitudinal/panel data > Contemporaneous correlation > GLS regression with correlated errors

Syntax

```
xtgls depvar [indepvars] [if] [in] [weight] [, options]
```

<i>options</i>	Description
Model	
<code>noconstant</code>	suppress constant term
<code>panels(iid)</code>	use i.i.d. error structure
<code>panels(heteroskedastic)</code>	use heteroskedastic but uncorrelated error structure
<code>panels(correlated)</code>	use heteroskedastic and correlated error structure
<code>corr(independent)</code>	use independent autocorrelation structure
<code>corr(ar1)</code>	use AR1 autocorrelation structure
<code>corr(psar1)</code>	use panel-specific AR1 autocorrelation structure
<code>rhotype(<i>calc</i>)</code>	specify method to compute autocorrelation parameter; see Options for details; seldom used
<code>igls</code>	use iterated GLS estimator instead of two-step GLS estimator
<code>force</code>	estimate even if observations unequally spaced in time
SE	
<code>nmk</code>	normalize standard error by $N - k$ instead of N
Reporting	
<code>level(#)</code>	set confidence level; default is level(95)
<code>display_options</code>	control columns and column formats, row spacing, line width, display of omitted variables and base and empty cells, and factor-variable labeling
Optimization	
<code>optimize_options</code>	control the optimization process; seldom used
<code>coeflegend</code>	display legend instead of statistics

A panel variable must be specified. For correlation structures other than independent, a time variable must be specified. A time variable must also be specified if `panels(correlated)` is specified. Use `xtset`; see [\[XT\] xtset](#).

indepvars may contain factor variables; see [\[U\] 11.4.3 Factor variables](#).

depvar and *indepvars* may contain time-series operators; see [\[U\] 11.4.4 Time-series varlists](#).

`by`, `collect`, and `statsby` are allowed; see [\[U\] 11.1.10 Prefix commands](#).

`aweight`s are allowed; see [\[U\] 11.1.6 weight](#).

`coeflegend` does not appear in the dialog box.

See [\[U\] 20 Estimation and postestimation commands](#) for more capabilities of estimation commands.

Options

Model
<code>noconstant</code> ; see [R] Estimation options .
<code>panels(<i>pdist</i>)</code> specifies the error structure across panels.
<code>panels(iid)</code> specifies a homoskedastic error structure with no cross-sectional correlation. This is the default.

`panels(heteroskedastic)` specifies a heteroskedastic error structure with no cross-sectional correlation.

`panels(correlated)` specifies a heteroskedastic error structure with cross-sectional correlation. If `p(c)` is specified, you must also specify a time variable (use `xtset`). The results will be based on a generalized inverse of a singular matrix unless $T \geq m$ (the number of periods is greater than or equal to the number of panels).

`corr(corr)` specifies the assumed autocorrelation within panels.

`corr(independent)` specifies that there is no autocorrelation. This is the default.

`corr(ar1)` specifies that, within panels, there is AR(1) autocorrelation and that the coefficient of the AR(1) process is common to all the panels. If `c(ar1)` is specified, you must also specify a time variable (use `xtset`).

`corr(psar1)` specifies that, within panels, there is AR(1) autocorrelation and that the coefficient of the AR(1) process is specific to each panel. `psar1` stands for panel-specific AR(1). If `c(psar1)` is specified, a time variable must also be specified; use `xtset`.

`rhotype(calc)` specifies the method to be used to calculate the autocorrelation parameter:

<code>regress</code>	regression using lags; the default
<code>dw</code>	Durbin–Watson calculation
<code>freg</code>	regression using leads
<code>nagar</code>	Nagar calculation
<code>theil</code>	Theil calculation
<code>tscorr</code>	time-series autocorrelation calculation

All the calculations are asymptotically equivalent and consistent; this is a rarely used option.

`igls` requests an iterated GLS estimator instead of the two-step GLS estimator for a nonautocorrelated model or instead of the three-step GLS estimator for an autocorrelated model. The iterated GLS estimator converges to the MLE for the `corr(independent)` models but does not for the other `corr()` models.

`force` specifies that estimation be forced even though the time variable is not equally spaced. This is relevant only for correlation structures that require knowledge of the time variable. These correlation structures require that observations be equally spaced so that calculations based on lags correspond to a constant time change. If you specify a time variable indicating that observations are not equally spaced, the (time dependent) model will not be fit. If you also specify `force`, the model will be fit, and it will be assumed that the lags based on the data ordered by the time variable are appropriate.

SE

`nmk` specifies that standard errors be normalized by $N - k$, where k is the number of parameters estimated, rather than N , the number of observations. Different authors have used one or the other normalization. [Greene \(2018, 313\)](#) remarks that whether a degree-of-freedom correction improves the small-sample properties is an open question.

Reporting

`level(#)`; see [\[R\] Estimation options](#).

`display_options`: `noci`, `nopvalues`, `noomitted`, `vsquish`, `noemptycells`, `baselevels`, `allbaselevels`, `nofvlabel`, `fvwrap(#)`, `fvwrapon(style)`, `cformat(%fmt)`, `pformat(%fmt)`, `sformat(%fmt)`, and `nolstretch`; see [R] [Estimation options](#).

Optimization

`optimize_options` control the iterative optimization process. These options are seldom used.

`iterate(#)` specifies the maximum number of iterations. When the number of iterations equals `#`, the optimization stops and presents the current results, even if convergence has not been reached. The default is `iterate(100)`.

`tolerance(#)` specifies the tolerance for the coefficient vector. When the relative change in the coefficient vector from one iteration to the next is less than or equal to `#`, the optimization process is stopped. `tolerance(1e-7)` is the default.

`log` and `nolog` specify whether to display the iteration log. The iteration log is displayed by default unless you used `set iterlog off` to suppress it; see `set iterlog` in [R] [set iter](#).

The following option is available with `xtgls` but is not shown in the dialog box:

`coeflegend`; see [R] [Estimation options](#).

Remarks and examples

Remarks are presented under the following headings:

[Introduction](#)
[Heteroskedasticity across panels](#)
[Correlation across panels \(cross-sectional correlation\)](#)
[Autocorrelation within panels](#)

Introduction

Information on GLS can be found in [Greene \(2018\)](#), [Maddala and Lahiri \(2006\)](#), [Davidson and MacKinnon \(1993\)](#), and [Judge et al. \(1985\)](#).

If you have many panels relative to periods, see [XT] [xtreg](#) and [XT] [xtgee](#). `xtgee`, in particular, provides capabilities similar to those of `xtgls` but does not allow cross-sectional correlation. On the other hand, `xtgee` allows a richer description of the correlation within panels as long as the same correlations apply to all panels. `xtgls` provides two unique features:

1. Cross-sectional correlation may be modeled (`panels(correlated)`).
2. Within panels, the AR(1) correlation coefficient may be unique (`corr(psar1)`).

`xtgls` allows models with heteroskedasticity and no cross-sectional correlation, but, strictly speaking, `xtgee` does not. `xtgee` with the `vce(robust)` option relaxes the assumption of equal variances, at least as far as the standard error calculation is concerned.

Also, `xtgls, panels(iid) corr(independent) nmk` is equivalent to `regress`.

The `nmk` option uses $n - k$ rather than n to normalize the variance calculation.

To fit a model with autocorrelated errors (`corr(ar1)` or `corr(psar1)`), the data must be equally spaced in time. To fit a model with cross-sectional correlation (`panels(correlated)`), panels must have the same number of observations (be balanced).

The equation from which the models are developed is given by

$$y_{it} = \mathbf{x}_{it}\boldsymbol{\beta} + \epsilon_{it}$$

where $i = 1, \dots, m$ is the number of units (or panels) and $t = 1, \dots, T_i$ is the number of observations for panel i . This model can equally be written as

$$\begin{bmatrix} \mathbf{y}_1 \\ \mathbf{y}_2 \\ \vdots \\ \mathbf{y}_m \end{bmatrix} = \begin{bmatrix} \mathbf{X}_1 \\ \mathbf{X}_2 \\ \vdots \\ \mathbf{X}_m \end{bmatrix} \boldsymbol{\beta} + \begin{bmatrix} \boldsymbol{\epsilon}_1 \\ \boldsymbol{\epsilon}_2 \\ \vdots \\ \boldsymbol{\epsilon}_m \end{bmatrix}$$

The variance matrix of the disturbance terms can be written as

$$E[\boldsymbol{\epsilon}\boldsymbol{\epsilon}'] = \boldsymbol{\Omega} = \begin{bmatrix} \sigma_{1,1}\boldsymbol{\Omega}_{1,1} & \sigma_{1,2}\boldsymbol{\Omega}_{1,2} & \cdots & \sigma_{1,m}\boldsymbol{\Omega}_{1,m} \\ \sigma_{2,1}\boldsymbol{\Omega}_{2,1} & \sigma_{2,2}\boldsymbol{\Omega}_{2,2} & \cdots & \sigma_{2,m}\boldsymbol{\Omega}_{2,m} \\ \vdots & \vdots & \ddots & \vdots \\ \sigma_{m,1}\boldsymbol{\Omega}_{m,1} & \sigma_{m,2}\boldsymbol{\Omega}_{m,2} & \cdots & \sigma_{m,m}\boldsymbol{\Omega}_{m,m} \end{bmatrix}$$

For the $\boldsymbol{\Omega}_{i,j}$ matrices to be parameterized to model cross-sectional correlation, they must be square (balanced panels).

In these models, we assume that the coefficient vector $\boldsymbol{\beta}$ is the same for all panels and consider a variety of models by changing the assumptions on the structure of $\boldsymbol{\Omega}$.

For the classic OLS regression model, we have

$$\begin{aligned} E[\epsilon_{i,t}] &= 0 \\ \text{Var}[\epsilon_{i,t}] &= \sigma^2 \\ \text{Cov}[\epsilon_{i,t}, \epsilon_{j,s}] &= 0 \quad \text{if } t \neq s \text{ or } i \neq j \end{aligned}$$

This amounts to assuming that $\boldsymbol{\Omega}$ has the structure given by

$$\boldsymbol{\Omega} = \begin{bmatrix} \sigma^2 \mathbf{I} & \mathbf{0} & \cdots & \mathbf{0} \\ \mathbf{0} & \sigma^2 \mathbf{I} & \cdots & \mathbf{0} \\ \vdots & \vdots & \ddots & \vdots \\ \mathbf{0} & \mathbf{0} & \cdots & \sigma^2 \mathbf{I} \end{bmatrix}$$

whether or not the panels are balanced (the $\mathbf{0}$ matrices may be rectangular). The classic OLS assumptions are the default `panels(iid)` and `corr(independent)` options for this command.

Heteroskedasticity across panels

In many cross-sectional datasets, the variance for each of the panels differs. It is common to have data on countries, states, or other units that have variation of scale. The heteroskedastic model is specified by including the `panels(heteroskedastic)` option, which assumes that

$$\boldsymbol{\Omega} = \begin{bmatrix} \sigma_1^2 \mathbf{I} & \mathbf{0} & \cdots & \mathbf{0} \\ \mathbf{0} & \sigma_2^2 \mathbf{I} & \cdots & \mathbf{0} \\ \vdots & \vdots & \ddots & \vdots \\ \mathbf{0} & \mathbf{0} & \cdots & \sigma_m^2 \mathbf{I} \end{bmatrix}$$

► Example 1

Greene (2012, 1112) reprints data in a classic study of investment demand by Grunfeld and Griliches (1960). Below we allow the variances to differ for each of the five companies.

```
. use https://www.stata-press.com/data/r19/invest2
. xtgls invest market stock, panels(hetero)
Cross-sectional time-series FGLS regression
Coefficients: generalized least squares
Panels:      heteroskedastic
Correlation: no autocorrelation

Estimated covariances   =          5      Number of obs      =      100
Estimated autocorrelations =          0      Number of groups   =       5
Estimated coefficients   =          3      Time periods        =      20
                                   Wald chi2(2)      =      865.38
                                   Prob > chi2       =      0.0000
```

invest	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
market	.0949905	.007409	12.82	0.000	.0804692	.1095118
stock	.3378129	.0302254	11.18	0.000	.2785722	.3970535
_cons	-36.2537	6.124363	-5.92	0.000	-48.25723	-24.25017



Correlation across panels (cross-sectional correlation)

We may wish to assume that the error terms of panels are correlated, in addition to having different scale variances. The variance structure is specified by including the `panels(correlated)` option and is given by

$$\Omega = \begin{bmatrix} \sigma_1^2 \mathbf{I} & \sigma_{1,2} \mathbf{I} & \cdots & \sigma_{1,m} \mathbf{I} \\ \sigma_{2,1} \mathbf{I} & \sigma_2^2 \mathbf{I} & \cdots & \sigma_{2,m} \mathbf{I} \\ \vdots & \vdots & \ddots & \vdots \\ \sigma_{m,1} \mathbf{I} & \sigma_{m,2} \mathbf{I} & \cdots & \sigma_m^2 \mathbf{I} \end{bmatrix}$$

Because we must estimate cross-sectional correlation in this model, the panels must be balanced (and $T \geq m$ for valid results). A time variable must also be specified so that `xtgls` knows how the observations within panels are ordered. `xtset` shows us that this is true.

► Example 2

```
. xtset
Panel variable: company (strongly balanced)
Time variable: time, 1 to 20
Delta: 1 unit

. xtgls invest market stock, panels(correlated)
Cross-sectional time-series FGLS regression
Coefficients: generalized least squares
Panels:      heteroskedastic with cross-sectional correlation
Correlation: no autocorrelation

Estimated covariances      =      15      Number of obs      =      100
Estimated autocorrelations =      0      Number of groups   =      5
Estimated coefficients     =      3      Time periods       =      20
                                Wald chi2(2)      =      1285.19
                                Prob > chi2       =      0.0000
```

invest	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
market	.0961894	.0054752	17.57	0.000	.0854583	.1069206
stock	.3095321	.0179851	17.21	0.000	.2742819	.3447822
_cons	-38.36128	5.344871	-7.18	0.000	-48.83703	-27.88552

The estimated cross-sectional covariances are stored in `e(Sigma)`.

```
. matrix list e(Sigma)
symmetric e(Sigma) [5,5]
      _ee      _ee2      _ee3      _ee4      _ee5
 _ee  9410.9061
 _ee2 -168.04631  755.85077
 _ee3 -1915.9538 -4163.3434  34288.49
 _ee4 -1129.2896 -80.381742  2259.3242  633.42367
 _ee5  258.50132  4035.872 -27898.235 -1170.6801  33455.511
```



► Example 3

We can obtain the MLE results by specifying the `igls` option, which iterates the GLS estimation technique to convergence:

```
. xtgls invest market stock, panels(correlated) igls
Iteration 1: Tolerance = .2127384
Iteration 2: Tolerance = .22817
(output omitted)
Iteration 1046: Tolerance = 1.000e-07
Cross-sectional time-series FGLS regression
Coefficients: generalized least squares
Panels:      heteroskedastic with cross-sectional correlation
Correlation: no autocorrelation

Estimated covariances      =          15      Number of obs      =          100
Estimated autocorrelations =           0      Number of groups   =           5
Estimated coefficients     =           3      Time periods        =          20
                                   Wald chi2(2)      =        558.51
Log likelihood              = -515.4222        Prob > chi2          =         0.0000
```

invest	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
market	.023631	.004291	5.51	0.000	.0152207	.0320413
stock	.1709472	.0152526	11.21	0.000	.1410526	.2008417
_cons	-2.216508	1.958845	-1.13	0.258	-6.055774	1.622759

Here the log likelihood is reported in the header of the output.



Autocorrelation within panels

The individual identity matrices along the diagonal of Ω may be replaced with more general structures to allow for serial correlation. `xtgls` allows three options so that you may assume a structure with `corr(independent)` (no autocorrelation); `corr(ar1)` (serial correlation where the correlation parameter is common for all panels); or `corr(pсар1)` (serial correlation where the correlation parameter is unique for each panel).

The restriction of a common autocorrelation parameter is reasonable when the individual correlations are nearly equal and the time series are short.

If the restriction of a common autocorrelation parameter is reasonable, this allows us to use more information in estimating the autocorrelation parameter to produce a more reasonable estimate of the regression coefficients.

When you specify `corr(ar1)` or `corr(pсар1)`, the iterated GLS estimator does not converge to the MLE.

► Example 4

If `corr(ar1)` is specified, each group is assumed to have errors that follow the same AR(1) process; that is, the autocorrelation parameter is the same for all groups.

```
. xtgls invest market stock, panels(hetero) corr(ar1)
Cross-sectional time-series FGLS regression
Coefficients: generalized least squares
Panels:      heteroskedastic
Correlation: common AR(1) coefficient for all panels (0.8651)

Estimated covariances      =          5      Number of obs      =       100
Estimated autocorrelations =          1      Number of groups   =        5
Estimated coefficients     =          3      Time periods        =       20
                                           Wald chi2(2)         =    119.69
                                           Prob > chi2          =     0.0000
```

invest	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
market	.0744315	.0097937	7.60	0.000	.0552362	.0936268
stock	.2874294	.0475391	6.05	0.000	.1942545	.3806043
_cons	-18.96238	17.64943	-1.07	0.283	-53.55464	15.62987



► Example 5

If `corr(psar1)` is specified, each group is assumed to have errors that follow a different AR(1) process.

```
. xtgls invest market stock, panels(iid) corr(psar1)
Cross-sectional time-series FGLS regression
Coefficients: generalized least squares
Panels:      homoskedastic
Correlation: panel-specific AR(1)

Estimated covariances      =          1      Number of obs      =       100
Estimated autocorrelations =          5      Number of groups   =        5
Estimated coefficients     =          3      Time periods        =       20
                                           Wald chi2(2)         =    252.93
                                           Prob > chi2          =     0.0000
```

invest	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
market	.0934343	.0097783	9.56	0.000	.0742693	.1125993
stock	.3838814	.0416775	9.21	0.000	.302195	.4655677
_cons	-10.1246	34.06675	-0.30	0.766	-76.8942	56.64499



Stored results

xtgls stores the following in `e()`:

Scalars

<code>e(N)</code>	number of observations
<code>e(N_ic)</code>	number of observations used to compute information criteria
<code>e(N_g)</code>	number of groups
<code>e(N_t)</code>	number of periods
<code>e(N_miss)</code>	number of missing observations
<code>e(n_cf)</code>	number of estimated coefficients
<code>e(n_cv)</code>	number of estimated covariances
<code>e(n_cr)</code>	number of estimated correlations
<code>e(df)</code>	degrees of freedom
<code>e(df_pear)</code>	degrees of freedom for Pearson χ^2
<code>e(df_ic)</code>	degrees of freedom for information criteria
<code>e(ll)</code>	log likelihood
<code>e(chi2)</code>	χ^2
<code>e(g_min)</code>	smallest group size
<code>e(g_avg)</code>	average group size
<code>e(g_max)</code>	largest group size
<code>e(rank)</code>	rank of $e(V)$
<code>e(rc)</code>	return code

Macros

<code>e(cmd)</code>	xtgls
<code>e(cmdline)</code>	command as typed
<code>e(depvar)</code>	name of dependent variable
<code>e(ivar)</code>	variable denoting groups
<code>e(tvar)</code>	variable denoting time within groups
<code>e(coefftype)</code>	estimation scheme
<code>e(corr)</code>	correlation structure
<code>e(vt)</code>	panel option
<code>e(rhotype)</code>	type of estimated correlation
<code>e(wtype)</code>	weight type
<code>e(wexp)</code>	weight expression
<code>e(title)</code>	title in estimation output
<code>e(chi2type)</code>	Wald; type of model χ^2 test
<code>e(rho)</code>	ρ
<code>e(properties)</code>	b V
<code>e(predict)</code>	program used to implement predict
<code>e(asbalanced)</code>	factor variables fvset as asbalanced
<code>e(asobserved)</code>	factor variables fvset as asobserved

Matrices

<code>e(b)</code>	coefficient vector
<code>e(Sigma)</code>	$\widehat{\Sigma}$ matrix
<code>e(V)</code>	variance–covariance matrix of the estimators

Functions

<code>e(sample)</code>	marks estimation sample
------------------------	-------------------------

In addition to the above, the following is stored in `r()`:

Matrices

<code>r(table)</code>	matrix containing the coefficients with their standard errors, test statistics, p -values, and confidence intervals
-----------------------	---

Note that results stored in `r()` are updated when the command is replayed and will be replaced when any r-class command is run after the estimation command.

Methods and formulas

The GLS results are given by

$$\begin{aligned}\widehat{\beta}_{\text{GLS}} &= (\mathbf{X}'\widehat{\Omega}^{-1}\mathbf{X})^{-1}\mathbf{X}'\widehat{\Omega}^{-1}\mathbf{y} \\ \widehat{\text{Var}}(\widehat{\beta}_{\text{GLS}}) &= (\mathbf{X}'\widehat{\Omega}^{-1}\mathbf{X})^{-1}\end{aligned}$$

For all our models, the Ω matrix may be written in terms of the Kronecker product:

$$\Omega = \Sigma_{m \times m} \otimes \mathbf{I}_{T_i \times T_i}$$

The estimated variance matrix is obtained by substituting the estimator $\widehat{\Sigma}$ for Σ , where

$$\widehat{\Sigma}_{i,j} = \frac{\widehat{\epsilon}_i' \widehat{\epsilon}_j}{T}$$

The residuals used in estimating Σ are first obtained from OLS regression. If the estimation is iterated, residuals are obtained from the last fitted model.

Maximum likelihood estimates may be obtained by iterating the FGLS estimates to convergence for models with no autocorrelation, `corr(independent)`.

The GLS estimates and their associated standard errors are calculated using $\widehat{\Sigma}^{-1}$. As Beck and Katz (1995) point out, the Σ matrix is of rank at most $\min(T, m)$ when you use the `panels(correlated)` option. For the GLS results to be valid (not based on a generalized inverse), T must be at least as large as m , as you need at least as many period observations as there are panels.

Beck and Katz (1995) suggest using OLS parameter estimates with asymptotic standard errors that are corrected for correlation between the panels. This estimation can be performed with the `xtpcse` command; see [XT] [xtpcse](#).

References

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Also see

- [XT] [xtgls postestimation](#) — Postestimation tools for xtgls
- [XT] [xtpcse](#) — Linear regression with panel-corrected standard errors
- [XT] [xtreg](#) — Linear models for panel data
- [XT] [xtregar](#) — Fixed- and random-effects linear models with an AR(1) disturbance
- [XT] [xtset](#) — Declare data to be panel data
- [R] [regress](#) — Linear regression
- [TS] [newey](#) — Regression with Newey–West standard errors
- [TS] [prais](#) — Prais–Winsten and Cochrane–Orcutt regression
- [U] [20 Estimation and postestimation commands](#)

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