

Outline

- Approaches to subset selection
- Regularization methods: ridge regression and lasso
- Choosing the penalty parameter

When Do Economists Care About Prediction? RCTs as a Case Study

THE CONVERSATION
Academic rigour, journalistic flair

How randomised trials became big in development economics

Published: December 9, 2019 8.18am GMT



A photograph showing two economists, Esther Duflo and Abhijit Banerjee, standing in front of a red sign for the 'Nobel Prize Press Conference' at the 'INSTITUTE OF ECONOMICS'. Duflo is on the left, wearing a grey blazer over a dark top, and Banerjee is on the right, wearing a yellow cardigan over a yellow shirt. They are both smiling. In the background, there's a blue table with water bottles and a whiteboard. A hand is visible in the foreground pointing towards the sign.

Esther Duflo (L) and Abhijit Banerjee, who, along with Michael Kremer (not pictured), won the 2019 Nobel Prize in Economic Sciences. EP/CJ Gunther

Source: *The Conversation* (2019)

When Do Economists Care About Prediction? RCTs as a Case Study

- Statistical power in a randomized trial depends on residual variance in the outcome
(Power is the probability of finding an impact – i.e. rejecting H_0 – if there is one.)
 - ▶ A typical RCT regression equation: $Y_{1,i} = \alpha + \beta D_i + \delta X_{0,i} + \gamma Y_{0,i} + \varepsilon_i$
 - ▶ $Y_{1,i}$ is the outcome, measure after the intervention
 - ▶ D_i is a dummy for being randomly assigned to the treatment group
 - ▶ $Y_{0,i}$ is the baseline value of the outcome
 - ▶ $X_{0,i}$ is a set of other baseline covariates that (one hopes) predict $Y_{1,i}$
 - ▶ The **minimum detectable effect** that a researcher can expect to measure through an RCT is proportional to the standard deviation of the residuals, i.e. to the unexplained variation in Y
- Economists running RCTs choose which covariates to measure, and pay for data collection
 - ▶ We want to measure X s that predict Y , and we don't want to throw money away
(by measuring a large number of baseline covariates that do not predict variation in Y)

Choosing Covariates in an RCT: The EMERGE Project



Encouraging
Multilingual
Early
Reading as the
Groundwork for
Education

Choosing Covariates in an RCT: The EMERGE Project

EMERGE was a cluster-randomized evaluation of an early literacy program in rural Kenya

- Intervention involved mother tongue storybooks and parent education
- Key child development outcomes of interest: literacy and vocabulary
- Large research team including me, Prof. Ozier, and two public health collaborators
- We designed survey instruments, and had to choose which variables to measure at baseline

Child development and educational outcomes tend to have high serial correlation

- Individual ability at time $t - 1$ is a strong predictor of individual ability at time t
- The right set of covariates can substantially increase effective sample size
- Measuring child development is costly in terms of time/money because each variable is constructed from multiple survey questions, and modules are administered one-on-one

Best Subset Selection

A **best subset selection** algorithm:

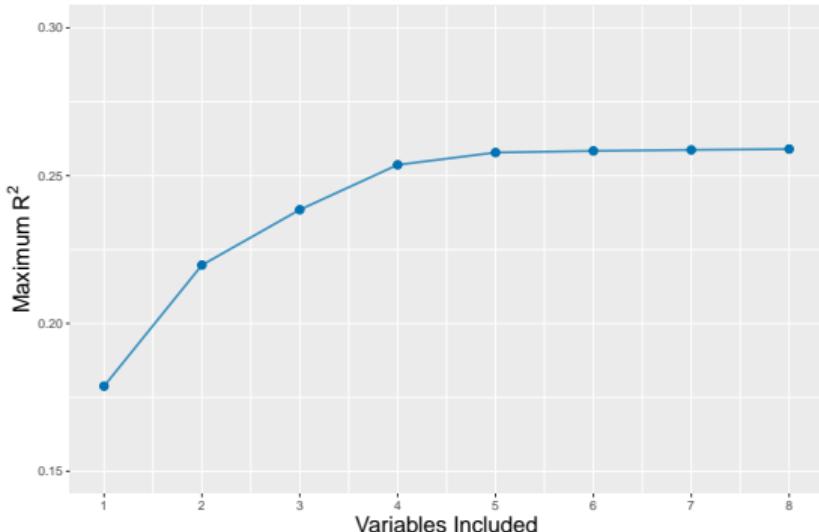
- For each number of possible covariates $k = 1, 2, \dots, p$,
 - ▶ Fit all models containing exactly k covariates
 - ▶ Identify the “best” in terms of R^2
- Choose the best subset using cross-validation or an alternative approach
 - ▶ Need to address the fact that R^2 always increases with k

Best Subset Selection Example: EMERGE

Use $N = 1,000$ data set on child development outcomes from EMERGE project

- literacy: measure of early literacy based on Early Grade Reading Assessment
- age_months: child age in months at time of survey
- male: dummy for boys
- haz: height-for-age z-score, measure of nutritional status
- receptive: receptive vocabulary, i.e. the ability to understand words (z-score)
- expressive: expressive vocabulary, i.e. the ability to produce words (z-score)
- fine_motor: fine motor skills (z-score)
- hh_size: household size
- mom_educ: mother's years of schooling

R^2 Is Increasing in the Number of Covariates



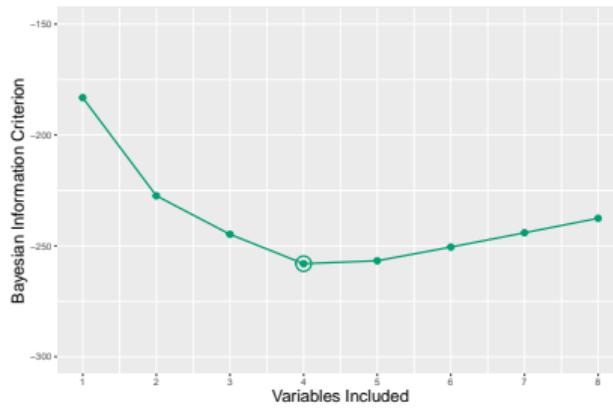
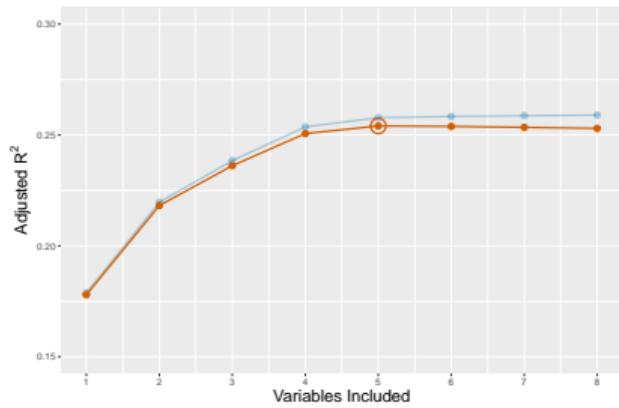
Variable	Number of Covariates							
	1	2	3	4	5	6	7	8
age_months					X	X	X	
male	X	X	X	X	X	X	X	X
haz		X	X	X	X	X	X	X
receptive				X	X	X	X	
expressive	X	X	X	X	X	X	X	X
fine_motor								X
hh_size						X	X	
mom_educ		X	X	X	X	X		

Choosing the Number of Covariates: Alternatives to Cross-Validation

Three alternatives to R^2 that adjust for the number of covariates in the specification, d

- Adjusted R^2 : $1 - \frac{RSS(n-d-1)}{TSS(n-1)}$ (seek to maximize)
- Akaike Information Criterion (AIC): $(RSS + 2d\hat{\sigma}^2) / n$ (seek to minimize)
- Bayesian Information Criterion (BIC): $(RSS + \ln(d)\hat{\sigma}^2) / n$ (seek to minimize)

Choosing the Number of Covariates: Alternatives to Cross-Validation



Best Subset Selection Is an Extension to OLS

In OLS, we seek to minimize:

$$\sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2$$

Best subset selection can be expressed as: choose β to minimize

$$\sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2 \text{ subject to } \sum_{j=1}^p I(\beta_j \neq 0) \leq s$$

where s is the number of regressors/predictors/features/covariates

Best Subset Selection Is Not Feasible with Many Covariates

Best subset selection is an extension to OLS that is solved algorithmically, not analytically

- When p is large, finding the best subset is computationally impossible ($2^P - 1$ regressions)
 - ▶ With 8 possible covariates: 255 regressions
 - ▶ With 20 possible covariates: over one million regressions
- Best subset selection makes sense when you can narrow the set of potential controls
 - ▶ Surveys often contain hundreds of questions
- Less computationally-intensive alternatives (forward and backward stepwise selection) exist but they are not robust to all patterns of correlation among potential covariates
 - ▶ Stepwise approaches involve adding (forward selection) or dropping (backward selection) the variable that gives the largest increase (forward) or smallest decrease (backward) in R^2

Shrinkage Operators: Machine Learning Extensions to OLS



Machine learning **shrinkage operators** (ridge regression, lasso) extend OLS to better predict Y

- Basic idea is to fully “kitchen sink” our regressions while proactively correcting for potential over-fitting, allowing us to leverage info from more covariates effectively

Lasso is attractive because it identifies a subset of X s that are most effective predictors of Y

Can We Improve on OLS?

A standard linear model may not be the best way to predict Y :

$$Y = \beta_0 + \beta_1 X_1 + \dots + \beta_p X_p + \varepsilon$$

Can we improve on OLS?

- When p is large relative to N , OLS is prone to over-fitting
- OLS explains both structural and spurious relationships in data

Like best subset selection, shrinkage operators minimize RSS subject to an additional constraint

$$\min_{\beta} \sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2 \text{ subject to } f(\beta) \leq s$$

Ridge Regression

Ridge regression solves the minimization problem:

$$\min_{\beta} \sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2 \text{ subject to } \sum_{j=1}^p \beta_j^2 \leq s$$

or, equivalently,

$$\min_{\beta} \sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2 + \lambda \sum_{j=1}^p \beta_j^2$$

for some **tuning parameter** $\lambda \geq 0$

Ridge regression shrinks OLS coefficients toward zero

- Shrinkage is more or less proportional, so ridge regression does not identify a subset of regressors to include in the regression model (it just down-weights some relative to others)

Shrinkage Operators: What's in a Name?

Like OLS, ridge regression has an analytical solution, as we can see in the $p = 1$ case:

$$\hat{\beta}_{OLS} = \frac{\sum_{i=1}^n x_i y_i - N\bar{X}\bar{Y}}{\sum_{i=1}^n x_i^2 - N\bar{X}^2} > \frac{\sum_{i=1}^n x_i y_i - N\bar{X}\bar{Y}}{\sum_{i=1}^n x_i^2 - N\bar{X}^2 + 2\lambda} = \hat{\beta}_{ridge}$$

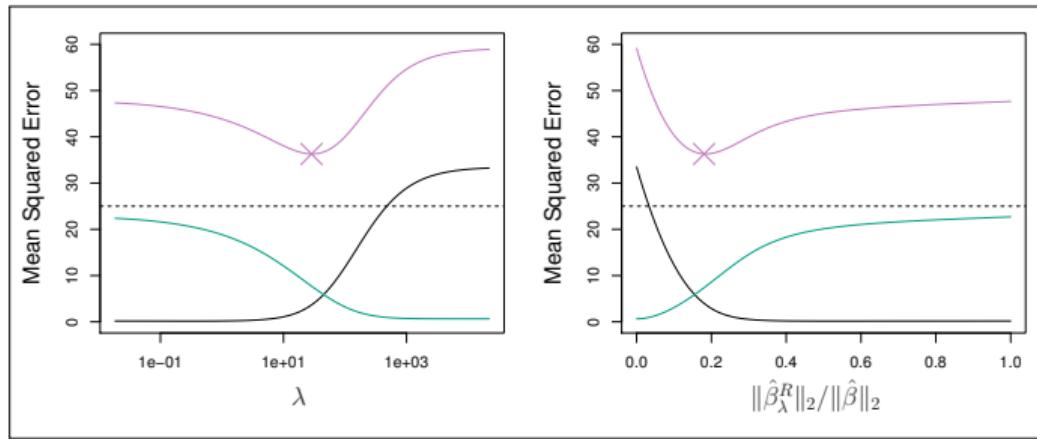
The (bivariate) ridge regression coefficient is smaller than the (bivariate) OLS coefficient

- When λ is close to 0, $\hat{\beta}_{ridge}$ is similar to $\hat{\beta}_{OLS}$
- $\hat{\beta}_{ridge}$ approaches 0 as λ gets large

With more than one independent variable, some ridge regression coefficients may be larger than OLS counterparts, and the coefficient on a specific X_k need not decline monotonically with λ

- Shrinkage is more or less proportional, so ridge regression does not identify a subset of regressors to include in the regression model (it just down-weights some relative to others)

OLS is BLUE, But Ridge Regression (Sometimes) Has Lower MSE



Source: James et al. (2021)

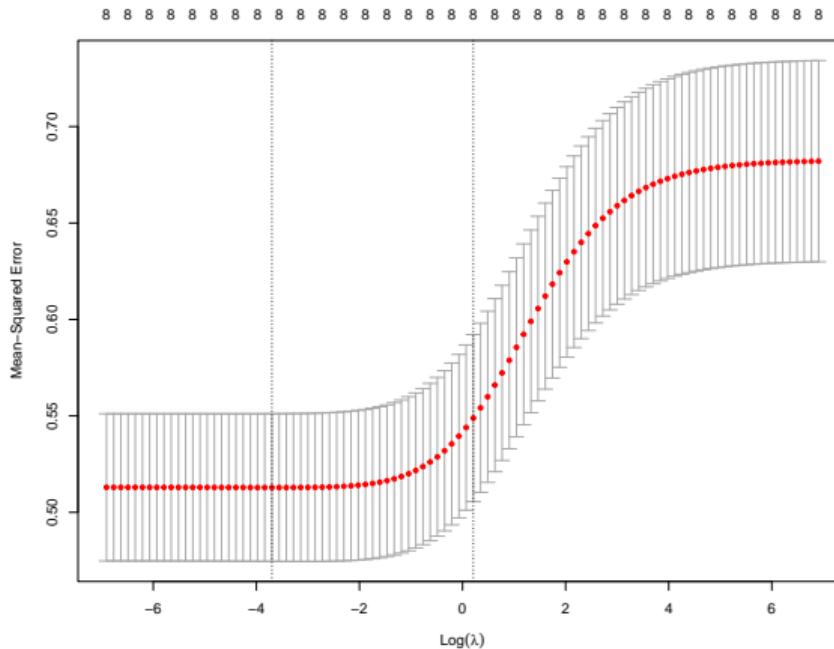
Gauss-Markov Theorem: OLS is the **best linear unbiased estimator** (BLUE) of Y

- Ridge regression is biased (black line), but has lower variance relative to the true underlying β (green line) and can therefore achieve lower MSE (pink line) for some λ s

Ridge Regression in Practice

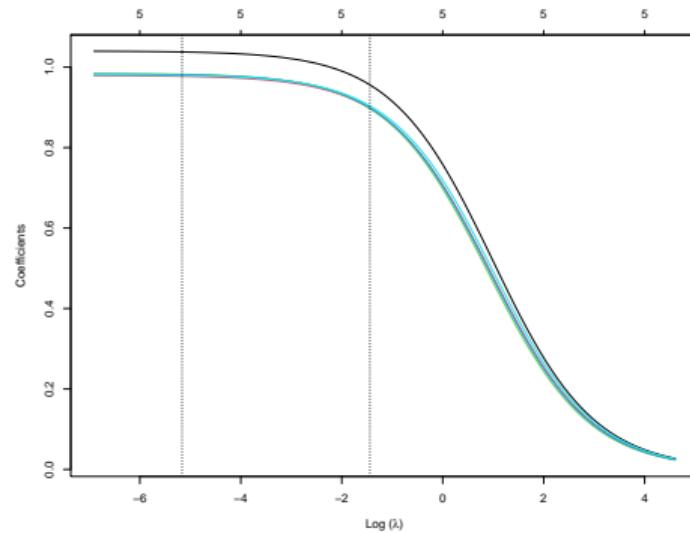
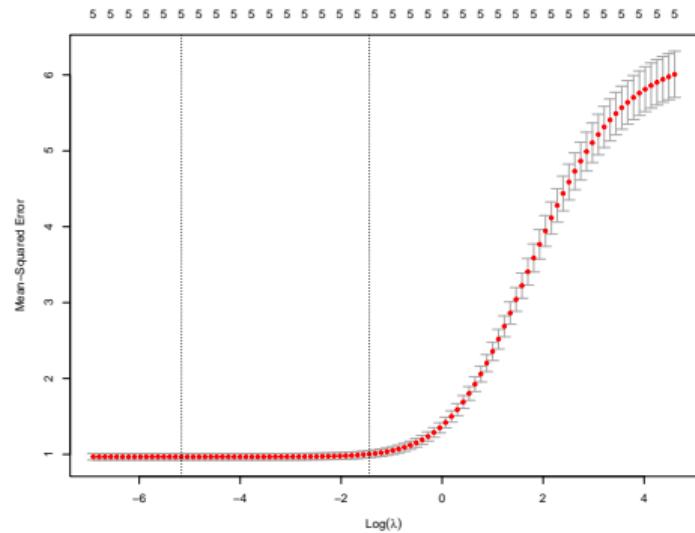
Variable	OLS (1)	Ridge Regression		
		$\lambda = 10^{-2}$ (2)	$\lambda = 10$ (3)	$\lambda = 10^4$ (4)
expressive	0.2543	0.2498	0.0247	0.0003
male	-0.3152	-0.3106	-0.0203	-0.0002
haz	0.0847	0.0844	0.0130	0.0002
mom_educ	0.0439	0.0436	0.0051	0.0001
receptive	0.0651	0.0671	0.0195	0.0002
age_months	0.0024	0.0024	0.0002	0.0000
hh_size	-0.0085	-0.0084	-0.0011	0.0000
fine_motor	0.0257	0.0269	0.0160	0.0002

Choosing the Penalty Parameter to Minimize Test MSE: EMERGE Data



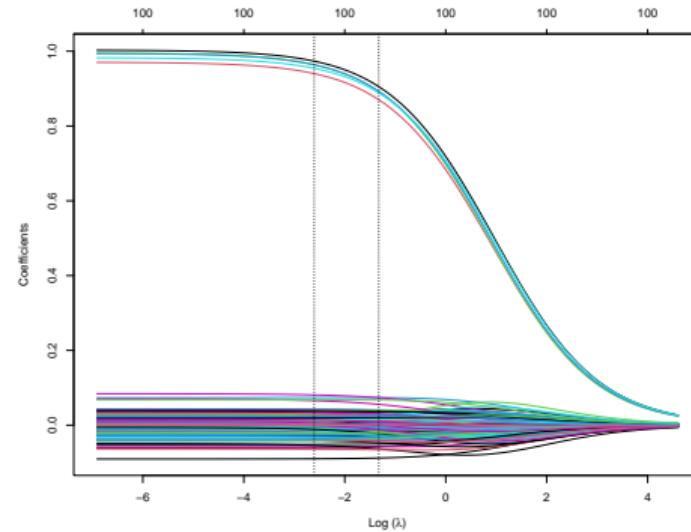
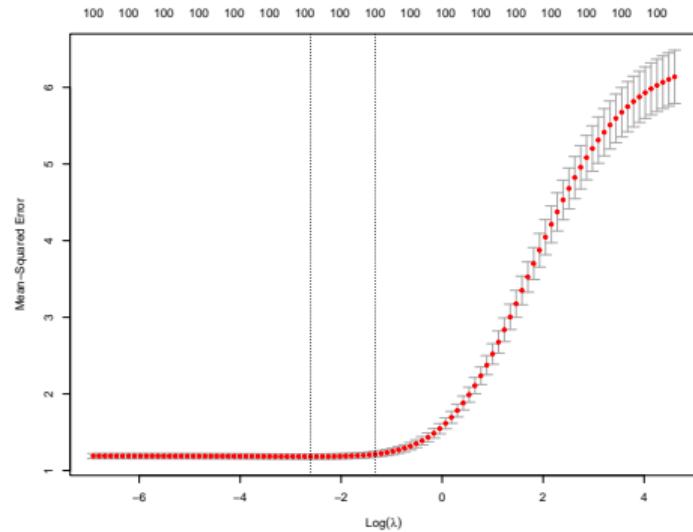
data source: EMERGE project (Jakiela, Ozier, Fernald, and Knauer 2021)

Ridge Regression in Simulated Data: $N = 1000, K = 5$



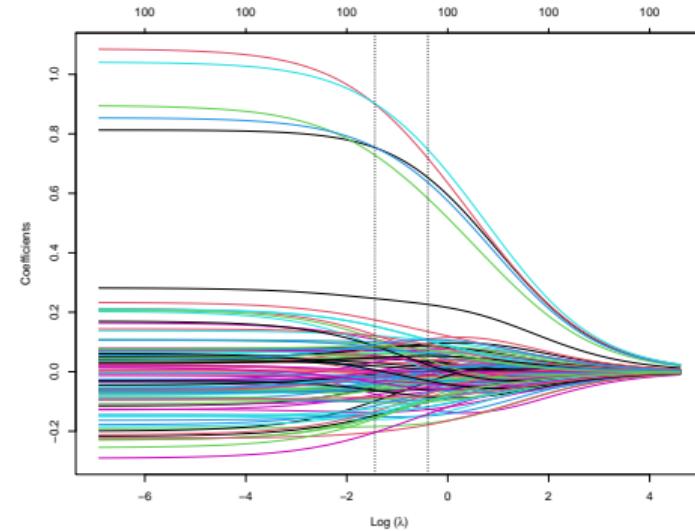
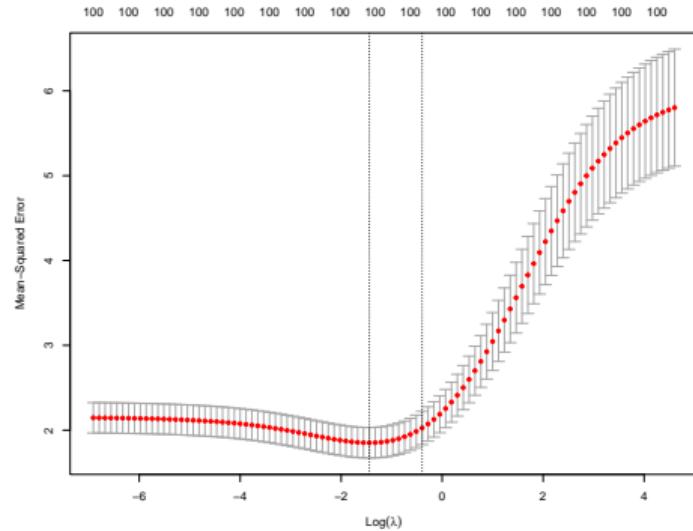
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 5$, $\varepsilon \sim N(0, 1)$, $N = 1000$, $K = 5$

Ridge Regression in Simulated Data: $N = 1000$, $K = 100$



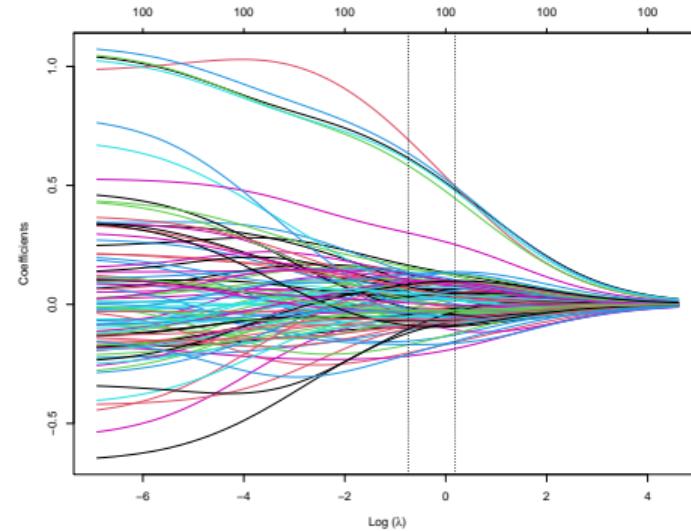
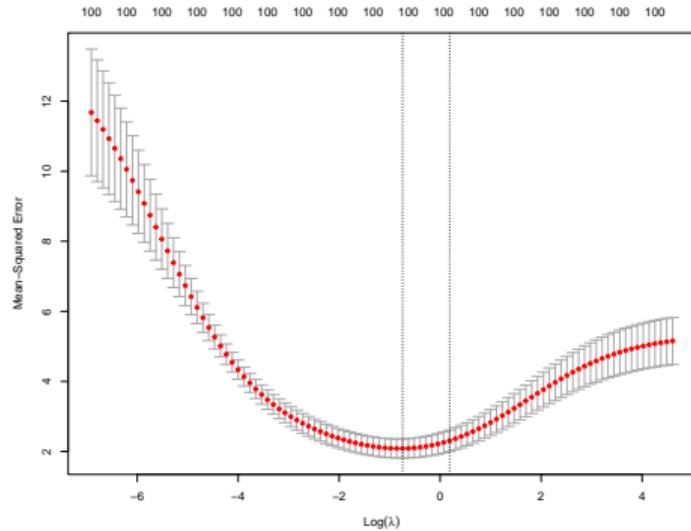
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 1000$, $K = 100$

Ridge Regression in Simulated Data: $N = 200$, $K = 100$



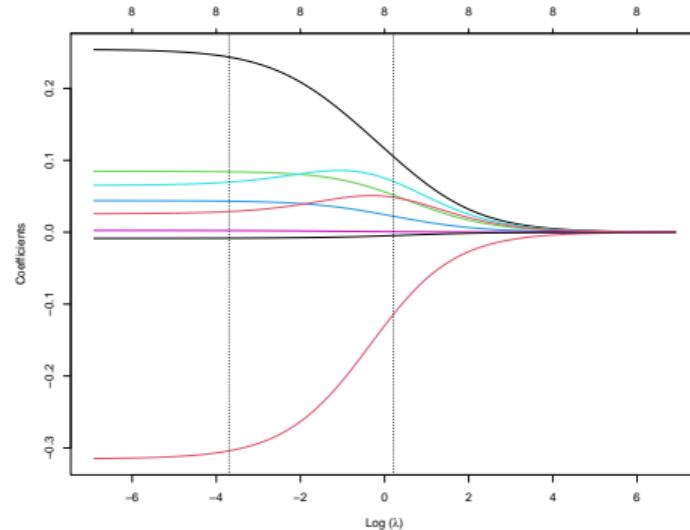
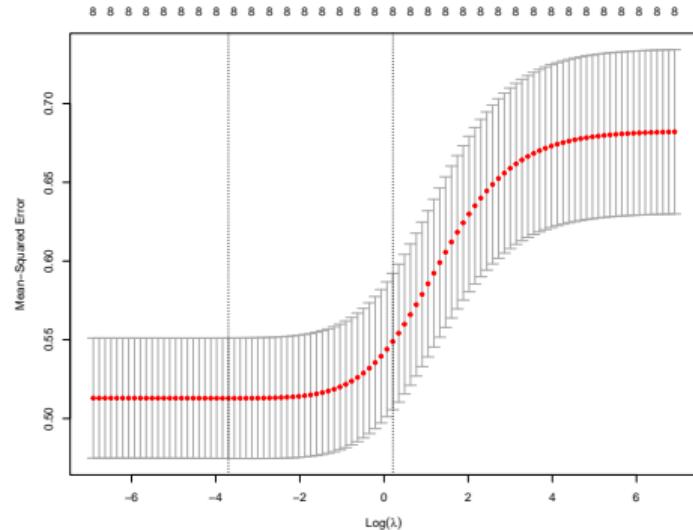
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 200$, $K = 100$

Ridge Regression in Simulated Data: $N = 120, K = 100$



data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 120$, $K = 100$

Ridge Regression in the EMERGE Data



data source: EMERGE project (Jakiela, Ozier, Fernald, and Knauer 2021)

Ridge Regression in Practice: Comparing MSEs in EMERGE Data

Splitting the data into a training data set and a test data set, we see that ridge reduces the MSE in the test data as expected, but not by much (relative to the SD of the outcome, 0.8258)

OLS	λ^*	λ^{1SE}
0.4928	0.4899	0.5930

λ^* is the λ that minimizes test MSE in cross-validation, λ^{1SE} is 1 SE higher than λ^*

Shrinkage Operators: Lasso

Lasso (Least Absolute Shrinkage and Selection Operator) seeks to minimize:

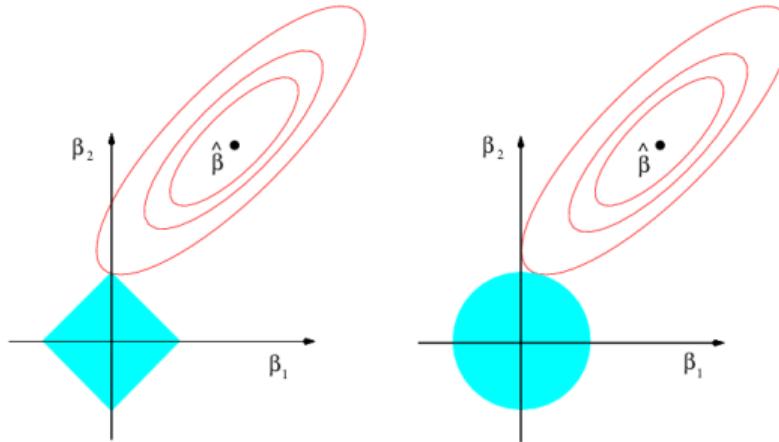
$$\min_{\beta} \sum_{i=1}^n \left(y_i - \beta_0 - \sum_{j=1}^p \beta_j x_{ij} \right)^2 + \lambda \sum_{j=1}^p |\beta_j|$$

for some **tuning parameter** $\lambda \geq 0$

Lasso combines benefits of subset selection, ridge regression; useful for choosing covariates

- Less computationally intensive than subset selection
- Sets some coefficients to 0 → identifies parsimonious model

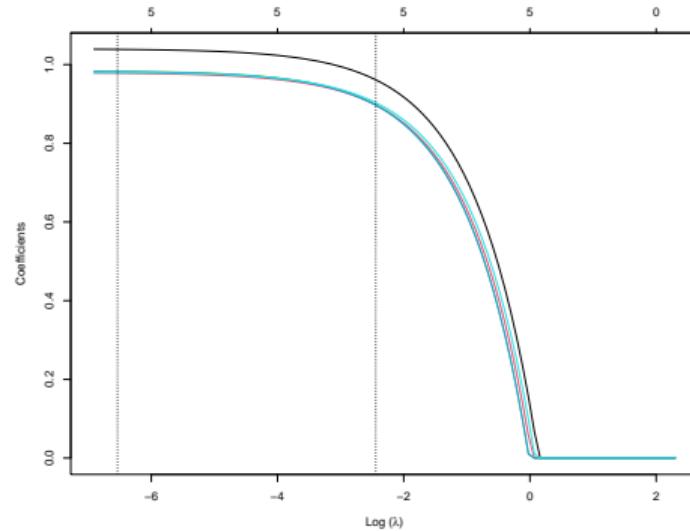
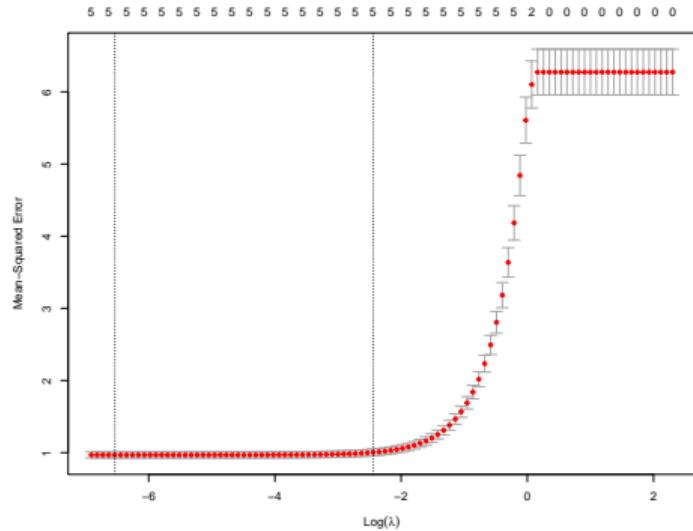
Lasso Sets Some Coefficients to Zero



Source: James et al. (2021)

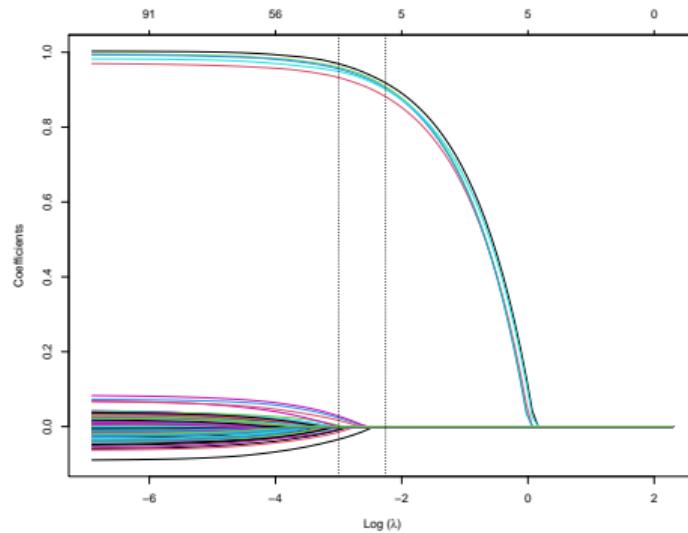
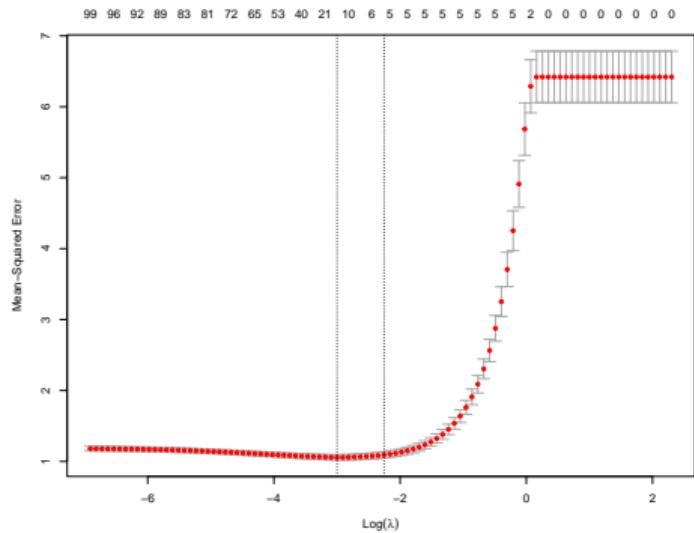
The lasso constraint region has sharp corners \Rightarrow some coefficients set to 0

Lasso in Simulated Data: $N = 1000$, $K = 5$



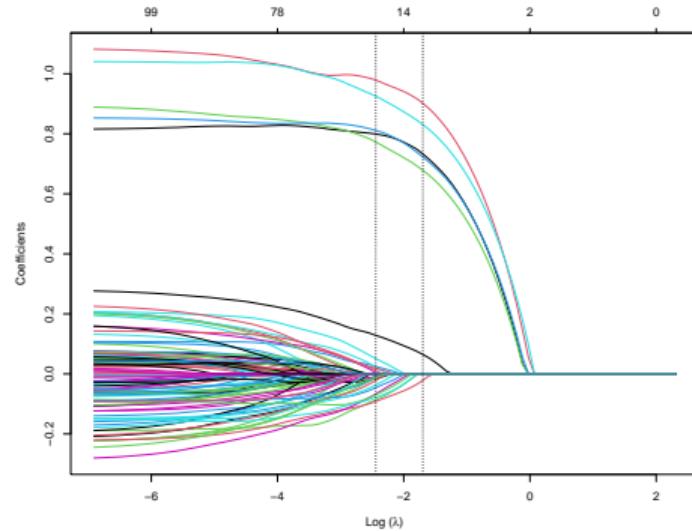
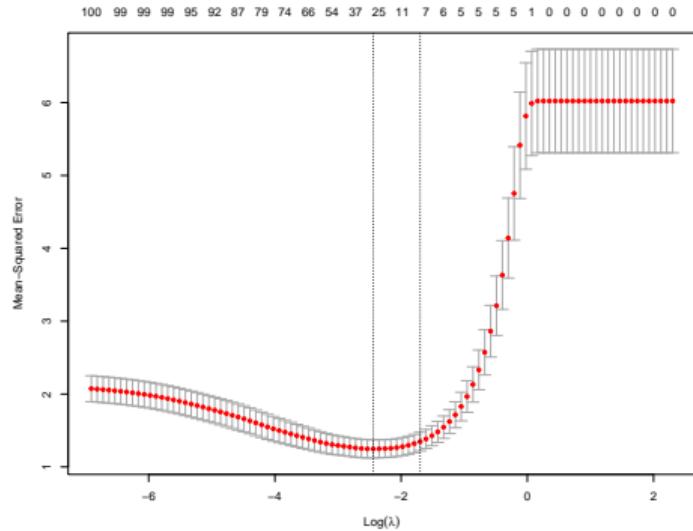
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 5$, $\varepsilon \sim N(0, 1)$, $N = 1000$, $K = 5$

Lasso in Simulated Data: $N = 1000$, $K = 100$



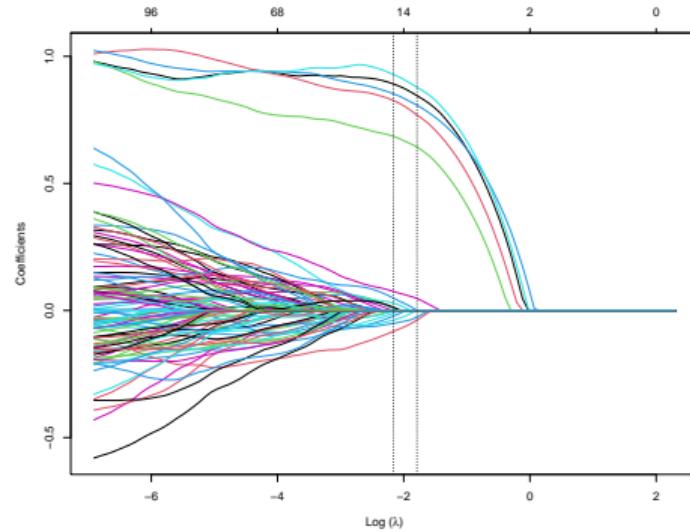
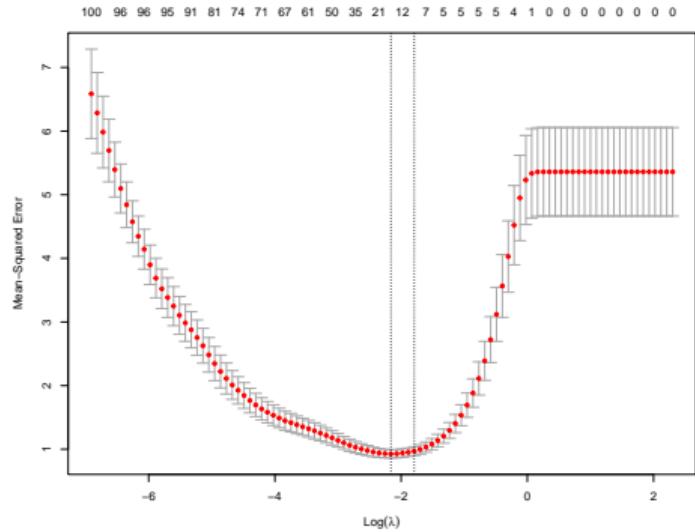
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 1000$, $K = 100$

Lasso in Simulated Data: $N = 200$, $K = 100$



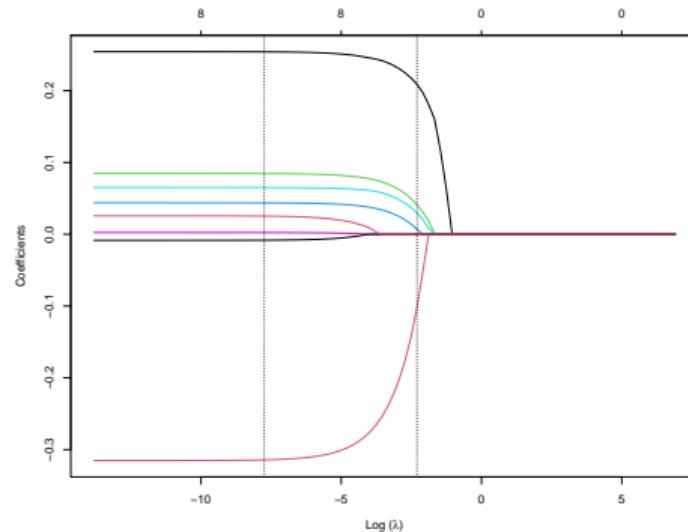
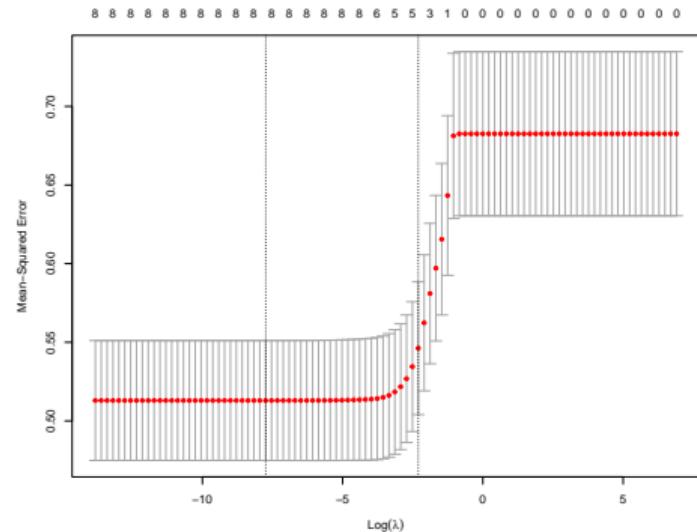
data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 200$, $K = 100$

Lasso in Simulated Data: $N = 120$, $K = 100$



data-generating process: $Y = \sum_{k=1}^5 X_k + \varepsilon$ where $X_k \sim N(0, 1)$ for $k = 1, \dots, 100$, $\varepsilon \sim N(0, 1)$, $N = 120$, $K = 100$

Lasso in Practice: EMERGE Data



data source: EMERGE project (Jakiela, Ozier, Fernald, and Knauer 2021)

Alternative “Data-Driven” Approach to Choosing λ

Belloni and Chernozhukov (2011), Belloni et al. (2012): alternative approach to choosing λ

- Chooses λ iteratively based on data, penalties vary across variables
- Errs on the side of choosing fewer controls to avoid over-fitting
- Allows for heteroskedasticity
- Designed to allow for valid post-selection lasso estimation (within a single data set)

Approaches may generate different sets of controls

- Costs of too many/too few may vary across empirical contexts

In $N = 120$, $K = 100$ simulated data, data-driven lasso $\Rightarrow 9 X$ variables selected

Comparing Approaches to Choosing Covariates via Lasso

Variable	OLS	λ^*	λ^{1SE}	λ^{DD}
age_months	X	X		
male	X	X	X	X
haz	X	X	X	X
receptive	X	X	X	X
expressive	X	X	X	X
fine_motor	X	X		X
hh_size	X	X		
mom_educ	X	X	X	X

Summary

Best subset selection, ridge regression, and lasso are constrained extensions to OLS

- Ridge and lasso are regularized: coefficients are shrunk toward zero to reduce over-fitting
- Best subset selection and lasso are useful for model selection (i.e. choosing covariates)

Lasso is now widely used by economists to choose a subset of (many) controls to include in OLS

- Number of controls selected depends on the penalty (or tuning) parameter
 - ▶ Cross-validation is optimizing prediction, leads to the inclusion of more controls
 - ▶ Data-driven approach of Belloni et al. (2012) or 1 SE rule typically better heuristics
- Desired number of controls may also depend on the cost of adding/including a variable
 - ▶ Expressive vocabulary, male dummy both predict emergent literacy in EMERGE data, but measuring expressive vocabulary probably costs thousands of times more per child

Epilogue: Covariate Selection via Post-Double-Selection (PDS) Lasso

EMERGE example focuses on choosing covariates **in advance**, deciding what to measure

- The error terms in the covariate selection process are independent of eventual analysis

Model selection – choosing covariates using your analysis sample – is more complicated

- Conducting statistical inference after covariate selection within a sample is complicated
- Post-double selection lasso is a widely used approach that addresses inference concerns
 - ▶ Step 1: run lasso to choose covariates that predict outcome Y
 - ▶ Step 2: run lasso to choose covariates that predict treatment of interest T
 - ▶ Step 3: run OLS including all covariates identified in first two steps, conduct inference

Lab # 6

Objective: use lasso to choose covariates in the EMERGE data

- ECON370-lab6-data.csv contains 200 observations and 25 variables
- You will create dummies for survey enumerator and randomization strata

Two approaches to choosing tuning/penalty parameter: CV and “data-driven” approach

- CV possible in both R and Python, “data-driven” approach only possible in R (?)
- Use CV to identify both MSE-minimizing tuning parameter and 1 SE alternative
 - ▶ 1 SE rule: find the SE of the MSE (across k folds) at the minimum, then find largest value of the tuning parameter that yields a test MSE below the sum of the minimum MSE + the SE
- Set of selected covariates will depend (a lot!) on your choice of tuning parameter