Spring 2023 Data C100/C200 Final Reference Sheet

Pandas

Suppose df is a DataFrame; s is a Series. import pandas as pd

| Function | Description |
|--|---|
| df[col] | Returns the column labeled col from df as a Series. |
| df[[col1, col2]] | Returns a DataFrame containing the columns labeled col1 and col2. |
| <pre>s.loc[rows] / df.loc[rows, cols]</pre> | Returns a Series/DataFrame with rows (and columns) selected by their index values. |
| <pre>s.iloc[rows] / df.iloc[rows, cols]</pre> | Returns a Series/DataFrame with rows (and columns) selected by their positions. |
| <pre>s.isnull() / df.isnull()</pre> | Returns boolean Series/DataFrame identifying missing values |
| <pre>s.fillna(value) / df.fillna(value)</pre> | Returns a Series/DataFrame where missing values are replaced by value |
| <pre>s.isin(values) / df.isin(values)</pre> | Returns a Series/DataFrame of booleans indicating if each element is in values. |
| df.drop(labels, axis) | Returns a DataFrame without the rows or columns named labels along axis (either 0 or 1) |
| <pre>df.rename(index=None, columns=None)</pre> | Returns a DataFrame with renamed columns from a dictionary index and/or columns |
| df.sort_values(by, ascending=True) | Returns a DataFrame where rows are sorted by the values in columns by |
| <pre>s.sort_values(ascending=True)</pre> | Returns a sorted Series. |
| <pre>s.unique()</pre> | Returns a NumPy array of the unique values |
| <pre>s.value_counts()</pre> | Returns the number of times each unique value appears in a Series |
| <pre>pd.merge(left, right, how='inner', on='a')</pre> | Returns a DataFrame joining left and right on the column labeled a; the join is of type inner |
| <pre>left.merge(right, left_on=col1, right_on=col2)</pre> | Returns a DataFrame joining left and right on columns labeled coll and col2. |
| df.pivot_table(index, columns, values=None, aggfunc='mean') | Returns a DataFrame pivot table where columns are unique values from columns (column name or list), and rows are unique values from index (column name or list); cells are collected values using aggfunc. If values is not provided, cells are collected for each remaining column with multi-level column indexing. |
| df.set_index(col) | Returns a DataFrame that uses the values in the column labeled col as the row index. |
| df.reset_index() | Returns a DataFrame that has row index 0, 1, etc., and adds the current index as a column. |

Let grouped = df.groupby(by) where by can be a column label or a list of labels.

| For a block | Provide the second s | |
|---|--|--|
| Function | Description | |
| <pre>grouped.count()</pre> | Return a Series containing the size of each group, excluding missing values | |
| <pre>grouped.size()</pre> | Return a Series containing size of each group, including missing values | |
| <pre>grouped.mean()/.min()/.max()</pre> | Return a Series/DataFrame containing mean/min/max of each group for each column, excluding missing values | |
| <pre>grouped.filter(f) grouped.agg(f)</pre> | Filters or aggregates using the given function f | |
| Function | Description | |
| s.str.len() | Returns a Series containing length of each string | |
| s.str[a:b] | Returns a Series where each element is a slice of the corresponding string indexed from a (inclusive, optional) to b (non-inclusive, optional) | |
| <pre>s.str.lower()/s.str.upper()</pre> | Returns a Series of lowercase/uppercase versions of each string | |
| .str.replace(pat, repl) | Returns a Series that replaces occurences of substrings matching the regex pat with string repl | |
| .str.contains(pat) | Returns a boolean Series indicating if a substring matching the regex pat is contained in each string | |
| s.str.extract(pat) | Returns a Series of the first subsequence of each string that matches the regex pat. If pat contains on group, then only the substring matching the group is extracted | |

Visualization

Function

plt.plot(x, y)

plt.scatter(x, y)

plt.bar(x, height)

plt.hist(x, bins=None)

Matplotlib: x and y are sequences of values. import matplotlib.pyplot as plt

Tukey-Mosteller Bulge Diagram.

| Description | Y ³ ,Y ² |
|--|--------------------------------|
| Creates a line plot of x against y | $\gamma \mid \chi$ |
| Creates a scatter plot of x against y | $\log X \frac{1}{\sqrt{2}}$ |
| Creates a histogram of x; bins can be an integer or a sequence | |
| Creates a bar plot of categories x and corresponding heights | $\lambda \mid \lambda$ |
| height | \sqrt{Y} , log Y |

Seaborn: x and y are column names in a DataFrame data. import seaborn as sns

| Function | Description |
|--|--|
| <pre>sns.countplot(data, x)</pre> | Create a barplot of value counts of variable x from data |
| <pre>sns.histplot(data, x, stat='count', kde=False) sns.displot(x, data, stat='count', rug=True, kde=True)</pre> | Creates a histogram of x from data, where bin statistics stat is one of 'count', 'frequency', 'probability', 'percent', and 'density'; optionally overlay a kernel density estimator. displot is similar but can optionally overlay a rug plot |
| sns.rugplot(data, x) | Adds a rug plot on the x-axis of variable x from data |
| sns.boxplot(data, x=None, y) sns.violinplot(data, x=None, y) | Create a boxplot of y, optionally factoring by categorical x, from data. violinplot is similar but also draws a kernel density estimator of y |
| sns.scatterplot(data, x, y) | Create a scatterplot of x versus y from data |
| <pre>sns.lmplot(x, y, data, fit_reg=True)</pre> | Create a scatterplot of ${\sf x}$ versus ${\sf y}$ from data, and by default overlay a least-squares regression line |
| sns.jointplot(x, y, data, kind) | Combine a bivariate scatterplot of x versus y from data, with univariate density plots of each variable overlaid on the axes; kind determines the visualization type for the distribution plot, can be scatter, kde or hist |

Regular Expressions

| Description | Operator | Description |
|--|---|---|
| Matches any character except \n | * | Matches preceding character/group zero or more times |
| Escapes metacharacters | ? | Matches preceding character/group zero or one times |
| Matches expression on either side of expression; has lowest priority of any operator | + | Matches preceding character/group one or more times |
| Predefined character group of digits (0-9), alphanumerics (a-z, A-Z, 0-9, and underscore), or whitespace, respectively | ^, \$ | Matches the beginning and end of the line, respectively |
| Inverse sets of d, w, s , respectively | () | Capturing group used to create a sub-expression |
| Matches preceding character/group exactly m times | [] | Character class used to match any of the specified characters or range (e.g. [abcde] is equivalent to [a-e]) |
| most ${\bf n}$ times. If either ${\bf m}$ or ${\bf n}$ are omitted, set lower/upper | [^] | Invert character class; e.g. [^a–c] matches all characters except a, b, c |
| | Matches any character except \n Escapes metacharacters Matches expression on either side of expression; has lowest priority of any operator Predefined character group of digits (0-9), alphanumerics (a-z, A-Z, 0-9, and underscore), or whitespace, respectively Inverse sets of \d, \w, \s, respectively Matches preceding character/group exactly m times Matches preceding character/group at least m times and at most n times. If either m or n are omitted, set lower/upper | Matches any character except \n * Escapes metacharacters ? Matches expression on either side of expression; has lowest priority of any operator + Predefined character group of digits (0-9), alphanumerics ^, \$ (a-z, A-Z, 0-9, and underscore), or whitespace, respectively ^, \$ Inverse sets of \d, \w, \s, respectively () Matches preceding character/group exactly m times [] |

Modified lecture example for capture groups:

import re

lines = '169.237.46.168 - - [26/Jan/2014:10:47:58 -0800] "GET ... HTTP/1.1"' re.findall(r'\[\d+\/(\w+)\/\d+:\d+:\d+ .+\]', line) # returns ['Jan']

| Function | Description |
|--|--|
| <pre>re.match(pattern, string)</pre> | Returns a match if zero or more characters at beginning of string matches pattern, else None |
| <pre>re.search(pattern, string)</pre> | Returns a match if zero or more characters anywhere in string matches pattern, else None |
| <pre>re.findall(pattern, string)</pre> | Returns a list of all non-overlapping matches of pattern in string (if none, returns empty list) |
| <pre>re.sub(pattern, repl, string)</pre> | Returns string after replacing all occurrences of pattern with repl |

Modeling

| Concept | Formula | Concept | Formula |
|------------------------|---|--|---|
| Variance, σ_x^2 | $\frac{1}{n}\sum_{i=1}^n (x_i-\bar{x})^2$ | Correlation <i>r</i> | $r=rac{1}{n}\sum_{i=1}^n rac{x_i-ar{x}}{\sigma_x}rac{y_i-ar{y}}{\sigma_y}$ |
| L_1 loss | $L_1(y,\hat{y}) = \mid y - \hat{y} \mid$ | Linear regression estimate of \boldsymbol{y} | $\hat{y}=	heta_0+	heta_1 x$ |
| $L_2 \log$ | $L_2(y,\hat{y})=(y-\hat{y})^2$ | Least squares linear regression | $\hat{	heta}_0 = ar{y} - \hat{	heta}_1 ar{x} \qquad \hat{	heta}_1 = r rac{\sigma_y}{\sigma_x}$ |

Empirical risk with loss ${\cal L}$

$$R(heta) = rac{1}{n}\sum_{i=1}^n L(y_i, \hat{y_i})$$

Ordinary Least Squares

Multiple Linear Regression Model: $\hat{\mathbb{Y}} = \mathbb{X}\theta$ with design matrix \mathbb{X} , response vector \mathbb{Y} , and predicted vector $\hat{\mathbb{Y}}$. If there are p features plus a bias/intercept, then the vector of parameters $\theta = [\theta_0, \theta_1, \dots, \theta_p]^T \in \mathbb{R}^{p+1}$. The vector of estimates $\hat{\theta}$ is obtained from fitting the model to the sample (\mathbb{X}, \mathbb{Y}) .

| Concept | Formula | Concept | Formula |
|---|--|---|---|
| Mean squared error | $R(heta) = rac{1}{n} \mathbb{Y} - \mathbb{X}	heta _2^2$ | Normal equation | $\mathbb{X}^T\mathbb{X}\hat{\theta}=\mathbb{X}^T\mathbb{Y}$ |
| Least squares estimate, if $\ensuremath{\mathbb{X}}$ is full rank | $\hat{	heta} = (\mathbb{X}^T \mathbb{X})^{-1} \mathbb{X}^T \mathbb{Y}$ | Residual vector, e | $e=\mathbb{Y}-\hat{\mathbb{Y}}$ |
| | | Multiple R^2 (coefficient of determination) | $R^2 = rac{	ext{variance of fitted values}}{	ext{variance of } y}$ |
| Ridge Regression L2 Regularization | $rac{1}{n} \mathbb{Y}-\mathbb{X}	heta _2^2+lpha 	heta _2^2$ | Squared L2 Norm of $	heta \in \mathbb{R}^d$ | $ 	heta _2^2 = \sum_{j=1}^d 	heta_j^2$ |
| Ridge regression estimate (closed form) | $\hat{	heta}_{	ext{ridge}} = (\mathbb{X}^T \mathbb{X} + nlpha I)^{-1} \mathbb{X}^T \mathbb{Y}$ | | |
| LASSO Regression L1 Regularization | $rac{1}{n} \mathbb{Y}-\mathbb{X}	heta _2^2+lpha 	heta _1$ | L1 Norm of $	heta \in \mathbb{R}^d$ | $ 	heta _1 = \sum_{j=1}^d 	heta_j $ |

Scikit-Learn

Package: sklearn.linear_model

| Linear Regression | Logistic Regression | Function(s) | Description |
|--------------------------|------------------------|---|---|
| \checkmark | - | LinearRegression(fit_intercept=True) | Returns an ordinary least squares Linear Regression model. |
| - | \checkmark | <pre>LogisticRegression(fit_intercept=True, penalty='l2', C=1.0)</pre> | Returns an ordinary least squares Linear Regression model. Hyperparameter C is inverse of regularization parameter, C = 1/ λ . |
| \checkmark | - | LassoCV(),RidgeCV() | Returns a Lasso (L1 Regularization) or Ridge (L2 regularization) linear model, respectively, and picks the best model by cross validation. |
| \checkmark | \checkmark | <pre>model.fit(X, y)</pre> | Fits the scikit-learn model to the provided X and y. |
| \checkmark | \checkmark | <pre>model.predict(X)</pre> | Returns predictions for the X passed in according to the fitted model. |
| \checkmark | \checkmark | <pre>model.predict_proba(X)</pre> | Returns predicted probabilities for the X passed in according to the fitted model. If binary classes, will return probabilities for both class 0 and 1. |
| \checkmark | \checkmark | model.coef_ | Estimated coefficients for the linear model, not including the intercept term. |
| \checkmark | \checkmark | model.intercept_ | Bias/intercept term of the linear model. Set to 0.0 if fit_intercept=False. |
| Package: <mark>sk</mark> | learn.model | _selection | |

Description

Function

train_test_split(*arrays, test_size=0.2) Returns two random subsets of each array passed in, with 0.8 of the array in the first subset and 0.2 in the second subset.

Probability

Let X have a discrete probability distribution P(X = x). X has expectation $\mathbb{E}[X] = \sum_x x P(X = x)$ over all possible values x, variance $\operatorname{Var}(X) = \mathbb{E}[(X - \mathbb{E}[X])^2]$, and standard deviation $\operatorname{SD}(X) = \sqrt{\operatorname{Var}(X)}$.

The covariance of two random variables X and Y is $\mathbb{E}[(X - \mathbb{E}[X])(Y - \mathbb{E}[Y])]$. If X and Y are independent, then Cov(X, Y) = 0.

| Notes | Property of Expectation | Property of Variance |
|--|---|--|
| X is a random variable. | | $\operatorname{Var}(X) = E[X^2] - (E[X])^2$ |
| X is a random variable, $a,b\in\mathbb{R}$ are scalars. | $\mathbb{E}[aX+b] = a\mathbb{E}[X] + b$ | $\operatorname{Var}(aX+b)=a^2\operatorname{Var}(X)$ |
| X,Y are random variables. | $\mathbb{E}[X+Y] = \mathbb{E}[X] + \mathbb{E}[Y]$ | $\operatorname{Var}(X+Y) = \operatorname{Var}(X) + \operatorname{Var}(Y) + 2\operatorname{Cov}(X,Y)$ |
| X is a Bernoulli random variable that takes on value 1 with probability p and 0 otherwise. | $\mathbb{E}[X]=p$ | $\operatorname{Var}(X) = p(1-p)$ |

Central Limit Theorem

Let (X_1,\ldots,X_n) be a sample of independent and identically distributed random variables drawn from a population with mean μ and standard deviation σ . The sample mean $\overline{X}_n = \sum_{i=1}^n X_i$ is normally distributed, where $\mathbb{E}[\overline{X}_n] = \mu$ and $\mathrm{SD}(\overline{X}_n) = \sigma/\sqrt{n}$.

Parameter Estimation and Gradient Descent

Parameter Estimation

Suppose for each individual with fixed input x, we observe a random response $Y = g(x) + \epsilon$, where g is the true relationship and ϵ is random noise with zero mean and variance σ^2 .

For a new individual with fixed input x, define our random prediction $\hat{Y}(x)$ based on a model fit to our observed sample (\mathbb{X}, \mathbb{Y}) . The model risk is the mean squared prediction error between Y and $\hat{Y}(x)$: $\mathbb{E}[(Y - \hat{Y}(x))^2] = \sigma^2 + \left(\mathbb{E}[\hat{Y}(x)] - g(x)\right)^2 + \operatorname{Var}(\hat{Y}(x))$.

Suppose that input x has p features and the true relationship g is linear with parameter $\theta \in \mathbb{R}^{p+1}$. Then $Y = f_{\theta}(x) = \theta_0 + \sum_{j=1}^p \theta_j x_j + \epsilon$ and $\hat{Y} = f_{\hat{\theta}}(x)$ for an estimate $\hat{\theta}$ fit to the observed sample (\mathbb{X}, \mathbb{Y}) .

Gradient Descent

Let $L(\theta, \mathbb{X}, \mathbb{Y})$ be an objective function to minimize over θ , with some optimal $\hat{\theta}$. Suppose $\theta^{(0)}$ is some starting estimate at t = 0, and $\theta^{(t)}$ is the estimate at step t. Then for a learning rate α , the gradient update step to compute $\theta^{(t+1)}$ is $\theta^{(t+1)} = \theta^{(t)} - \alpha \nabla_{\theta} L(\theta^{(t)}, \mathbb{X}, \mathbb{Y})$, where $\nabla_{\theta} L(\theta^{(t)}, \mathbb{X}, \mathbb{Y})$ is the partial derivative/gradient of L with respect to θ , evaluated at $\theta^{(t)}$. SQL

SQLite syntax:

SELECT [DISTINCT]
 {* | expr [[AS] c_alias]
 {,expr [[AS] c_alias] ...}}
FROM tableref {, tableref}
[[INNER | LEFT] JOIN table_name
 ON qualification_list]
[WHERE search_condition]
[GROUP BY colname {,colname...}]
[HAVING search_condition]
[ORDER BY column_list]
[LIMIT number]
[OFFSET number of rows];

| Syntax | Description | |
|---------------------------------------|--|--|
| SELECT column_expression_list | List is comma-separated. Column expressions may include aggregation functions (MAX, FIRST, COUNT, etc). AS renames columns. DISTINCT selects only unique rows. | |
| FROM s INNER JOIN t ON cond | Inner join tables s and t using cond to filter rows; the INNER keyword is optional. | |
| FROM s LEFT JOIN t ON cond | Left outer join of tables s and t using cond to filter rows. | |
| FROM s, t | Cross join of tables ${\sf s}$ and ${\sf t}$: all pairs of a row from ${\sf s}$ and a row from ${\sf t}$ | |
| WHERE a IN cons_list | Select rows for which the value in column a is among the values in a cons_list. | |
| ORDER BY RANDOM LIMIT n | Draw a simple random sample of n rows. | |
| ORDER BY a, b DESC | Order by column a (ascending by default) , then b (descending). | |
| CASE WHEN pred THEN cons ELSE alt END | Evaluates to cons if pred is true and alt otherwise. Multiple WHEN/THEN pairs can be included, and ELSE is optional. | |
| WHERE s.a LIKE 'p' | Matches each entry in the column a of table s to the text pattern p. The wildcard $\%$ matches at least zero characters. | |
| LIMIT number | Keep only the first number rows in the return result. | |
| OFFSET number | Skip the first number rows in the return result. | |

Principal Component Analysis (PCA)

The *i*-th Principal Component of the matrix X is defined as the *i*-th column of $U\Sigma$ defined by Singular Value Decomposition (SVD).

 $X = U\Sigma V^T$ is the SVD of X if U and V^T are matrices with orthonormal columns and Σ is a diagonal matrix. The diagonal entries of Σ , $[s_1, \ldots, s_r, 0, \ldots, 0]$, are known as singular values of X, where $s_i > s_j$ for i < j and $r = \operatorname{rank}(X)$.

Define the design matrix $X \in \mathbb{R}^{n \times p}$. Define the total variance of X as the sum of individual variances of the p features. The amount of variance captured by the *i*-th principal component is equivalent to s_i^2/n_i , where n is the number of datapoints.

| Syntax | Description |
|---|---|
| <pre>np.linalg.svd(X, full_matrices = True)</pre> | SVD of X with shape (M, N) that returns u, s, vt, where s is a 1D array of X's singular values. If full_matrices=True, u and vt have shapes (M, M) and (N, N) respectively; otherwise shapes are (M, K) and (K, N) , respectively, where K = min(M, N). |
| Classification and Lagistic Degrapsion | |

Classification and Logistic Regression

Confusion Matrix

Columns are the predicted values \hat{y} and rows are the actual classes y.

| | $\hat{y}=0$ | $\hat{y}=1$ |
|-------|---------------------|---------------------|
| y = 0 | True negative (TN) | False Positive (FP) |
| y = 1 | False negative (FN) | True Positive (TP) |

Classification Performance

Suppose you predict n datapoints.

| Metric | Formula | Other Names |
|------------|--------------------|----------------------------------|
| Accuracy | $\frac{TP+TN}{n}$ | |
| Precision | $\frac{TP}{TP+FP}$ | |
| Recall/TPR | $\frac{TP}{TP+FN}$ | True Positive Rate, Sensitivity |
| FPR | $\frac{FP}{FP+TN}$ | False Positive Rate, Specificity |

An ROC curve visualizes TPR vs. FPR for different thresholds T.

Logistic Regression Model: For input feature vector x, $\hat{P}_{\theta}(Y = 1|x) = \sigma(x^T\theta)$, where $\sigma(z) = 1/(1 + e^{-z})$. The estimate $\hat{\theta}$ is the parameter θ that minimizes the average cross-entropy loss on training data. For a single datapoint, define cross-entropy loss as $-[y \log(p) + (1 - y) \log(1 - p)]$, where p is the probability that the response is 1.

Logistic Regression Classifier: For a given input x and trained logistic regression model with parameter θ , compute $p = \hat{P}(Y = 1|x) = \sigma(x^T\theta)$. predict response \hat{y} with classification threshold T as follows:

$$\hat{y} = ext{classify}(x) = egin{cases} 1 & p \geq T \ 0 & ext{otherwise} \end{cases}$$

Clustering

K-Means Clustering: Pick an arbitrary k, and randomly place k "centers", each a different color. Then repeat until convergence:

- 1. Color points according to the closest center (defined as squared distance).
- 2. Move center for each color to center of points with that color.

K-Means minimizes inertia, defined as the sum of squared distances from each datapoint to its center.

Agglomerative Clustering: Assign each datapoint to its own cluster. Then, recursively merge pairs of clusters together until there are k clusters remaining.

A datapoint's silhouette score S is defined as $S = (B - A) / \max(A, B)$, where A is the mean distance to other points in its cluster, and B is the mean distance to points in its closest cluster.