Group 24 – House Prices Prediction

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1. Project Task

The goal of this project is to build Machine Learning models to predict the sale prices of houses in Ames, Iowa given the data sets (year 2006-2010) from Kaggle.

The difficulties of this project include cleaning 80 original variables, building effective algorithms by ourselves, finding best parameters for models, and ensembling models to reach final high score standing in Kaggle Leaderboard.

2. Dataset, Metric & Feature Engineering

Training set 1460 rows & Testing set 1459 rows.



Find potential important features, Delete highly correlated X's; Delete outliers Fill missing value: add 0, "None", or mode case by case

Continuous: Log transformation

Ordinal: Label encoding; Categorical: One hot encoding (dummies)

Dual Form:

Training set 1456 rows * (202 features + 1) columns

Training	Validation	Testing	Total
1092	364	1459	2915

Metric: 4-fold cross validation - root mean square error (RMSE) between log(predicted sale price) and log(observed sale price)

3. Approach

CV RMSE

0.118954

0.121036

0.126783

0.128861

0.131965

0.137224

0.142892

0.223191

0.826657

0.826657

Minimize cost function:

1. Linear Regression:

$$J(\theta) = \frac{1}{2m} \sum_{i=1}^{m} (h_{\theta}(x^{(i)}) - y^{(i)})^{2}$$

- Method 1: Gradient Descent
- Method 2: Normal Equation
 - $\theta = (X^T X)^{-1} X^T y$
- Method 3: scikit-learn package

2. Ridge Regression:

Minimize cost function:

$$J(\theta) = \frac{1}{2m} \left[\sum_{i=1}^{m} (h_{\theta}(x^{(i)}) - y^{(i)})^{2} + \lambda \sum_{j=1}^{n} \theta_{j}^{2} \right]$$

- Method 1: Gradient Descent
- Method 2: Normal Equation:

$$\theta = (X^T X + \lambda I')^{-1} X^T y, \lambda > 0, I' = diag(0,1,1,...,1)$$

- Method 3: scikit-learn package

3. Lasso Regression:

Minimize cost function:

$$J(\theta) = \frac{1}{2m} \left[\sum_{i=1}^{m} (h_{\theta}(x^{(i)}) - y^{(i)})^{2} + \lambda \sum_{j=1}^{n} |\theta_{j}| \right]$$

- Method 1: Gradient Descent
- Method 2: scikit-learn package

Model

XGBoost Reg (xgboost)

Support Vector Reg (our)

Elastic Net Reg (sklearn)

Ridge Reg (our normal eqtn.)

Random Forest Reg (sklearn)

Linear Reg (our grad. descent)

Lasso Reg (our grad. descent)

Ridge Reg (sklearn)

Linear Reg (sklearn)

Lasso Reg (sklearn)

4. Results

	<u> </u>	,			
	Reg (our grad. de	0.826686			
Kaggle Leader Score					
#	Team	Score	Entries		
1	Alex plus BG que Lucas	0.0000	1		
2	Lucas le bg	0.03176	6 4		
3	Zheng Pan	0.0802	1 1		
803	Our Best Stacking	0.11864	4 15		
1389	Our Best Single Model XGBoost	0.12543	5		

4. Elastic Net Regression:

Minimize cost function:

$$J(\theta) = \frac{1}{2m} \left[\frac{\sum_{i=1}^{m} (h_{\theta}(x^{(i)}) - y^{(i)})^{2}}{+\lambda_{1} \sum_{j=1}^{n} \theta_{j}^{2} + \lambda_{2} \sum_{j=1}^{n} |\theta_{j}|} \right]$$

- Method 1: Gradient Descent
- Method 2: scikit-learn package Figure 1: Difference between Ridge, Lasso and Elastic Net Regression

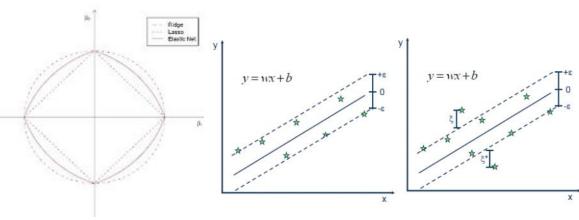
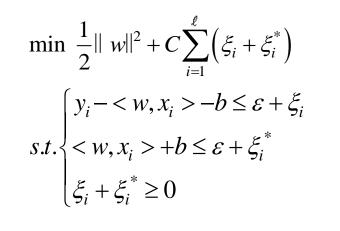


Figure 2: Comparison between SVM (left) and SVR (right)

6. XGBoost Regression 5. Support Vector Regression:



- Compared to other methods, Gradient Descent can't make model to fit Y with low CV RMSE.
- Normal equation for Ridge Regression can work almost as good as Ridge Regression in sklearn, even better in Stacking.
- Normal equations (Lasso & Elastic Net) are hard to compute and invertible singular matrix issue often occurs.
- Lasso and Elastic Net can't make CV RMSE better, but useful for feature selection.
- Some important features are: Above Ground Living Area, Size of Garage in Cars Capacity, Original Construction Date, and Zoning Classification of the sale.

Prediction

 $s.t. \sum_{n=1}^{N} (\alpha_n - \hat{\alpha}_n) = 0,$

 $L(a, \hat{a}) = -\frac{1}{2} \sum_{n=1}^{N} \sum_{n=1}^{N} (a_n - \hat{a}_n) (a_m - \hat{a}_m) k(x_n, x_m)$

 $-\varepsilon \sum_{n=1}^{N} (a_n + \hat{a}_n) + \sum_{n=1}^{N} (a_n - \hat{a}_n) t_n$

 $y(\mathbf{x}) = \sum_{n=0}^{\infty} (a_n - \hat{a}_n) k(\mathbf{x}, \mathbf{x}_n) + b$

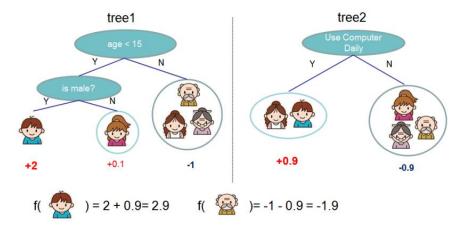


Figure 3: Tree Ensemble Model.

$L(\phi) = \sum_{i} l(\hat{y}_i, y_i) + \sum_{i} \Omega(f_k) \Omega(f) = \gamma T + \frac{1}{2} \lambda ||w||^2$ $\hat{y}_{i}^{(t)} = \sum_{k}^{t} f_{k}(x_{i}) = \hat{y}_{i}^{(t-1)} + f_{t}(x_{i})$

Figure 4: Structure Score Calculation

$$y_{i}^{(t)} = \sum_{k=1}^{n} f_{k}(\mathbf{x}_{i}) = y_{i}^{(t-1)} + f_{t}(\mathbf{x}_{i})$$

$$L^{(t)} \square \sum_{i=1}^{n} \left[l(\mathbf{y}_{i}, \hat{\mathbf{y}}^{(t-1)}) + g_{i}f_{t}(\mathbf{x}_{i}) + \frac{1}{2}h_{i}f_{t}^{2}(\mathbf{x}_{i}) \right] + \Omega(f_{t})$$

$$L^{(t)} = \sum_{i=1}^{n} \left[g_{i}f_{t}(\mathbf{x}_{i}) + \frac{1}{2}h_{i}f_{t}^{2}(\mathbf{x}_{i}) \right] + \gamma T + \frac{1}{2}\lambda \sum_{j=1}^{T} w_{j}^{2}$$

$$L^{(t)}(q) = -\frac{1}{2} \sum_{j=1}^{T} \frac{\left(\sum_{i \in I_{j}} g_{i}\right)^{2}}{\sum_{i \in I_{i}} h_{i} + \lambda} + \gamma T$$

7. Random Forest Regression

scikit-learn package, try to find best # of tree depth and best # of trees in the forest

8. Stacking:

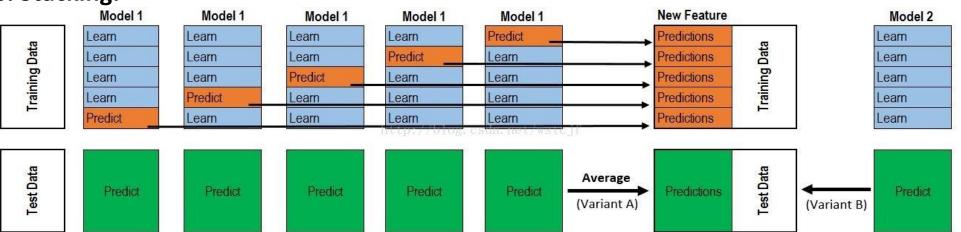


Figure 5: Stacking example flow chart

- The more models combined in the first level of the stacking, the better the performance of ensembled model is. (Ridge + Lasso + ElaNet + SVR + XGB + Random)
- The "simpler" the model in the second level of the stacking, the better the performance of ensembled model is. (Linear Reg works best)
- Max_depth and min_weight parameters has very great influence on the final result and should be optimized through Grid Search first. Regularization parameter gamma determines the complexity of the tree and should be optimized then. After fixing these parameters, subsample and colsample_bytree needs furthur optimization. XGBoost model has some insensitive parameters which can be determined manually instead of time-consuming fine-grained Grid Search.

5. Future Works

- Compute normal equations for Lasso and Ela Net.
- Program for decision trees and random forests
- Learn better algorithms from open source code in library sklearn and XGB.
- Try other model ensembling methods.