

**mepoisson** — Multilevel mixed-effects Poisson regression

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## Description

`mepoisson` fits mixed-effects models for count responses. The conditional distribution of the response given the random effects is assumed to be Poisson.

## Quick start

### *Without weights*

Two-level Poisson regression of  $y$  on  $x$  with random intercepts by `lev2`

```
mepoisson y x || lev2:
```

Add `evar` measuring exposure

```
mepoisson y x, exposure(evar) || lev2:
```

Same as above, but report incidence-rate ratios

```
mepoisson y x, exposure(evar) || lev2:, irr
```

Add [indicators](#) for levels of categorical variable `a` and random coefficients on  $x$

```
mepoisson y x i.a || lev2: x, irr
```

Three-level random-intercept model of  $y$  on  $x$  with `lev2` nested within `lev3`

```
mepoisson y x || lev3: || lev2:
```

### *With weights*

Two-level Poisson regression of  $y$  on  $x$  with random intercepts by `lev2` and observation-level frequency weights `wvar1`

```
mepoisson y x [fweight=wvar1] || lev2:
```

Two-level random-intercept model from a two-stage sampling design with PSUs identified by `psu` using PSU-level and observation-level sampling weights `wvar2` and `wvar1`, respectively

```
mepoisson y x [pweight=wvar1] || psu:, pweight(wvar2)
```

Add secondary sampling stage with units identified by `ssu` having weights `wvar2` and PSU-level weights `wvar3` for a three-level random-intercept model

```
mepoisson y x [pw=wvar1] || psu:, pw(wvar3) || ssu:, pw(wvar2)
```

Same as above, but `svyset` data first

```
svyset psu, weight(wvar3) || ssu, weight(wvar2) || _n, weight(wvar1)
svy: mepoisson y x || psu: || ssu:
```

## Menu

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## Syntax

```
mepoisson depvar fe_equation [ || re_equation ] [ || re_equation ... ] [ , options ]
```

where the syntax of *fe\_equation* is

```
[ indepvars ] [ if ] [ in ] [ weight ] [ , fe_options ]
```

and the syntax of *re\_equation* is one of the following:

for random coefficients and intercepts

```
levelvar: [ varlist ] [ , re_options ]
```

for random effects among the values of a factor variable in a crossed-effects model

```
levelvar: R.varname
```

*levelvar* is a variable identifying the group structure for the random effects at that level or is `_all` representing one group comprising all observations.

<i>fe_options</i>	Description
<code>Model</code>	
<code>noconstant</code>	suppress the constant term from the fixed-effects equation
<code>exposure(<i>varname</i><sub>e</sub>)</code>	include $\ln(\text{varname}_e)$ in model with coefficient constrained to 1
<code>offset(<i>varname</i><sub>o</sub>)</code>	include <i>varname</i> <sub>o</sub> in model with coefficient constrained to 1

<i>re_options</i>	Description
<code>Model</code>	
<code>covariance(<i>vartype</i>)</code>	variance–covariance structure of the random effects
<code>noconstant</code>	suppress constant term from the random-effects equation
<code>fweight(<i>varname</i>)</code>	frequency weights at higher levels
<code>iweight(<i>varname</i>)</code>	importance weights at higher levels
<code>pweight(<i>varname</i>)</code>	sampling weights at higher levels

<i>options</i>	Description
Model	
<code>constraints</code> ( <i>constraints</i> )	apply specified linear constraints
SE/Robust	
<code>vce</code> ( <i>vcetype</i> )	<i>vcetype</i> may be <code>oim</code> , <code>opg</code> , <code>robust</code> , or <code>cluster</code> <i>clustvar</i>
Reporting	
<code>level</code> (#)	set confidence level; default is <code>level(95)</code>
<code>irr</code>	report fixed-effects coefficients as incidence-rate ratios
<code>nocnsreport</code>	do not display constraints
<code>notable</code>	suppress coefficient table
<code>noheader</code>	suppress output header
<code>nogroup</code>	suppress table summarizing groups
<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
Integration	
<code>intmethod</code> ( <i>intmethod</i> )	integration method
<code>intpoints</code> (#)	set the number of integration (quadrature) points for all levels; default is <code>intpoints(7)</code>
Maximization	
<code>maximize_options</code>	control the maximization process; seldom used
<code>startvalues</code> ( <i>svmethod</i> )	method for obtaining starting values
<code>startgrid</code> [ ( <i>gridspec</i> ) ]	perform a grid search to improve starting values
<code>noestimate</code>	do not fit the model; show starting values instead
<code>dnnumerical</code>	use numerical derivative techniques
<code>collinear</code>	keep collinear variables
<code>coeflegend</code>	display legend instead of statistics
<hr/>	
<i>vartype</i>	Description
<code>independent</code>	one unique variance parameter per random effect and all covariances 0; the default unless the <code>R.</code> notation is used
<code>exchangeable</code>	equal variances for random effects and one common pairwise covariance
<code>identity</code>	equal variances for random effects and all covariances 0; the default if the <code>R.</code> notation is used
<code>unstructured</code>	all variances and covariances to be distinctly estimated
<code>fixed</code> ( <i>matname</i> )	user-selected variances and covariances constrained to specified values; the remaining variances and covariances unrestricted
<code>pattern</code> ( <i>matname</i> )	user-selected variances and covariances constrained to be equal; the remaining variances and covariances unrestricted

<i>intmethod</i>	Description
<u>mv</u> aghermite	mean–variance adaptive Gauss–Hermite quadrature; the default unless a crossed random-effects model is fit
<u>mc</u> aghermite	mode-curvature adaptive Gauss–Hermite quadrature
<u>pc</u> aghermite	Pinheiro–Chao mode-curvature adaptive Gauss–Hermite quadrature
<u>g</u> hermite	nonadaptive Gauss–Hermite quadrature
<u>l</u> aplace	Laplacian approximation; the default for crossed random-effects models
<u>pc</u> laplace	Pinheiro–Chao Laplacian approximation

*indepvars* and *varlist* may contain factor variables; see [U] 11.4.3 **Factor variables**.

*depvar*, *indepvars*, and *varlist* may contain time-series operators; see [U] 11.4.4 **Time-series varlists**.

*bayes*, *by*, *collect*, and *svy* are allowed; see [U] 11.1.10 **Prefix commands**. For more details, see [BAYES] **bayes: mepoisson**.

*vce()* and *weights* are not allowed with the *svy* prefix; see [SVY] **svy**.

*fweights*, *iwweights*, and *pweights* are allowed; see [U] 11.1.6 **weight**. Only one type of weight may be specified. Weights are not supported under the Laplacian approximation or for crossed models.

*startvalues()*, *startgrid*, *noestimate*, *dnumerical*, *collinear*, and *coeflegend* do not appear in the dialog box.

See [U] 20 **Estimation and postestimation commands** for more capabilities of estimation commands.

## Options

### Model

*noconstant* suppresses the constant (intercept) term and may be specified for the fixed-effects equation and for any of or all the random-effects equations.

*exposure*(*varname<sub>e</sub>*) specifies a variable that reflects the amount of exposure over which the *depvar* events were observed for each observation;  $\ln(\text{varname}_e)$  is included in the fixed-effects portion of the model with the coefficient constrained to be 1.

*offset*(*varname<sub>o</sub>*) specifies that *varname<sub>o</sub>* be included in the fixed-effects portion of the model with the coefficient constrained to be 1.

*covariance*(*vartype*) specifies the structure of the covariance matrix for the random effects and may be specified for each random-effects equation. *vartype* is one of the following: *independent*, *exchangeable*, *identity*, *unstructured*, *fixed*(*matname*), or *pattern*(*matname*).

*covariance*(*independent*) covariance structure allows for a distinct variance for each random effect within a random-effects equation and assumes that all covariances are 0. The default is *covariance*(*independent*) unless a crossed random-effects model is fit, in which case the default is *covariance*(*identity*).

*covariance*(*exchangeable*) structure specifies one common variance for all random effects and one common pairwise covariance.

*covariance*(*identity*) is short for “multiple of the identity”; that is, all variances are equal and all covariances are 0.

*covariance*(*unstructured*) allows for all variances and covariances to be distinct. If an equation consists of *p* random-effects terms, the unstructured covariance matrix will have  $p(p + 1)/2$  unique parameters.

`covariance(fixed(matname))` and `covariance(pattern(matname))` covariance structures provide a convenient way to impose constraints on variances and covariances of random effects. Each specification requires a *matname* that defines the restrictions placed on variances and covariances. Only elements in the lower triangle of *matname* are used, and row and column names of *matname* are ignored. A missing value in *matname* means that a given element is unrestricted. In a `fixed(matname)` covariance structure, (co)variance ( $i, j$ ) is constrained to equal the value specified in the  $i, j$ th entry of *matname*. In a `pattern(matname)` covariance structure, (co)variances ( $i, j$ ) and ( $k, l$ ) are constrained to be equal if  $matname[i, j] = matname[k, l]$ .

`fweight(varname)` specifies frequency weights at higher levels in a multilevel model, whereas frequency weights at the first level (the observation level) are specified in the usual manner, for example, `[fw=fwtvar1]`. *varname* can be any valid Stata variable name, and you can specify `fweight()` at levels two and higher of a multilevel model. For example, in the two-level model

```
. mecmd fixed_portion [fw = wt1] || school: ... , fweight(wt2) ...
```

the variable `wt1` would hold the first-level (the observation-level) frequency weights, and `wt2` would hold the second-level (the school-level) frequency weights.

`iweight(varname)` specifies importance weights at higher levels in a multilevel model, whereas importance weights at the first level (the observation level) are specified in the usual manner, for example, `[iw=iwtvar1]`. *varname* can be any valid Stata variable name, and you can specify `iweight()` at levels two and higher of a multilevel model. For example, in the two-level model

```
. mecmd fixed_portion [iw = wt1] || school: ... , iweight(wt2) ...
```

the variable `wt1` would hold the first-level (the observation-level) importance weights, and `wt2` would hold the second-level (the school-level) importance weights.

`pweight(varname)` specifies sampling weights at higher levels in a multilevel model, whereas sampling weights at the first level (the observation level) are specified in the usual manner, for example, `[pw=pwtvar1]`. *varname* can be any valid Stata variable name, and you can specify `pweight()` at levels two and higher of a multilevel model. For example, in the two-level model

```
. mecmd fixed_portion [pw = wt1] || school: ... , pweight(wt2) ...
```

variable `wt1` would hold the first-level (the observation-level) sampling weights, and `wt2` would hold the second-level (the school-level) sampling weights.

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

#### SE/Robust

`vce(vcetype)` specifies the type of standard error reported, which includes types that are derived from asymptotic theory (`oim`, `opg`), that are robust to some kinds of misspecification (`robust`), and that allow for intragroup correlation (`cluster clustvar`); see [\[R\] vce\\_option](#). If `vce(robust)` is specified, robust variances are clustered at the highest level in the multilevel model.

#### Reporting

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

`irr` reports estimated fixed-effects coefficients transformed to incidence-rate ratios, that is,  $\exp(\beta)$  rather than  $\beta$ . Standard errors and confidence intervals are similarly transformed. This option affects how results are displayed, not how they are estimated or stored. `irr` may be specified either at estimation or upon replay.

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

`notable` suppresses the estimation table, either at estimation or upon replay.

`noheader` suppresses the output header, either at estimation or upon replay.

`nogroup` suppresses the display of group summary information (number of groups, average group size, minimum, and maximum) from the output header.

*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](#).

#### Integration

`intmethod(intmethod)` specifies the integration method to be used for the random-effects model. `mvaghermite` performs mean–variance adaptive Gauss–Hermite quadrature; `mcaghermite` and `pcaghermite` perform mode-curvature adaptive Gauss–Hermite quadrature; `ghermite` performs nonadaptive Gauss–Hermite quadrature; and `laplace` and `pclaplace` perform the Laplacian approximation, equivalent to mode-curvature adaptive Gaussian quadrature with one integration point. Techniques `pcaghermite` and `pclaplace` obtain the random-effects mode and curvature using the efficient hierarchical decomposition algorithm described in [Pinheiro and Chao \(2006\)](#). For hierarchical models, this algorithm takes advantage of the design structure to minimize memory use and utilizes a series of orthogonal triangulations to compute the factored random-effects Hessian indirectly, avoiding the sparse full Hessian. Techniques `mcaghermite` and `laplace` use Cholesky factorization on the full Hessian. For four- and higher-level hierarchical designs, there can be dramatic computation-time differences.

The default integration method is `mvaghermite` unless a crossed random-effects model is fit, in which case the default integration method is `laplace`. The Laplacian approximation has been known to produce biased parameter estimates; however, the bias tends to be more prominent in the estimates of the variance components rather than in the estimates of the fixed effects.

For crossed random-effects models, estimation with more than one quadrature point may be prohibitively intensive even for a small number of levels. For this reason, the integration method defaults to the Laplacian approximation. You may override this behavior by specifying a different integration method.

`intpoints(#)` sets the number of integration points for quadrature. The default is `intpoints(7)`, which means that seven quadrature points are used for each level of random effects. This option is not allowed with `intmethod(laplace)`.

The more integration points, the more accurate the approximation to the log likelihood. However, computation time increases as a function of the number of quadrature points raised to a power equaling the dimension of the random-effects specification. In crossed random-effects models and in models with many levels or many random coefficients, this increase can be substantial.

#### Maximization

*maximize\_options*: `difficult`, `technique(algorithm_spec)`, `iterate(#)`, `[no]log`, `trace`, `gradient`, `showstep`, `hessian`, `showtolerance`, `tolerance(#)`, `ltolerance(#)`, `nrtolerance(#)`, `nonrtolerance`, and `from(init_specs)`; see [R] [Maximize](#). Those that require special mention for `mepoisson` are listed below.

`from()` accepts a properly labeled vector of initial values or a list of coefficient names with values. A list of values is not allowed.

The following options are available with `mepoisson` but are not shown in the dialog box:

`startvalues(svmethod)`, `startgrid[(gridspec)]`, `noestimate`, and `dnnumerical`; see [ME] [meglm](#).

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

## Remarks and examples

[stata.com](#)

For a general introduction to `me` commands, see [ME] [me](#).

Remarks are presented under the following headings:

[Introduction](#)

[Two-level models](#)

[Higher-level models](#)

## Introduction

Mixed-effects Poisson regression is Poisson regression containing both fixed effects and random effects. In longitudinal data and panel data, random effects are useful for modeling intracluster correlation; that is, observations in the same cluster are correlated because they share common cluster-level random effects.

`mepoisson` allows for many levels of random effects. However, for simplicity, for now we consider the two-level model, where for a series of  $M$  independent clusters, and conditional on a set of random effects  $\mathbf{u}_j$ ,

$$\Pr(y_{ij} = y | \mathbf{x}_{ij}, \mathbf{u}_j) = \exp(-\mu_{ij}) \mu_{ij}^y / y! \quad (1)$$

for  $\mu_{ij} = \exp(\mathbf{x}_{ij}\boldsymbol{\beta} + \mathbf{z}_{ij}\mathbf{u}_j)$ ,  $j = 1, \dots, M$  clusters, with cluster  $j$  consisting of  $i = 1, \dots, n_j$  observations. The responses are counts  $y_{ij}$ . The  $1 \times p$  row vector  $\mathbf{x}_{ij}$  are the covariates for the fixed effects, analogous to the covariates you would find in a standard Poisson regression model, with regression coefficients (fixed effects)  $\boldsymbol{\beta}$ . For notational convenience here and throughout this manual entry, we suppress the dependence of  $y_{ij}$  on  $\mathbf{x}_{ij}$ .

The  $1 \times q$  vector  $\mathbf{z}_{ij}$  are the covariates corresponding to the random effects and can be used to represent both random intercepts and random coefficients. For example, in a random-intercept model,  $\mathbf{z}_{ij}$  is simply the scalar 1. The random effects  $\mathbf{u}_j$  are  $M$  realizations from a multivariate normal distribution with mean  $\mathbf{0}$  and  $q \times q$  variance matrix  $\boldsymbol{\Sigma}$ . The random effects are not directly estimated as model parameters but are instead summarized according to the unique elements of  $\boldsymbol{\Sigma}$ , known as variance components. One special case of (1) places  $\mathbf{z}_{ij} = \mathbf{x}_{ij}$  so that all covariate effects are essentially random and distributed as multivariate normal with mean  $\boldsymbol{\beta}$  and variance  $\boldsymbol{\Sigma}$ .

As noted in section 13.7 of [Rabe-Hesketh and Skrondal \(2022\)](#), the inclusion of a random intercept causes the marginal variance of  $y_{ij}$  to be greater than the marginal mean, provided the variance of the random intercept is not 0. Thus the random intercept in a mixed-effects Poisson model produces overdispersion, a measure of variability above and beyond that allowed by a Poisson process; see [R] [nbreg](#) and [ME] [menbreg](#).

Below we present examples of mixed-effects Poisson regression; refer to [ME] [me](#) and [ME] [meglm](#) for additional examples including crossed random-effects models. A two-level Poisson model can also be fit using `xtpoisson` with the `re` option; see [XT] [xtpoisson](#). In the absence of random effects, mixed-effects Poisson regression reduces to standard Poisson regression; see [R] [poisson](#).

## Two-level models

### ▷ Example 1: Two-level random-intercept model

Breslow and Clayton (1993) fit a mixed-effects Poisson model to data from a randomized trial of the drug progabide for the treatment of epilepsy.

```
. use https://www.stata-press.com/data/r18/epilepsy
(Epilepsy data; progabide drug treatment)
. describe
Contains data from https://www.stata-press.com/data/r18/epilepsy.dta
Observations:      236      Epilepsy data; progabide drug
                        treatment
Variables:          8      31 May 2022 14:09
                        (_dta has notes)
```

Variable name	Storage type	Display format	Value label	Variable label
subject	byte	%9.0g		Subject ID: 1-59
seizures	int	%9.0g		No. of seizures
treat	byte	%9.0g	treat	Treatment
visit	float	%9.0g		Doctor's visit
lage	float	%9.0g		log(age), mean-centered
lbas	float	%9.0g		log(0.25*baseline seizures), mean-centered
lbas_trt	float	%9.0g		lbas/treat interaction
v4	byte	%8.0g		Fourth visit indicator

Sorted by: subject

Originally from [Thall and Vail \(1990\)](#), data were collected on 59 subjects (31 on progabide, 28 on placebo). The number of epileptic seizures (`seizures`) was recorded during the two weeks prior to each of four doctor visits (`visit`). The treatment group is identified by the indicator variable `treat`. Data were also collected on the logarithm of age (`lage`) and the logarithm of one-quarter the number of seizures during the eight weeks prior to the study (`lbas`). The variable `lbas_trt` represents the interaction between `lbas` and treatment. `lage`, `lbas`, and `lbas_trt` are mean centered. Because the study originally noted a substantial decrease in seizures prior to the fourth doctor visit, an indicator, `v4`, for the fourth visit was also recorded.

[Breslow and Clayton \(1993\)](#) fit a random-effects Poisson model for the number of observed seizures

$$\log(\mu_{ij}) = \beta_0 + \beta_1 \text{treat}_{ij} + \beta_2 \text{lbas}_{ij} + \beta_3 \text{lbas\_trt}_{ij} + \beta_4 \text{lage}_{ij} + \beta_5 \text{v4}_{ij} + u_j$$

for  $j = 1, \dots, 59$  subjects and  $i = 1, \dots, 4$  visits. The random effects  $u_j$  are assumed to be normally distributed with mean 0 and variance  $\sigma_u^2$ .



```

. mepoisson seizures treat lbas lbas_trt lage v4 || subject:
Fitting fixed-effects model:
Iteration 0:  Log likelihood = -1016.4106
Iteration 1:  Log likelihood = -819.20112
Iteration 2:  Log likelihood = -817.66006
Iteration 3:  Log likelihood = -817.65925
Iteration 4:  Log likelihood = -817.65925
Refining starting values:
Grid node 0:  Log likelihood = -680.40523
Refining starting values (unscaled likelihoods):
Grid node 0:  Log likelihood = -680.40523
Fitting full model:
Iteration 0:  Log likelihood = -680.40523 (not concave)
Iteration 1:  Log likelihood = -672.95766 (not concave)
Iteration 2:  Log likelihood = -667.14039
Iteration 3:  Log likelihood = -665.51823
Iteration 4:  Log likelihood = -665.29165
Iteration 5:  Log likelihood = -665.29067
Iteration 6:  Log likelihood = -665.29067
Mixed-effects Poisson regression          Number of obs    =      236
Group variable: subject                   Number of groups =       59
                                           Obs per group:
                                           min =           4
                                           avg =          4.0
                                           max =           4
Integration method: mvaghermite           Integration pts. =        7
                                           Wald chi2(5)     =      121.70
                                           Prob > chi2      =       0.0000
Log likelihood = -665.29067

```

seizures	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
treat	-.9330306	.4007512	-2.33	0.020	-1.718489	-.1475727
lbas	.8844225	.1312033	6.74	0.000	.6272689	1.141576
lbas_trt	.3382561	.2033021	1.66	0.096	-.0602087	.736721
lage	.4842226	.3471905	1.39	0.163	-.1962582	1.164703
v4	-.1610871	.0545758	-2.95	0.003	-.2680536	-.0541206
_cons	2.154578	.2199928	9.79	0.000	1.7234	2.585756
subject						
var(_cons)	.2528664	.0589844			.1600801	.399434

LR test vs. Poisson model: chibar2(01) = 304.74      Prob >= chibar2 = 0.0000

The number of seizures before the fourth visit does exhibit a significant drop, and the patients on progabide demonstrate a decrease in frequency of seizures compared with the placebo group. The subject-specific random effects also appear significant:  $\hat{\sigma}_u^2 = 0.25$  with standard error 0.06.

Because this is a simple random-intercept model, you can obtain equivalent results by using `xtpoisson` with the `re` and `normal` options.



### ► Example 2: Two-level random-slope model

In their study of PQL, [Breslow and Clayton \(1993\)](#) also fit a model where they dropped the fixed effect on `v4` and replaced it with a random subject-specific linear trend over the four doctor visits. The model they fit is

$$\log(\mu_{ij}) = \beta_0 + \beta_1 \text{treat}_{ij} + \beta_2 \text{lbas}_{ij} + \beta_3 \text{lbas\_trt}_{ij} + \beta_4 \text{lage}_{ij} + \beta_5 \text{visit}_{ij} + u_j + v_j \text{visit}_{ij}$$

where  $(u_j, v_j)$  are bivariate normal with 0 mean and variance–covariance matrix

$$\Sigma = \text{Var} \begin{bmatrix} u_j \\ v_j \end{bmatrix} = \begin{bmatrix} \sigma_u^2 & \sigma_{uv} \\ \sigma_{uv} & \sigma_v^2 \end{bmatrix}$$

```
. mepoisson seizures treat lbas lbas_trt lage visit || subject: visit,
> covariance(unstructured) intpoints(9) nolog
Mixed-effects Poisson regression      Number of obs      =      236
Group variable: subject                Number of groups   =      59
                                       Obs per group:
                                       min =            4
                                       avg =            4.0
                                       max =            4
Integration method: mvaghermite        Integration pts.   =      9
Wald chi2(5)                          =     115.56
Prob > chi2                             =      0.0000
Log likelihood = -655.68103
```

seizures	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
treat	-.9286592	.4021715	-2.31	0.021	-1.716901	-.1404175
lbas	.8849762	.1312535	6.74	0.000	.627724	1.142228
lbas_trt	.3379759	.2044471	1.65	0.098	-.062733	.7386849
lage	.4767192	.3536276	1.35	0.178	-.2163781	1.169817
visit	-.2664098	.1647098	-1.62	0.106	-.5892352	.0564156
_cons	2.099555	.2203749	9.53	0.000	1.667629	2.531482
subject						
var(visit)	.5314803	.229385			.2280928	1.238405
var(_cons)	.2514923	.0587902			.1590534	.3976549
subject						
cov(visit, _cons)	.0028715	.0887037	0.03	0.974	-.1709846	.1767276

```
LR test vs. Poisson model: chi2(3) = 324.54      Prob > chi2 = 0.0000
```

Note: LR test is conservative and provided only for reference.

In the above, we specified the `covariance(unstructured)` option to allow correlation between  $u_j$  and  $v_j$ , although on the basis of the above output it probably was not necessary—the default independent structure would have sufficed. In the interest of getting more accurate estimates, we also increased the number of quadrature points to nine, although the estimates do not change much when compared with estimates based on the default seven quadrature points.

The essence of the above-fitted model is that after adjusting for other covariates, the log trend in seizures is modeled as a random subject-specific line, with intercept distributed as  $N(\beta_0, \sigma_u^2)$  and slope distributed as  $N(\beta_5, \sigma_v^2)$ . From the above output,  $\hat{\beta}_0 = 2.10$ ,  $\hat{\sigma}_u^2 = 0.25$ ,  $\hat{\beta}_5 = -0.27$ , and  $\hat{\sigma}_v^2 = 0.53$ .

You can predict the random effects  $u_j$  and  $v_j$  by using `predict` after `mepoisson`; see [ME] [mepoisson postestimation](#). Better still, you can obtain a predicted number of seizures that takes these random effects into account.

## Higher-level models

### ▷ Example 3: Three- and four-level random-intercept model

Rabe-Hesketh and Skrondal (2022, exercise 13.7) describe data from the *Atlas of Cancer Mortality in the European Economic Community* (EEC) (Smans, Mair, and Boyle 1993). The data were analyzed in Langford, Bentham, and McDonald (1998) and record the number of deaths among males due to malignant melanoma during 1971–1980.

```
. use https://www.stata-press.com/data/r18/melanoma
(Skin cancer (melanoma) data)
. describe
Contains data from https://www.stata-press.com/data/r18/melanoma.dta
Observations:      354      Skin cancer (melanoma) data
Variables:         6       30 May 2022 17:10
                        (_dta has notes)
```

Variable name	Storage type	Display format	Value label	Variable label
nation	byte	%11.0g	n	Nation ID
region	byte	%9.0g		Region ID: EEC level-I areas
county	int	%9.0g		County ID: EEC level-II/level-III areas
deaths	int	%9.0g		No. deaths during 1971-1980
expected	float	%9.0g		No. expected deaths
uv	float	%9.0g		UV dose, mean-centered

Sorted by:

Nine European nations (variable `nation`) are represented, and data were collected over geographical regions defined by EEC statistical services as level I areas (variable `region`), with deaths being recorded for each of 354 counties, which are level II or level III EEC-defined areas (variable `county`, which identifies the observations). Counties are nested within regions, and regions are nested within nations.

The variable `deaths` records the number of deaths for each county, and `expected` records the expected number of deaths (the exposure) on the basis of crude rates for the combined countries. Finally, the variable `uv` is a measure of exposure to ultraviolet (UV) radiation.

In modeling the number of deaths, one possibility is to include dummy variables for the nine nations as fixed effects. Another is to treat these as random effects and fit the three-level random-intercept Poisson model,

$$\log(\mu_{ijk}) = \log(\text{expected}_{ijk}) + \beta_0 + \beta_1 \text{uv}_{ijk} + u_k + v_{jk}$$

for nation  $k$ , region  $j$ , and county  $i$ . The model includes an exposure term for expected deaths.

```
. mepoisson deaths uv, exposure(expected) || nation: || region:
Fitting fixed-effects model:
Iteration 0:  Log likelihood = -2136.5847
Iteration 1:  Log likelihood = -1723.8955
Iteration 2:  Log likelihood = -1723.7727
Iteration 3:  Log likelihood = -1723.7727
Refining starting values:
Grid node 0:  Log likelihood = -1166.6536
Refining starting values (unscaled likelihoods):
Grid node 0:  Log likelihood = -1166.6536
Fitting full model:
Iteration 0:  Log likelihood = -1166.6536 (not concave)
Iteration 1:  Log likelihood = -1152.2741 (not concave)
Iteration 2:  Log likelihood = -1146.3094 (not concave)
Iteration 3:  Log likelihood = -1119.8479 (not concave)
Iteration 4:  Log likelihood = -1108.0129 (not concave)
Iteration 5:  Log likelihood = -1098.8067
Iteration 6:  Log likelihood = -1095.7563
Iteration 7:  Log likelihood = -1095.3164
Iteration 8:  Log likelihood = -1095.31
Iteration 9:  Log likelihood = -1095.31
Mixed-effects Poisson regression          Number of obs    =        354
```

Grouping information

Group variable	No. of groups	Observations per group		
		Minimum	Average	Maximum
nation	9	3	39.3	95
region	78	1	4.5	13

```
Integration method: mvaghermite          Integration pts. =        7
Wald chi2(1) =        6.12
Prob > chi2 =        0.0134
Log likelihood = -1095.31
```

deaths	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
uv	-.0282041	.0113998	-2.47	0.013	-.0505473	-.0058608
_cons	-.0639672	.1335515	-0.48	0.632	-.3257234	.197789
ln(expected)	1	(exposure)				
nation						
var(_cons)	.1371732	.0723303			.048802	.3855676
nation>						
region						
var(_cons)	.0483483	.0109079			.0310699	.0752353

```
LR test vs. Poisson model: chi2(2) = 1256.93          Prob > chi2 = 0.0000
```

Note: LR test is conservative and provided only for reference.

By including an exposure variable that is an expected rate, we are in effect specifying a linear model for the log of the standardized mortality ratio, the ratio of observed deaths to expected deaths that is based on a reference population. Here the reference population is all nine nations.

Looking at the estimated variance components, we can see there is more unobserved variability between nations than between regions within each nation. This may be due to, for example, country-specific informational campaigns on the risks of sun exposure.



```

Integration method: mcaghermite                    Integration pts. =          7
Log likelihood = -1086.6754                        Wald chi2(1)      =          8.62
                                                    Prob > chi2       =          0.0033

```

deaths	Coefficient	Std. err.	z	P> z	[95% conf. interval]	
uv	-.0334702	.0113968	-2.94	0.003	-.0558075	-.0111329
_cons	-.0864583	.1299275	-0.67	0.506	-.3411115	.168195
ln(expected)	1	(exposure)				
nation var(_cons)	.1288627	.0681643			.0456949	.3634011
nation> region var(_cons)	.0406279	.0105154			.0244633	.0674735
nation> region> county var(_cons)	.0146672	.0050979			.0074215	.0289867

```
LR test vs. Poisson model: chi2(3) = 1274.19          Prob > chi2 = 0.0000
```

Note: LR test is conservative and provided only for reference.

In the above, we used `intmethod(mcaghermite)`, which is not only faster but also produces estimates that closely agree with those obtained with the default `mvaghermite` integration method.

◀

## Stored results

`mepoisson` stores the following in `e()`:

### Scalars

<code>e(N)</code>	number of observations
<code>e(k)</code>	number of parameters
<code>e(k_dv)</code>	number of dependent variables
<code>e(k_eq)</code>	number of equations in <code>e(b)</code>
<code>e(k_eq_model)</code>	number of equations in overall model test
<code>e(k_f)</code>	number of fixed-effects parameters
<code>e(k_r)</code>	number of random-effects parameters
<code>e(k_rs)</code>	number of variances
<code>e(k_rc)</code>	number of covariances
<code>e(df_m)</code>	model degrees of freedom
<code>e(ll)</code>	log likelihood
<code>e(N_clust)</code>	number of clusters
<code>e(chi2)</code>	$\chi^2$
<code>e(p)</code>	<i>p</i> -value for model test
<code>e(ll_c)</code>	log likelihood, comparison model
<code>e(chi2_c)</code>	$\chi^2$ , comparison test
<code>e(df_c)</code>	degrees of freedom, comparison test
<code>e(p_c)</code>	<i>p</i> -value for comparison test
<code>e(rank)</code>	rank of <code>e(V)</code>
<code>e(ic)</code>	number of iterations
<code>e(rc)</code>	return code
<code>e(converged)</code>	1 if converged, 0 otherwise

### Macros

<code>e(cmd)</code>	<code>meglm</code>
<code>e(cmd2)</code>	<code>mepoisson</code>

<code>e(cmdline)</code>	command as typed
<code>e(depvar)</code>	name of dependent variable
<code>e(wtype)</code>	weight type
<code>e(wexp)</code>	weight expression (first-level weights)
<code>e(fweightk)</code>	<code>fweight</code> variable for <i>k</i> th highest level, if specified
<code>e(iweightk)</code>	<code>iweight</code> variable for <i>k</i> th highest level, if specified
<code>e(pweightk)</code>	<code>pweight</code> variable for <i>k</i> th highest level, if specified
<code>e(covariates)</code>	list of covariates
<code>e(ivals)</code>	grouping variables
<code>e(model)</code>	poisson
<code>e(title)</code>	title in estimation output
<code>e(link)</code>	log
<code>e(family)</code>	poisson
<code>e(clustvar)</code>	name of cluster variable
<code>e(offset)</code>	offset
<code>e(intmethod)</code>	integration method
<code>e(n_quad)</code>	number of integration points
<code>e(chi2type)</code>	Wald; type of model $\chi^2$
<code>e(vce)</code>	<code>vce</code> type specified in <code>vce()</code>
<code>e(vcetype)</code>	title used to label Std. err.
<code>e(opt)</code>	type of optimization
<code>e(which)</code>	max or min; whether optimizer is to perform maximization or minimization
<code>e(ml_method)</code>	type of ml method
<code>e(user)</code>	name of likelihood-evaluator program
<code>e(technique)</code>	maximization technique
<code>e(datasignature)</code>	the checksum
<code>e(datasignaturevars)</code>	variables used in calculation of checksum
<code>e(properties)</code>	b V
<code>e(estat_cmd)</code>	program used to implement <code>estat</code>
<code>e(predict)</code>	program used to implement <code>predict</code>
<code>e(marginsnotok)</code>	predictions disallowed by <code>margins</code>
<code>e(marginswtype)</code>	weight type for <code>margins</code>
<code>e(marginswexp)</code>	weight expression for <code>margins</code>
<code>e(asbalanced)</code>	factor variables <code>fvset</code> as <code>asbalanced</code>
<code>e(asobserved)</code>	factor variables <code>fvset</code> as <code>asobserved</code>
Matrices	
<code>e(b)</code>	coefficient vector
<code>e(Cns)</code>	constraints matrix
<code>e(iolog)</code>	iteration log (up to 20 iterations)
<code>e(gradient)</code>	gradient vector
<code>e(N_g)</code>	group counts
<code>e(g_min)</code>	group-size minimums
<code>e(g_avg)</code>	group-size averages
<code>e(g_max)</code>	group-size maximums
<code>e(V)</code>	variance-covariance matrix of the estimators
<code>e(V_modelbased)</code>	model-based variance
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, <i>p</i> -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

`mepoisson` is a convenience command for `meglm` with a `log` link and an `poisson` family; see [ME] `meglm`.

In a two-level Poisson model, for cluster  $j$ ,  $j = 1, \dots, M$ , the conditional distribution of  $\mathbf{y}_j = (y_{j1}, \dots, y_{jn_j})'$ , given a set of cluster-level random effects  $\mathbf{u}_j$ , is

$$\begin{aligned} f(\mathbf{y}_j | \mathbf{u}_j) &= \prod_{i=1}^{n_j} [\{\exp(\mathbf{x}_{ij}\boldsymbol{\beta} + \mathbf{z}_{ij}\mathbf{u}_j)\}^{y_{ij}} \exp\{-\exp(\mathbf{x}_{ij}\boldsymbol{\beta} + \mathbf{z}_{ij}\mathbf{u}_j)\} / y_{ij}!] \\ &= \exp \left[ \sum_{i=1}^{n_j} \{y_{ij}(\mathbf{x}_{ij}\boldsymbol{\beta} + \mathbf{z}_{ij}\mathbf{u}_j) - \exp(\mathbf{x}_{ij}\boldsymbol{\beta} + \mathbf{z}_{ij}\mathbf{u}_j) - \log(y_{ij}!)\} \right] \end{aligned}$$

Defining  $c(\mathbf{y}_j) = \sum_{i=1}^{n_j} \log(y_{ij}!)$ , where  $c(\mathbf{y}_j)$  does not depend on the model parameters, we can express the above compactly in matrix notation,

$$f(\mathbf{y}_j | \mathbf{u}_j) = \exp \{ \mathbf{y}'_j (\mathbf{X}_j \boldsymbol{\beta} + \mathbf{Z}_j \mathbf{u}_j) - \mathbf{1}' \exp(\mathbf{X}_j \boldsymbol{\beta} + \mathbf{Z}_j \mathbf{u}_j) - c(\mathbf{y}_j) \}$$

where  $\mathbf{X}_j$  is formed by stacking the row vectors  $\mathbf{x}_{ij}$  and  $\mathbf{Z}_j$  is formed by stacking the row vectors  $\mathbf{z}_{ij}$ . We extend the definition of  $\exp(\cdot)$  to be a vector function where necessary.

Because the prior distribution of  $\mathbf{u}_j$  is multivariate normal with mean  $\mathbf{0}$  and  $q \times q$  variance matrix  $\boldsymbol{\Sigma}$ , the likelihood contribution for the  $j$ th cluster is obtained by integrating  $\mathbf{u}_j$  out of the joint density  $f(\mathbf{y}_j, \mathbf{u}_j)$ ,

$$\begin{aligned} \mathcal{L}_j(\boldsymbol{\beta}, \boldsymbol{\Sigma}) &= (2\pi)^{-q/2} |\boldsymbol{\Sigma}|^{-1/2} \int f(\mathbf{y}_j | \mathbf{u}_j) \exp(-\mathbf{u}'_j \boldsymbol{\Sigma}^{-1} \mathbf{u}_j / 2) d\mathbf{u}_j \\ &= \exp\{-c(\mathbf{y}_j)\} (2\pi)^{-q/2} |\boldsymbol{\Sigma}|^{-1/2} \int \exp\{h(\boldsymbol{\beta}, \boldsymbol{\Sigma}, \mathbf{u}_j)\} d\mathbf{u}_j \end{aligned} \quad (2)$$

where

$$h(\boldsymbol{\beta}, \boldsymbol{\Sigma}, \mathbf{u}_j) = \mathbf{y}'_j (\mathbf{X}_j \boldsymbol{\beta} + \mathbf{Z}_j \mathbf{u}_j) - \mathbf{1}' \exp(\mathbf{X}_j \boldsymbol{\beta} + \mathbf{Z}_j \mathbf{u}_j) - \mathbf{u}'_j \boldsymbol{\Sigma}^{-1} \mathbf{u}_j / 2$$

and for convenience, in the arguments of  $h(\cdot)$  we suppress the dependence on the observable data  $(\mathbf{y}_j, \mathbf{X}_j, \mathbf{Z}_j)$ .

The integration in (2) has no closed form and thus must be approximated; see *Methods and formulas* in [ME] `meglm` for details.

`mepoisson` supports multilevel weights and survey data; see *Methods and formulas* in [ME] `meglm` for details.

## References

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## Also see

- [ME] **mepoisson postestimation** — Postestimation tools for mepoisson
- [ME] **menbreg** — Multilevel mixed-effects negative binomial regression
- [ME] **me** — Introduction to multilevel mixed-effects models
- [BAYES] **bayes: mepoisson** — Bayesian multilevel Poisson regression
- [SEM] **Intro 5** — Tour of models (*Multilevel mixed-effects models*)
- [SVY] **svy estimation** — Estimation commands for survey data
- [XT] **xtpoisson** — Fixed-effects, random-effects, and population-averaged Poisson models
- [U] **20 Estimation and postestimation commands**

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