CS 6140 Machine Learning: Assignment - 3 (Total Points: 100)

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Q1. Implementing Decision Tree from Scratch (60 points)

You will implement the decision tree algorithm for the mushroom dataset. The dataset is provided to you with this assignment (mushroom.csv). The objective is to classify whether a mushroom is poisonous or edible based on various features. You will write a program that uses some algorithm to construct a decision tree, and use the tree to classify new instances. Your implementation should include the following functions:

- 1. EDA on dataset
- 2. Calculate the entropy of a dataset
- 3. Split a dataset based on an attribute and its value
- 4. Select the attribute to split the dataset on
- 5. Build a decision tree
- 6. Classify an instance using the decision tree
- 7. Calculate testing accuracy.

Note: Perform EDA at steps like splitting data etc.

EDA on dataset

```
In [1]: import logging
   import pandas as pd
   import seaborn as sb
   import sklearn as sk
   import numpy as np
   import matplotlib.pyplot as plt

from collections import Counter
   from sklearn.preprocessing import LabelEncoder
```

```
from sklearn.model_selection import train_test_split
from sklearn.metrics import accuracy_score
from sklearn.linear_model import LinearRegression
from sklearn.tree import DecisionTreeClassifier
```

```
In [2]: def stack_bar_break_out_by_edibility(dataframe, attribute: str):
            Drawing a stack bar breakout by editibility. This would help us analyze what
            try:
                if attribute == 'class':
                    sb.set(font scale=1.4)
                    dataframe[f'{attribute}'].value_counts().plot(kind='bar', figsize=(7, 6), rot=0, color=['blue', 'red']);
                    plt.title("Count of Poisonsous vs Edible mushrooms", y=1.02);
                else:
                    dataframe.groupby([f'{attribute}', 'class']).size().unstack().plot(kind='bar',
                                                                                        stacked=True.
                                                                                        figsize=(7, 6),
                                                                                        rot=0,
                                                                                        color=['blue', 'red'])
                    plt.title(f'Count {str(attribute).upper()} break out by Edibility');
                plt.show()
            except Exception as err:
                logging.error(f'Fail with error: {err}', exc_info = True)
                raise err
        def read csv():
            Read mushrooms.csv
            read file = pd.read csv(r'C:\Users\Leaksmy Heng\Documents\GitHub\cs6140\HW3\mushrooms.csv')
            return read_file
        def conduct eda():
            Read csv and conduct the exploratory data analysis.
            @param file path: file path that we want to read
            @type file_path: string
            @return read file
            @rtype: dataframe
            read_file = read_csv()
```

```
# understand the shape of our data
shape_of_data = read_file.shape
total_number_of_rows = shape_of_data[0]
total_number_of_columns = shape_of_data[1]
print(shape of data)
# understand the data, and its metrics
describe data = read file.describe().T
print(describe_data)
# understand the data type
datatype = read_file.info()
print(datatype)
# display a frew row
sample = read file.head()
print(sample)
# examine count of labelling data and build a stack graph
list_of_columns = read_file.columns
for row in list of columns:
    stack_bar_break_out_by_edibility(read_file, row)
# based on the stack bar, there are only 1 veil-type. And the veil type is both edible and poison
# hence we could drop that
read_file = read_file.drop('veil-type', axis = 1)
# new data frame after dropping
df = read_file.columns
# encode the column
encoder = LabelEncoder()
for column in df:
    read_file[column] = encoder.fit_transform(read_file[column])
# check if there are any null value
null_val = read_file.isnull().sum()
print(null_val)
# checking correlation now that our data is encoded
corr = read file.corr()
sb.heatmap(corr, square=True)
return read_file
```

Build a decision tree to include:

```
In [3]: class DecisionTree:
            Creating a decision tree using entropy.
            def __init__(self, max_depth=None, min_samples_split=2, min_impurity=1e-7):
                Create a constructor for Decision Tree using CART classification and regression trees.
                 @param max depth: max depthh hyperparameter that is used to stop the recursive when splitting data
                 @param min samples split: another hyperparameter that is used to stop the recursive when splitting data
                 @param min impurity: another hyperparameter that is used to stop the recursive when splitting data,
                 default value of 0.0000001
                self.max depth = max depth
                 self.min samples split = min samples split
                 self.min impurity = min impurity
                 self.tree = {}
            def split_data(self, x_training, feature, threshold):
                Splitting the training set into 2 subsets using a feature and a threshold.
                This will be used recursively to split the sub-subsets.
                 @param x training: training data set
                 @param feature: feature
                 @param threshold: threshold. Example is the mushroom eatable?
                left = np.where(x training[:, feature] <= threshold)</pre>
                 right = np.where(x training[:, feature] > threshold)
                 return left[0], right[0]
            def entropy(self, y):
                 calculating the entropy of the labeling data.
                 @param y: labeling data
                 label, each_label_counts = np.unique(y, return_counts=True)
                 probability = each label counts / np.sum(each label counts)
                 entropy = -np.sum(probability * np.log2(probability))
```

```
return entropy
def information gain(self, y, left y, right y):
    calculating information gain.
   @param y: labeling data
    @param left y: labeling data on the left child node after splitting
    @param right y: labeling data on the right child node after splitting
   left probability = len(left y) / len(y)
    right probability = len(right y) / len(y)
    entropy_parent = self.entropy(y)
    entropy left = self.entropy(left y)
    entropy right = self.entropy(right y)
    information_gain = entropy_parent - left_probability*entropy_left - right_probability*entropy_right
    return information gain
def select_attribute(self, x_training, y_training):
   Selecting the pair with best feature & best threshold (k, tk) that produce the purest subsets.
   To select the pair, we need to use an information_gain to check.
   @param x_training: training data set
    @param y training: label training data set
   best feature = None
   best threshold = None
   max info gain = -np.inf
   features = x_training.shape[1]
    # check each feature to find the best one with its threshold
   for feature in range(features):
        each feature = x training[:, feature]
        thresholds = np.unique(each_feature)
        for threshold in thresholds:
            left_node, right_node = self.split_data(x_training, feature, threshold)
            if len(left node) == 0 or len(right node) == 0:
                continue
            # calculating information gain
            info_gain = self.information_gain(y_training, y_training[left_node], y_training[right_node])
```

```
if info gain > max info gain:
                best_feature = feature
                best threshold = threshold
                max_info_gain = info_gain
    if max_info_gain < self.min_impurity:</pre>
        best feature = None
        best threshold = None
    return best feature, best threshold
def majority class(self, y):
    Determine the majority class in a given set of labels.
    @param y training: label training data set
   counter = Counter(y)
    majority_class = counter.most_common(1)[0][0]
    return majority_class
def fit(self, x_training, y_training, depth = 0):
   Training the dataset.
   @param x_training: training data set that exclude the labeling data
   @param y training: labeling training data set
   training_data, training_features = x_training.shape
    classes = len(np.unique(y training))
   # stop training the data if the depth of the tree is equal to the maximum depth
   # or if the training data set is less than the sample data for splitting
    # or if the number of classes (edible vs poisonous) is == 1
    if (depth == self.max_depth) or (training_data < self.min_samples_split) or (classes == 1):</pre>
        self.tree = self.majority class(y training)
        return self.tree
    best feature, best threshold = self.select attribute(x training, y training)
    if best feature is None or best threshold is None:
        self.tree = self.majority class(y training)
        return self.tree
    # recursively splitting the training data and increment the depth of the tree by 1
   left_tree, right_tree = self.split_data(x_training, best_feature, best_threshold)
```

```
left_node = self.fit(x_training[left_tree], y_training[left_tree], depth + 1)
   right_node = self.fit(x_training[right_tree], y_training[right_tree], depth + 1)
   # store the tree in the dictionary
   self.tree = {'feature': best feature,
                 'threshold': best_threshold,
                 'left_node': left_node,
                 'right node': right node
    return self.tree
def traverse_tree(self, sample):
   Traversing the tree.
   @sample: sample data set
   node = self.tree
   while isinstance(node, dict):
       if sample[node['feature']] <= node['threshold']:</pre>
            node = node['left_node']
        else:
            node = node['right_node']
    return node
def predict(self, x testing):
   Predicting algorithm. This algorithm requires traversing the tree from root to leave.
   @param x testing: testing data set that you use to predict
   y pred = np.zeros(x testing.shape[0])
   for i, tree in enumerate(x_testing):
       y_pred[i] = self.traverse_tree(tree)
   return y_pred
```

```
In [4]: data = conduct_eda()
```

(8124, 23)

	count	unique	top	freq
class	8124	2	е	4208
cap-shape	8124	6	Х	3656
cap-surface	8124	4	У	3244
cap-color	8124	10	n	2284
bruises	8124	2	f	4748
odor	8124	9	n	3528
gill-attachment	8124	2	f	7914
gill-spacing	8124	2	С	6812
gill-size	8124	2	b	5612
gill-color	8124	12	b	1728
stalk-shape	8124	2	t	4608
stalk-root	8124	5	b	3776
stalk-surface-above-ring	8124	4	S	5176
stalk-surface-below-ring	8124	4	S	4936
stalk-color-above-ring	8124	9	W	4464
stalk-color-below-ring	8124	9	W	4384
veil-type	8124	1	р	8124
veil-color	8124	4	W	7924
ring-number	8124	3	0	7488
ring-type	8124	5	р	3968
spore-print-color	8124	9	W	2388
population	8124	6	V	4040
habitat	8124	7	d	3148
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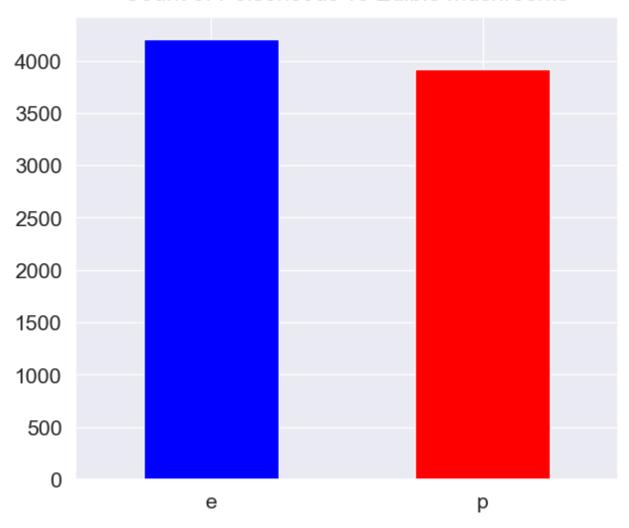
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 8124 entries, 0 to 8123

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Data	columns (total 23 columns):	
#	Column	Non-Null Count	Dtype
0	class	8124 non-null	object
1	cap-shape	8124 non-null	object
2	cap-surface	8124 non-null	object
3	cap-color	8124 non-null	object
4	bruises	8124 non-null	object
5	odor	8124 non-null	object
6	gill-attachment	8124 non-null	object
7	gill-spacing	8124 non-null	object
8	gill-size	8124 non-null	object
9	gill-color	8124 non-null	object
10	stalk-shape	8124 non-null	object
11	stalk-root	8124 non-null	object
12	stalk-surface-above-ring	8124 non-null	object
13	stalk-surface-below-ring	8124 non-null	object
14	stalk-color-above-ring	8124 non-null	object

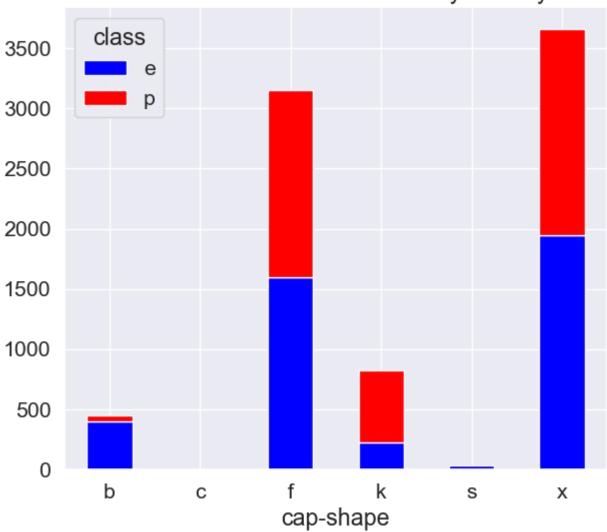
```
15 stalk-color-below-ring
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                               8124 non-null
                                              object
16 veil-type
                               8124 non-null
17 veil-color
                                               object
                               8124 non-null
18 ring-number
                              8124 non-null
                                              object
19 ring-type
                               8124 non-null
                                               object
20 spore-print-color
                                              object
                              8124 non-null
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[5 rows x 23 columns]

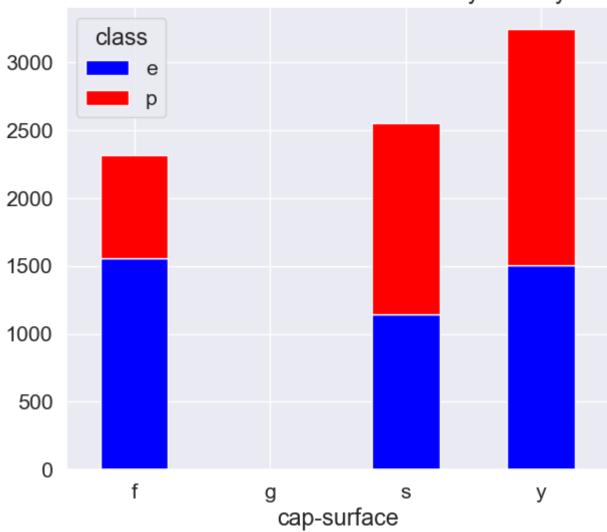
Count of Poisonsous vs Edible mushrooms



Count CAP-SHAPE break out by Edibility



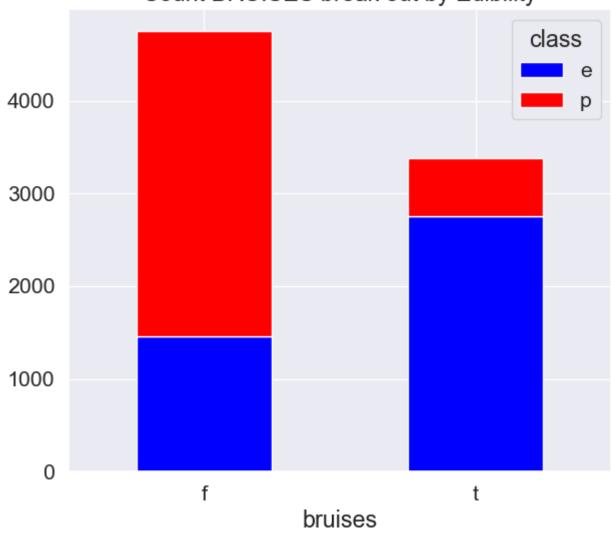
Count CAP-SURFACE break out by Edibility



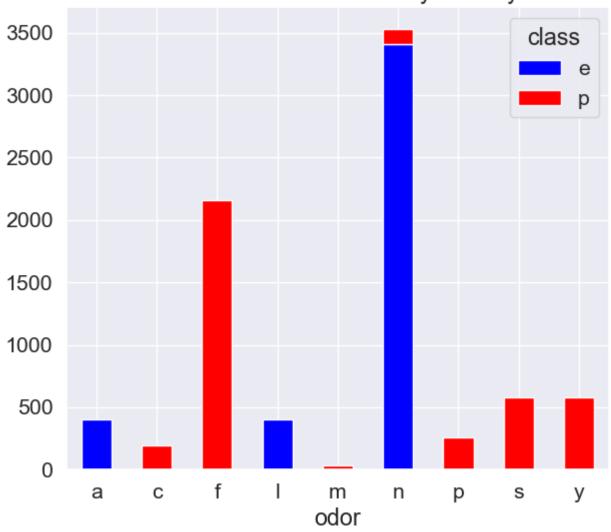
Count CAP-COLOR break out by Edibility



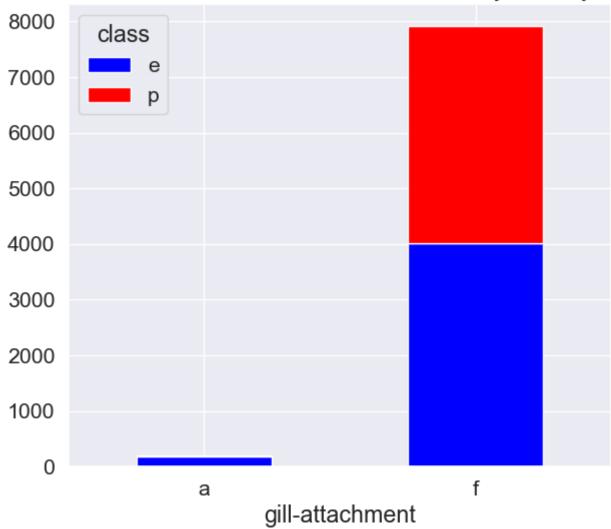
Count BRUISES break out by Edibility



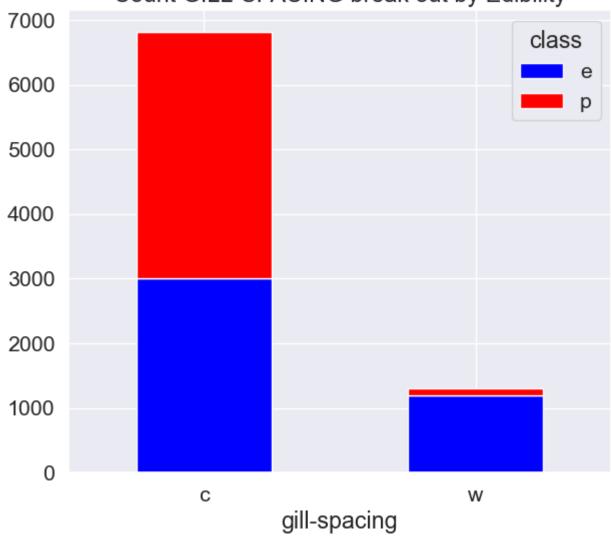


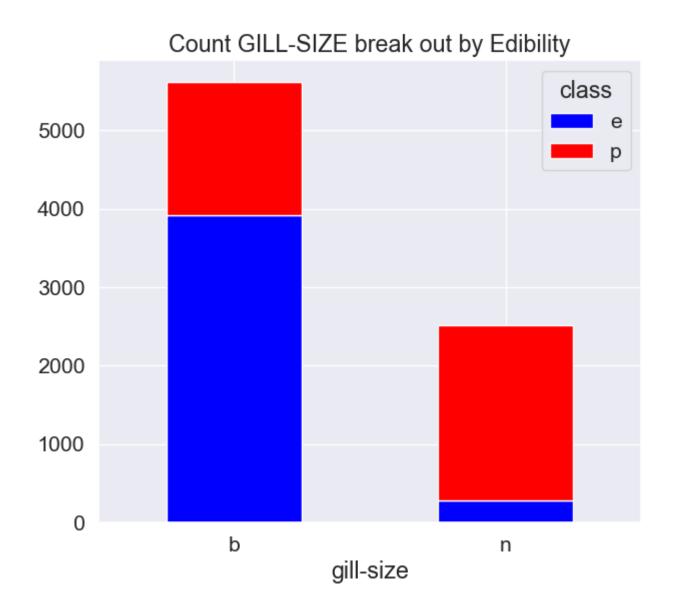


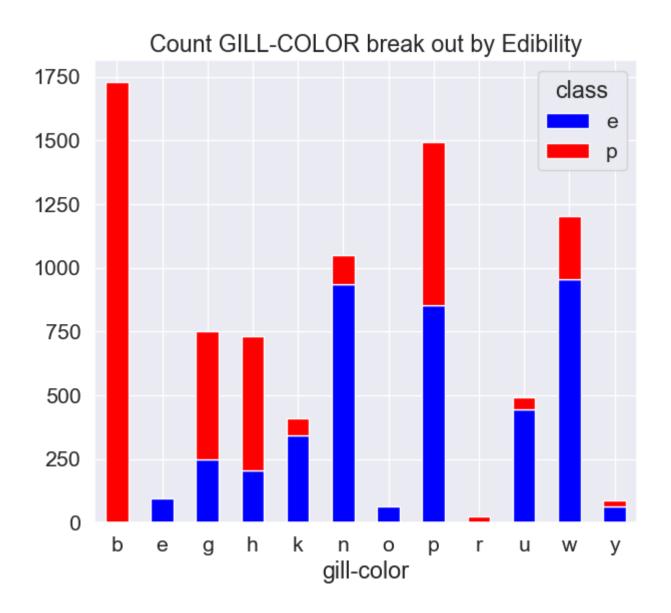
Count GILL-ATTACHMENT break out by Edibility



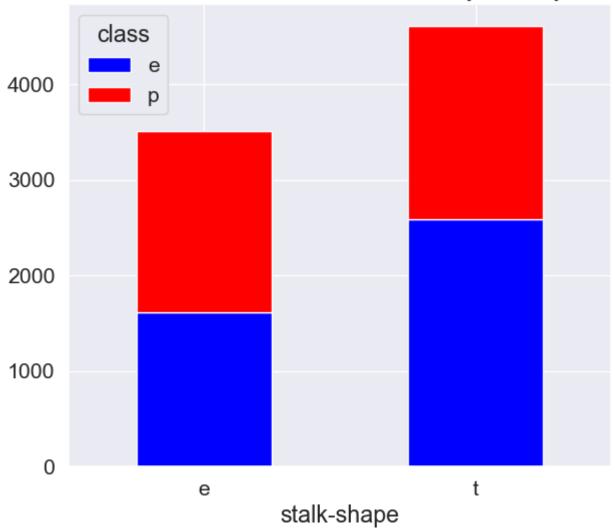
Count GILL-SPACING break out by Edibility



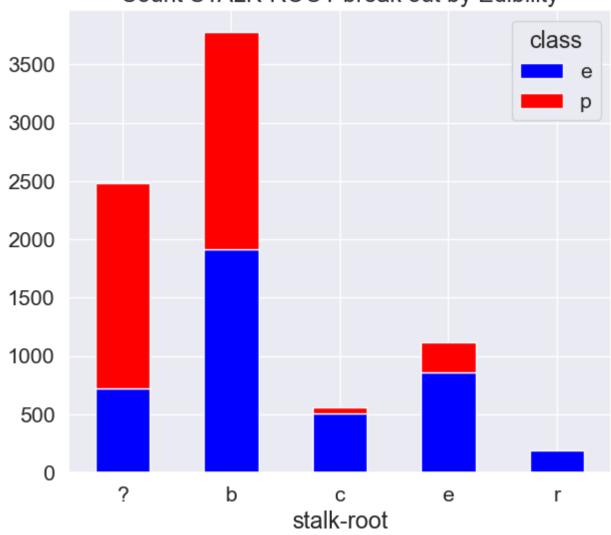




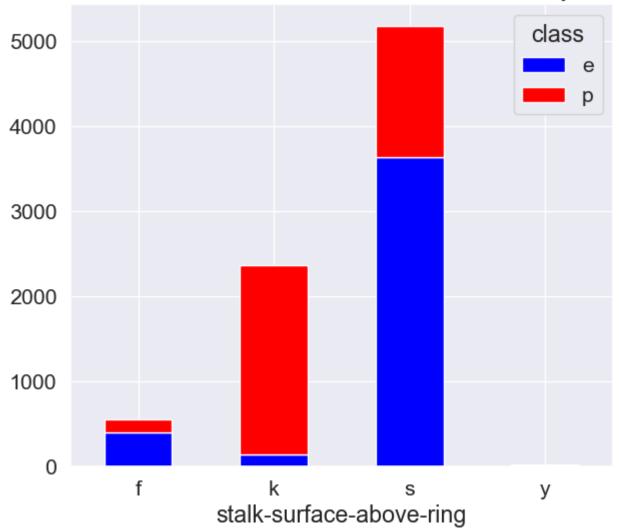
Count STALK-SHAPE break out by Edibility



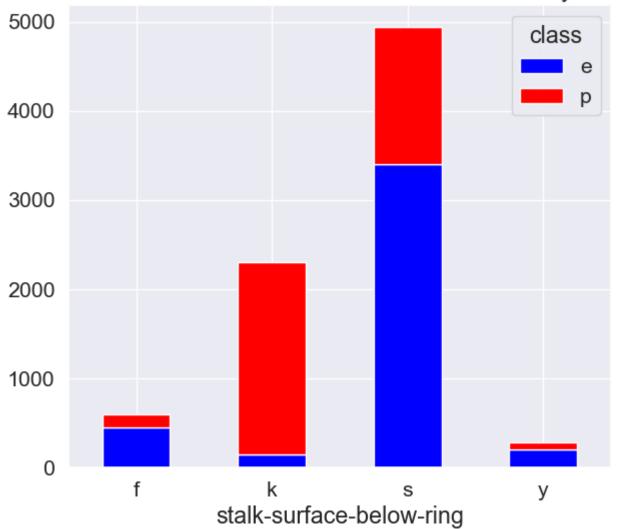




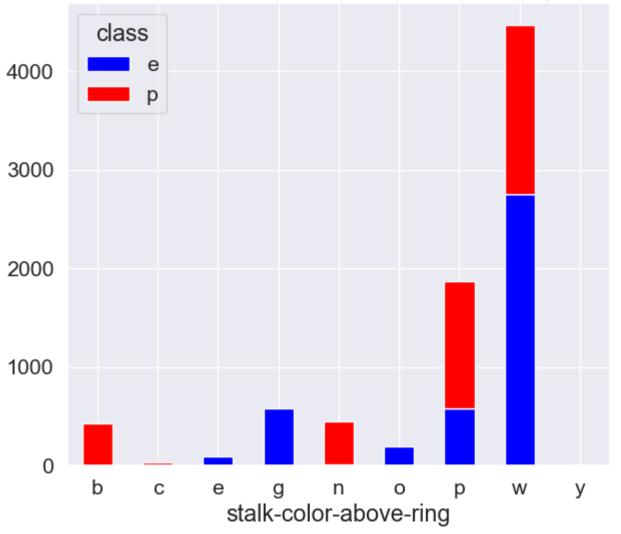
Count STALK-SURFACE-ABOVE-RING break out by Edibility



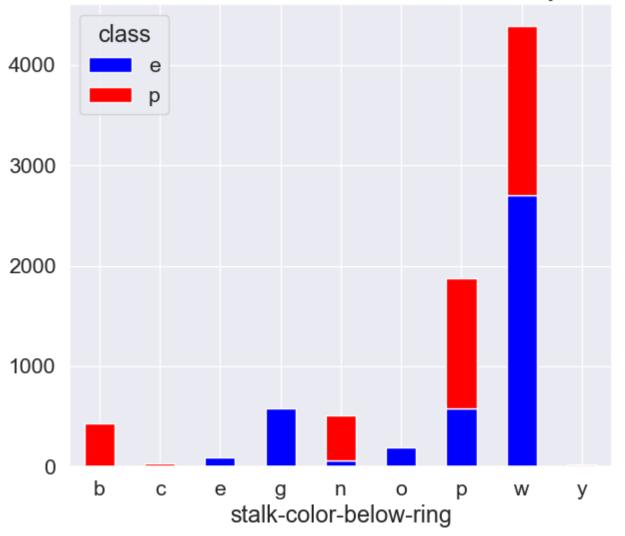
Count STALK-SURFACE-BELOW-RING break out by Edibility



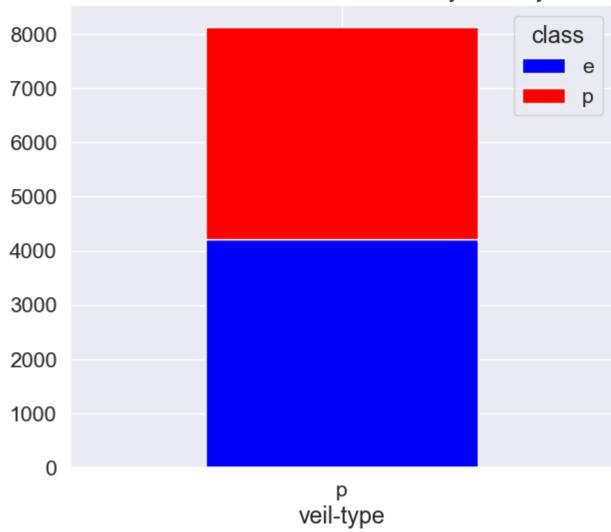
Count STALK-COLOR-ABOVE-RING break out by Edibility



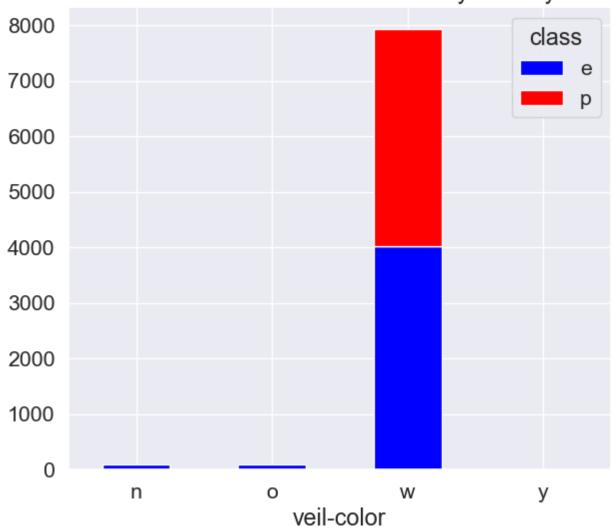
Count STALK-COLOR-BELOW-RING break out by Edibility



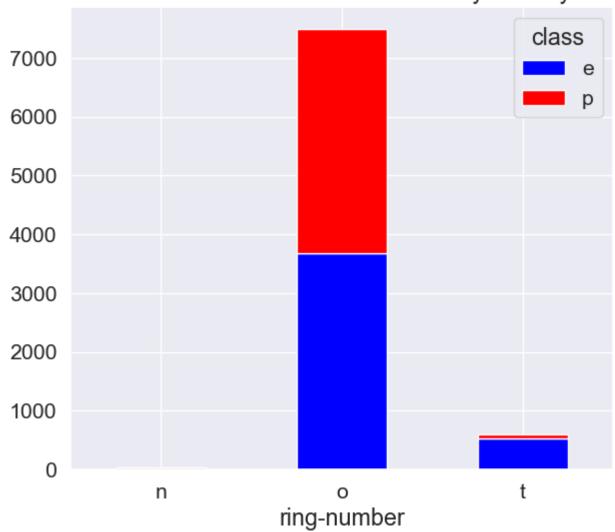
Count VEIL-TYPE break out by Edibility



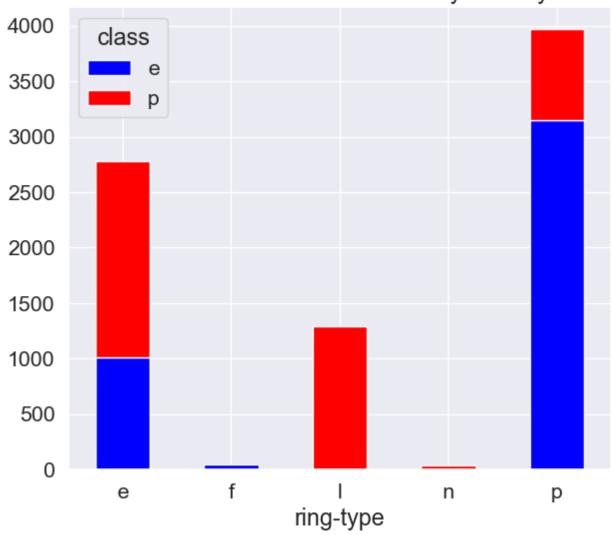
Count VEIL-COLOR break out by Edibility

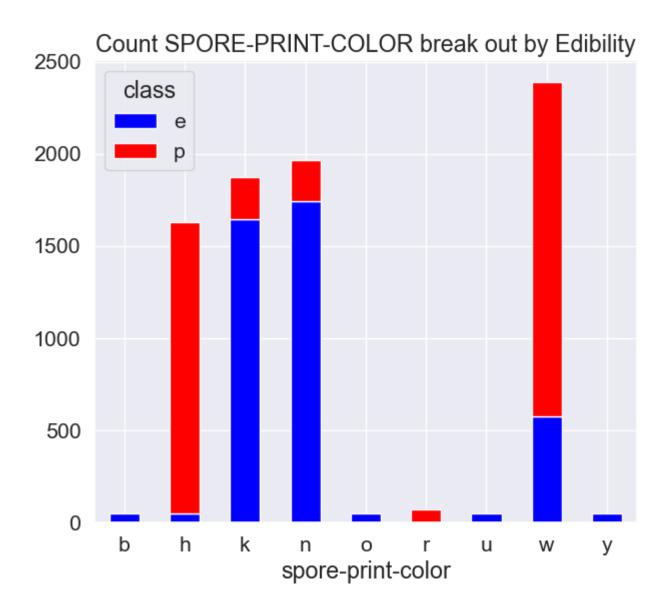


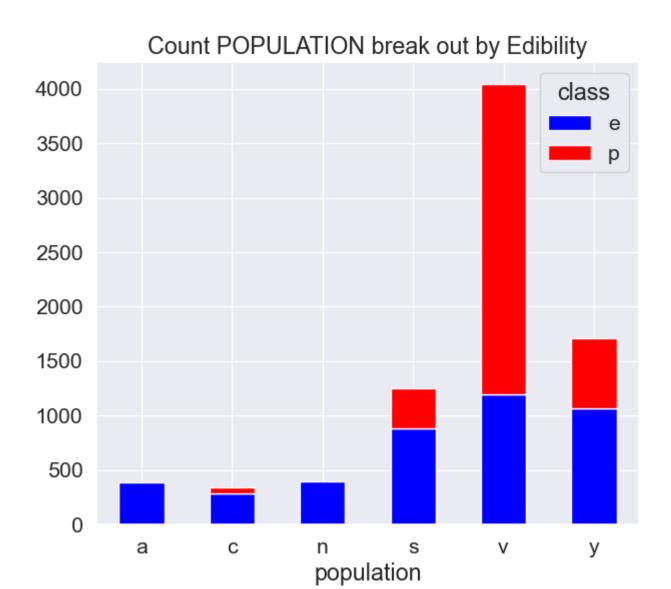
Count RING-NUMBER break out by Edibility

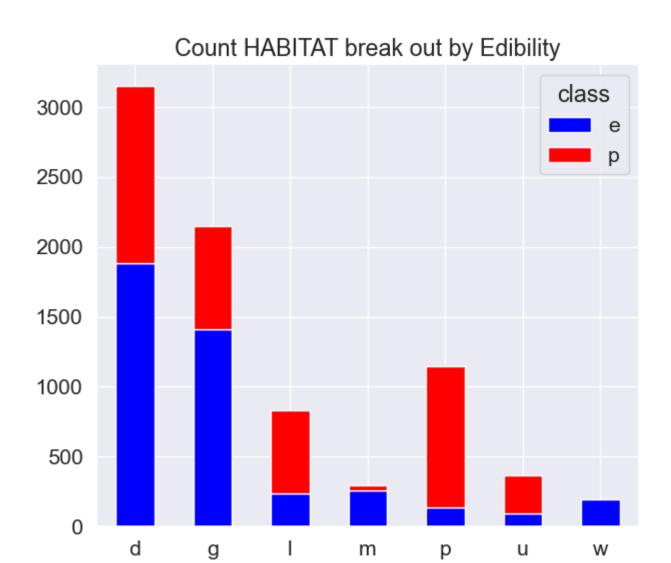


Count RING-TYPE break out by Edibility



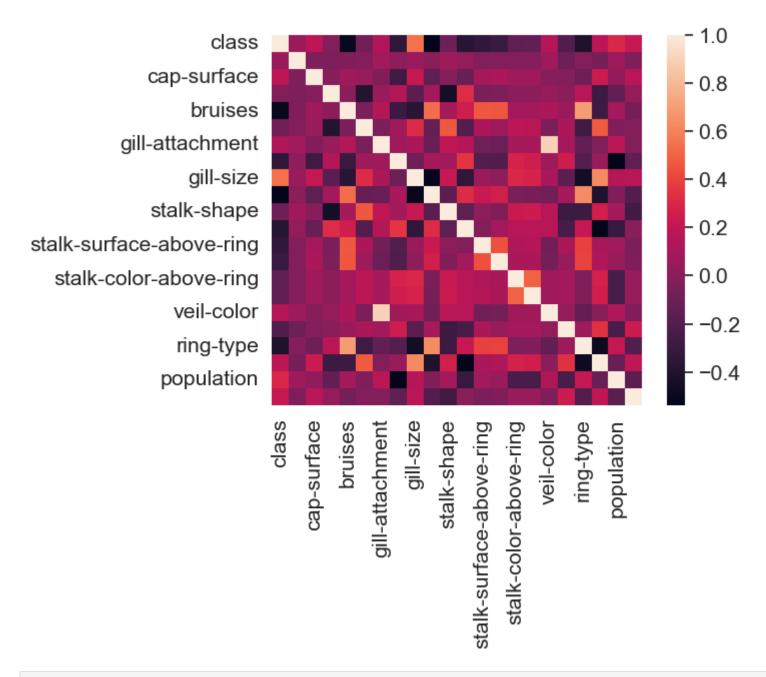






habitat

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```
In [5]: # split data into training and testing set
    x = data.iloc[:, :-1]
    y = data.iloc[:, 0]
    x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=0.07, random_state=0)
```

```
# Train the decision tree
tree = DecisionTree()
tree.fit(x_train.values, y_train.values)

# Test the decision tree
y_pred = tree.predict(x_test.values)

# Test the decision tree
y_pred = tree.predict(x_test.values)

In [6]:

# checking the acuracy of the data
accuracy = accuracy_score(y_test, y_pred)
print(f"Accuracy: {accuracy:.2f}")
```

Accuracy: 1.00

Q.2 Implementing Decision Tree using Sklearn (10 points)

- You will implement the decision tree algorithm using the scikit-learn library.
- You will write a program that trains and tests a decision tree model on the mushroom dataset.
- Your implementation should include the following steps:
- 1. Load the mushroom dataset

tree_clf.fit(x_train, y_train)

- 2. Split the dataset into training and testing sets
- 3. Train a decision tree model on the training set
- 4. Test the model on the testing set and calculate the accuracy

tree_clf = DecisionTreeClassifier(criterion='entropy')

```
In [7]: # already load the mushroom data set through calling conducting EDA on the top (data = conduct_eda())
# Hence we will just used that to split the data here

# split data into training and testing set
x = data.iloc[:, :-1]
y = data.iloc[:, 0]
x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=0.07, random_state=0)
In [8]: # Train a decision tree
```

```
Out[8]: DecisionTreeClassifier(criterion='entropy')

In [9]: # Test the model on the testing set and calculate the accuracy
    y_pred = tree_clf.predict(x_test)
    accuracy = accuracy_score(y_test, y_pred)
    print(f"Accuracy: {accuracy:.2f}")
```

Accuracy: 1.00

Q.3 Comparison between the results from Scratch and Sklearn (5 points)

• You will compare the results obtained from your implementation from scratch with the results obtained from the scikit-learn library.

The results obtain by my implementation and scikit-learn are the same. Both results are at 1.00 or 100% accuracy.

Q3. Neural Networks (25 Points)

- 1. **Forward Propagation**: Calculate the value of L (7 points). Use the weights and the input values to calculate the value of L using forward propagation.
- 2. **Backward Propagation**: Compute the derivatives (18 points). Use backward propagation to find the derivatives of the following variables with respect to each input and weight: (A1,A2,B1,B2,C1,C2,D). Use the chain rule to find the derivatives of each variable with respect to the inputs and weights.

