Is the test error unbiased for these programs?

```
# Given dataset of 1000-by-50 feature
# matrix X, and 1000-by-1 labels vector
mu = np.mean(X, axis=0)
X = X - mu

idx = np.random.permutation(1000)
TRAIN = idx[0:900]
TEST = idx[900::]

ytrain = y[TRAIN]
Xtrain = X[TRAIN,:]

# Solve for argmin_w ||Xtrain*w - ytrain||_2
w = np.linalg.solve( np.dot(Xtrain.T, Xtrain),
                     np.dot(Xtrain.T, ytrain) )
b = np.mean(ytrain)

ytest = y[TEST]
Xtest = X[TEST,:]

train_error = np.dot( np.dot(Xtrain, w)+b - ytrain,
                      np.dot(Xtrain, w)+b - ytrain )/len(TRAIN)
test_error = np.dot( np.dot(Xtest, w)+b - ytest,
                     np.dot(Xtest, w)+b - ytest )/len(TEST)

print('Train error = ',train_error)
print('Test error = ',test_error)
```

Xtest is bleeding into w through μ

NO

```
# Given dataset of 1000-by-50 feature
# matrix X, and 1000-by-1 labels vector
idx = np.random.permutation(1000)
TRAIN = idx[0:900]
TEST = idx[900::]

ytrain = y[TRAIN]
Xtrain = X[TRAIN,:]
Xtrain_avg = np.mean(Xtrain, axis=0)
Xtrain = Xtrain - Xtrain_avg

# Solve for argmin_w ||Xtrain*w - ytrain||_2
w = np.linalg.solve( np.dot(Xtrain.T, Xtrain),
                     np.dot(Xtrain.T, ytrain) )
b = np.mean(ytrain)

ytest = y[TEST]
Xtest = X[TEST,:]
Xtest_avg = np.mean(Xtest, axis=0)
Xtest = Xtest - Xtest_avg
                Xtrain_avg
train_error = np.dot( np.dot(Xtrain, w)+b - ytrain,
                      np.dot(Xtrain, w)+b - ytrain )/len(TRAIN)
test_error = np.dot( np.dot(Xtest, w)+b - ytest,
                     np.dot(Xtest, w)+b - ytest )/len(TEST)

print('Train error = ',train_error)
print('Test error = ',test_error)
```

Is the test error unbiased for this program?

```
# Given dataset of 1000-by-50 feature
# matrix X, and 1000-by-1 labels vector
idx = np.random.permutation(1000)
TRAIN = idx[0:800]
VAL = idx[800:900]
TEST = idx[900::]

ytrain = y[TRAIN]
Xtrain = X[TRAIN,:]
yval = y[VAL]
Xval = X[VAL,:]

err = np.zeros(50)
for d in range(1,51):
    w, b = fit(Xtrain[:,0:d], ytrain)
    yval_hat = predict(w, b, Xval[:,0:d])
    err[d-1] = np.mean((yval_hat-yval)**2)
d_best = np.argmin(err)+1

Xtot = np.concatenate((Xtrain, Xval), axis=0)
ytot = np.concatenate((ytrain, yval), axis=0)
w, b = fit(Xtot[:,0:d_best], ytot)

ytest = y[TEST]
Xtest = X[TEST,:]

ytot_hat = predict(w, b, Xtot[:,0:d_best])
tot_train_error = np.mean((ytot_hat-ytot)**2)
ytest_hat = predict(w, b, Xtest[:,0:d_best])
test_error = np.mean((ytest_hat-ytest)**2)

print('Train error = ',train_error)
print('Test error = ',test_error)
```

```
def fit(Xin, Yin):
    mu = np.mean(Xin, axis=0)
    Xin = Xin - mu
    w = np.linalg.solve( np.dot(Xin.T, Xin),
                         np.dot(Xin.T, Yin) )
    b = np.mean(Yin) - np.dot(w, mu)
    return w, b

def predict(w, b, Xin):
    return np.dot(Xin, w)+b
```

Cross-Validation

How… How… How???????

- How do we pick the regularization constant λ…
- How do we pick the number of basis functions…

- We could use the test data, but…

How… How… How???????

- How do we pick the regularization constant λ…
- How do we pick the number of basis functions…

- We could use the test data, but…

  - Never ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever ever train on the test data

(LOO) Leave-one-out cross validation

- **Consider a validation set with 1 example:**
  - **D – training data** $D = \{(x_i, y_i)\}_{i=1}^{n}$
  - **D\j – training data with j th data point (x_j ,y_j) moved to validation set** $D/j = \{(x_1, y_1) \ldots, (x_{j-1}, y_{j-1}), (x_{j+1}, y_{j+1}) \ldots (x_n)$
- **Learn classifier $f_{D\setminus j}$ with $D\setminus j$ dataset**
- **Estimate true error** as squared error on predicting $\mathbf{y_j}$:
  - **Unbiased estimate of $error_{true}(f_{D\setminus j})$!**

- **LOO cross validation**: Average over all data points *j*:
  - **For each data point you leave out, learn a new classifier $f_{D\setminus j}$**
  - **Estimate error** as:

(LOO) Leave-one-out cross validation

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  - **D – training data**
  - **D\j – training data with j th data point (x_j ,y_j) moved to validation set**
- **Learn classifier $f_{D\setminus j}$ with D\j dataset**
- **Estimate true error** as squared error on predicting $\mathbf{y_j}$:
  - **Unbiased estimate of $\text{error}_{true}(f_{D\setminus j})$!**

- **LOO cross validation**: Average over all data points *j*:
  - **For each data point you leave out, learn a new classifier $f_{D\setminus j}$**
  - **Estimate error** as:

$$\text{error}_{LOO} = \frac{1}{n}\sum_{j=1}^{n}(y_j - f_{\mathcal{D}\setminus j}(x_j))^2$$

LOO cross validation is (almost) unbiased estimate!

- When computing LOOCV error, we only use *N-1* data points
  - So it's not estimate of true error of learning with *N* data points
  - Usually pessimistic, though – learning with less data typically gives worse answer

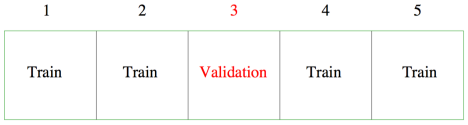
- LOO is almost unbiased! Use LOO error for model selection!!!
  - E.g., picking λ

Computational cost of LOO

- Suppose you have 100,000 data points
- You implemented a great version of your learning algorithm
  - Learns in only 1 second
- Computing LOO will take about 1 day!!!
  -

Use *k*-fold cross validation

- Randomly divide training data into *k* equal parts
  - $D_1,\ldots,D_k$
- For each *i*
  - Learn classifier $f_{D\setminus Di}$ using data point not in D_i
  - Estimate error of $f_{D\setminus Di}$ on validation set D_i:



$$\text{error}_{\mathcal{D}_i} = \frac{1}{|\mathcal{D}_i|} \sum_{(x_j,y_j)\in\mathcal{D}_i} (y_j - f_{\mathcal{D}\setminus\mathcal{D}_i}(x_j))^2$$

Use *k*-fold cross validation

- **Randomly divide training data into *k* equal parts**
  - **$D_1,\ldots,D_k$**
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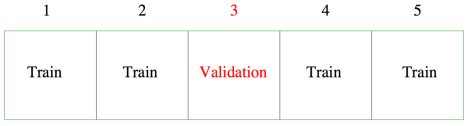


$$\text{error}_{\mathcal{D}_i} = \frac{1}{|\mathcal{D}_i|} \sum_{(x_j,y_j)\in\mathcal{D}_i} (y_j - f_{\mathcal{D}\backslash\mathcal{D}_i}(x_j))^2$$

- **k-fold cross validation error is average** over data splits:

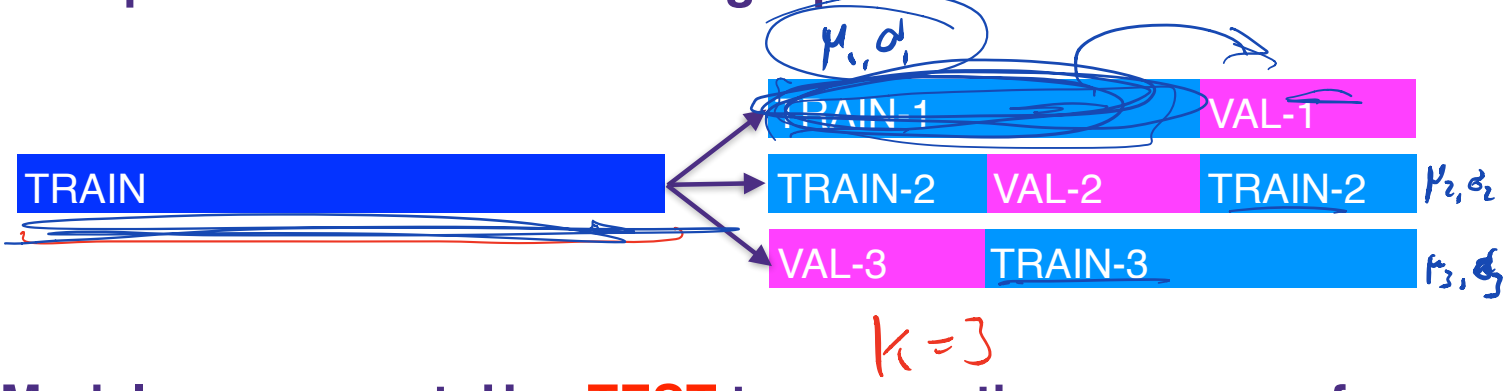
$$error_{k-fold} = \frac{1}{k}\sum_{i=1}^{k} error_{\mathcal{D}_i}$$

- **k-fold cross validation properties:**
  - **Much faster to compute** than LOO
  - **More (pessimistically) biased** – using much less data, only *n(k-1)/k*
  - **Usually, k = 10**

Recap

- Given a dataset, begin by splitting into

- Model selection: Use k-fold cross-validation on **TRAIN** to train predictor and choose magic parameters such as λ



- Model assessment: Use **TEST** to assess the accuracy of the model you output
  - **Never ever ever ever ever train or choose parameters based on the test data**

Example 1

- You wish to predict the stock price of zoom.us given historical stock price data
- You use all daily stock price up to Jan 1, 2020 as **TRAIN** and Jan 2, 2020 - April 13, 2020 as **TEST**
- What's wrong with this procedure?

#1 Rule of ML:
Train and Test set are drawn i.i.d.

$x_{t,i}$ = ith stock price at time t.

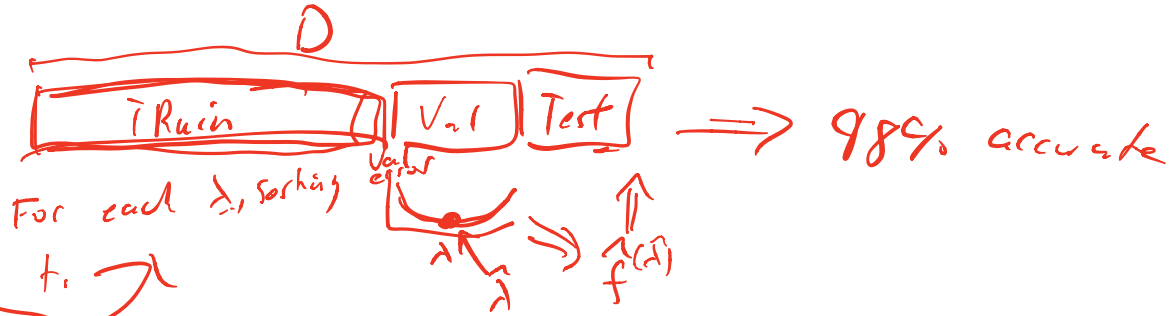
$$x_{t+1,i} = \sum_{s=0}^{k} a_s x_{t-s,i} + \underbrace{\varepsilon_{t+1,i}}_{\mathcal{N}(0,1)}$$

Example 2

- **Given 10,000-dimensional data and n examples, we pick a subset of 50 dimensions that have the highest correlation with labels in the training set:**

50 indices j that have largest $\frac{|\sum_{i=1}^{n} x_{i,j} y_i|}{\sqrt{\sum_{i=1}^{n} x_{i,j}^2}}$

- **After picking our 50 features, we then use CV with the training set to train ridge regression with regularization λ**
- **What's wrong with this procedure?**



Recap

- Learning is…
  - Collect some data
    - E.g., housing info and sale price
  - Randomly split dataset into TRAIN, VAL, and TEST
    - E.g., 80%, 10%, and 10%, respectively
  - Choose a hypothesis class or model
    - E.g., linear with non-linear transformations
  - Choose a loss function
    - E.g., least squares with ridge regression penalty on TRAIN
  - Choose an optimization procedure
    - E.g., set derivative to zero to obtain estimator, cross-validation on VAL to pick num. features and amount of regularization
  - Justifying the accuracy of the estimate
    - E.g., report TEST error

Simple Variable Selection
LASSO: Sparse Regression

Sparsity

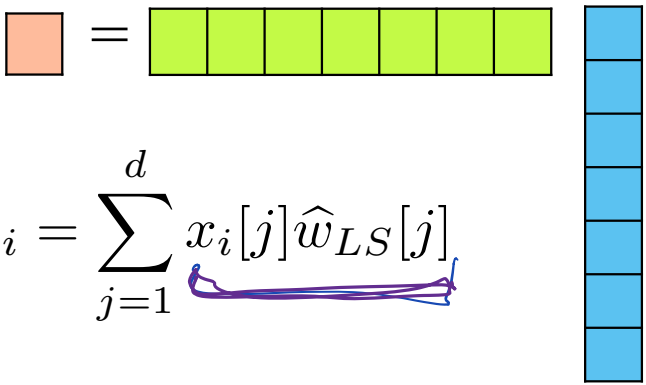
$$\widehat{w}_{LS} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2$$

- **Vector w is sparse, if many entries are zero**

Sparsity

$$\widehat{w}_{LS} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2$$

- **Vector w is sparse, if many entries are zero**
  - **Efficiency**: If size(**w**) = 100 Billion, each prediction is expensive:
    - **If w is sparse, prediction computation only depends on number of non-zeros**



$$\widehat{y}_i = \widehat{w}_{LS}^{\top} x_i = \sum_{j=1}^{d} x_i[j]\widehat{w}_{LS}[j]$$

Sparsity

$$\widehat{w}_{LS} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2$$

- **Vector w is sparse, if many entries are zero**
  - **Interpretability**: What are the relevant dimension to make a prediction?

- How do we find "best" subset among all possible?

Lot size
Single Family
Year built
Last sold price
Last sale price/sqft
Finished sqft
Unfinished sqft
Finished basement sqft
floors
Flooring types
Parking type
Parking amount
Cooling
Heating
Exterior materials
Roof type
Structure style

Dishwasher
Garbage disposal
Microwave
Range / Oven
Refrigerator
Washer
Dryer
Laundry location
Heating type
Jetted Tub
Deck
Fenced Yard
Lawn
Garden
Sprinkler System

Finding best subset: Exhaustive

- Try all subsets of size 1, 2, 3, … and one that minimizes validation error
- Problem?

Finding best subset: Greedy

Forward stepwise:
Starting from simple model and iteratively add features most useful to fit

Backward stepwise:
Start with full model and iteratively remove features least useful to fit

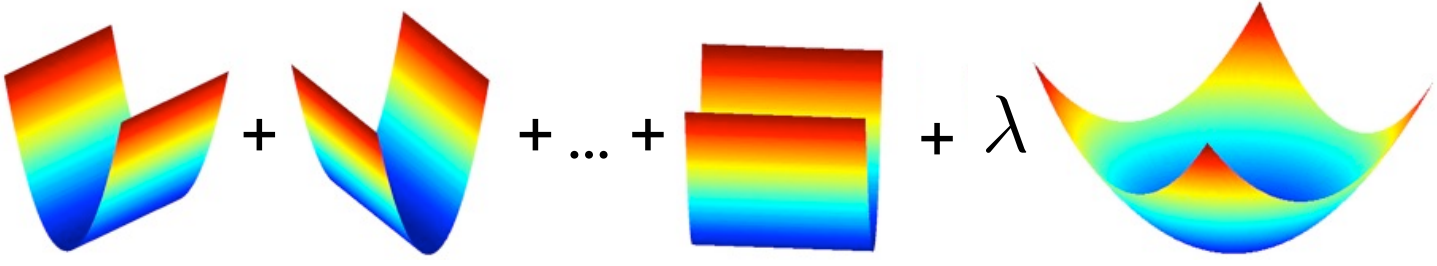
Combining forward and backward steps:
In forward algorithm, insert steps to remove features no longer as important

Lots of other variants, too.

Finding best subset: Regularize

Ridge regression makes coefficients small

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$



Finding best subset: Regularize

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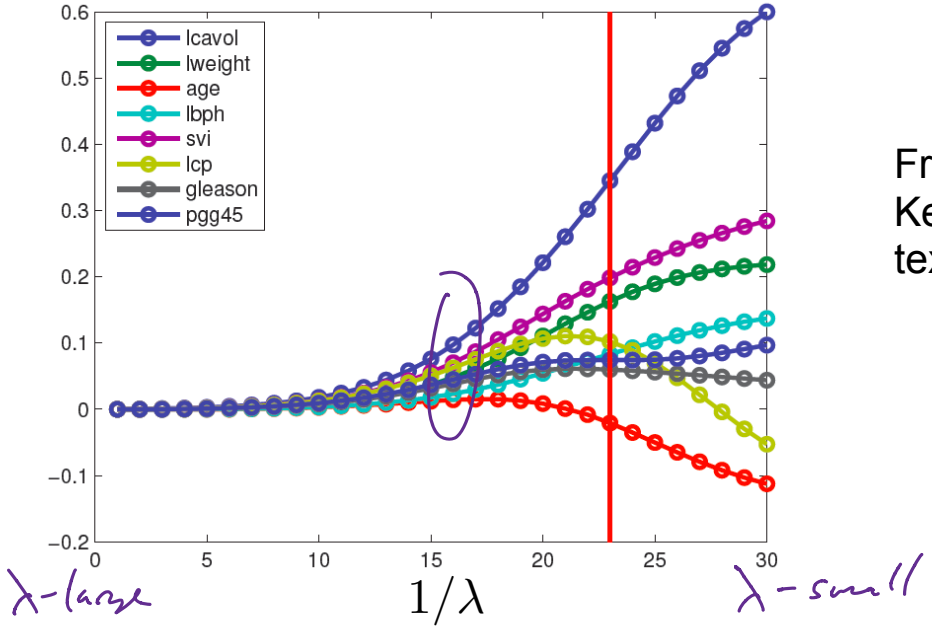
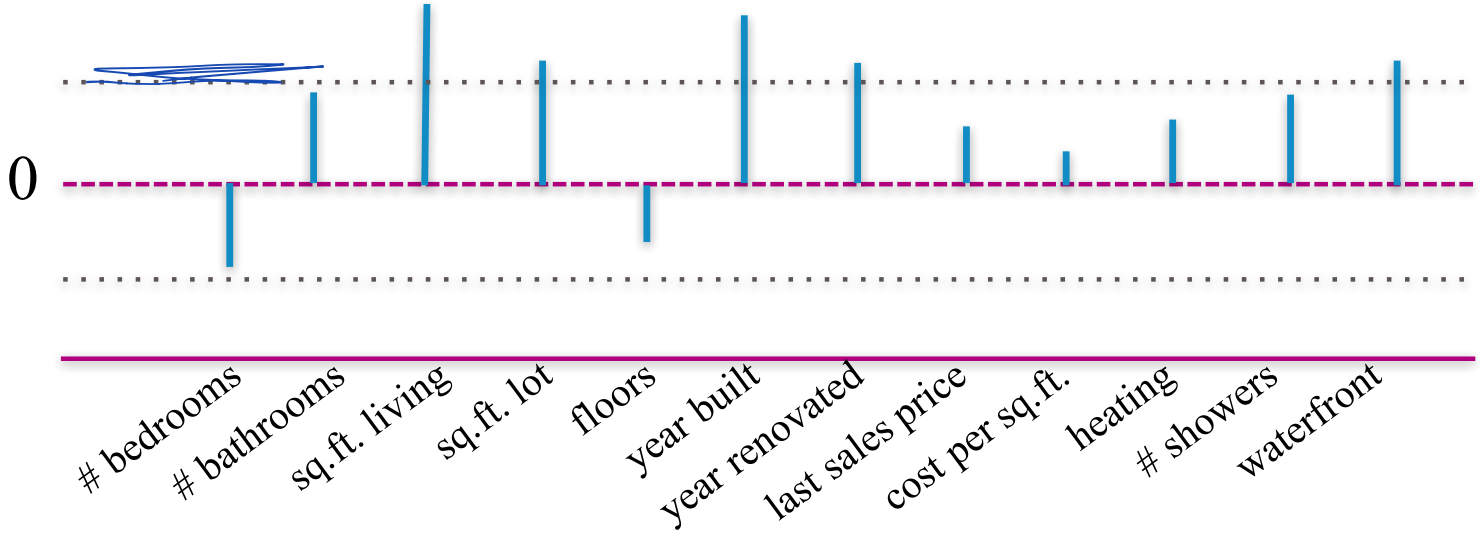


From
Kevin Murphy
textbook

Thresholded Ridge Regression

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$

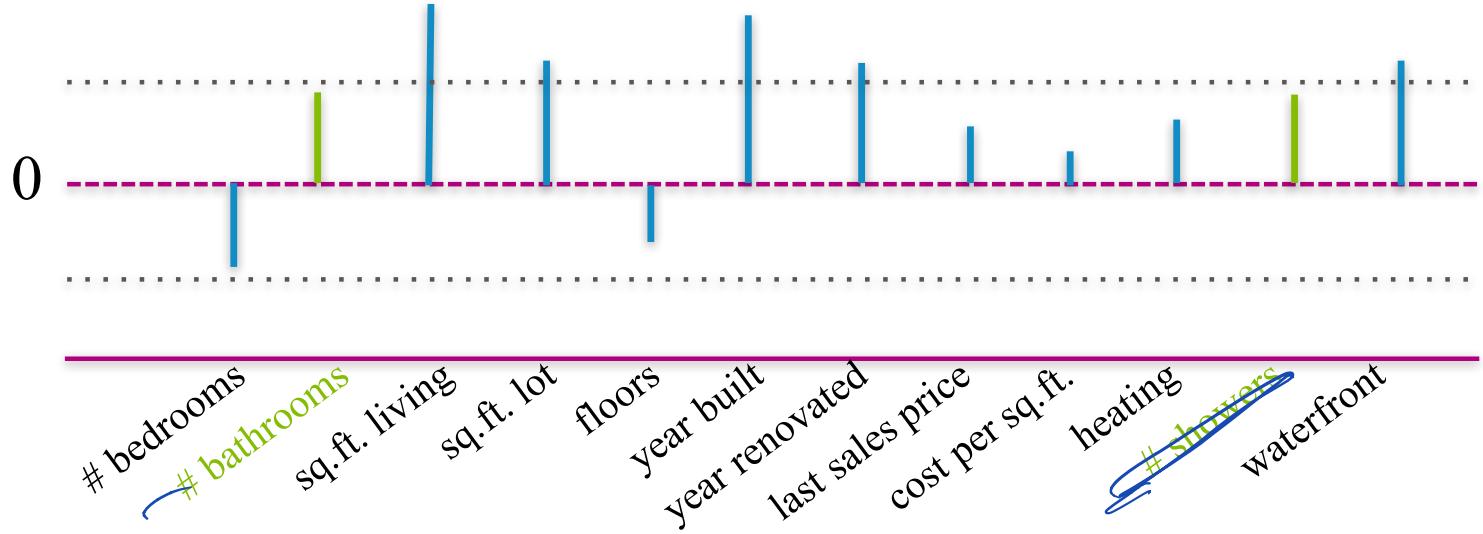
Why don't we just set **small** ridge coefficients to 0?



Thresholded Ridge Regression

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$

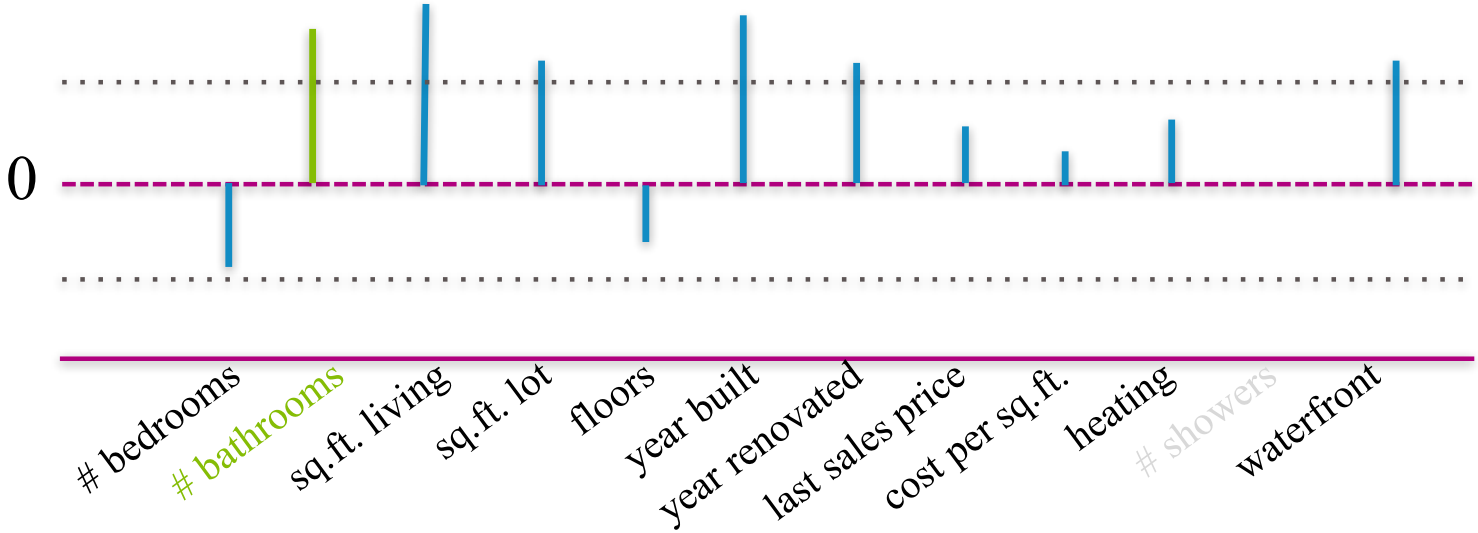
Consider two related features (bathrooms, showers)



Thresholded Ridge Regression

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$

What if we **didn't** include showers? Weight on bathrooms increases!

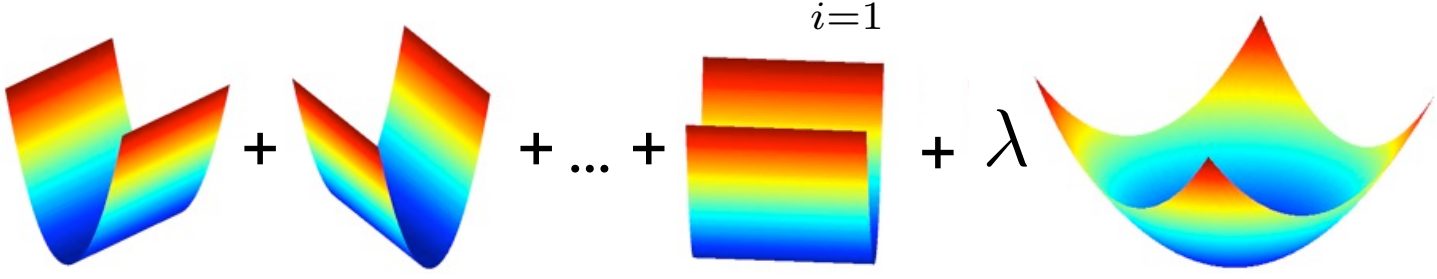


Can another regularizer perform selection automatically?

Recall Ridge Regression

- Ridge Regression objective:

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$



$$\|w\|_p = \left(\sum_{i=1}^{d} |w|^p\right)^{1/p}$$

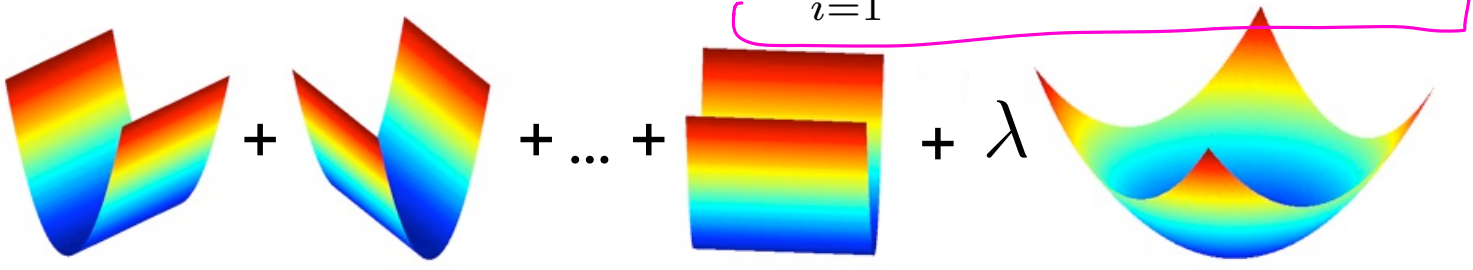
$p = 2$ $\left(\sum_{i=1}^{d} w_i^2\right)^{1/2}$

$p = 1$ $\sum_{i=1}^{d} |w_i|$

Ridge vs. Lasso Regression

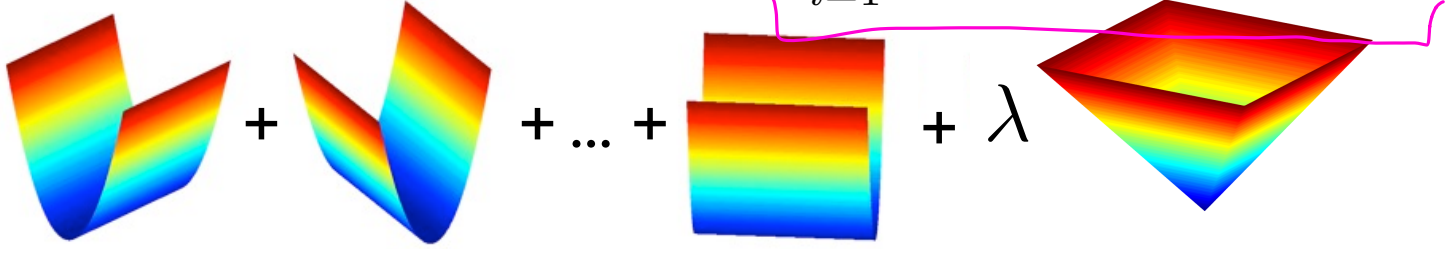
- Ridge Regression objective:

$$\widehat{w}_{ridge} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_2^2$$



- Lasso objective:

$$\widehat{w}_{lasso} = \arg\min_{w} \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda ||w||_1$$



Penalized Least Squares

$$\text{Ridge} : r(w) = ||w||_2^2 \qquad \text{Lasso} : r(w) = ||w||_1$$

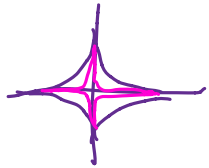
$$\widehat{w}_r = \arg\min_w \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda r(w)$$

regularizer

ℓ_0-norm: $\quad ||w||_0 = \sum_{i=1}^{d} \mathbb{1}\{w_i \neq 0\}$

$= \#\{\text{non-zeros in } w\}$

$= \lim_{p \to 0} \left(\sum_{i=1}^{d} |w_i|^p \right)^{1/p}$



Penalized Least Squares

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$$\widehat{w}_r = \arg\min_w \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 + \lambda r(w)$$

For any $\lambda \geq 0$ for which $\widehat{w}_r$ achieves the minimum, there exists a $\nu \geq 0$ such that

$$\widehat{w}_r = \arg\min_w \sum_{i=1}^{n} \left(y_i - x_i^T w\right)^2 \qquad \text{subject to } r(w) \leq \nu$$

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