

Building Better Board Game

September 7, 2016

1 Building a Better Board Game

1.1 Setup

1.1.1 Import Libraries

```
In [1]: import warnings
import numpy as np
import pandas as pd

with warnings.catch_warnings():
    warnings.simplefilter("ignore")
    %matplotlib inline
    import matplotlib.pyplot as plt
    import matplotlib

print("Matplotlib version {} is installed.".format(matplotlib.__version__))

# Style 'ggplot' makes prettier plots.
# (plt.style isn't supported before matplotlib v1.4)
# Comment the following line, and this notebook will
# still run on matplotlib <1.4, but with default plot styling.
plt.style.use('ggplot')
```

Matplotlib version 1.5.1 is installed.

1.1.2 Set some global variables for this notebook

```
In [2]: # Set this to the number of CPUs you're comfortable
# having this notebook use during model creation.
cpus = 3

# This is used throughout the notebook for consistent results.
seed = 12
```

1.2 Data Exploration / Data Preprocessing

Read all of our data in as a pandas data frame

```
In [3]: all_data = pd.read_csv("../Data/CSV/games.csv")
```

How many observations (games) are in our set?

```
In [4]: print("The data set contains {} games.".format(len(all_data)))
```

The data set contains 84593 games.

Get an idea of the range of values for each feature in the data

```
In [5]: with warnings.catch_warnings():
        warnings.simplefilter("ignore")
        print(all_data.describe())
```

id	year	ratingScore	ratingCount	ratingStdDev	\
count	84593.000000	84593.000000	84593.000000	84593.000000	84593.000000
mean	80013.233152	1807.385008	4.279174	110.736988	0.774242
std	63960.226532	588.475029	3.039157	929.482652	0.814887
min	1.000000	-3500.000000	0.000000	0.000000	0.000000
25%	23001.000000	1985.000000	0.000000	0.000000	0.000000
50%	60049.000000	2004.000000	5.333330	2.000000	0.695211
75%	139950.000000	2011.000000	6.716100	15.000000	1.419710
max	202858.000000	2018.000000	10.000000	59423.000000	4.500000

	weightAvg	weightLightPct	weightMediumLightPct	weightMediumPct	\
count	84593.000000	71714.000000	71714.000000	71714.000000	
mean	0.876848	19.957703	16.550744	10.183793	
std	1.160127	35.385330	29.738281	23.392949	
min	0.000000	0.000000	0.000000	0.000000	
25%	0.000000	0.000000	0.000000	0.000000	
50%	0.000000	0.000000	0.000000	0.000000	
75%	1.785700	25.000000	25.000000	0.000000	
max	5.000000	100.000000	100.000000	100.000000	

	weightMediumHeavyPct	weightHeavyPct	playerAgeMin	playtimeMin	\
count	71714.000000	71714.000000	84593.000000	84593.000000	
mean	3.200719	1.401069	7.019162	47.198243	
std	12.743502	8.428136	6.808049	327.225953	
min	0.000000	0.000000	0.000000	0.000000	
25%	0.000000	0.000000	0.000000	6.000000	
50%	0.000000	0.000000	8.000000	30.000000	
75%	0.000000	0.000000	12.000000	60.000000	
max	100.000000	100.000000	1333.000000	60120.000000	

	playtimeMax	playersStatedMin	playersStatedMax	playersBestMin	\
count	84593.000000	84593.000000	84593.000000	84593.000000	
mean	51.168395	1.98825	5.682491	2.219664	
std	341.891784	0.92677	55.144456	1.210552	
min	0.000000	0.000000	0.000000	0.000000	
25%	5.000000	2.000000	2.000000	2.000000	
50%	30.000000	2.000000	4.000000	2.000000	
75%	60.000000	2.000000	6.000000	2.000000	
max	60120.000000	99.000000	11299.000000	99.000000	

	playersBestMax	priceAverage	priceStdDev	
count	84593.000000	19461.000000	19461.000000	...
mean	5.263934	24.716603	7.954304	...
std	54.600688	33.740133	16.690697	...
min	0.000000	0.010000	0.000000	...
25%	2.000000	9.950000	0.000000	...
50%	4.000000	16.663333	3.925003	...

```

75%          6.000000      29.084300      9.479418 ...
max         11299.000000    1300.000000    581.085046 ...

```

[8 rows x 163 columns]

It seems that most features have values for most games, with the exception of priceAverage and priceStdDev.

```

In [6]: feature = "priceAverage"
        percentage = len(all_data[feature].dropna())/float(len(all_data)) * 100
        print("{:.3f}% of games have a value for the '{}' feature.".format(percentage, feature))

```

23.005% of games have a value for the 'priceAverage' feature.

How many games have values for all features in the set?

```

In [7]: all_without_na = all_data.dropna()
        percentage = len(all_without_na)/float(len(all_data)) * 100
        print("{:.3f}% of games have a values for all features.".format(percentage, feature))
        print("{} games remain once filtering out those with null feature values.".format(len(all_without_na)))

```

22.573% of games have a values for all features.

19095 games remain once filtering out those with null feature values.

Some games have feature values that we might choose to consider as outliers.

```

In [8]: all_without_na.year.describe()

```

```

Out[8]: count      19095.000000
        mean        1980.718408
        std         209.432133
        min         -3500.000000
        25%         1994.000000
        50%         2006.000000
        75%         2012.000000
        max          2016.000000
        Name: year, dtype: float64

```

As you can see above, some games were published as early as 3500 B.C. Perhaps we should limit the year of publication to exclude ancient Egyptian games.

```

In [9]: since_1950 = all_without_na.query("year >= 1950")
        print("{} games remain once filtering out those published prior to 1950.".format(len(since_1950)))

```

18799 games remain once filtering out those published prior to 1950.

Ratings are crowd-sourced. Ratings may be less reliable if there were too few people contributing to the rating.

How many games have at least three votes cast toward their rating?

```

In [10]: at_least_three_raters = since_1950.query("ratingCount >= 5")
         print("{} games remain once filtering out those rated by fewer than 5 people").format(len(at_least_three_raters))

```

17622 games remain once filtering out those rated by fewer than 5 people

Our target variable is going to be 'ratingScore'.

Let's see how the ratingScore value is distributed.

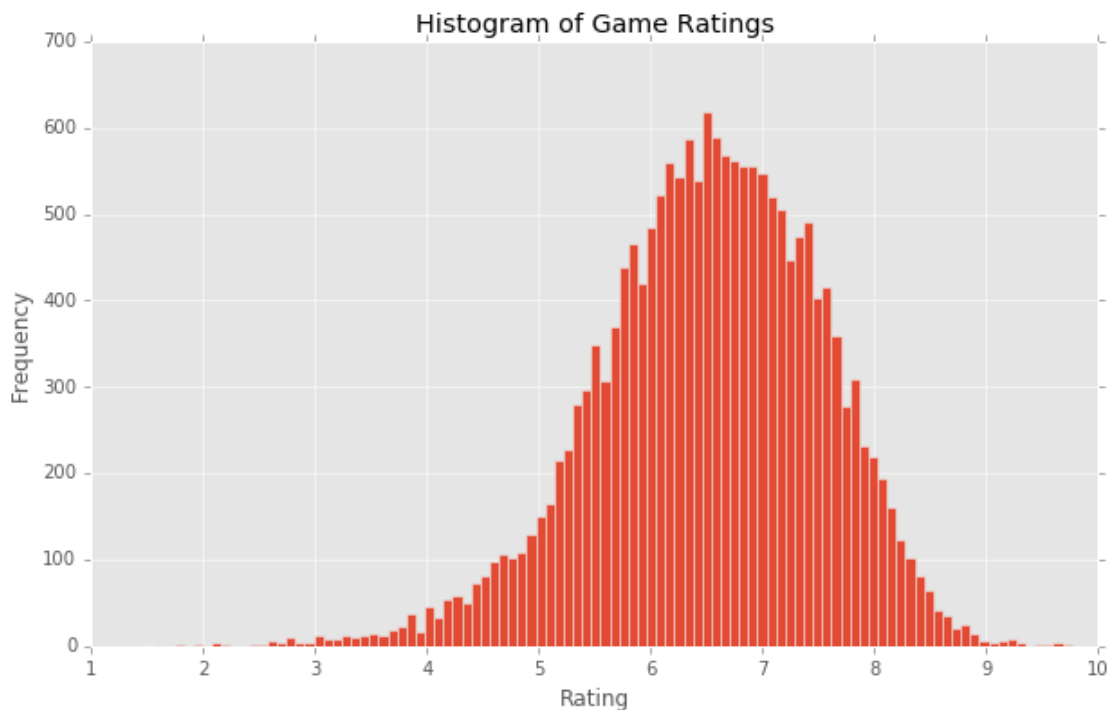
```
In [11]: df = at_least_three_raters
df.ratingScore.describe()

Out[11]: count      17622.000000
         mean         6.511106
         std         0.999621
         min         1.500000
         25%         5.875985
         50%         6.554890
         75%         7.221110
         max         9.764290
         Name: ratingScore, dtype: float64
```

It seems as if 1.0 is the minimum allowable score, and 10.0 is the max.
A histogram will probably help us see how these ratings are distributed.

1.3 Exploratory Visualization

```
In [12]: plt.figure(figsize=(10,6))
plt.hist(df.ratingScore, 100)
plt.xlim(xmin=1, xmax=10)
plt.xlabel("Rating")
plt.ylabel("Frequency")
plt.title("Histogram of Game Ratings")
plt.show()
```



Well, that looks like a skewed Gaussian distribution. Ratings of 1.0 and 10.0 are quite uncommon.
The mean of our subset of the data is shifted toward higher ratings.

```
In [13]: print "The mean rating for our subset of games is {:.3f}.".format(np.mean(df.ratingScore))
```

The mean rating for our subset of games is 6.511.

Now let's separate the columns that we intend to use as features from that which is the target.

```
In [14]: # Why are we excluding features?
# 'id' holds no predictive value
# 'name', and 'url' are strings, and cannot be features in their current form
# 'ratingCount' and 'ratingStdDev' will likely be unavailable features during prediction
# 'ratingScore' is the target variable, and should not be used to train a model
excluded = ["id", "name", "url", "ratingScore", "ratingCount", "ratingStdDev"]
feature_cols = [col for col in df.columns if col not in excluded]

# This is our target
target_col = "ratingScore"

# Create an n*m data frame where n is the number of games, and m is the number of features
X_all = df[feature_cols]
# Create an n*1 data frame where n is the number of games, each with the single target feature
y_all = df[target_col]
```

1.4 Implementation

We need to split this data into (at least) two portions. * A training set - 80% of total * We will use 5-fold cross-validation to get predictions on this portion. * A held-out test (or validation) set - 20% of total * We will use this set to prove the performance of the final model.

```
In [15]: # Select features (X) and corresponding labels (y) for the trainingCV and holdout sets
# train_test_split() shuffles data randomly. We will set the random state for consistency
from sklearn.cross_validation import train_test_split
# X/y_train are 80% of data, for use in training/testing models.
# X/y_holdout are 20% of data, for use in proving model performance at the end.
X_train, X_holdout, y_train, y_holdout = train_test_split(
    np.array(X_all), np.array(y_all),
    train_size=0.8,
    random_state=seed)

from sklearn.cross_validation import KFold
# We will split X_train into train and test 5 times via KFold
kf = KFold(X_train.shape[0], n_folds=5, shuffle=True, random_state=seed)

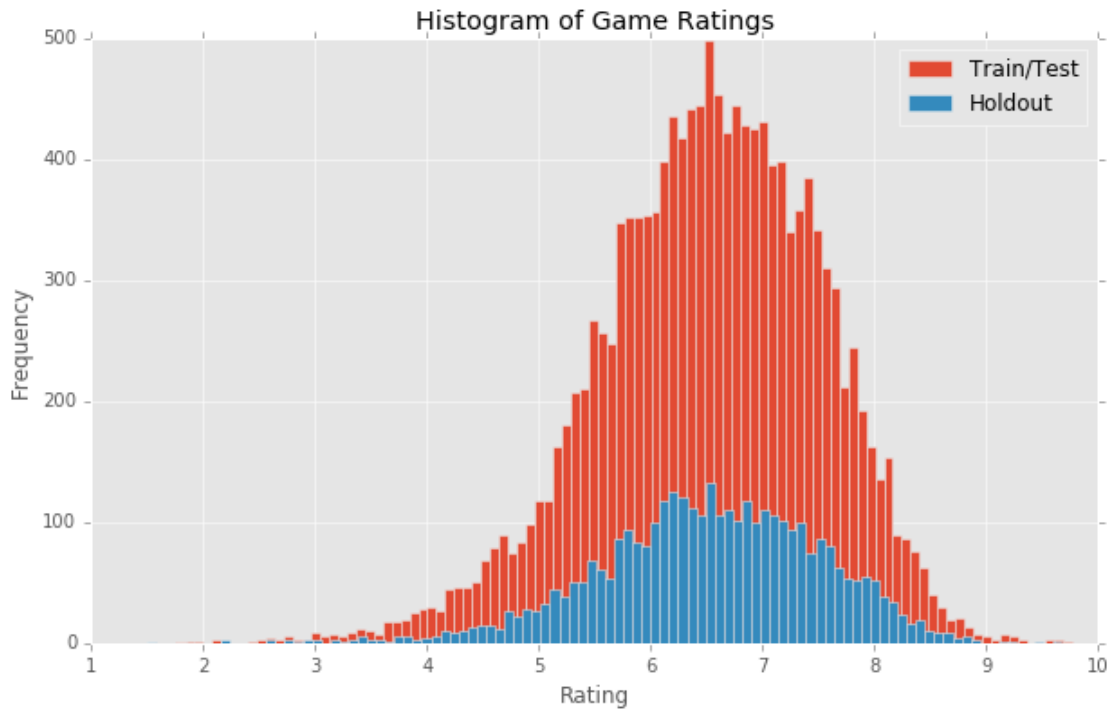
print "Cross-validated Training set includes {} samples".format(len(X_train))
print "Holdout/Validation set includes      {} samples".format(len(X_holdout))
```

```
Cross-validated Training set includes 14097 samples
Holdout/Validation set includes      3525 samples
```

With random sampling, we would anticipate the histogram of the target variable in the training and test sets to be similarly distributed.

```
In [16]: plt.figure(figsize=(10,6))
plt.hist(y_train, 100, label="Train/Test")
plt.hist(y_holdout, 100, label="Holdout")
plt.legend(loc='upper right')
plt.xlim(xmin=1, xmax=10)
plt.xlabel("Rating")
plt.ylabel("Frequency")
```

```
plt.title("Histogram of Game Ratings")
plt.show()
```



It's useful to time training jobs. Let's set up some train and predict wrapper functions to do so.

```
In [17]: # Helper functions to train and predict
import time

def train(learner, X_train, y_train):
    print "Training {}...".format(learner.__class__.__name__)
    start = time.time()
    learner.fit(X_train, y_train)
    end = time.time()
    print "Done.\nTraining time (secs): {:.3f}".format(end - start)

def predict(learner, features):
    print "Predicting target using {}...".format(learner.__class__.__name__)
    start = time.time()
    y_pred = learner.predict(features)
    end = time.time()
    print "Done.\nPrediction time (secs): {:.3f}".format(end - start)
    return pd.Series(y_pred)
```

1.4.1 Random Forest regressor

A tree-based learner The first child learner we wish to try is Random Forest. Since there are some parameters that might need to be tuned, we'll make use of GridSearchCV to learn what the best choices for those parameters might be.

```

In [18]: # Random Forest is a tree-based learner.
from sklearn.ensemble import RandomForestRegressor
random_forest_reg = RandomForestRegressor(
    n_estimators=2000, # More is better (but slower)
    oob_score=True,   #
    random_state=seed)

# We'll mess with some parameters in our grid search.
# A smart choice for max_features is sqrt(num_features).
# For our 159 features, that would be 12.6, so I'm starting in that range.
parameters = {
    'max_features': [10,20,30,40,50,60,80],
    'min_samples_split': [2,3,4,6]
}

# We'll need these to do a grid search.
from sklearn.metrics import make_scorer
from sklearn.grid_search import GridSearchCV

# We'll use MSE as our function to determine the winner in the grid search
from sklearn.metrics import mean_squared_error
mse_scorer = make_scorer(mean_squared_error, greater_is_better=False)

# Set up the grid search, using training data in 5 folds
rf_grid = GridSearchCV(random_forest_reg, parameters, scoring=mse_scorer, cv=kf, n_jobs=cpus)

```

Perform the grid search.

```

In [19]: # Fit model to training data - this will take a while.
train(rf_grid, X_train, y_train)
# Let's print which parameter(s) were chosen as the best in the grid
print "\nBest model: {}".format(rf_grid.best_estimator_)

```

Training GridSearchCV...

Done.

Training time (secs): 6991.263

```

Best model: RandomForestRegressor(bootstrap=True, criterion='mse', max_depth=None,
    max_features=40, max_leaf_nodes=None, min_samples_leaf=1,
    min_samples_split=3, min_weight_fraction_leaf=0.0,
    n_estimators=2000, n_jobs=1, oob_score=True, random_state=12,
    verbose=0, warm_start=False)

```

It might help to visualize the value from the scorer function across the parameter grid.

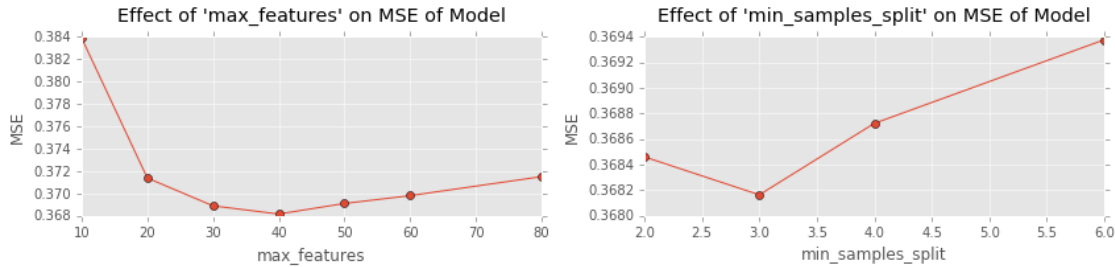
```

In [20]: # Subset the grid_scores_ list to isolate one dependent variable
f, (max_features, min_samples_split) = plt.subplots(1, 2, figsize=(12, 3))

# First plot 'max_features'
models = zip(*[(abs(mean), std, p['max_features'])
    for p, mean, std in [x for x in rf_grid.grid_scores_ if x[0]['min_samples_split'] == 3]])
max_features.plot(models[2], models[0], 'o-')
max_features.set_title("Effect of 'max_features' on MSE of Model", y=1.05)
max_features.set_xlabel("max_features")
max_features.set_ylabel("MSE")

```

```
# then plot 'min_samples_split'
models = zip(*[(abs(mean), std, p['min_samples_split'])
               for p, mean, std in [x for x in rf_grid.grid_scores_ if x[0]['max_features'] == 40]])
min_samples_split.plot(models[2], models[0], 'o-')
min_samples_split.set_title("Effect of 'min_samples_split' on MSE of Model", y=1.05)
min_samples_split.set_xlabel("min_samples_split")
min_samples_split.set_ylabel("MSE")
f.tight_layout(h_pad=10.05)
plt.show()
```



- 'max_features' seems to produce a models with the lowest mean squared error around 40 features.
- 'min_samples_split' clearly seems to perform best with a value of 3.

Get Random Forest predictions on all the training data now that we've determined good parameter values.

This is scoring the training set in 5 folds, and collecting the predictions from the held out portion at each fold, to ensure no bias in predictions on the training set.

```
In [21]: from sklearn.cross_validation import cross_val_predict
         from sklearn.metrics import mean_squared_error
         from scipy.stats import pearsonr
```

```
# Helper utility to show cross-validated performance statistics on a training set.
def show_training_performance(estimator, X, y, cv=5):
    # Compute predictions on the entire set by
    # collecting the predictions on the held-out set from each fold
    y_pred = cross_val_predict(estimator, X, y, cv=cv)
    print "Mean Squared Error: {}".format(mean_squared_error(y, y_pred))
    print "Pearson Correlation: {}".format(pearsonr(y, y_pred)[0])
    return y_pred
```

For the metrics we've decided to use for model comparison, how does this model perform?

```
In [22]: y_rf_pred = show_training_performance(rf_grid.best_estimator_, X_train, y_train, cv=kf)
```

Mean Squared Error: 0.3681646466

Pearson Correlation: 0.79626430548

Then let's visualize the actual vs predicted ratingScore, to see how close to the diagonal they are.

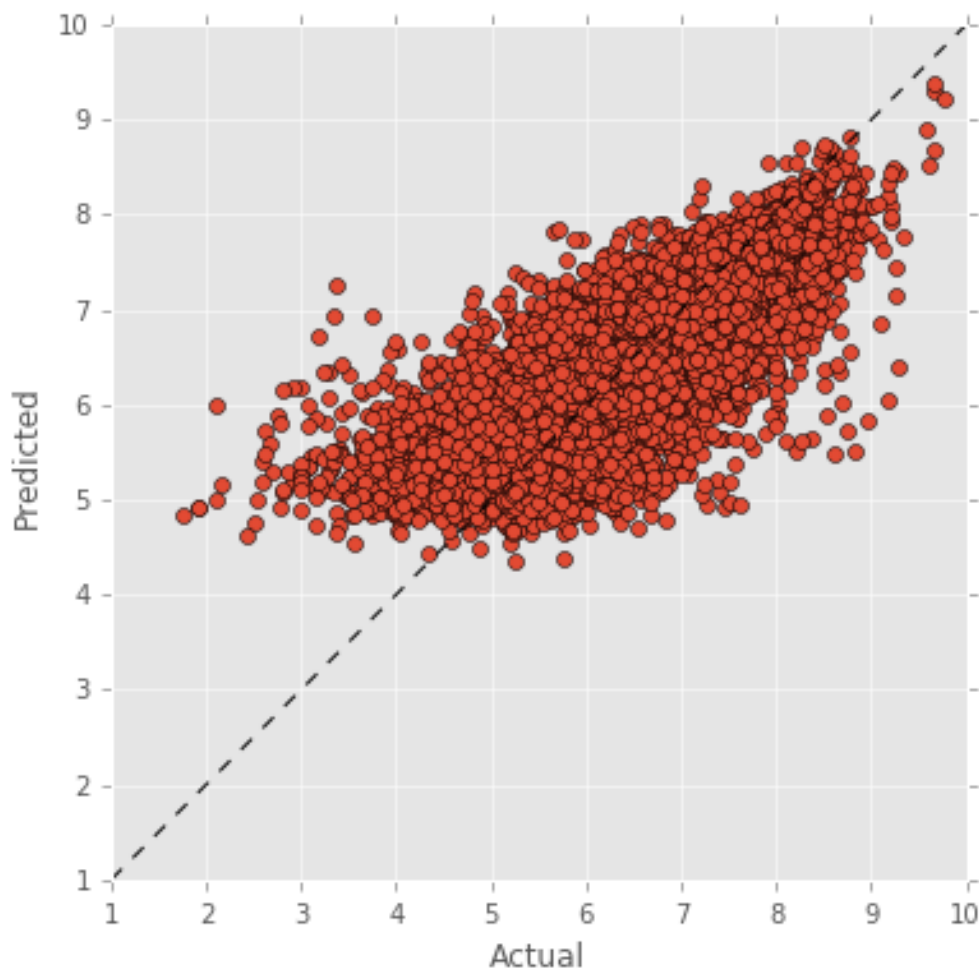
```
In [23]: # Helper utility to create actual vs. predicted plots
         def plot_actual_v_predicted(title, actual, predicted):
```



```
plt.figure(figsize=(6,6))
plt.plot(actual, predicted, 'o')
plt.plot([1, 10], [1, 10], 'k--')
plt.xlabel("Actual")
plt.ylabel("Predicted")
plt.title("Actual vs. Predicted values for ratingScore - {}".format(title), y=1.05)
plt.xlim(1, 10)
plt.ylim(1, 10)
plt.gca().set_aspect('equal', adjustable='box')
plt.show()
```

In [24]: `plot_actual_v_predicted("RandomForest", y_train, y_rf_pred)`

Actual vs. Predicted values for ratingScore - RandomForest



This model seems hesitant to predict values below 4. The largest errors are thus on the lowest-scoring games.

Now that we've determined good parameter values, let's store them for later use.

In [25]: `# Lock in our best-performing parameters for this estimator`
`random_forest_reg.set_params(**rf_grid.best_params_)`

```
Out [25]: RandomForestRegressor(bootstrap=True, criterion='mse', max_depth=None,
                                max_features=40, max_leaf_nodes=None, min_samples_leaf=1,
                                min_samples_split=3, min_weight_fraction_leaf=0.0,
                                n_estimators=2000, n_jobs=1, oob_score=True, random_state=12,
                                verbose=0, warm_start=False)
```

1.4.2 K-Nearest-Neighbors

An incredibly simple distance-based 'learner'

```
In [26]: from sklearn.neighbors import KNeighborsRegressor
         k_neighbors_reg = KNeighborsRegressor( weights='distance' )
         # We'll mess with some parameters in our grid search.
         # n_neighbors is now many neighbors to consider for each point
         parameters = {
             'n_neighbors': np.arange(2,32,2)
         }
         # Set up the grid search, using "base" data in 10 folds
         knn_grid = GridSearchCV(k_neighbors_reg, parameters, scoring=mse_scorer, cv=kf, n_jobs=cpus)
```

What were the best values for the hyper-parameter(s) within our grid?

```
In [27]: # KNN models use Euclidian distance, so relative scales of features matter.
         # Here we are standardizing the feature values.
         from sklearn import preprocessing
         scaler = preprocessing.StandardScaler().fit(X_all)
         X_train_scaled = scaler.transform(X_train)
         # Fit model to training data
         train(knn_grid, X_train_scaled, y_train)
         # Let's print which parameter(s) were chosen as the best in the grid
         print "\nBest model: {}".format(knn_grid.best_estimator_)
```

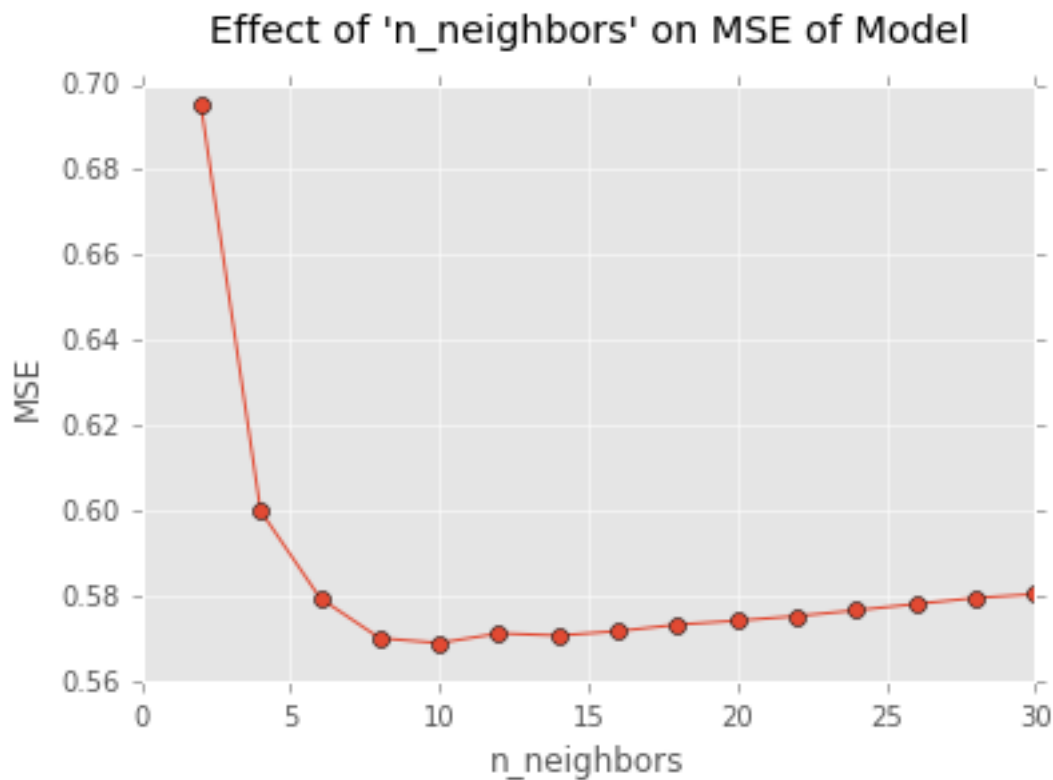
Training GridSearchCV...

Done.

Training time (secs): 346.834

```
Best model: KNeighborsRegressor(algorithm='auto', leaf_size=30, metric='minkowski',
                                metric_params=None, n_jobs=1, n_neighbors=10, p=2,
                                weights='distance')
```

```
In [28]: # Plot MSE vs 'n_neighbors'
         models = zip(*[(abs(mean), std, p['n_neighbors'])
                        for p, mean, std in [x for x in knn_grid.grid_scores_]])
         plt.plot(models[2],models[0], 'o-')
         plt.title("Effect of 'n_neighbors' on MSE of Model", y=1.05)
         plt.xlabel("n_neighbors")
         plt.ylabel("MSE")
         plt.show()
```



- `n_neighbors` had an clear impact the resulting model's mean squared error. **10 Seemed to be a good value** (though not by much)

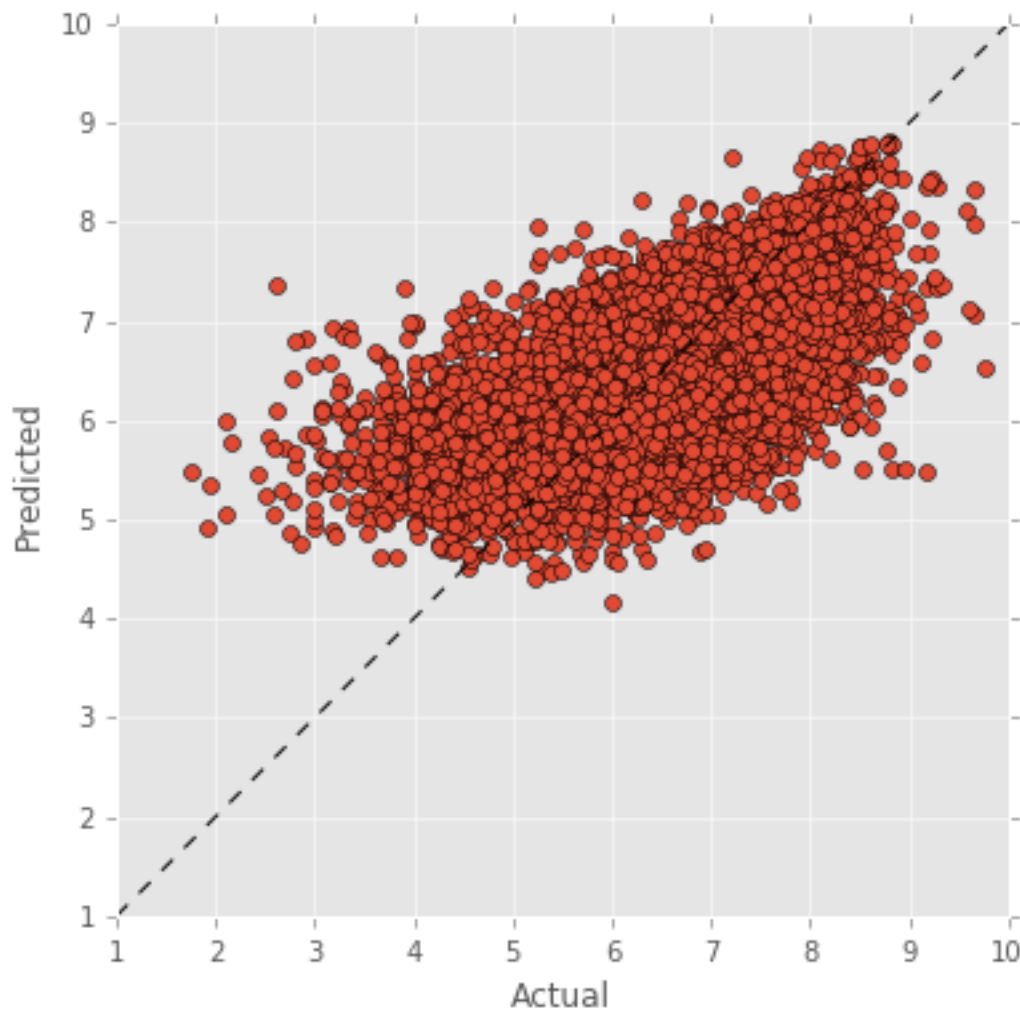
Get K Neighbor predictions on all the training data now that we've determined good parameter values.

```
In [29]: y_knn_pred = show_training_performance(knn_grid.best_estimator_, X_train_scaled, y_train, cv=kf)
          plot_actual_v_predicted("KNeighbors", y_train, y_knn_pred)
```

Mean Squared Error: 0.568845614465

Pearson Correlation: 0.660520144685

Actual vs. Predicted values for ratingScore - KNeighbors



Interestingly, this model does predict values below 4 at times, though only when the games scored above a 4.

```
In [30]: # Lock in our best-performing parameters for this estimator
         k_neighbors_reg.set_params(**knn_grid.best_params_)
```

```
Out[30]: KNeighborsRegressor(algorithm='auto', leaf_size=30, metric='minkowski',
                             metric_params=None, n_jobs=1, n_neighbors=10, p=2,
                             weights='distance')
```

1.4.3 Support Vector regressor

A slow learner, with many important (and dependent) parameters I initially attempted to test multiple parameters in this grid. After **9 days**, modeling had not finished. I was eventually forced to try small sets of parameters and values so that I could achieve results at all.

- I first varied the value of 'C', before settling on a best value of 10.

- Then I varied the value of ‘gamma’, with ‘C’ fixed, and settled on a value of 0.001.

I am aware that this is not an ideal method, as these two parameters are likely dependent, and if I had time for a more exhaustive grid search, the winning pair of these features might indeed be different for this feature set.

```
In [31]: # The Support Vector Machine (SVM) as a regression is referred to as SVR.
from sklearn.svm import SVR
support_vector_reg = SVR()
# We'll mess with some parameters in our grid search.
# C - Penalty parameter (1/lambda)
# kernel - method of mapping multi-dimensional feature space into lower dimensions to fit a pl
parameters = {
    'C': [1,10,100], # Determined through prior grid searches on only this parameter
    'gamma': [0.0001,0.001,0.01], # Subsequent single-parameter grid search
}
# Set up the GridSearch with cross-validation
svr_grid = GridSearchCV(support_vector_reg, parameters, scoring=mse_scorer, cv=kf, n_jobs=cpus)

In [32]: # Fit model to training data (with standardized data)
train(svr_grid, X_train_scaled, y_train)
# Let's print which parameter(s) were chosen as the best in the grid
print "Best model: {}".format(svr_grid.best_estimator_)
```

Training GridSearchCV...

Done.

Training time (secs): 1673.969

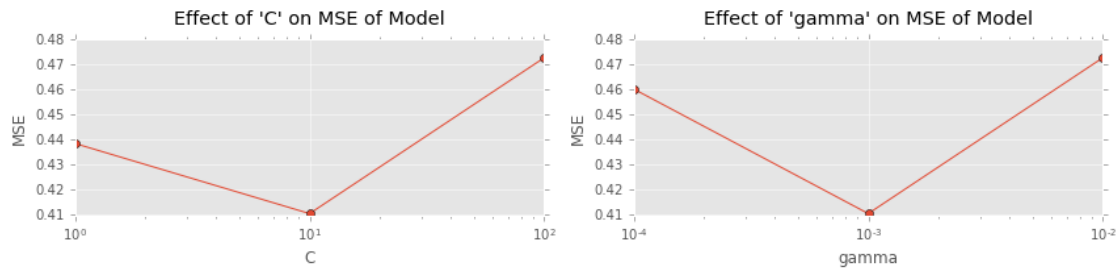
Best model: SVR(C=10, cache_size=200, coef0=0.0, degree=3, epsilon=0.1, gamma=0.001, kernel='rbf', max_iter=-1, shrinking=True, tol=0.001, verbose=False)

What were the best values for the hyper-parameter(s) within our grid?

```
In [33]: # Subset the grid_scores_ list to isolate one dependent variable
f, (c, gamma) = plt.subplots(1, 2, figsize=(12, 3))

# First plot 'max_features'
models = zip(*[(abs(mean), std, p['C'])
               for p, mean, std in [x for x in svr_grid.grid_scores_ if x[0]['gamma'] == 0.001]])
c.plot(models[2],models[0], 'o-')
c.set_title("Effect of 'C' on MSE of Model", y=1.05)
c.set_xlabel("C")
c.set_xscale("log")
c.set_ylabel("MSE")

# then plot 'min_samples_split'
models = zip(*[(abs(mean), std, p['gamma'])
               for p, mean, std in [x for x in svr_grid.grid_scores_ if x[0]['C'] == 10]])
gamma.plot(models[2],models[0], 'o-')
gamma.set_title("Effect of 'gamma' on MSE of Model", y=1.05)
gamma.set_xlabel("gamma")
gamma.set_xscale("log")
gamma.set_ylabel("MSE")
f.tight_layout(h_pad=10.05)
plt.show()
```

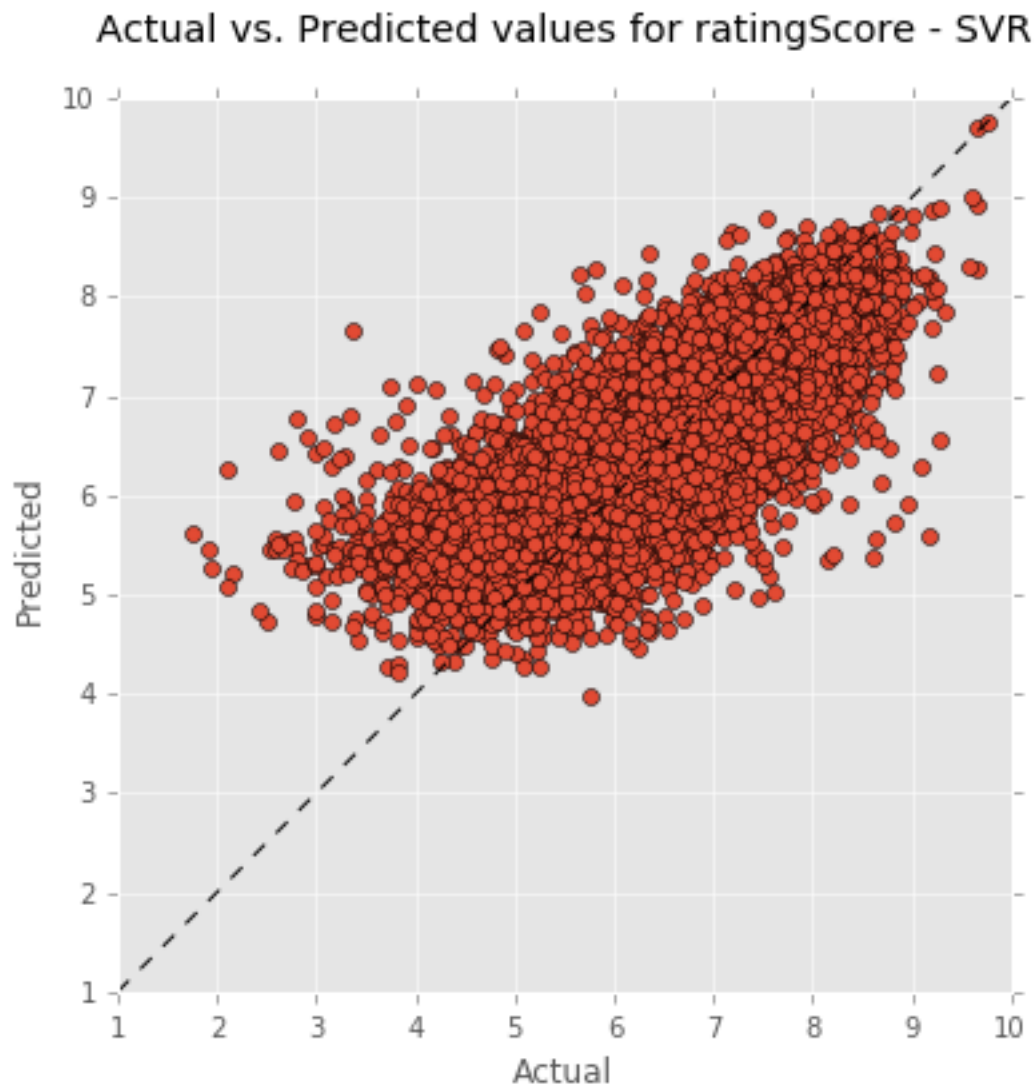


Let's again visualize the actual vs predicted ratingScore

```
In [34]: y_svr_pred = show_training_performance(svr_grid.best_estimator_, X_train_scaled, y_train, cv=kf)
         plot_actual_v_predicted("SVR", y_train, y_svr_pred)
```

Mean Squared Error: 0.410589573021

Pearson Correlation: 0.768832708798



```
In [35]: # Lock in our best-performing parameters for this estimator
support_vector_reg.set_params(**svr_grid.best_params_)
```

```
Out[35]: SVR(C=10, cache_size=200, coef0=0.0, degree=3, epsilon=0.1, gamma=0.001,
           kernel='rbf', max_iter=-1, shrinking=True, tol=0.001, verbose=False)
```

1.4.4 Gradient Boosting Regressor

Another ensemble model What if we stacked these two learners together to create another model?

```
In [36]: from sklearn.ensemble import GradientBoostingRegressor
gradient_boost_reg = GradientBoostingRegressor(
    max_features=40,      # Choosing (for better or worse) to use the optimal value from the RF
    n_estimators=1000,    # More is better (but slower)
    random_state=seed)
# We'll mess with one parameter in our grid search. (GBR takes a long time)
# max_depth - How far to grow each of the trees
parameters = {
    'max_depth': [2, 4, 6, 8, 10, 15, 20],
}
# Set up the GridSearch with cross-validation
gbr_grid = GridSearchCV(gradient_boost_reg, parameters, scoring=mse_scorer, cv=kf, n_jobs=cpus)
```

What were the best values for the hyper-parameter(s) within our grid?

```
In [37]: # Fit model to training data - not nearly as long-running as RandomForest
train(gbr_grid, X_train, y_train)
print "Best model: {}".format(gbr_grid.best_estimator_)
```

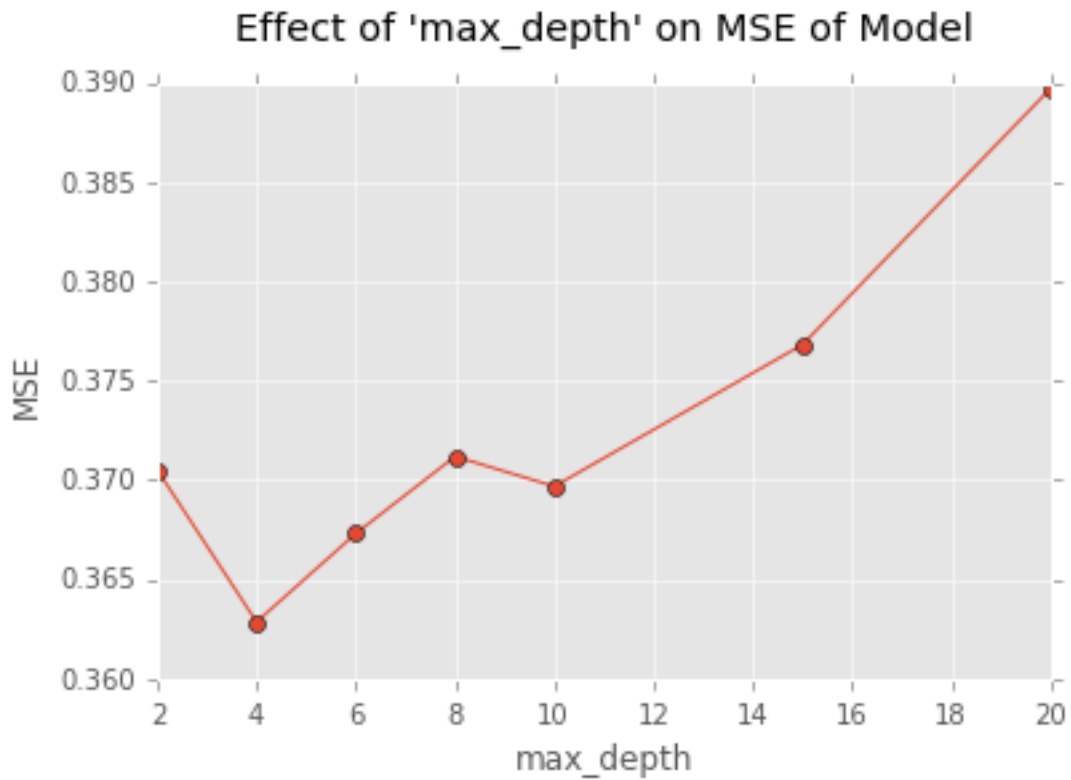
Training GridSearchCV...

Done.

Training time (secs): 2688.459

```
Best model: GradientBoostingRegressor(alpha=0.9, init=None, learning_rate=0.1, loss='ls',
    max_depth=4, max_features=40, max_leaf_nodes=None,
    min_samples_leaf=1, min_samples_split=2,
    min_weight_fraction_leaf=0.0, n_estimators=1000,
    presort='auto', random_state=12, subsample=1.0, verbose=0,
    warm_start=False)
```

```
In [38]: # Plot 'max_depth' vs MSE
models = zip(*[(abs(mean), std, p['max_depth']) for p, mean, std in gbr_grid.grid_scores_])
plt.plot(models[2], models[0], 'o-')
plt.title("Effect of 'max_depth' on MSE of Model", y=1.05)
plt.xlabel("max_depth")
plt.ylabel("MSE")
plt.show()
```



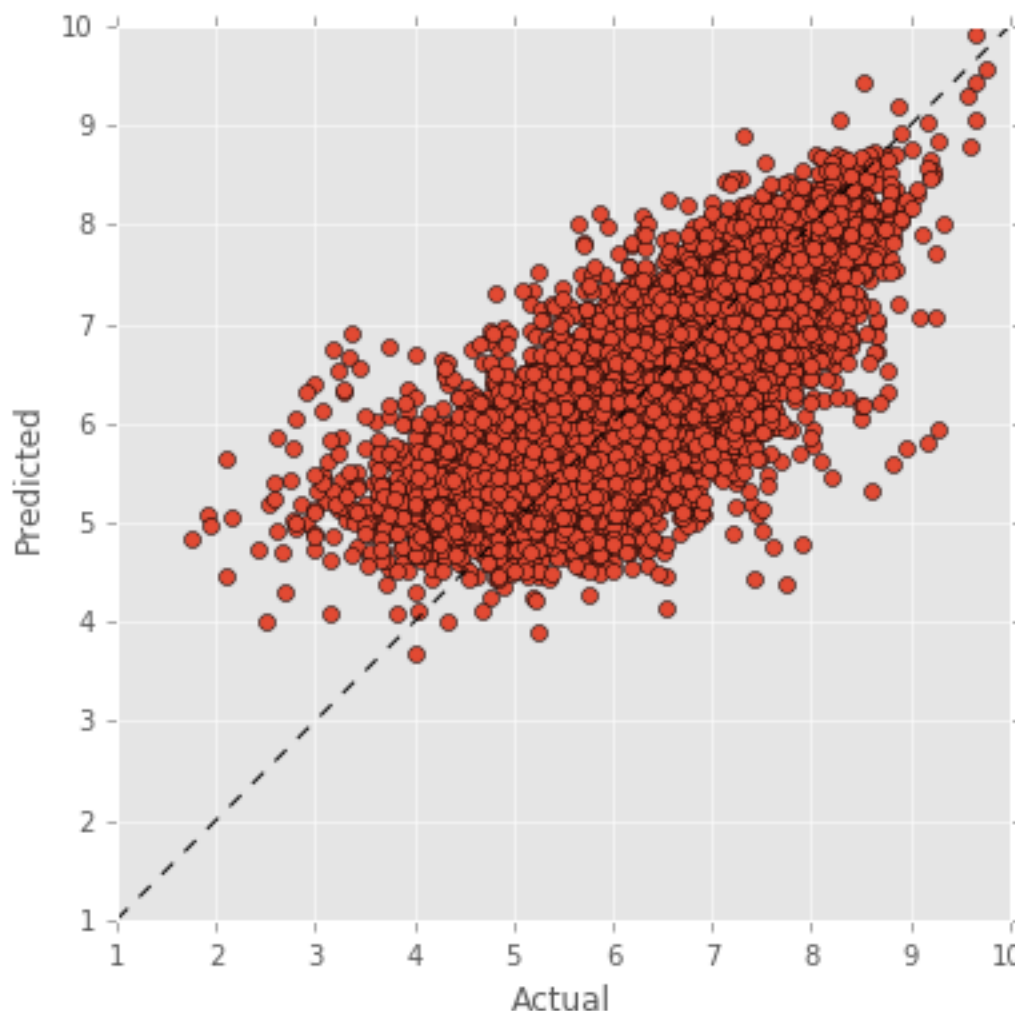
Let's again visualize the actual vs predicted ratingScore

```
In [39]: y_gbr_pred = show_training_performance(gbr_grid.best_estimator_, X_train, y_train, cv=kf)
         plot_actual_v_predicted("GBR", y_train, y_gbr_pred)
```

Mean Squared Error: 0.362856147435

Pearson Correlation: 0.798432186811

Actual vs. Predicted values for ratingScore - GBR



```
In [40]: # Lock in our best-performing parameters for this estimator
         gradient_boost_reg.set_params(**gbr_grid.best_params_)
```

```
Out[40]: GradientBoostingRegressor(alpha=0.9, init=None, learning_rate=0.1, loss='ls',
                                     max_depth=4, max_features=40, max_leaf_nodes=None,
                                     min_samples_leaf=1, min_samples_split=2,
                                     min_weight_fraction_leaf=0.0, n_estimators=1000,
                                     presort='auto', random_state=12, subsample=1.0, verbose=0,
                                     warm_start=False)
```

1.5 Creating a Pipeline for Stacking

I couldn't seem to find a standard implementation for this.

It's a bit tedious to train all of the child learners, then stack them manually. Why don't we make an estimator that accepts an array of child estimators, and a meta (stacking) estimator, and does the work for us?

```

In [41]: from sklearn.externals.joblib import Parallel, delayed
         from sklearn.metrics import make_scorer, mean_squared_error
         from scipy.stats import pearsonr
         from sklearn.cross_validation import cross_val_predict

def _fit_estimator(estimator, X, y, scaler=False, cv=10, verbose=0, random_state=None):
    """Private function used to fit with a single estimator in parallel."""
    # If requested for this estimator, scale the features
    if(scaler):
        X = scaler.transform(X)
    # Get cross-validated predictions on all of the data - to feed to stacker
    y_pred = cross_val_predict(estimator, X, y, cv=10)
    if verbose > 0:
        print "Estimator: {}".format(estimator.__class__.__name__)
        print "  Mean Squared Error: {}".format(mean_squared_error(y, y_pred))
        print "  Pearson Correlation: {}".format(pearsonr(y, y_pred)[0])
    # Train the final estimator on all data
    estimator.fit(X, y)
    # Return the cross-validated predictions for this child estimator
    return y_pred

def _predict_estimator(estimator, X, scaler=False, verbose=0):
    """Private function used to predict with a single estimator in parallel."""
    # If requested for this estimator, scale the features
    if(scaler):
        X = scaler.transform(X)
    # Return the cross-validated predictions for this child estimator
    return estimator.predict(X)

class StackedRegression:

    def __init__(self, estimators, meta_estimator, cv=None, random_state=None, n_jobs=1, verbose=0):
        self.__class__.__name__ = 'StackedRegression'
        self.estimators = estimators
        self.meta_estimator = meta_estimator
        self.cv = cv
        self.random_state = random_state
        self.n_jobs = n_jobs
        self.verbose = verbose

    def fit(self, X, y):
        """Fit the model according to the given training data."""
        # Parallel loop: fit each estimator, storing cross-validated predictions
        self.child_predictions_ = np.transpose(np.array(
            Parallel(n_jobs=self.n_jobs, verbose=self.verbose, backend="threading")(
                delayed(_fit_estimator)(
                    est['estimator'], X, y, scaler=est['scaler'] if 'scaler' in est else False,
                    verbose=self.verbose, random_state=self.random_state)
                for est in self.estimators)))
        # fit the meta-estimator on the predictions from the child estimators
        self.stack_pred_ = _fit_estimator(self.meta_estimator['estimator'],
            self.child_predictions_, y, scaler=False,
            cv=self.cv, verbose=self.verbose, random_state=self.random_state)
        return self

```

```

def predict(self, X):
    """Predict class for X."""
    # First predict with all child estimators
    # Parallel loop: predict with each child estimator
    child_predictions = np.transpose(np.array(
        Parallel(n_jobs=self.n_jobs, verbose=self.verbose, backend="threading")(
            delayed(_predict_estimator)(est['estimator'], X, scaler=est['scaler'] if 'scaler' in est else None,
                                       verbose=self.verbose)
            for est in self.estimators)))
    # fit the meta-estimator on the predictions from the child estimators
    stack_predictions = _predict_estimator(self.meta_estimator['estimator'], child_predictions,
                                          scaler=False,
                                          verbose=self.verbose)

    return stack_predictions

```

Now let's instantiate a meta-estimator, and the StackedRegressor.

We'll be stacking the predictions of: * Random Forest * K-Nearest Neighbors * Support Vector Regression
 * Gradient Boosting Regression

With a meta-estimator (stacker) of: * LassoCV - a linear modeler

NOTE: We stored the best_params_ of each grid search above, so each child estimator will only attempt to make one model, with those best parameters this time around.

```

In [42]: # I'll use a linear model (Lasso) as the "meta-estimator" or stacker
from sklearn.linear_model import LassoCV
lasso_cv_reg = LassoCV()

# Provide standard scaling for some learners
from sklearn import preprocessing
standard_scaler = preprocessing.StandardScaler().fit(X_all)

# Create the stacking regression learner
stacker = StackedRegression(
    # Child estimators
    [{'estimator': random_forest_reg},
     {'estimator': k_neighbors_reg, 'scaler': standard_scaler},
     {'estimator': support_vector_reg, 'scaler': standard_scaler},
     {'estimator': gradient_boost_reg}],
    # Stacking meta-estimator
    {'estimator': lasso_cv_reg},
    # Additional settings, passed to lower-level estimators
    cv=kf,
    #verbose=1,
    random_state=seed,
    n_jobs=cpus)

```

Let's give it a whirl!

```

In [43]: # Now train it
train(stacker, X_train, y_train)

```

Training StackedRegression...

Done.

Training time (secs): 2012.814

1.6 Each Estimator's Performance on The Held-out Set

```
In [83]: # Show holdout performance
def show_holdout_performance(estimator, X, actual_y):
    # Compute predictions on the held out set
    predicted_y = estimator.predict(X)
    print(estimator.__class__.__name__)
    mse = mean_squared_error(actual_y, predicted_y)
    corr = pearsonr(actual_y, predicted_y)[0]
    print(" Mean Squared Error: {}".format(mse))
    print(" Pearson Correlation: {}".format(corr))
    return (mse, corr, predicted_y)

# Iterate over each estimator, scale features as necessary, score the held-out set
performance = pd.DataFrame()
for est in ([{'estimator': random_forest_reg},
             {'estimator': k_neighbors_reg, 'scaler': scaler},
             {'estimator': support_vector_reg, 'scaler': scaler},
             {'estimator': gradient_boost_reg}]):
    if('scaler' in est):
        X = est['scaler'].transform(X_holdout)
    else:
        X = X_holdout
    (mse, corr, predictions) = show_holdout_performance(est['estimator'], X, y_holdout)
    performance = performance.append(pd.Series([est['estimator'].__class__.__name__, mse, corr]))
```

```
RandomForestRegressor
Mean Squared Error: 0.369919958452
Pearson Correlation: 0.794390683847
KNeighborsRegressor
Mean Squared Error: 0.535479812911
Pearson Correlation: 0.683169963154
SVR
Mean Squared Error: 0.408616531813
Pearson Correlation: 0.769350107821
GradientBoostingRegressor
Mean Squared Error: 0.358261623101
Pearson Correlation: 0.800556639349
```

It'd be interesting to note how the linear stacker weighted the importance of each of the child learners.

```
In [45]: pd.Series(stacker.meta_estimator['estimator'].coef_,
                  index=[e['estimator'].__class__.__name__ for e in stacker.estimators])
```

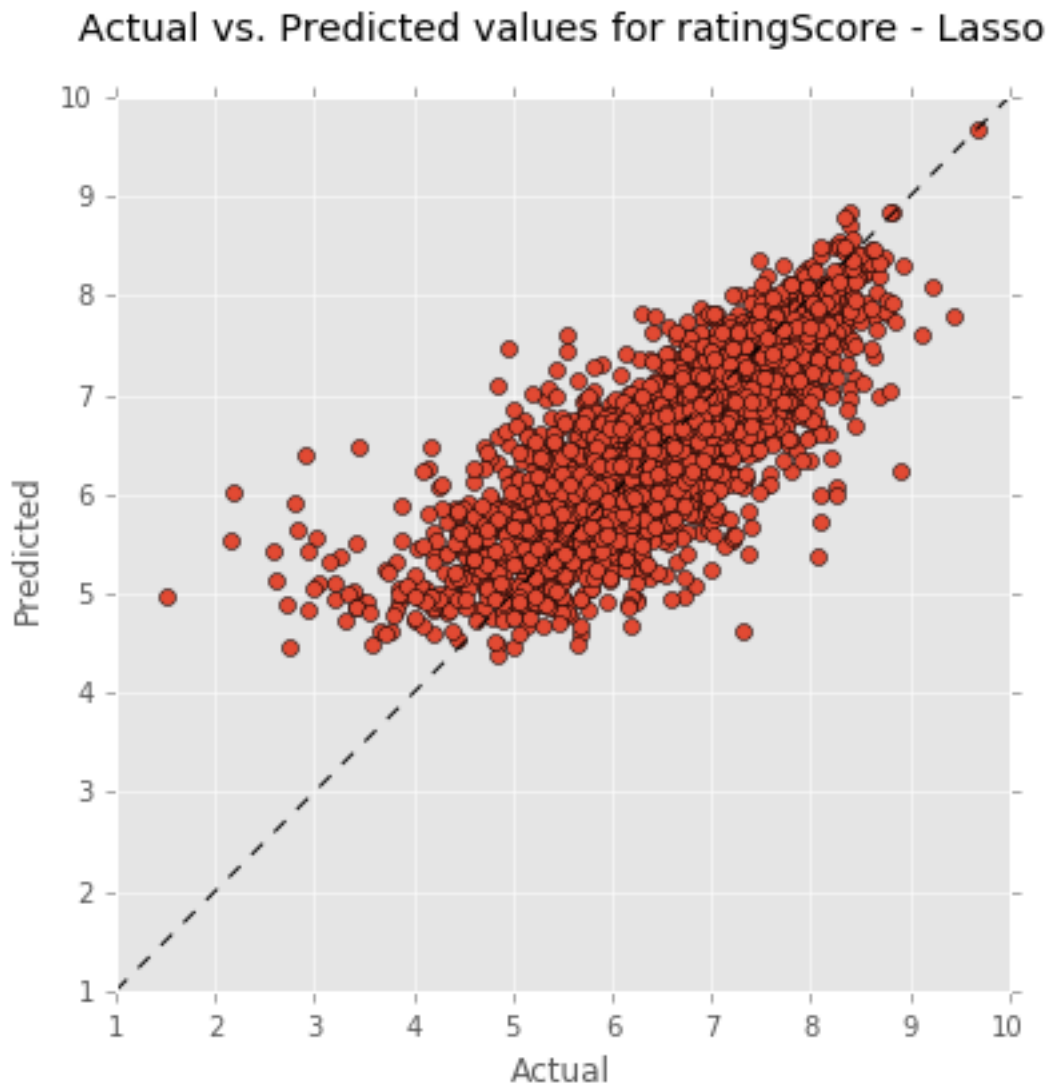
```
Out[45]: RandomForestRegressor    0.450035
KNeighborsRegressor              0.082644
SVR                             0.114073
GradientBoostingRegressor        0.423877
dtype: float64
```

```
In [84]: (mse, corr, final_predictions) = show_holdout_performance(stacker, X_holdout, y_holdout)
performance = performance.append(pd.Series([stacker.__class__.__name__, mse, corr]), ignore_in
```

```
StackedRegression
Mean Squared Error: 0.344842206614
Pearson Correlation: 0.808836787278
```

It looks as if we've been able to achieve a higher correlation via stacking than from any learner individually (barely).

```
In [60]: plt.figure(figsize=(6,6))
plt.plot(y_holdout, final_predictions, 'o')
plt.plot([1, 10], [1, 10], 'k--')
plt.xlabel("Actual")
plt.ylabel("Predicted")
plt.title("Actual vs. Predicted values for ratingScore - Lasso", y=1.05)
plt.xlim(1, 10)
plt.ylim(1, 10)
plt.gca().set_aspect('equal', adjustable='box')
plt.show()
```



1.7 An Attempt at Improvement

With 84593 games, we were forced to shrink our data to 17622 examples to ensure that all features had values.

What if we dropped `priceAverage`, and `priceStdDev`, since the majority of the games did not have values for these features?

This should help to answer which is more important to the creation of a strong model:

- The inclusion of `priceAverage` and `priceStdDev`, potentially predictive features
- The inclusion of more training data to these learners

Repeating the subsetting of data from above, but after removing two columns from the feature vector

```
In [48]: # Drop the two price-based features
no_prices = all_data.drop(['priceAverage', 'priceStdDev'], axis=1)
print("Games initially: {}".format(len(no_prices)))
# Drop all rows with NA in any feature column
no_prices_without_na = no_prices.dropna()
print("Games without NAs: {}".format(len(no_prices_without_na)))
# Filter with the same assumptions as previously
df_no_prices = no_prices_without_na.query("year >= 1950 & ratingCount >= 5")
print("Games after filter: {}".format(len(df_no_prices)))

# Split into features and target
excluded = ["id", "name", "url", "ratingScore", "ratingCount", "ratingStdDev"]
feature_cols = [col for col in df_no_prices.columns if col not in excluded]
X_all_no_prices = df_no_prices[feature_cols]
y_all_no_prices = df_no_prices[target_col]

# Again, an 80/20 split, but from a larger data pool
X_train_no_prices, X_holdout_no_prices, y_train_no_prices, y_holdout_no_prices = train_test_split(
    np.array(X_all_no_prices), np.array(y_all_no_prices),
    train_size=0.8,
    random_state=seed)
```

Games initially: 84593

Games without NAs: 71714

Games after filter: 32240

That's a shame. It seems that the same games that don't have price history information are often the ones with ratings provided by fewer than five raters. I suppose that would make sense, but it also limits our usable data.

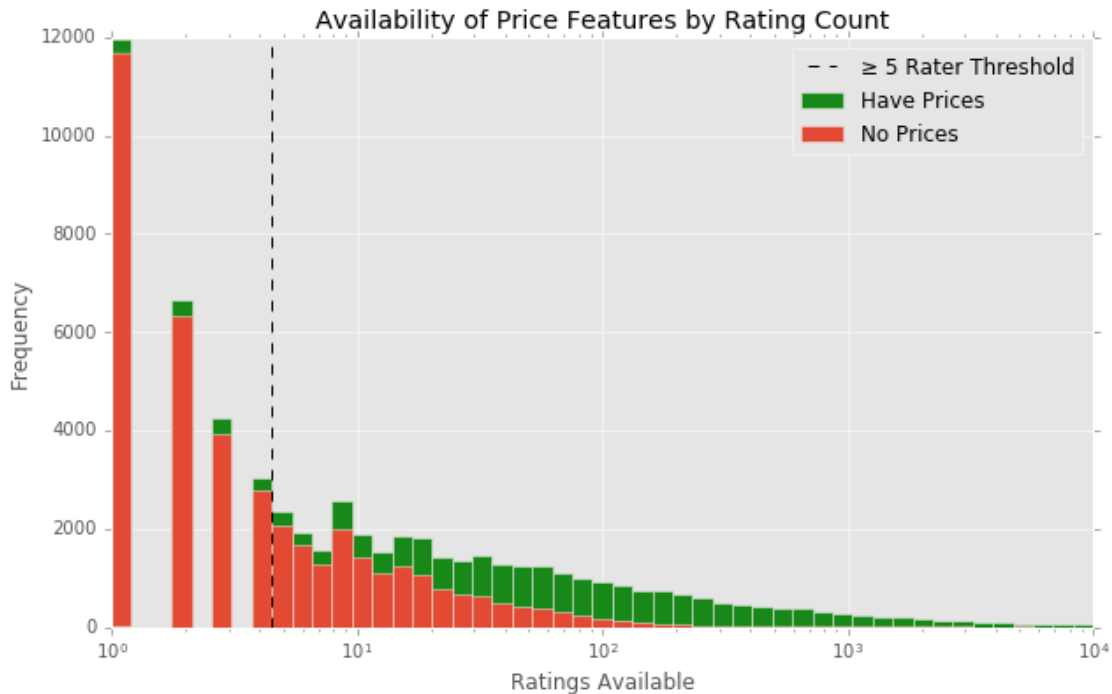
Let's produce a plot to see if what we're inferring is correct.

```
In [49]: # For Unicode '≥' symbol
from __future__ import unicode_literals

plt.figure(figsize=(10,6))
plt.hist(all_data['ratingCount'],
         bins=np.logspace(0, 4, 50), color="#188918", label="Have Prices")
plt.hist(all_data.query("priceAverage != priceAverage")['ratingCount'], # See citation
         bins=np.logspace(0, 4, 50), color="#e24a33", label="No Prices")
# Line at the threshold of 5 raters
plt.plot([4.5, 4.5], [0, 12000], 'k--', label="≥ 5 Rater Threshold")
plt.gca().set_xscale('log')
plt.ylabel("Frequency")
```

```
plt.xlabel("Ratings Available")
plt.legend(loc='upper right')
plt.title("Availability of Price Features by Rating Count")
plt.show()
```

Citation - Hack to use pandas query() to find rows with null value
<http://stackoverflow.com/questions/26535563/querying-for-nan-and-other-names-in-pandas>



Yup. The first time around: - We removed the observations (i.e games or rows) with no price data - the red portion in the chart. - Then we removed the remaining (green) observations with fewer than five raters. This removed a negligible amount of data.

However, the second time around: - We left in the observations with price data (red) - Then again removed the observations with fewer than five raters (red + green to the left of the threshold), which resulted in a large amount of lost data.

It is pretty clear from this chart that once approximately one-hundred raters have participated, the game is nearly guaranteed to have available price data.

Now we're ready to train the child learners, and the stacker.

Note:, I have chosen not to perform a new grid search for new parameter values. This seems a reasonable choice, as the number (and values) of features are nearly the same, and this will save a good deal of time.

```
In [50]: # We should re-calculate the standard scaler with this data (it shouldn't differ by much)
         standard_scaler_no_prices = preprocessing.StandardScaler().fit(X_all_no_prices)

         # So we don't overwrite our models from before.
         from sklearn.base import clone

         # Create the stacking regression learner
         stacker_no_prices = StackedRegression(
             # Child estimators - cloning the previous ones here to not overwrite the original models
```

```

[{'estimator': clone(random_forest_reg)},
 {'estimator': clone(k_neighbors_reg), 'scaler': standard_scaler_no_prices},
 {'estimator': clone(support_vector_reg), 'scaler': standard_scaler_no_prices},
 {'estimator': clone(gradient_boost_reg)}],
# Stacking meta-estimator
{'estimator': clone(lasso_cv_reg)},
# Additional settings, passed to lower-level estimators
cv=kf,
#verbose=1,
random_state=seed,
n_jobs=cpus)

# Now train it
train(stacker_no_prices, X_train_no_prices, y_train_no_prices)

# How were the child learners weighted?
pd.Series(stacker_no_prices.meta_estimator['estimator'].coef_,
          index=[e['estimator'].__class__.__name__ for e in stacker_no_prices.estimators])

# Predict on the holdout set
final_predictions_no_prices = show_holdout_performance(stacker_no_prices, X_holdout_no_prices,

```

Training StackedRegression...

Done.

Training time (secs): 4513.931

StackedRegression

Mean Squared Error: 0.557246071492

Pearson Correlation: 0.765398399292

The correlation of 0.7653 is lower than that of 0.8088 from the model that included price-based features.

It seems that in this case, price-based features provided more predictive power than nearly doubling the size of the training set.

1.8 Playing at Board Game Design

Now that we have a model that predicts crowd-sourced game ratings with reasonable accuracy, can we use it to help us design a best-selling game?

Let's explore the 157-dimensional feature space by creating 100,000 reasonable feature vectors, creating predicted ratings for each, and evaluating the vectors that score well.

```

In [51]: # We want to get consistent results
np.random.seed(seed)
# Number of proposed game feature vectors to create
games_to_create = 100000

# Pre-allocate a pandas DataFrame of n x m (games_to_create observations, by 157 features)
excluded = ["id", "name", "url", "ratingScore", "ratingCount", "ratingStdDev"]
feature_cols = [col for col in all_data.columns if col not in excluded]
proposed_games = pd.DataFrame(0, index=np.arange(games_to_create), columns=feature_cols)
for feature in feature_cols:
    # Randomly choose from all seen values for this feature, 100,000 times
    proposed_games[feature] = np.random.choice(all_data[feature].dropna(), games_to_create)

# Enforce some data rules
# 1. All weight*Pct features for a given row should sum to 100.

```



```

import re
weight_features = filter(re.compile('^weight.*Pct$').match, feature_cols)
proposed_games[weight_features] = np.random.dirichlet(
    np.ones(len(weight_features))/4,
    size=games_to_create) * 100
#temp_weights = pd.DataFrame(np.random.rand(games_to_create, len(weight_features)))
#proposed_games[weight_features] = temp_weights.div(pd.Series(temp_weights.sum(axis=1)), axis=

# 2. weightAvg should reflect the percentages previously computed.
proposed_games['weightAvg'] = (
    (proposed_games['weightLightPct'] / 100 * 1) +
    (proposed_games['weightMediumLightPct'] / 100 * 2) +
    (proposed_games['weightMediumPct'] / 100 * 3) +
    (proposed_games['weightMediumHeavyPct'] / 100 * 4) +
    (proposed_games['weightHeavyPct'] / 100 * 5))

# 3. All subdomain* features for a given row should sum to 100.
subdomain_features = filter(re.compile('^subdomain.*').match, feature_cols)
proposed_games[subdomain_features] = np.random.dirichlet( # See citation
    np.ones(len(subdomain_features))/4,
    size=games_to_create) * 100
#temp_weights = pd.DataFrame(np.random.rand(games_to_create, len(subdomain_features)))
#proposed_games[subdomain_features] = temp_weights.div(pd.Series(temp_weights.sum(axis=1)), ax

# 4. Let's assume that our potential game will be published in 2017.
proposed_games['year'] = np.repeat(2017, games_to_create)

# 5. Let's fix the number of players at between 2 and 4.
proposed_games['playersStatedMin'] = np.repeat(2, games_to_create)
proposed_games['playersStatedMax'] = np.repeat(4, games_to_create)
proposed_games['playersBestMin'] = np.repeat(2, games_to_create)
proposed_games['playersBestMax'] = np.repeat(4, games_to_create)

# 6. Play time should be between 30 and 60 minutes.
proposed_games['playtimeMin'] = np.repeat(30, games_to_create)
proposed_games['playtimeMax'] = np.repeat(60, games_to_create)

# http://stackoverflow.com/questions/18659858/generating-a-list-of-random-numbers-summing-to-1

```

Now we can use this set of 100000 prospective game designs, and see if any will likely receive high scores.

```
In [52]: predicted_ratings = predict(stacker, proposed_games)
```

Predicting target using StackedRegression...

Done.

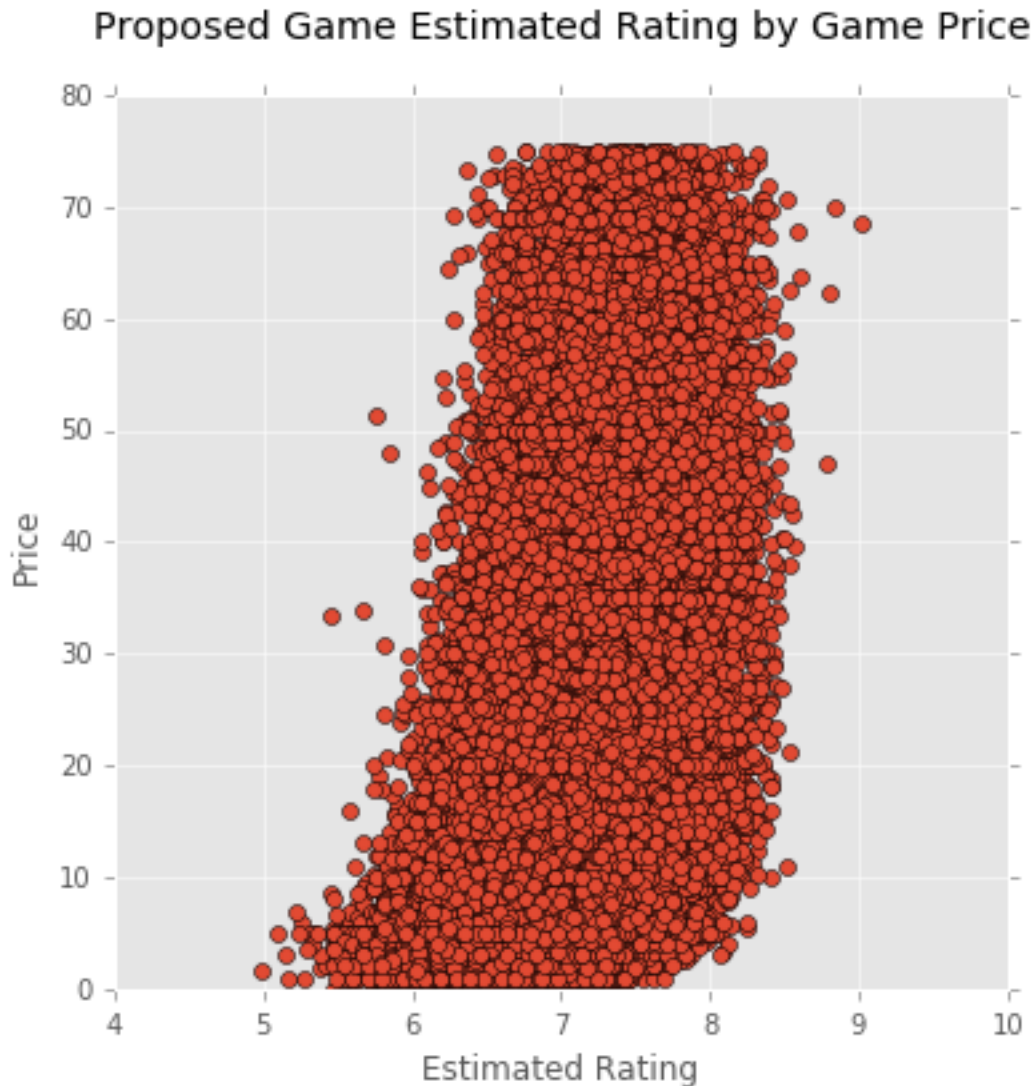
Prediction time (secs): 549.008

Let's plot a comparison of the ratings of these potential games, and their prices, to see if these seem to be related.

```
In [107]: proposed_games['predicted_rating'] = predicted_ratings
proposed_games.sort(columns='predicted_rating', ascending=False)
```

```
# Let's limit our potential games to those under £75, and see how the ratings fall by price.
plt.figure(figsize=(6,6))
```

```
plt.plot(proposed_games.query("priceAverage < 75")['predicted_rating'],
         proposed_games.query("priceAverage < 75")['priceAverage'],
         'o')
plt.xlabel("Estimated Rating")
plt.ylabel("Price")
plt.title("Proposed Game Estimated Rating by Game Price", y=1.05)
plt.show()
```



Price doesn't seem to correlate strongly with rating, until the price drops below \$10.

Let's see if we can isolate some common game categories or mechanics among the best (and worst) rated games.

```
In [151]: # Get some aggregate information about the entire game set
all_mechanics = proposed_games[filter(re.compile('^mechanic.*').match,
                                     feature_cols)].sum(axis=0).order(ascending=False)
all_categories = proposed_games[filter(re.compile('^categor.*').match,
                                     feature_cols)].sum(axis=0).order(ascending=False)
```

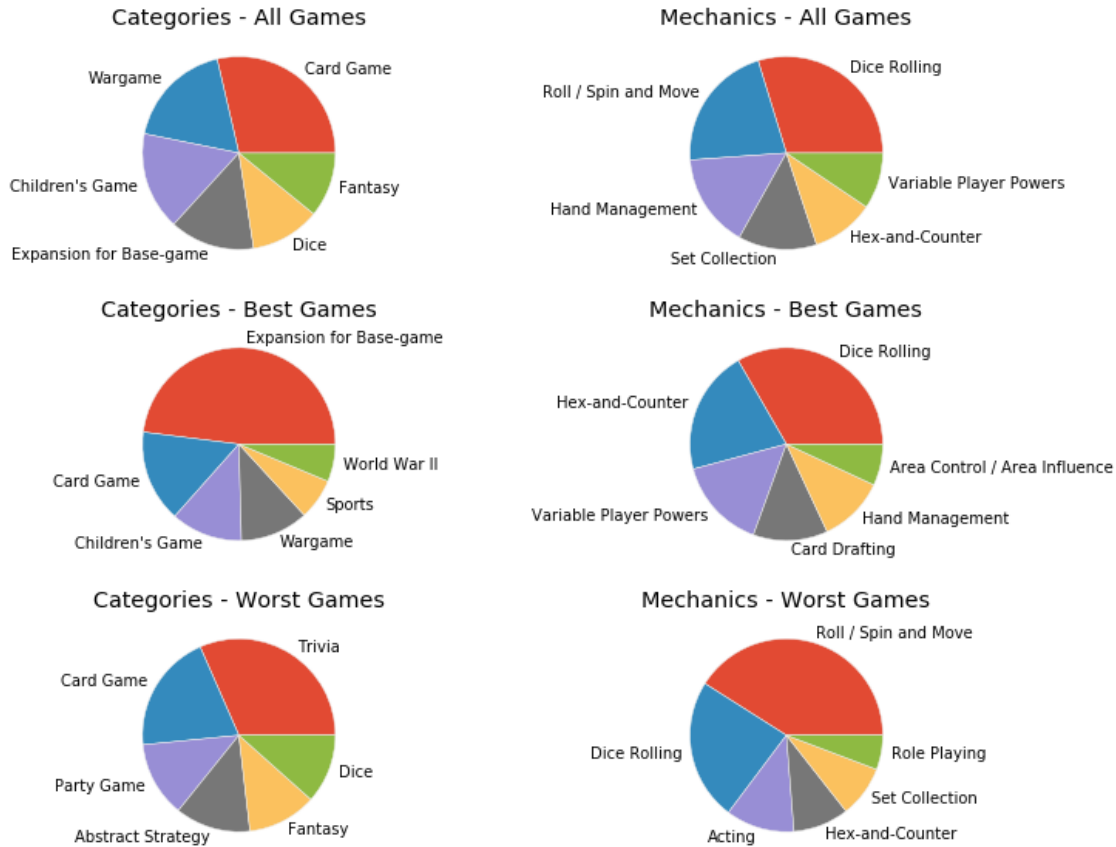
```

# Get some aggregate information about the best-rated X games in this set
limit = 100
best_games = proposed_games.sort(columns='predicted_rating', ascending=False).iloc[0:limit+1]
best_mechanics = best_games[filter(re.compile('^mechanic.*').match,
                                   feature_cols)].sum(axis=0).order(ascending=False)
best_categories = best_games[filter(re.compile('^category.*').match,
                                   feature_cols)].sum(axis=0).order(ascending=False)

# The same for the worst X games
worst_games = proposed_games.sort(columns='predicted_rating', ascending=True).iloc[0:limit+1]
worst_mechanics = worst_games[filter(re.compile('^mechanic.*').match,
                                   feature_cols)].sum(axis=0).order(ascending=False)
worst_categories = worst_games[filter(re.compile('^category.*').match,
                                   feature_cols)].sum(axis=0).order(ascending=False)

# Create some pie charts
r = re.compile('(mechanic|category):')
colors = plt.rcParams['axes.color_cycle']
f, axes = plt.subplots(3, 2, figsize=(10, 8))
f.tight_layout()
for (subplot, frequencies, title) in zip(axes.flat,
    (all_categories, all_mechanics,
     best_categories, best_mechanics,
     worst_categories, worst_mechanics),
    ('Categories - All Games', 'Mechanics - All Games',
     'Categories - Best Games', 'Mechanics - Best Games',
     'Categories - Worst Games', 'Mechanics - Worst Games')):
    # Create a pair of plots for each tuple
    subplot.set_aspect('equal', adjustable='box')
    subplot.pie(frequencies.iloc[0:6],
        labels=[r.sub('', 1) for 1 in frequencies.iloc[0:6].index.tolist()],
        colors=colors)
    subplot.set_title(title)
#Show them all - this might be messy.
plt.show()

```



1.8.1 Common threads

- Card Games as a category are common among all games, and similarly represented in both best, and worst rated games
- Dice rolling as a mechanic is also similarly represented in both best and worst games

1.8.2 Best performer attributes

- The best performing games are disproportionally expansions for base games - perhaps just because hit games are more likely to spawn expansions in the first place
- Children's games and war games are frequently categories for well-ranked games that do not often appear in poorly ranked games
- 'Hex-and-counter', 'variable player powers', and 'card drafting' are game mechanics found more often in well-ranked games

1.8.3 Worst performer attributes

- Trivia and party games seem to be more likely to fare poorly in ratings
- Games with 'roll/spin and move', and 'acting' mechanics are to be avoided.

In [174]: *#Exploring individual games*
Uncomment the next line to see the non-truncated list of all the features for a given game
#pd.options.display.max_rows = 999
`print(worst_games.iloc[1])`

```

year                2017.000000
weightAvg           1.950217
weightLightPct      70.273723
weightMediumLightPct 4.658333
weightMediumPct     3.606147
weightMediumHeavyPct 2.696095
weightHeavyPct      18.765702
playerAgeMin        12.000000
playtimeMin         30.000000
playtimeMax         60.000000
playersStatedMin    2.000000
playersStatedMax    4.000000
playersBestMin      2.000000
playersBestMax      4.000000
priceAverage        5.000000
...
mechanic:Trading    0.000000
mechanic:Trick-taking 0.000000
mechanic:Variable Phase Order 0.000000
mechanic:Variable Player Powers 0.000000
mechanic:Voting     1.000000
mechanic:Worker Placement 0.000000
subdomain:Abstract Strategy Games 4.564371
subdomain:Children's Games 32.107203
subdomain:Customizable Games 0.068859
subdomain:Family Games 13.344570
subdomain:Party Games 0.522216
subdomain:Strategy Games 3.300906
subdomain:Thematic Games 33.762077
subdomain:Wargames 12.329797
predicted_rating    5.097614
Name: 77933, Length: 160, dtype: float64

```

Best game suggestion - though potential outlier: An expansion for a strategy word game with an American civil war theme, priced at \$63. (I'd play that)

A few other reasonable candidates:

- A Napoleonic wargame with a hex-and-counter mechanic for \$70
- A complex adventure/horror card game aimed at families for \$39
- An action/dexterity dice-rolling game with a set-collecting mechanic for \$38
- A children's animal-themed simulation party-game for \$18

Designs to avoid:

- Don't try to market a customizable children's game in both the 'ancient' and 'fighting' categories, involving a stock-trading mechanic
- Please don't. Any any price.
- A thematic children's trivia game that includes dice, voting, and negotiation - for the suggested price of \$5

1.9 Conclusion

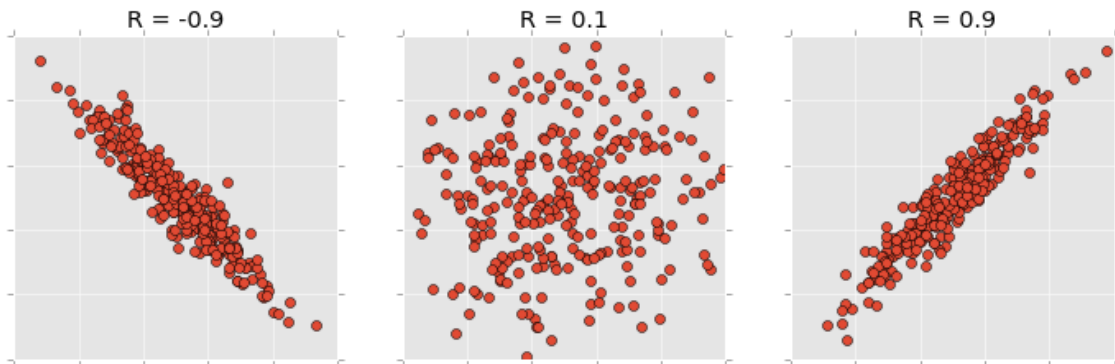
1. I was able to produce a model that achieved a higher Pearson correlation when stacking learners than I was from any learner individually.

2. I was able to surpass the intended threshold of a model whose predictions correlate with those of humans at a value of 0.8 or better.
3. In attempting to train the model with more data, but two fewer features, the model produced fared worse than the original, leading me to the conclusion that the price features have power when predicting the rating of an individual board game.

1.10 Illustrative Plots for the Paper

1.10.1 Correlation Coefficients

```
In [54]: # Settings that apply to all plots
f, (neg, zero, pos) = plt.subplots(1, 3, figsize=(12, 4))
plots = [
    (neg, [[.03, -.03], [.001, .001]]), # Negative correlation
    (zero, [[.05, 0], [0, .05]]),      # Random (near-zero) correlation
    (pos, [[.03, .03], [-.001, .001]]) # Positive correlation
for (subplot, covariance) in (plots):
    # Distributions centered around the midpoint between 0 and 1
    mean = [0.5, 0.5]
    x, y = np.random.multivariate_normal(mean, covariance, 300).T
    subplot.set_aspect('equal', adjustable='box')
    subplot.plot(x, y, 'o')
    subplot.set_xlim(0,1)
    subplot.set_ylim(0,1)
    subplot.axes.get_xaxis().set_ticklabels([])
    subplot.axes.get_yaxis().set_ticklabels([])
    subplot.set_title("R = {0:.1f}".format(pearsonr(x, y)[0]))
#Show them all
plt.show()
```



1.10.2 Comparative Model Performance

```
In [104]: # Provide meaningful column names
performance=performance.rename(columns={0:'Estimator',
                                         1:'Mean Squared Error',
                                         2:'Pearson Correlation'})

# Sort by correlation (ascending)
performance = performance.sort('Pearson Correlation')
print(performance)
```

Estimator	Mean Squared Error	Pearson Correlation	
1	KNeighborsRegressor	0.535480	0.683170
2	SVR	0.408617	0.769350
0	RandomForestRegressor	0.369920	0.794391
3	GradientBoostingRegressor	0.358262	0.800557
4	StackedRegression	0.344842	0.808837

[5 rows x 3 columns]

```
In [103]: index = np.arange(5)
          bar_width = 0.4
          opacity = 0.4
          plt.figure(figsize=(6,6))

          bars1 = plt.bar(index, performance['Mean Squared Error'], bar_width,
                           alpha=opacity,
                           color='b',
                           label='Mean-Squared Error')

          bars2 = plt.bar(index + bar_width, performance['Pearson Correlation'], bar_width,
                           alpha=opacity,
                           color='g',
                           label='Correlation')

          plt.xlabel('Estimator')
          plt.ylabel('Metric Value')
          plt.title('Comparative Model Performance\nMSE and Correlation', y=1.05)
          plt.ylim(0.2,1)
          plt.xticks(index + bar_width, ('KNN', 'SVR', 'RF', 'GBR', 'Stacker'))
          plt.legend()
          plt.show()
```

Comparative Model Performance
MSE and Correlation

