Lecture 5 LASSO Regularization

EE-UY 4563/EL-GY 9123: INTRODUCTION TO MACHINE LEARNING

PROF. SUNDEEP RANGAN





Learning Objectives

- ☐ Formulate a linear estimation problem with a regularization
- □ Compute an L1-regularized estimate (LASSO) using sklearn tools
- ☐ Compute the optimal regularization level using cross validation
- ☐ Interpret results from a LASSO path
- ☐ Set regularizer based on a probabilistic prior





Outline

Motivating Example: Predicting prostate cancer from a PSA test

- ☐ Model selection from LASSO regularization
- ☐ Probabilistic interpretation

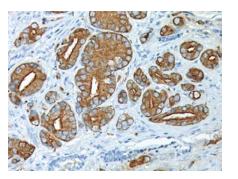


Prostate Specific Antigen Testing

- ☐ PSA levels easily tested
- ☐ High PSA believed to be associated with prostate cancer
 - Potential tool for screening
- ☐ Classic 1989 study by Thomas et al:
 - Measured PSA level of 102 men prior to prostate removal
 - Measured characteristics of prostate from samples
 - Characteristics include cancer volume, weight, ...
- ☐ Data analysis:
 - What characteristics predict PSA?

Stamey, Thomas A., et al. "Prostate specific antigen in the diagnosis and treatment of adenocarcinoma of the prostate. II. Radical prostatectomy treated patients." The Journal of urology 141.5 (1989): 1076-1083.









Data

- ☐ Prostate dataset widely-used in ML classes
- ☐ Can be downloaded from many sites
- ☐Samples = 97 patients
- ■8 features of the prostate
- ☐ Target variable = lpsa (log PSA)

```
# Get data
url = 'https://web.stanford.edu/~hastie/ElemStatLearn/datasets/prostate.data'
df = pd.read_csv(url, sep='\t', header=0)
df = df.drop('Unnamed: 0', axis=1) # skip the column of indices
```

```
The data frame has the following components:
lcavol
      log(cancer volume)
lweight
     log(prostate weight)
1bph
     log(benign prostatic hyperplasia amount)
svi
      seminal vesicle invasion
1cp
      log(capsular penetration)
gleason
      Gleason score
pgg45
     percentage Gleason scores 4 or 5
lpsa
     log(prostate specific antigen)
```





First Try: Linear Model

☐ Simple idea: Use linear regression

$$y \approx \hat{y} = \beta_0 + \beta_1 x_1 + \dots + \beta_d x_d$$

- y = lpsa (target PSA level)
- $x_1, \dots, x_d = \text{prostate features } (d = 8)$
- ■Why linear regression?
 - Easy to compute / interpret
 - Coefficients are easy to interpret
 - Larger coefficients ⇒ larger influence of feature on PSA

```
ns_train = nsamp // 2
ns_test = nsamp - ns_train
X_tr = X[:ns_train,:]  # Gets the first ns_train rows of X
y_tr = y[:ns_train]  # Gets the correspoinding rows of y

print("num samples train = %d, test = %d" % (ns_train, ns_test))
num samples train = 48, test = 49

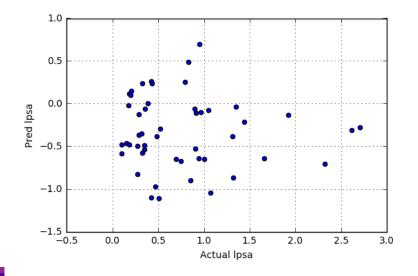
regr = linear_model.LinearRegression()
regr.fit(X_tr,y_tr)
```

Model Does Not Generalize

- Evaluate model with cross validation
 - Train on 48 samples
 - Measure RSS on 49 samples
- ☐ Test RSS is very high
- ☐ Scatter plot shows no predictive ability
- ■What happened?
- □ Can we do a better model?

```
X_ts = X[ns_train:,:]
y_ts = y[ns_train:]
y_ts_pred = regr.predict(X_ts)
RSS_rel_ts = np.mean((y_ts_pred-y_ts)**2)/(np.std(y_ts)**2)
print("Normalized test RSS = {0:f}".format(RSS_rel_ts))
```

Normalized test RSS = 4.539225





Outline

☐ Motivating Example: Predicting prostate cancer from a PSA test

Model selection from LASSO regularization

☐ Probabilistic interpretation



Intuition

- ■We know from last lecture:
 - Too many parameters ⇒ Large generalization error
- ☐ In this data set, only a few factors are likely significant
- ☐But, we don't know which one
- □Can we automatically identify them?
- □ Idea: Fit model under constraint:
 - Force only a few parameters to be non-zero
- ☐General idea of regularization:
 - Constrain the parameters with prior knowledge

```
The data frame has the following components:
lcavol
      log(cancer volume)
lweight
      log(prostate weight)
age
1bph
      log(benign prostatic hyperplasia amount)
svi
      seminal vesicle invasion
1cp
      log(capsular penetration)
gleason
      Gleason score
pgg45
      percentage Gleason scores 4 or 5
lpsa
      log(prostate specific antigen)
```





Regularized LS Estimation

☐ Standard least squares estimation (from Lecture 3):

$$\hat{\beta} = \arg\min_{\beta} RSS(\beta), \qquad RSS(\beta) = \sum_{i=1}^{n} (y_i - \hat{y}_i)^2$$

☐ Regularized estimator:

$$\hat{\beta} = \arg\min_{\beta} J(\beta), \qquad J(\beta) = RSS(\beta) + \phi(\beta)$$

- $RSS(\beta)$ = prediction error from before
- $\phi(\beta)$ = regularizing function.
- \square Concept: Regularizer penalizes β that are "unlikely"
 - Constrains estimate to smaller set of parameters



Two Common Regularizers

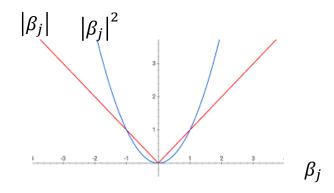
☐ Ridge regression (called L2)

$$\phi(\beta) = \alpha \sum_{j=1}^{d} |\beta_j|^2$$

□ LASSO regression (called L1)

$$\phi(\beta) = \alpha \sum_{j=1}^{d} |\beta_j|$$

- \square Both penalize large β_i
- lacksquare Level of regularization controlled by lpha



L1 and L2 Norm

- ☐ Ridge and LASSO regularized estimated are often written with vector norms
- ☐ Ridge cost function:

$$J(\beta) = \sum_{i=1}^{n} (y_i - \hat{y}_i)^2 + \alpha \sum_{j=1}^{d} |\beta_j|^2 = ||y - A\beta||^2 + \alpha ||\beta||^2$$

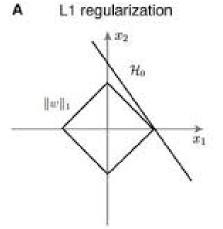
□LASSO cost function:

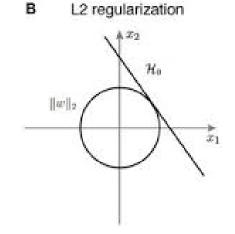
$$J(\beta) = \sum_{i=1}^{n} (y_i - \hat{y}_i)^2 + \alpha \sum_{j=1}^{d} |\beta_j| = ||y - A\beta||^2 + \alpha ||\beta||_1$$

• $\|\beta\|_1$ = L1 norm (pronounced ell-1)

Ridge vs LASSO

- □Optimization can be easily performed for L1 and L2 regularizers
 - Regularizer is convex
 - More on this later
- □L1 tends to lead to more sparse solutions
 - More coefficients are zero
 - Will focus this lecture on L1





Selecting Regularization Level

- \square How do we select regularization level α ?
 - Higher $\alpha \Rightarrow$ More constrained / simpler model
 - ∘ Lower α ⇒ More complex model
- □ Similar to inverse of model order
- \square Find α via cross-validation





Computing LASSO in python

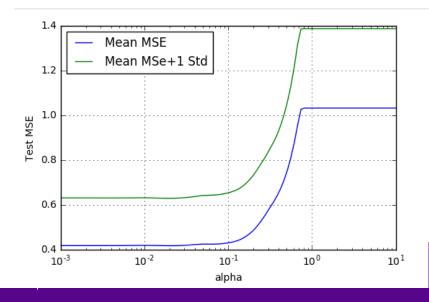
☐ Use sklearn Lasso method

Cross validation loop

Outer loop: Loop over folds

 \circ Inner loop: Loop over α

Measure mean and std deviation of MSE



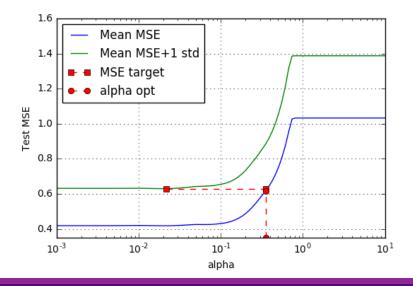
```
: # Create a k-fold cross validation object
  kf = sklearn.model selection.KFold(n splits=nfold,shuffle=True)
  # Create the LASSO model. We use the `warm start` parameter so
  # This speeds up the fitting.
  model = linear_model.Lasso(warm_start=True)
  # Regularization values to test
  nalpha = 100
  alphas = np.logspace(-3,1,nalpha)
  # MSE for each alpha and fold value
  mse = np.zeros((nalpha,nfold))
  for ifold, ind in enumerate(kf.split(X)):
      # Get the training data in the split
      Itr, Its = ind
      X_{tr} = X[Itr,:]
      y_{tr} = y[Itr]
      X_{ts} = X[Its,:]
      y_ts = y[Its]
      # Compute the lasso path for the split
      for ia, a in enumerate(alphas):
          # Fit the model on the training data
          model.alpha = a
          model.fit(X tr,y tr)
          # Compute the prediction error on the test data
          y_ts_pred = model.predict(X_ts)
          mse[ia,ifold] = np.mean((y_ts_pred-y_ts)**2)
```





Using One Standard Deviation Rule

- ☐ Use one standard deviation rule from before
 - \circ Find α_0 with minimum mean MSE, mean_mean
 - Set mse_tgt = mse_mean[α_0] + mse_std[α_0]
 - Find largest α where mse_mean[α] < mse_tgt



```
# Find the minimum MSE and MSE target
imin = np.argmin(mse_mean)
mse_tgt = mse_mean[imin] + mse_std[imin]
alpha_min = alphas[imin]

# Find the least complex model with mse_mean < mse_tgt
I = np.where(mse_mean < mse_tgt)[0]
iopt = I[-1]
alpha_opt = alphas[iopt]
print("Optimal alpha = %f" % alpha_opt)</pre>
```



Coefficients

- \square Select α via cross-validation
- ☐ Then, find coefficients using all training data.
- ☐ Final coefficients are sparse:
 - Only two factors are non-zeros
 - Lcavol: log cancer volume

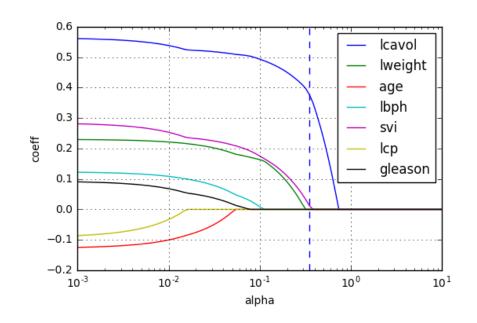
```
model.alpha = alpha_opt
model.fit(X,y)

# Print the coefficients
for i, c in enumerate(model.coef_):
    print("%8s %f" % (names_x[i], c))|

lcavol 0.376872
lweight 0.000000
    age 0.000000
    lbph 0.000000
    svi 0.012024
    lcp 0.000000
gleason 0.000000
```

LASSO path

- \Box Useful to plot coefficients as a function of α .
- ☐ Called the LASSO path
- ☐ Indicates relative importance of different factors
- ☐ For this data set:
 - lcavol most important
- □ Don't draw medical conclusions
 - Need more detailed significance testing
 - Complex subject for another class...



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Probabilistic interpretation



Bayesian Estimation

□ Suppose that true data generated from probabilistic model:

$$y = A\beta + w, \qquad w \sim N(0, \sigma^2 I)$$

 \square Consider maximum a posterior (MAP) estimator of β :

$$\hat{\beta} = \arg\max_{\beta} p(\beta|y, A)$$

 $\hat{\beta} = \text{Most likely parameter value given evidence } y, A$

□ Bayes Rule: $p(\beta|y, A) = p(y|A, \beta)p(\beta)/p(y|A)$

□ Hence: $\hat{\beta} = \arg \max_{\beta} p(y|A, \beta)p(\beta)$

 \circ Likelihood: $p(y|A,\beta)$ How well β matches data

 \circ Prior: p(eta): How well eta agrees with prior knowledge / constraints

☐ More in probability class...





Bayes Estimation with Logarithms

□Often easier to use logarithms:

$$\hat{\beta} = \arg \max_{\beta} p(y|A,\beta)p(\beta) = \arg \min_{\beta} [-\ln p(y|A,\beta)p(\beta)]$$

$$= \arg \min_{\beta} [-\ln p(y|A,\beta) - \ln p(\beta)]$$

- Gaussian density: $\ln p(y|A,\beta) = -\frac{1}{2\sigma^2} ||y A\beta||^2$
- Hence

$$\hat{\beta} = \arg\min_{\beta} \left[\frac{1}{2\sigma^2} \|y - A\beta\|^2 - \ln p(\beta) \right] = \arg\min_{\beta} [\|y - A\beta\|^2 + \phi(\beta)]$$

- \Box Conclusion: MAP estimate = regularized LS with $\phi(\beta) = -2\sigma^2 \ln p(\beta)$
 - \circ Penalize β proportional to $-\ln p(\beta)$
 - \circ Less likely β penalized more

