# Classification of breast cancer data [35 points]

You will implement a Support Vector Machine (SVM) model for the classification of breast cancer data and compare its performance with K-nearest Neighbors.

### Loading the data

```
In [1]: import pandas as pd
    import numpy as np
    from matplotlib import pyplot as plt
    import seaborn as sns

# added for final run of notebook to suppress deprecation warnings
    import warnings
    warnings.filterwarnings("ignore")

In [2]: ## loading the raw data using pandas
    bc = pd.read_csv('data.csv')
    bc = bc.drop(bc.columns[-1],axis=1)
    bc.head()
```

#### Out[2]:

	id	diagnosis	radius_mean	texture_mean	perimeter_mean	area_mean	smoothness_m
0	842302	М	17.99	10.38	122.80	1001.0	0.11
1	842517	М	20.57	17.77	132.90	1326.0	30.0
2	84300903	М	19.69	21.25	130.00	1203.0	0.10
3	84348301	М	11.42	20.38	77.58	386.1	0.14
4	84358402	М	20.29	14.34	135.10	1297.0	0.10
_							

5 rows × 32 columns

Pre-process and understand data (5 pts)

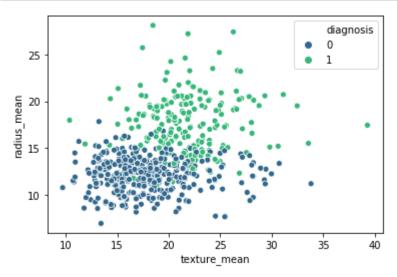
The data label is the diagnosis (M = malignant, B = benign).

Question (2pts): Convert diagnosis column to binary 0-1 label where 1 incidates malignancy.

**Question (3pts):** Plot the first 2 features against diagnosis. Is the data linearly separable and what happens if we fit a hard-margin SVM to the data?

**Answer:** No, the data is not perfectly linearly separable, so if we attempt to fit a hard-margin SVM to the data, the SVM will guit and no solution will be found.

```
In [4]: # plotting the the first 2 features against diagnosis
sns.scatterplot(x='texture_mean', y='radius_mean', data=bc, hue='diagnosis', p
alette='viridis');
```



#### Splitting the data: training and test

```
In [5]: from sklearn.model_selection import train_test_split

X = bc.drop(["id", "diagnosis"], axis=1)
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.25, rand om_state=0)
```

# 1. Support Vector Machine (SVM) (20 pts)

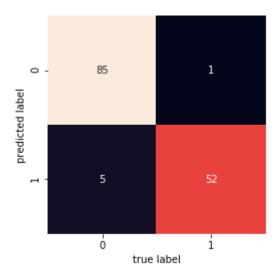
### Question (15pts): Train and test a SVM model with the following requirements:

- Use GridSearchCV to find the best combination of model hyper-parameters: kernel: {linear, rbf}; for both kernel, C:{1, 10, 100, 1000}; for rbf kernel 'gamma' = {1e-3, 1e-4, 1e-5}.
- Try both accuracy and recall metric as scoring function in GridSearchCV
- Report the best hyper-parameter setting for each metric
- · Report and plot the confusion matrix of the best estimator for each metric

```
In [6]: from sklearn.svm import SVC
        from sklearn.model selection import GridSearchCV
        from sklearn.metrics import confusion matrix
        svc = SVC()
        param_grid = [{'kernel': ['rbf'], 'gamma': [1e-3, 1e-4, 1e-5], 'C': [1, 10, 10]
        0, 1000]},
                       {'kernel': ['linear'], 'C': [1, 10, 100, 1000]}]
        scores = ['accuracy', 'recall']
        for sc in scores:
            print("# Tuning hyper-parameters for %s" % sc)
            svc = GridSearchCV(SVC(), param_grid, cv=5, scoring=sc)
            svc.fit(X train, y train)
            print()
            print('Best parameters for %s:' %sc)
            print(svc.best_params_)
            print()
            y pred svc = svc.predict(X test)
            # computing and plotting confusion matrix
            mat = confusion_matrix(y_test, y_pred_svc)
            sns.heatmap(mat.T, square=True, annot=True, fmt='d', cbar=False)
            plt.xlabel('true label')
            plt.ylabel('predicted label')
            plt.show()
```

# Tuning hyper-parameters for accuracy

```
Best parameters for accuracy:
{'C': 1, 'kernel': 'linear'}
```

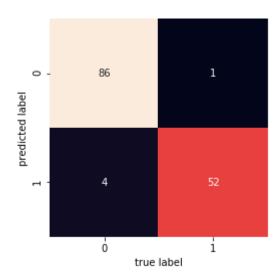


# Tuning hyper-parameters for recall

C:\Users\amand\Anaconda3\lib\site-packages\sklearn\model\_selection\\_search.p y:841: DeprecationWarning: The default of the `iid` parameter will change fro m True to False in version 0.22 and will be removed in 0.24. This will change numeric results when test-set sizes are unequal.

DeprecationWarning)

Best parameters for recall:
{'C': 10, 'kernel': 'linear'}



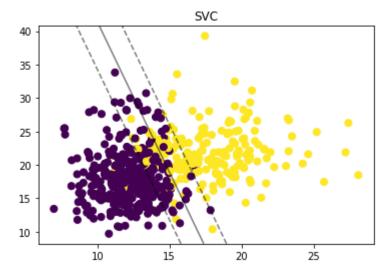
**Question (5pts):** Fit an SVM model to a slice of data that contains only the first 2 features and visualize decision bouldary based on these two features.

```
In [7]: # code credited to Recitation 9 notes
        def plot_svc_decision_function(model, ax=None, plot_support=True):
             """Plot the decision function for a 2D SVC"""
            if ax is None:
                ax = plt.gca()
            xlim = ax.get_xlim()
            ylim = ax.get ylim()
            # create grid to evaluate model
            x = np.linspace(xlim[0], xlim[1], 30)
            y = np.linspace(ylim[0], ylim[1], 30)
            Y, X = np.meshgrid(y, x)
            xy = np.vstack([X.ravel(), Y.ravel()]).T
            P = model.decision function(xy).reshape(X.shape)
            # plot decision boundary and margins
            ax.contour(X, Y, P, colors='k',
                       levels=[-1, 0, 1], alpha=0.5,
                       linestyles=['--', '--'])
            # plot support vectors
            if plot_support:
                ax.scatter(model.support vectors [:, 0],
                           model.support_vectors_[:, 1],
                            s=300, linewidth=3, facecolors='none');
            ax.set xlim(xlim)
            ax.set ylim(ylim)
```

```
In [8]: from sklearn import metrics

# for display purposes, we fit the model on the first two features
svc_best = SVC(kernel='linear', C=1)
svc_best.fit(X_train.loc[:,'radius_mean':'texture_mean'], y_train)

# Plotting the decision boundary for all data (both train and test)
plt.scatter(x='radius_mean', y='texture_mean', data=X, c=y, s=50, cmap='viridis')
plot_svc_decision_function(svc_best);
plt.title('SVC')
plt.show()
```



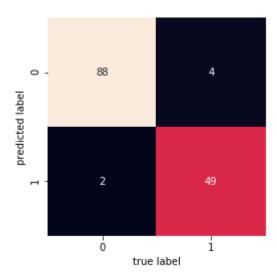
# 2. K-nearest Neighbors classifier (10pts)

**Question:** Repeat all the tasks for SVM to train a kNN classifier where the hyper-parameter is only k (number of neighbors) ranging from 1 to 50 in kNN. Compare the results with those of SVM, which classifier gives better results?

**Answer:** Of the classifiers trained, the linear-kernel SVM with C=10 did the best. Both the SVM and KNN models misclassified approximately the same total number of observations (5 for SVM and 6 for KNN), however the SVM model only had one false negative, whereas the KNN model had 4. In the case of breast cancer prediction, the false negative is the more potentially harmful case, so we would want to choose the model that minimizes false negatives over false positives.

```
In [13]: from sklearn.neighbors import KNeighborsClassifier
         knn = KNeighborsClassifier()
         k range = list(np.arange(1,51))
         param_grid = dict(n_neighbors=k_range)
         scores = ['accuracy', 'recall']
         for sc in scores:
             print("# Tuning hyper-parameters for %s" % sc)
             grid_knn = GridSearchCV(KNeighborsClassifier(), param_grid, cv=5, scoring=
         sc)
             grid_knn.fit(X_train,y_train)
             knn = grid_knn
             y pred knn = knn.fit(X train, y train)
             print('Best parameters for %s:' %sc)
             print(grid_knn.best_params_)
             print()
             y_pred_knn = knn.predict(X_test)
             mat = confusion_matrix(y_test, y_pred_knn)
             sns.heatmap(mat.T, square=True, annot=True, fmt='d', cbar=False)
             plt.xlabel('true label')
             plt.ylabel('predicted label')
             plt.show()
```

# Tuning hyper-parameters for accuracy
Best parameters for accuracy:
{'n\_neighbors': 9}



# Tuning hyper-parameters for recall

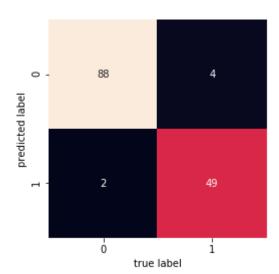
C:\Users\amand\Anaconda3\lib\site-packages\sklearn\model\_selection\\_search.p y:841: DeprecationWarning: The default of the `iid` parameter will change fro m True to False in version 0.22 and will be removed in 0.24. This will change numeric results when test-set sizes are unequal.

DeprecationWarning)

Best parameters for recall:
{'n\_neighbors': 9}

C:\Users\amand\Anaconda3\lib\site-packages\sklearn\model\_selection\\_search.p y:841: DeprecationWarning: The default of the `iid` parameter will change fro m True to False in version 0.22 and will be removed in 0.24. This will change numeric results when test-set sizes are unequal.

DeprecationWarning)



```
In [14]: # Plotting the decision boundary for all data (both train and test)
         from matplotlib.colors import ListedColormap
         # for display purposes, we fit the model on the first two features
         knn best = KNeighborsClassifier(n neighbors=9)
         knn_best.fit(X_train.loc[:,'radius_mean':'texture_mean'], y_train)
         # copying X test, y test to X set and y set. Remeber X has two columns
         X_set, y_set = X.loc[:,'radius_mean':'texture_mean'], np.array(y)
         # creating a meshgrid where each point is classifed using learned kNN
         X1, X2 = np.meshgrid(np.arange(start = X_set.iloc[:, 0].min() - 1, stop = X_se
         t.iloc[:, 0].max() + 1, step = 0.01),
                              np.arange(start = X set.iloc[:, 1].min() - 1, stop = X se
         t.iloc[:, 1].max() + 1, step = 0.01))
         plt.contourf(X1, X2, knn_best.predict(np.array([X1.ravel(), X2.ravel()]).T).re
         shape(X1.shape),
                      alpha = 0.25, cmap = ListedColormap(('purple', 'yellow')))
         plt.xlim(X1.min(), X1.max())
         plt.ylim(X2.min(), X2.max())
         plt.scatter(x='radius_mean', y='texture_mean', data=X, c=y, s=50, cmap='viridi
         s')
         plt.title('KNN')
         plt.show()
```

