	<pre>X_real array([0.</pre>
	0.70707071, 0.71717172, 0.72727273, 0.73737374, 0.74747475, 0.7575756, 0.767676777, 0.7777778, 0.78787879, 0.7979798, 0.80808081, 0.81818182, 0.82828283, 0.83838384, 0.84848485, 0.55858586, 0.86868687, 0.87878788, 0.8888889, 0.8989899, 0.90909091, 0.91919192, 0.9292293, 0.93939394, 0.94949495, 0.95959596, 0.96969697, 0.97979798, 0.98989899, 1.]) Y_real array([0.00000000e+00, 6.34239197e-02, 1.26592454e-01, 1.89251244e-01, 2.51147987e-01, 3.12033446e-01, 3.71662456e-01, 4.29794912e-01, 4.86196736e-01, 5.40640817e-01, 5.92907929e-01, 6.42787610e-01, 6.90079011e-01, 7.34591709e-01, 7.76146464e-01, 8.14575952e-01, 8.49725430e-01, 8.81453363e-01, 9.09631995e-01, 9.34147860e-01, 9.54902241e-01, 9.71811568e-01, 9.98684776e-01, 9.8388464e-01, 9.98867339e-01, 9.99874128e-01, 9.96854776e-01, 9.88821442e-01, 9.78802446e-01, 9.63842159e-01, 9.45000819e-01, 9.22354294e-01, 8.95993774e-01, 8.66025404e-01, 8.32569855e-01, 7.9761841e-01, 8.95993774e-01, 8.66025404e-01, 9.75761841e-01,
	7.55749574e-01, 7.12694171e-01, 6.66769001e-01, 4.5822652e-01, 4.09030535e-01, 3.42020143e-01, 2.81732557e-01, 2.20310533e-01, 1.58001396e-01, 9.5056043ae-02, 3.17279335e-02, -3.17279335e-02, -9.5056043ae-02, -1.58001396e-01, -2.20310533e-01, -2.81732557e-01, -3.42020143e-01, -4.0930535e-01, -4.58226522e-01, -5.13677392e-01, -5.67059864e-01, -6.18158986e-01, -6.66769001e-01, -7.12694171e-01, -7.55749574e-01, -9.2254294e-01, -9.45000819e-01, -9.63842159e-01, -9.78802446e-01, -9.88251442e-01, -9.96854776e-01, -9.9886739e-01, -9.9886739e-01, -9.9886739e-01, -9.9886739e-01, -9.9838464e-01, -9.84807753e-01, -9.7111568e-01, -9.94500241e-01, -9.34147860e-01, -9.09631995e-01, -8.4575952e-01, -7.76146464e-01, -7.34591709e-01, -6.09079011e-01, -6.42787610e-01, -5.92907929e-01, -5.40640817e-01, -2.51147987e-01, -1.89251244e-01, -1.26592454e-01, -6.34239197e-02, -2.44929360e-16]) import matplotlib.pyplot as plt
	<pre>import seaborn as sns //matplotlib inline sns.set_style('white') sns.set_context('talk') sns.set_palette('dark') # Plot of the noisy (sparse) ax = data.set_index('x')['y'].plot(ls='', marker='o', label='data') ax.plot(X_real, Y_real, ls='', marker=''', label='real function') ax.legend() ax.set(xlabel='x data', ylabel='y data');</pre>
4000	1.0 0.5 0.0 -0.5 -1.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0
\/\ \/\	 Using the PolynomialFeatures class from Scikit-learn's preprocessing library, create 20th order polynomial features. Fit this data using linear regression. Plot the resulting predicted value compared to the calculated data. Interest that PolynomialFeatures requires either a dataframe (with one column, not a Series) or a 2D array of dimension (X , 1), where X is the length. Interest that polynomial features requires either data frame with one column or a 2D array of dimension (X, 1), we want to ensure that we have a two-dimensional array, we can't just pass in just one column or the have to make it in the form of perhaps a data frame, as we see here, or a two-dimensional array. So we import polynomial features, our linear regression. From sklearn.preprocessing import PolynomialFeatures from sklearn.linear_model import LinearRegression # Setup the polynomial features degree = 20
	<pre>pf = PolynomialFeatures(degree) lr = LinearRegression() # use linear regression on that expanded version of X. # Extract the X- and Y- data from the dataframe X_data = data[['x']] # data frame version of that column [[]], if we did a single bracket, we would actually error out here. Y_data = data['y'] # Create the features and fit the model X_poly = pf.fit_transform(X_data) lr = lr.fit(X_poly, Y_data) Y_pred = lr.predict(X_poly) # Plot the result plt.plot(X_data, Y_data, marker='o', ls='', label='data', alpha=1) plt.plot(X_real, Y_real, ls='', label='real function') plt.plot(X_data, Y_pred, marker='o', alpha=.5, label='predictions w/ polynomial features')</pre>
***************************************	plt.legend() ax = plt.gca() ax.set(xlabel='x data', ylabel='y data'); 1.0 0.5 data
:	 -1.0 0.0 0.2 0.4 0.6 0.8 1.0 x data Performing the regression on using the data with polynomial features using ridge regression (α=0.001) and lasso regression (α=0.0001). Plotting the results, as was done in Question 1. Also plotting the magnitude of the coefficients obtained from these regressions, and compare them to those obtained from linear regression in the previous question. The linear regression coefficients will likely need separate plot (or their own y-axis) due to their large magnitude. # Mute the sklearn warning about regularization import warnings warnings.filterwarnings('ignore', module='sklearn')
	<pre>#We're going to say rr equal to our Ridge object, with an alpha = 0.001. #We're then going to fit that to our X_poly that we defined before. #So that's 20 degree polynomial as well as our Y_data. # The ridge regression model rr = Ridge(alpha=0.001) rr = rr.fit(X_poly, Y_data) Y_pred_rr = rr.predict(X_poly) # The lasso regression model lassor = Lasso(alpha=0.0001) lassor = Lasso(alpha=0.0001) lassor = lassor.fit(X_poly, Y_data) Y_pred_lr = lassor.predict(X_poly)</pre>
	# The plot of the predicted values plt.plot(X_data, Y_data, marker='o', ls='', label='data') plt.plot(X_real, Y_real, ls='', label='real function') plt.plot(X_data, Y_pred, label='linear regression', marker='^', alpha=.5) plt.plot(X_data, Y_pred_rr, label='ridge regression', marker='^', alpha=.5) plt.plot(X_data, Y_pred_lr, label='lasso regression', marker='^', alpha=.5) plt.legend() ax = plt.gca() ax.set(xlabel='x data', ylabel='y data');
;	0.5 0.0 -0.5 -1.0 data
c	coefficients = pd.DataFrame() coefficients['linear regression'] = lr.coefravel() coefficients['linear regression'] = rr.coefravel() coefficients['lasso regression'] = lassor.coefravel() coefficients = coefficients.applymap(abs) coefficients.describe() # Huge difference in scale between non-regularized vs regularized regression linear regression ridge regression lasso regression count 2.100000e+01 21.000000 mean 5.777551e+13 2.169397 2.167284 std 6.031941e+13 2.900278 4.706731 min 1.619371e+07 0.000000 0.000000
1 r 1 d	25% 3.416390e+12 0.467578 0.00000 50% 3.621188e+13 1.017272 0.252181 75% 1.070540e+14 2.883507 1.641353 max 1.655347e+14 12.429635 20.176708 (coefficients>0).sum() Linear regression 20 20 120 120 120 120 120 120 120 120 1
	<pre># Setup the dual y-axes ax1 = plt.axes() ax2 = ax1.twinx() # Plot the linear regression data ax1.plot(lr.coefravel(),</pre>
	ax1.set_ylim(-2e14, 2e14) ax2.set_ylim(-25, 25) # Combine the legends h1, l1 = ax1.get_legend_handles_labels() h2, l2 = ax2.get_legend_handles_labels() ax1.legend(h1+h2, l1+l2) ax1.set(xlabel='coefficients',ylabel='linear regression') ax2.set(ylabel='ridge and lasso regression') ax1.set_xticks(range(len(lr.coef_))); 2 1e14 -20 .50
۸۷(linear regression 1
8	 Import the data with Pandas, remove any null values, and one hot encode categoricals. Either Scikit-learn's feature encoders or Pandas get_dummies method can be used. Split the data into train and test sets. Log transform skewed features. Scaling can be attempted, although it can be interesting to see how well regularization works without scaling features. filepath = os.sep.join(data_path + ['Ames_Housing_Sales.csv']) data = pd.read_csv("C:\\Users\\rsnen\\Desktop\\IBM\\Ames_Housing_Sales.csv", sep=',') len(data.columns) data.head()
5 CI	1 1262.0 0.0 0.0 None 3 1Fam TA Gd 978.0 0.0 0.0 Pave 6 1262.0 AllPub 298.0 1976 2 920.0 866.0 0.0 None 3 1Fam TA Mn 486.0 0.0 0.0 Pave 6 920.0 AllPub 0.0 2001 3 961.0 756.0 0.0 None 3 1Fam Gd No 216.0 0.0 0.0 Pave 7 756.0 AllPub 0.0 1915 4 1145.0 1053.0 0.0 None 4 1Fam TA Av 655.0 0.0 0.0 Pave 9 1145.0 AllPub 192.0 2000 Trows × 80 columns Treate a list of categorial data and one-hot encode. Pandas one-hot encoder (get_dummies) works well with data that is defined as a categorical.
I	# Do the one hot encoding data = pd.get_dummies(data, columns=one_hot_encode_cols) data = pd.get_dummies(data, drop_first = True) data.columns Index(['1stFlrSF', '2ndFlrSF', '3SsnPorch', 'BedroomAbvGr', 'BsmtFinSF1',
Tł W	train, test = train_test_split(data, test_size=0.3, random_state=42) here are a number of columns that have skewed featuresa log transformation can be applied to them. Note that this includes the SalePrice, our predictor. However, let's keep that one as is. # Create a list of float colums to check for skewing mask = data.dtypes == np.float float_cols = data.columns[mask] We're then going to take that, filter down here to only those that are above 0.75, we're then going to sort those values from largest to smallest, we're then going to put it into a DataFrame. So from Pandas series, we have the dot translate it into a Pandas DataFrame. Since it's now a new Pandas DataFrame, we're then going to have a default of a column called zero. We're just going to rename that skew. skew_limit = 0.75 skew_vals = train[float_cols].skew() skew_cols = (skew_vals
	Skew_vals
	ScreenPorch 4.599803 BsmtFinSF2 4.466378 EnclosedPorch 3.218303 LotFrontage 3.138032 MasVnrArea 2.492814 OpenPorchSF 2.295489 SalePrice 2.106910 BsmtFinSF1 2.010766 TotalBsmtSF 1.979164 1stFIrSF 1.539692
Γr	GrLivArea 1.455564 WoodDeckSF 1.334388 BsmtUnfSF 0.900308 GarageArea 0.838422 2ndFIrSF 0.773655 ransform all the columns where the skew is greater than 0.75, excluding "SalePrice". # OPTIONAL: Let's look at what happens to one of these features, when we apply np.log1p visually. field = "BsmtFinSF1" fig, (ax_before, ax_after) = plt.subplots(1, 2, figsize=(10, 5)) train[field].hist(ax=ax_before)
	train[field].apply(np.log1p).hist(ax=ax_after) ax_before.set(title='before np.log1p', ylabel='frequency', xlabel='value') ax_after.set(title='after np.log1p', ylabel='frequency', xlabel='value') fig.suptitle('Field "{}".format(field)); # a little bit better before np.log1p
;	# Mute the setting with a copy warnings pd options mode chained_assignment = None
Se	<pre>for col in skew_cols.index.tolist(): if col == "SalePrice": continue train[col] = np.log1p(train[col]) test[col] = test[col].apply(np.log1p) # same thing eparate features from predictor. feature_cols = [x for x in train.columns if x != 'SalePrice'] X_train = train[feature_cols] y_train = train['SalePrice'] X_test = test[feature_cols] y_test = test['SalePrice'] * Writing a function</pre>
	<pre>from sklearn.metrics import mean_squared_error def rmse(ytrue, ypredicted): return np.sqrt(mean_squared_error(ytrue, ypredicted)) • Fit a basic linear regression model • print the root-mean-squared error for this model • plot the predicted vs actual sale price based on the model. from sklearn.linear_model import LinearRegression linearRegression = LinearRegression().fit(X_train, y_train)</pre>
6	<pre>linearRegression_rmse = rmse(y_test, linearRegression.predict(X_test)) print(linearRegression_rmse) 33477.1807710954 f = plt.figure(figsize=(6,6)) ax = plt.axes() ax.plot(y_test, linearRegression.predict(X_test),</pre>
Jistor Dries	ylim=lim, title='Linear Regression Results'); Linear Regression Results 500000 400000 200000
Ri	100000 100000 200000 300000 400000 500000 Actual Price regression uses L2 normalization to reduce the magnitude of the coefficients. This can be helpful in situations where there is high variance. The regularization functions in Scikit-learn each contain versions that bross-validation built in.
N	 Fit a regular (non-cross validated) Ridge model to a range of α values and plot the RMSE using the cross validated error function you created above. Use
1 M eli Si	ridgeCV_rmse = rmse(y_test, ridgeCV.predict(X_test)) print(ridgeCV.alpha_, ridgeCV_rmse) 15.0 32883.672192475 fluch like the RidgeCV function, there is also a LassoCV function that uses an L1 regularization function and cross-validation. L1 regularization will selectively shrink some coefficients, effectively performing feature limination. the LassoCV function does not allow the scoring function to be set. However, the custom error function (rmse) created above can be used to evaluate the error on the final model. imiliarly, there is also an elastic net function with cross validation, ElasticNetCV, which is a combination of L2 and L1 regularization. Fit a Lasso model using cross validation and determine the optimum value for \(\alpha \) and the RMSE using the function created above. Note that the magnitude of \(\alpha \) may be different from the Ridge model.
U:	 Repeat this with the Elastic net model. Compare the results via table and/or plot. Isset the following alphas: [1e-5, 5e-5, 0.0001, 0.0005] from sklearn.linear_model import LassoCV alphas2 = np.array([1e-5, 5e-5, 0.0001, 0.0005]) lassoCV = LassoCV(alphas=alphas2, max_iter=5e4, cv=3).fit(X_train, y_train) lassoCV_rmse = rmse(y_test, lassoCV.predict(X_test))
0 W	print(lassoCV.alpha_, lassoCV_rmse) # Lasso is slower 0.0005 43180.91016592929 We can determine how many of these features remain non-zero. print('Of {} coefficients, {} are non-zero with Lasso.'.format(len(lassoCV.coef_), len(lassoCV.coefnonzero()[0]))) Of 294 coefficients, 275 are non-zero with Lasso. low lets try the elastic net, with the same alphas as in Lasso, and l1_ratios between 0.1 and 0.9 from sklearn.linear_model import ElasticNetCV l1_ratios = np.linspace(0.1, 0.9, 9)
O C	<pre>elasticNetCV = ElasticNetCV(alphas=alphas2,</pre>
E	rmse_df
	<pre>models = [ridgeCV, lassoCV, elasticNetCV] for mod, lab in zip(models, labels): ax.plot(y_test, mod.predict(X_test),</pre>
Prodictor Dries	Linear Regression Results 500000 400000 300000 1000000 0 0
_€ R(-200000 Ridge Lasso ElasticNet 100000 200000 300000 400000 500000 600000 Actual Price et's explore Stochastic gradient descent in this exercise. tecall that Linear models in general are sensitive to scaling. However, SGD is <i>very</i> sensitive to scaling. foreover, a high value of learning rate can cause the algorithm to diverge, whereas a too low value may take too long to converge. Fit a stochastic gradient descent model without a regularization penalty (the relevant parameter is penalty).
	 Now fit stochastic gradient descent models with each of the three penalties (L2, L1, Elastic Net) using the parameter values determined by cross validation above. Do not scale the data before fitting the model. Compare the results to those obtained without using stochastic gradient descent. # Import SGDRegressor and prepare the parameters from sklearn.linear_model import SGDRegressor model_parameters_dict = {
	<pre>'alpha': elasticNetCV.alpha_, 'li_ratio': elasticNetCV.li_ratio_} } new_rmses = {} for modellabel, parameters in model_parameters_dict.items(): # following notation passes the dict items as arguments \$GD = \$GDRegressor(**parameters) \$GD.fit(X_train, y_train) new_rmses[modellabel] = rmse(y_test, \$GD.predict(X_test)) rmse_df['RMSE-SGD'] = pd.Series(new_rmses) rmse_df RMSE RMSE-SGD Linear 633477.180771 6.422430e+15</pre>
V	Ridge 32883.672192 5.236489e+15 Lasso 43180.910166 2.201001e+15 ElasticNet 35206.951603 9.431410e+15 Pass in eta0=1e-7 when creating the instance of SGDClassifier . Re-compute the errors for all the penalties and compare. # Import SGDRegressor and prepare the parameters from sklearn.linear_model import SGDRegressor
	<pre>model_parameters_dict = { 'Linear': {'penalty': 'none'}, 'Lasso': {'penalty': '12',</pre>
E	SGD.fit(X_train, y_train) new_rmses[modellabel] = rmse(y_test, SGD.predict(X_test)) rmse_df['RMSE-SGD-learningrate'] = pd.Series(new_rmses) rmse_df RMSE
	 Fit a MinMaxScaler to X_train create a variable X_train_scaled . Using the scaler, transform X_test and create a variable X_test_scaled . Apply the same versions of SGD to them and compare the results. Don't pass in a eta0 this time. from sklearn.preprocessing import MinMaxScaler scaler = MinMaxScaler() X_train_scaled = scaler.fit_transform(X_train) X_test_scaled = scaler.fit_scaled(X_train) X_test_scaled = scaler.transform(X_test) new_rmses = {} for modellabel, parameters in model_parameters_dict.items(): # following notation passes the dict items as arguments SGD = SGDRegressor(**parameters) SGD.fit(X_train_scaled, y_train) new_rmses[modellabel] = rmse(y_test, SGD.predict(X_test_scaled))
E	SGD.fit(X_train_scaled, y_train) new_rmses[modellabel] = rmse(y_test, SGD.predict(X_test_scaled)) rmse_df['RMSE-SGD-scaled'] = pd.Series(new_rmses) rmse_df RMSE
	<pre>scaler = MinMaxScaler() X_train_scaled = scaler.fit_transform(X_train) X_test_scaled = scaler.transform(X_test) new_rmses = {} for modellabel, parameters in model_parameters_dict.items(): # following notation passes the dict items as arguments SGD = SGDRegressor(**parameters) SGD fit(X_train_scaled, Y_train) new_rmses[modellabel] = rmse(y_test, SGD.predict(X_test_scaled)) rmse_df['RMSE-SGD-scaled'] = pd.Series(new_rmses) rmse_df</pre> RMSE_RMSE-SGD-RMSE-SGD-learningrate RMSE-SGD-scaled
	RMSE RMSE-SGD RMSE-SGD-learningrate RMSE-SGD-scaled Linear 633477.180771 6.422430e+15 72821.103928 33399.992821 Ridge 32883.672192 5.236489e+15 75389.918918 77772.380255 Lasso 43180.910166 2.201001e+15 75155.270525 33237.088478 ElasticNet 35206.951603 9.431410e+15 74218.759038 33437.882444