

Models for clustered data

Jon Zelner

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University of Michigan School of Public Health

jzelner@umich.edu

www.jonzelner.net

Agenda

- Re-intro to GLMs

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- Why?

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- Why?
- Radon!

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- Why?
- Radon!
- Self-assessments.

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- Gives flexibility in the types of **outcomes** and **inputs** we can model.
- Model **non-linear** response using **linear** predictors

- Represent hierarchical/multi-level structure

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- Model spatial autocorrelation

Advanced GLM tricks

- Represent hierarchical/multi-level structure
- Model spatial autocorrelation
- Allow variance parameters to change with individual-level covariates to accommodate heteroskedasticity

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- Outcome assumed to have distribution with mean determined by regression coefficients.
- Link function, $g()$, translates between linear predictor and the mean of the distribution function.

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- Conditional mean for individual i denoted by $\hat{y}_i = g(y_i^*)$

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And others:

- Gamma, Exponential, Negative Binomial...

GLM with Normally-distributed errors

GLM with normally distributed errors \approx OLS regression

What are assumptions of ordinary least squares regression?

GLM translates OLS into an explicitly **probabilistic** framework

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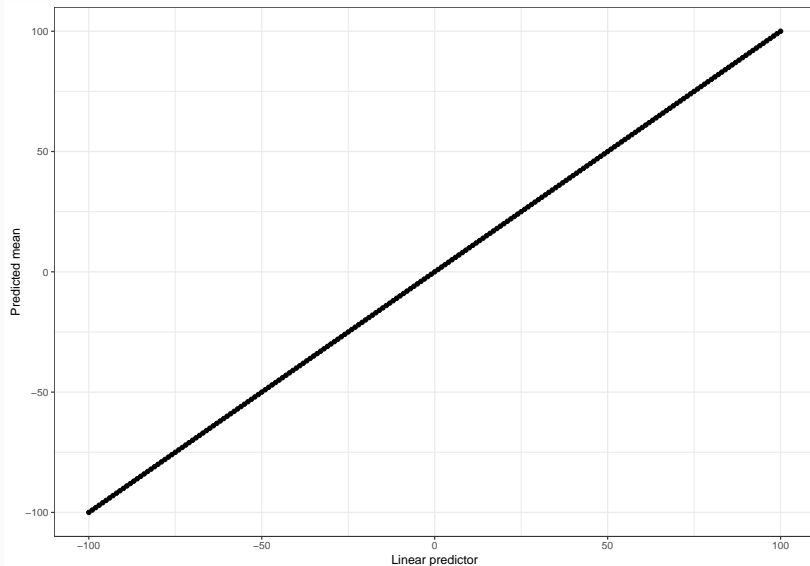
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- $\epsilon_i \sim \text{Normal}(0, \sigma^2)$

Identity link function maps real numbers to real numbers



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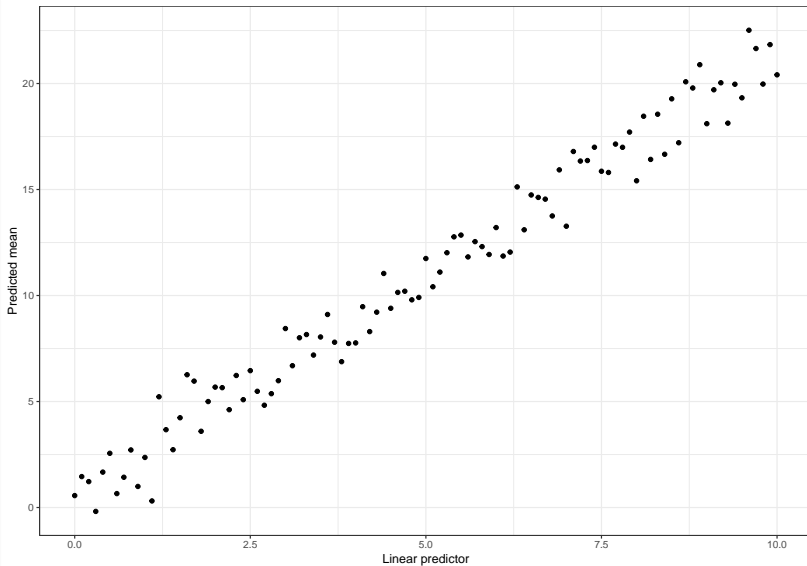
Interpreting results of Gaussian GLM

- α mean of y_i when $x_i = 0$
- β is change in mean of y_i for each one-unit change in x_i
- Just like in Merlo example, we also estimate the value of σ^2 that measures the amount of individual-level variability in outcomes

Generate some fake data in R

```
## Predictors
x <- seq(from = 0.0, to = 10.0, by = 0.1)
## Regression coefficient
b <- 2.1
## Intercept
a <- 0.4
## Variance
sd <- 1.0
## The Data
y <- a + b * x + rnorm(length(x), 0, sd)
```

The Data



Recover the parameters

```
m <- glm(y ~ x, data = df)
```

Parameter estimates

Table 1: Fitting generalized (gaussian/identity) linear model: $y \sim x$

	Estimate	Std. Error	t value	Pr(> t)
(Intercept)	0.55	0.2	2.7	0.0071
x	2.1	0.035	60	1.6e-79

Model Residuals

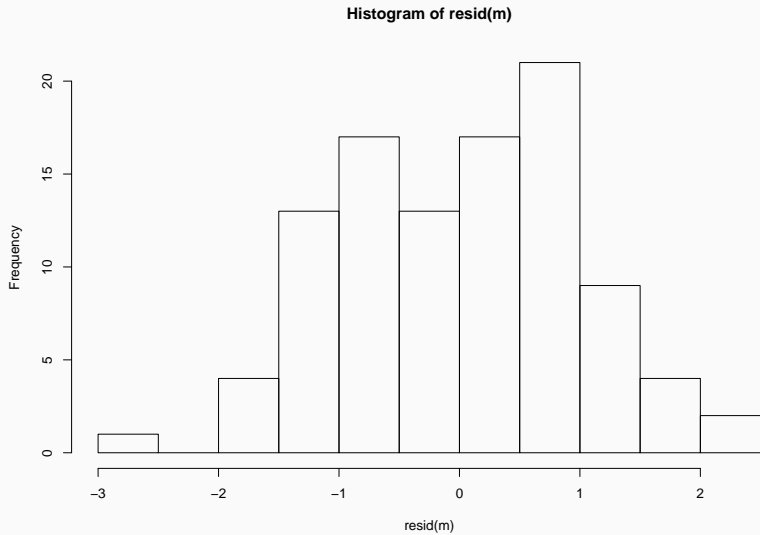


Table 2: Analysis of Variance Model

		Sum	Mean		
	Df	Sq	Sq	F value	Pr(>F)
x	1	3721	3721	3579	1.6e-79
Residuals	99	103	1	NA	NA

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- We'll talk about these soon enough, but first...
- Both of these necessitate calculating the **likelihood** of the data.

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For Gaussian GLM, f = Normal distribution, $\theta = \{\alpha, \beta, \sigma^2\}$

Motivations for Multilevel Modeling

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- When might you want them to both vary?

A re-introduction to Generalized Linear Models (GLMs) for clustered data

Notation for a classic un-clustered GLM

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- α is expected outcome when $x_i = 0$
- ϵ_i are independently and identically distributed (i.i.d.) errors

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In plain-ish English:

- Observation y_{ij} of individual i is a function of $\alpha + \beta x_i$ and normally distributed **errors** (ϵ_i) with mean zero and variance σ^2 .

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Another way of writing it:

- $y_i \sim N(\alpha + \beta x_i, \sigma^2)$

Three Approaches to Modeling Clustered Data



Which door will you choose?

Door #1: Ignore clustering and fit a normal GLM

- *Pool* data across all units, i.e. ignore clustering.
- i.e. fit model $y_{ij} = \alpha + \beta x_i + \epsilon_i$

Is this a good idea? Why or why not?

NO!



Complete pooling ignores potential sources of *observed* and *unobserved*. unit-level **confounding**.

Pooling clustered data violates assumption of independent errors

A **pooled** model:

$$y_i = \alpha + \beta x + \epsilon_i \quad (1)$$

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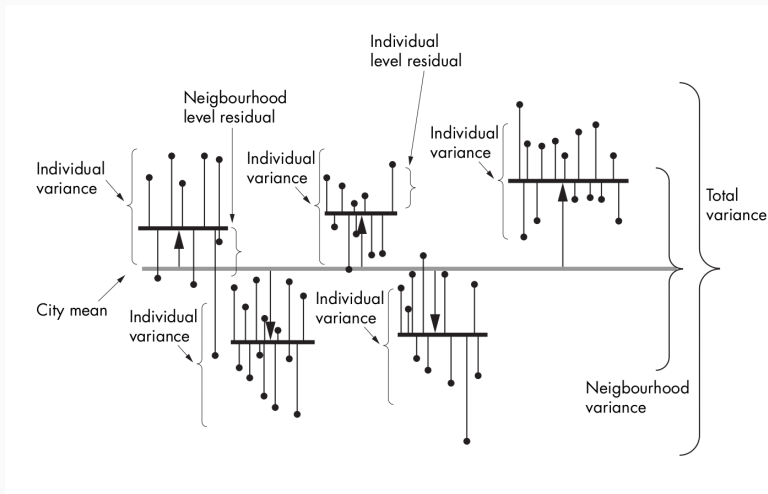
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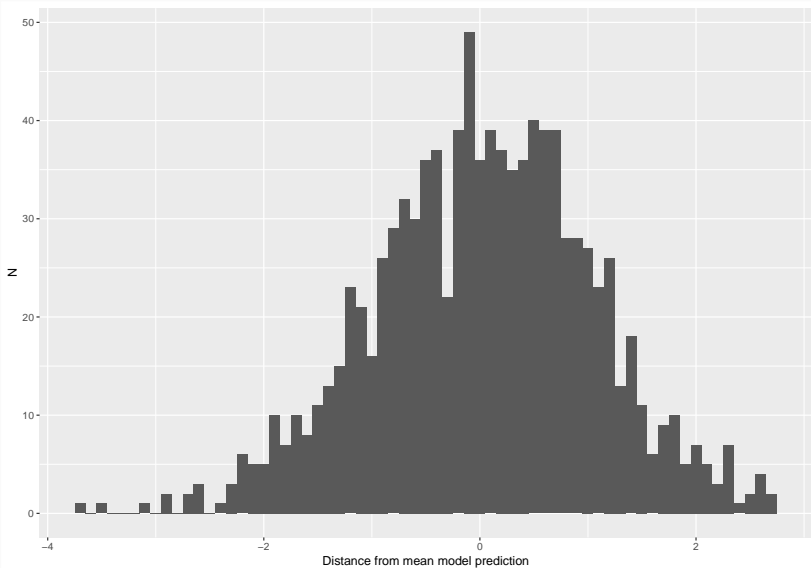
- y_i is a combination of systematic variation ($\alpha + \beta x$) and *uncorrelated* random noise (ϵ_i) where:

$$i.i.d. \epsilon \sim Normal(0, \sigma^2) \quad (2)$$

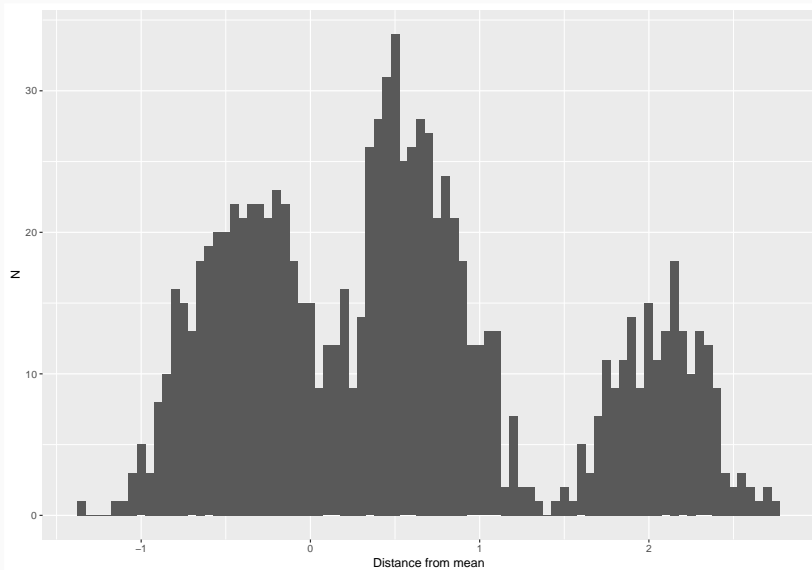
Clustering may result in correlation between average differences from mean



Your **residuals** should look like this



When you ignore clustering you may see something like:



Door #2: Fit a different model to each cluster

Fit *unpooled* model to each unit (j), assuming outcomes in each unit are independent:

- $y_{ij} = \alpha_j + \beta_j x_i + \epsilon_{ij}$
- $\epsilon_{ij} \sim N(0, \sigma_j^2)$

More danger!



Totally unpooled models run the risk of **overfitting** the data, particularly in small samples.

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- Some units (e.g. counties) may have few observations, making *unpooled* models impractical
- We may want to allow some effect of exposure (e.g. having a basement) to be consistent across counties.

Door #3: Partial Pooling!

- Allow effects to vary across clusters, but constrain them with a **prior** distribution.

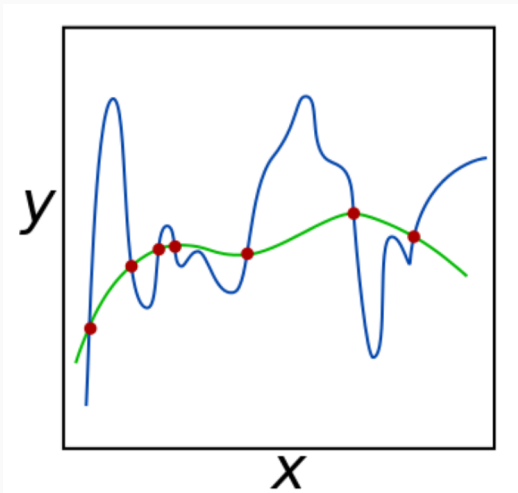
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- Allow effects to vary across clusters, but constrain them with a **prior** distribution.
- This approach accommodates variation across units without assuming they have no similarity.
- More likely to make accurate **out-of-sample** predictions than the fully-pooled or unpooled examples.

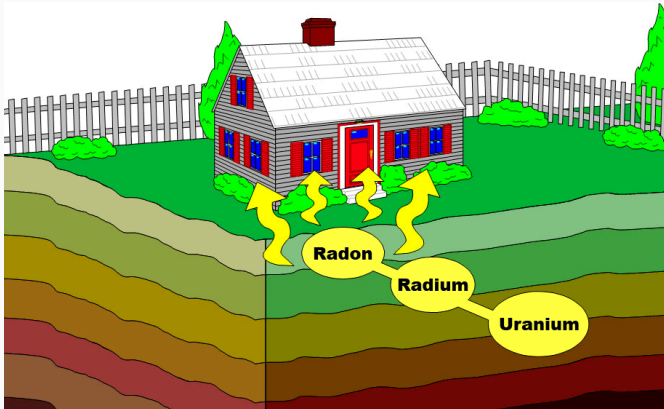
Partial pooling = Regularization



Both functions fit the data perfectly...which one should you prefer?

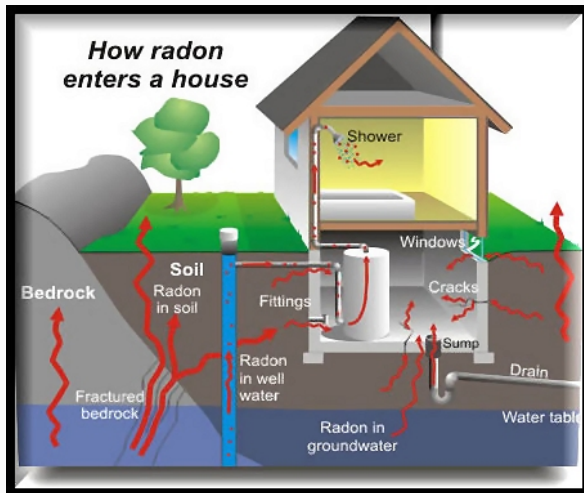
Radon Example

Radon is a carcinogenic gas



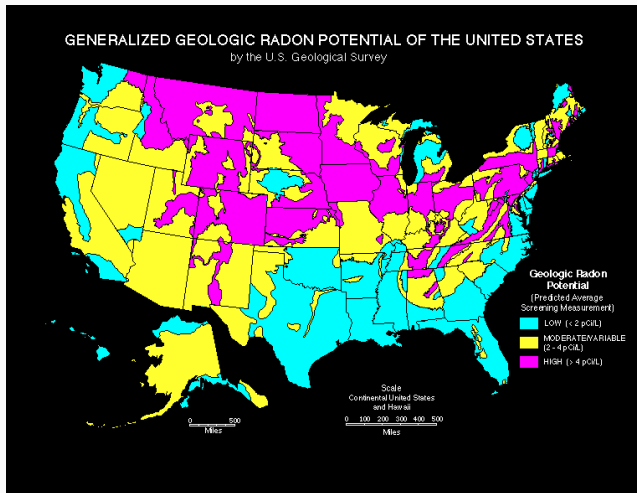
Radon is a byproduct of decaying soil uranium.

Radon enters a house more easily when it is built into the ground



Ann Arbor is a radon hotspot!

Considerable geographic variation in radon potential



Ann Arbor is a radon hotspot!

Trust me on this one...



My very own radon mitigation system.

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What should a model that accounts for important sources of variation in household radon potential include?

- County-level variation in soil uranium.
- Whether or not the radon measurement was taken in a basement.

Random intercepts account for county-level variation

Gelman [Gelman2006] proposes a multi-level model to measure household radon in household i in county j , y_{ij} :

$$\cdot y_{ij} \sim N(\alpha_j + \beta x_{ij}, \sigma_y^2), \text{ for } i = 1, \dots, n_j, j = 1, \dots, J$$

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Where:

- α_j is average, non-basement radon measure at county level
- β is fixed effect measuring average change in radon level in houses with a basement.
- σ_y^2 represents within-county variation in risk

Include predictors of county-level variation in second level

County-level random intercept is a function of county soil uranium measure, u_j :

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Where:

- γ_0 is expected household radon measure when $u_j = 0$
- γ_1 scales expected county-level uranium with u_j
- σ_α^2 is between-county variation in radon risk not measured by u_j .

Putting it all together

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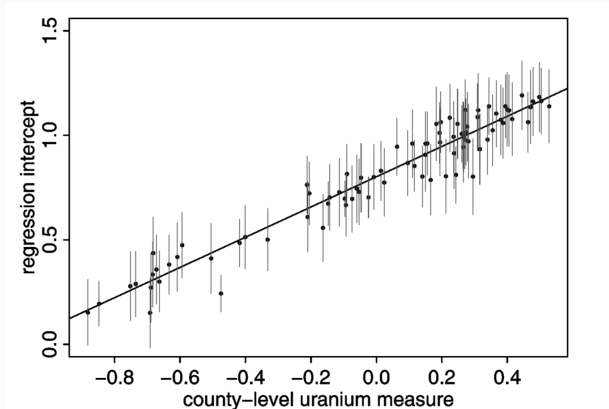
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Household-level radon measure is a function of having a basement and county-level intercept:

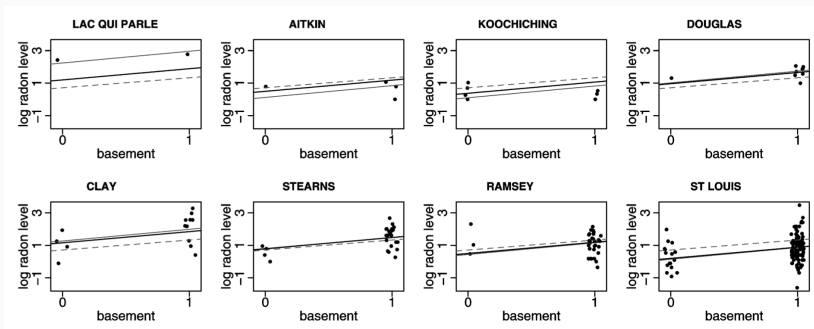
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County-level radon levels vary with soil uranium measures



County-level intercept, α_j , (± 1 standard error) as a function of county-level uranium.

Model predictions vs. radon measures by county



Multi-level regression line, $y = \alpha_j + \beta x$, from 8 Minnesota counties. Un-pooled estimates = light grey line; Totally pooled estimates = dashed grey line.

Next Time

- Hands-on with the Radon example

References
