

AP: Duflo (2001)

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Abstract

1. Robust Standard Errors

- A typical strategy in applied work is to report **heteroskedasticity-robust standard errors** for OLS regressions (“robust” option in Stata).
- Estimator for the variance of the asymptotic distribution of the OLS estimator that is **(asymptotically) consistent both under homo- and heteroskedasticity**.

8 Under homoskedasticity:

- OLS standard errors are not only **consistent**, but also the **best linear unbiased estimate** (BLUE, most efficient).
- Therefore, it is not recommended to use the “robust” ones.

12 Under heteroskedasticity:

- OLS estimator of standard errors is **inconsistent**, but the “robust” one is consistent.
- When the regression approximates a non-linear CEF, we have heteroskedastic errors in general, so better to use the robust one. Eg As soon as the dependent variable is not a continuous variable (e.g. binary), the CEF is not linear (easy proof). Then you will have, **by construction**, heteroskedastic errors.

- Since we do not know the CEF or form of heteroskedasticity, **in large samples it is safer to use robust standard errors**.
- But, both the OLS and the robust estimators are **biased in small samples**, and the robust ones can be **more biased**.
- **Conservative strategy:** use the larger of OLS (conventional) and robust standard errors, typically the robust ones.

26 2. Clustering and the Moulton Factor

- With large samples, we do not need to care too much about heteroskedasticity.
- However, **clustering (group structures) can affect standard errors substantially**, even with large samples.

- E.g., we randomize at the school level, but outcomes are measured at the student level.
- Clustering impacts the off-diagonal terms in the variance-covariance matrix of errors. Basically you have a block diagonal matrix: **correlation within groups**, but **no correlation across groups!** This affects the error term. Obs are no longer iid. You underestimate standard errors, you overestimate the precision of your estimator (the idea is that the true variability in real world is higher, and it is as if you have fewer than the number of observations to learn from the data)

- **Failure to control for within-cluster error correlation** can lead to:
 - Misleadingly small standard errors,
 - Inflated t-statistics,
 - Falsely small p-values.

- If we assume that **non-diagonal terms in the variance-covariance matrix are constant**, we can compute the exact bias and derive a correction: the **Moulton factor**.

Model for Clustered Observations

$$Y_{ig} = \beta_0 + \beta_1 X_g + e_{ig}$$

- $g = 1, \dots, G$ groups. Covariates X vary at group level.
- Test scores in the same school are correlated \Rightarrow errors e_{ig} are correlated.
- $\mathbb{E}(e_{ig} e_{jg}) = \rho_e \sigma^2 \neq 0$, about two observations i and j within the same group g.
- where considering a **Random effects model**: $e_{ig} = v_g + \eta_{ig}$, so we can decompose the observation i error as the sum of a group level error and an individual level error. The idea is putting all of the correlation among individuals belonging to the same cluster in this term and assume η_{ig} is homoskedastic and uncorrelated across individuals, if you do this you find β_1

Intraclass correlation coefficient:

$$\rho_e = \frac{\sigma_v^2}{\sigma_v^2 + \sigma_\eta^2}$$

This measure how similar are observations within a class.

Interpretation

- If $\sigma_v^2 = 0 \Rightarrow \rho_e = 0$
All variance is from individual noise \Rightarrow No within-group correlation.
- If $\sigma_\eta^2 = 0 \Rightarrow \rho_e = 1$
All individuals in a cluster have the same error \Rightarrow Perfect intra-cluster correlation.

Implications for Standard Errors

- When $\rho_e > 0$, individuals in the same group are not independent.
- If you ignore this and assume i.i.d. errors:
 - * Your standard errors will be too small.
 - * You overstate precision, leading to over-rejection of null hypotheses.

Moulton Factor

$$\sqrt{\frac{V(\hat{\beta}_1)}{V_c(\hat{\beta}_1)}} = \sqrt{1 + (n - 1)\rho_e}$$

- $V(\hat{\beta}_1)$: correct variance, $V_c(\hat{\beta}_1)$: conventional variance ignoring clustering.
- n : average group size.
- **Bias increases with group size n and ρ_e .**
- as soon as rho is greater than zero and group have more than 2 observations you have true standard errors > those assumed under homoskedasticity.
- if rho is equal to zero true se are equal to homoskedastic ones
- When $\rho_e = 1$, bias factor = \sqrt{n} . With rho = 1, **the bias is of size n**. No additional info from more observations in same group. So, n observations within a cluster are not adding any independence

information. But if you mistakenly treat them as n independent observations, you overestimate the information by a factor of n. All errors within group are exactly the same, thus their outcome should be the same, and more observations do not provide any additional information! you get no info by observing different units within the same group, you have actually n observations = n number of groups.

General Moulton Factor The covariates play a key role in the generalized Moulton correction because the bias from clustering depends not only on the correlation of the errors within groups, but also on the correlation of the covariates within groups.

$$\sqrt{\frac{V(\hat{\beta}_1)}{V_c(\hat{\beta}_1)}} = \sqrt{1 + \left(\frac{V(n_g)}{E(n_g)} + E(n_g) - 1 \right) \rho_e \rho_x}$$

- ρ_x : intraclass correlation of X .
- Formula adapts if n_g not constant (different group sizes).

Correcting Clustered SE: 5 Approaches

1. **Parametric:** use Moulton formula. Moulton correction is a parametric rule to correct for intraclass correlation. THE CORRECTION IS TRIVIAL, SVEGLIA GUARDA LA FORMULA SU ED APPLICALA!
2. **Cluster SE:** flexible. Consistent if G (number of groups) is large.
3. **Group Averages:** most conservative strategy: constraint myself to running a reg at the groups levels (throw away all individual level information). ISSUE, SCREENSHOT!
4. **GLS or MLE:** One would estimate the parameters of the error correlation model, and estimate the original model by feasible generalized least squares. Not robust to wrong specifications.
5. **Block Bootstrap:** resample clusters, compute $\hat{\beta}_1^{(b)}$, and estimate variance across B bootstrap samples. Need to do bootstrap with blocks to keep in all bootstrap samples the group structure!

When to Use Cluster Option in Stata?

- Must have a random sample of clusters from a super-population with large G (BASICALLY, LARGE NUMBER OF GROUPS).
- With regressor randomized **within clusters** \Rightarrow clustering not needed.
- With regressor randomized **at cluster level** \Rightarrow must cluster.

3. Bootstrapping

1. **Block (or pairs) Bootstrap:** re-sample entire clusters instead of individuals. Do the following steps B times (recommended $B \geq 400$):
 - 0.1 Form G clusters based on (y_i, x_i) by resampling with replacement G times from the original sample of clusters.
 - 0.2 Compute B estimates, one $\hat{\beta}_i$, for each bootstrap sample. Compute the variance of the B estimates.
2. **Problem:** does not have much better properties than cluster option. But, you can use it for methods when that option is not available.

Resampling technique that can be used as an alternative to inference based on asymptotic formulas.

Draw repeatedly from your sample of N observations, as if it were the population.

Get the bootstrap sampling distribution as the distribution of the estimator across the different draws.

The bootstrap can be useful when we do not have asymptotic formulas for the estimator, or for asymptotic refinements (i.e. reduction in finite sample bias for consistent estimators).

4. Serial Correlation

4.1. Serial Correlation in Panels and DID Models

Suppose the treatment d_{it} varies at the state level and over time, and we have individual-level data:

$$y_{its} = \gamma d_{ts} + x'_{its} \delta + \alpha_s + \delta_t + u_{its}$$

- **Key concern:** with serial correlation in errors over time within states, we must adjust standard errors for:

- group correlation (individuals in the same state),
- and serial correlation (within a state over time).

Cases

1. Serial correlation but no group correlation:

Use robust SEs or ARIMA-based methods to model serial correlation directly.

2. No serial correlation, but group correlation within states

(e.g., shocks independent over time but correlated within state): Cluster SEs at the group*time level.

3. Both serial correlation (within state) and group correlation:

Cluster at the **group level** (e.g., state), as a conservative and recommended approach. Because serial correlation cannot be ignored, you must collapse to the state level, which gives you fewer clusters (bigger clusters less fun).

Important: We need a *large number of groups* to estimate intra-cluster correlations over time reliably.

4.2. What to Cluster Over?

- Individuals often reside in subregions within larger regions.
- **Recommendation:** cluster at the **region** level (Cameron and Miller, 2015).
- Clustering at subregion level is incorrect if regressors or errors are correlated across subregions within the same region.
- **Practical strategy:** start clustering at the lowest level, try higher levels, and stop when standard errors stabilize.
- *Note:* larger clusters tend to have higher standard errors due to greater average n .

4.3. Small Number of Clusters

- Clustered SEs: t-statistic converges to $N(0, 1)$ only as $G \rightarrow \infty$.
- With finite G : inference should be based on a t-distribution with $G - 1$ degrees of freedom.
- **Donald and Lang (2007):** if there are K regressors invariant within cluster (take same values), use t-distribution with $G - K$ degrees of freedom.
- With few clusters, SEs are downward biased even with clustering \rightarrow over-rejection of the null.
- Empirical rule: clustering at group level is typically safe if $G \geq 50$ and cluster sizes are similar.

4.4. Fixes for Few Clusters

What if less than 50 clusters?

1. Bias-Corrected Cluster-Robust Variance

Cluster generalization of HC2/HC3 heteroskedasticity-robust formulas. Helps, but doesn't fully eliminate bias or over-rejection.

2. Cluster Bootstrap with Asymptotic Refinement (e.g., Wild Cluster Bootstrap)

1. Estimate unrestricted model, get $\hat{\beta}_k$ and t-stat.
2. Re-estimate under $\hat{\beta}_k = 0$ get restricted $\tilde{\beta}$, residuals \tilde{u} .
3. Do B bootstrap replications:

$$y_{ig}^b = X_{ig} \tilde{\beta} + \tilde{u} \cdot v_g^b$$

where $v_g^b = \pm 1$, randomly assigned per cluster. boh.

- 191 4. Estimate t-statistic for each bootstrap draw.
 192 5. Compute p-value = share of draws with bootstrapped t-stat >
 193 original.
 194 6. In Stata: use `cgmwildboot` or `boottest`.

195 ■ **Randomization Inference**

- 196 • Instead of drawing samples from a super population, we can
 197 draw from repeatedly sampling the treatment assignment.
 198 • Define your statistic (e.g. difference in means between treat-
 199 ment and control) and derive its exact distribution by calculat-
 200 ing the test statistic under each possible permutation of treat-
 201 ment allocation.
 202 • Assume you have a sample with 18 units, 9 assigned to treat-
 203 ment, you have $\binom{18}{9} = \frac{18!}{9! \cdot 9!} = 48,620$ different possible assign-
 204 ment vectors, get one statistic for each.
 205 • We ask how likely it is to observe a value of the test statistic as
 206 large as the one observed: count how many of the 48,620 test
 207 statistics have a value larger than the one observed.
 208 • Lets say it is true for 1,500 of them, then your Exact p-value
 209 is $\frac{1500}{48620} = 0.03$, and you can reject the null of no difference in
 210 means at the 97% level.
 211 • If the number of permutations is very large, you can calculate
 212 the statistic for a random sub-sample of permutations and get
 213 close to exact p-values.

214 **Stata Implementation: ritest**

- 215 • `ritest` (Hess, 2017 Stata Journal), `ssc install ritest`
 216 • `ritest treatment _b[treatment], reps(1000)`
 217 `seed(1111) cluster(id): reg Y treatment,`
 218 `cluster(id)`