Practical 4

```
import numpy as np
import cvxopt
from cvxopt import matrix as cvxopt_matrix
from matplotlib import pyplot as plt
from sklearn.svm import SVC
from sklearn.datasets import load_wine
from sklearn.tree import DecisionTreeClassifier
import pandas as pd
```

Task 0: The Data

We will work with the data from practical 3. Load the data and split it into a training and test set. You can re-use the data splitting function from Practical 2.

```
# def split_data(X, y, frac=0.3, seed=None):
      len_X = len(X)
      indices = np.random.permutation(len_X)
      indices_test, indices_train = indices[:int(len(indices)*frac)], indices[int(len(indice
s)*frac):]
     y_train, y_test = y[indices_train], y[indices_test]
     X_train, X_test = X[indices_train], X[indices_test]
      return y_train, y_test, X_train, X_test
# from Practical 2
def split_data(X, y, frac=0.3, max_samples=None, seed=4711):
    if seed is not None:
        np.random.seed(seed)
    indices = np.random.permutation(len(X))
    indices = indices[:max_samples]
    indices_test, indices_train = indices[:int(len(indices)*frac)], indices[int(len(indices)*
frac):]
    assert (len(set(indices))) == len(indices)
    assert len(set(indices_test).intersection(indices_train)) == 0
   X_train, X_test = X[indices_train], X[indices_test]
   y_train, y_test = y[indices_train], y[indices_test]
    return X_train, X_test, y_train, y_test
# X_train, X_test, y_train, y_test = split_data(X_2d, t_2d)
```

```
# Load data
X_2d, t_2d = np.load('data/nonlin_2d_data.npy')[:,:2], np.load('data/nonlin_2d_data.npy')[:,
2]
```

```
# split data
X_train, X_test, t_train, t_test = split_data(X_2d, t_2d, seed=1)
```

In SVMs, each data sample x_n has a corresponding lagrange multiplier α_n which indicates if x_n is a support vector. In the latter case $\alpha_n > 0$ holds. The goal of learning the SVM is to figure out which samples are support vectors by learning α . The dual SVM optimizes the following quadratic program.

$$\min rac{1}{2} \sum_{n=1}^N \sum_{m=1}^N lpha_n lpha_m t_n t_m k(\mathbf{x}_n, \mathbf{x}_m) - \sum_{n=1}^N lpha_n$$

subject to

$$0 < \alpha_n < C$$

$$\sum_{n=1}^N lpha_n t_n = 0$$

The quadratic program solver expects the following form:

$$\min rac{1}{2} lpha^T P lpha + \mathbf{q}^T lpha$$

subject to

$$A\alpha = b$$

$$G\alpha \le h$$

Here, A and G are matrices with one row per individual constraint. Similarly, b and h are vectors with one element per individual constraint.

Having trained the SVM, a prediction for an input \mathbf{x} is made by:

$$y = sign([\sum_{n}^{N} lpha_n t_n k(\mathbf{x}, \mathbf{x}_n)] + b)$$

Task 1.1

Use the code provided below as a basis to express the constrained optimization problem in terms of P,q,A,b,G and h and implement a function <code>fit_svm</code> which passes these variables to the provided QP solver. Fit a SVM on the training data and extract its parameters.

Hints: - The box constraint $0 \le \alpha_n \le C$ defines two constraints of the form $G\alpha_n \le h$ for each α_n . - The inequality $x \ge 0$ is equivalent to $-x \le 0$.

```
def rbf_kernel(x_n, x_m, gamma=1/5):
    tmp = np.square(x_n - x_m)
    tmp = -(1/gamma) * np.sum(tmp)
    return np.exp(tmp)
```

```
def accuracy(y_true, y_pred):
    n_c = np.sum((y_true == y_pred)) #number of correctly classified
    acc = n_c/len(y_true)
    return acc
```

```
def fit_svm(X, t, kernel, C=1.0):
    '''Fit SVM using data (X,t), specified kernel and parameter C.'''
   t = np.array([-1 if l == 0 else 1 for l in t])
    tmat = np.diag(t)
    Ptmp = np.array([kernel(xn,xm) for xn in X for xm in X]).reshape((len(X), len(X)))
    P = np.matmul(np.matmul(tmat, Ptmp), tmat)
   q = -1 * np.ones(len(X))
    A = np.array([float(k) for k in t]).reshape((1,len(X)))
   b = np.array([0.]).reshape((1,1))
   onev = np.array([-1, 1])
    G = np.kron(np.identity(len(X)), onev.reshape((2,1)))
    tmp = np.array([0,C])
   h = np.tile(tmp, len(X))
    assert P.shape == (len(X), len(X))
    assert len(q) == len(X)
    assert A.shape == (1, len(X)) and A.dtype == 'float'
    assert b.shape == (1, 1)
    assert len(G) == 2 * len(X)
    assert len(h) == 2 * len(X)
    return solve_quadratic_program(P, q, A, b, G, h)
def solve_quadratic_program(P, q, A, b, G, h):
    '''Uses cvxopt to solve the quadratic program.'''
   P, q, A, b, G, h = [cvxopt.matrix(var) for var in [P, q, A, b, G, h]]
    minimization = cvxopt.solvers.qp(P, q, G, h, A, b)
    lagr_mult = np.ravel(minimization['x'])
    return lagr_mult
def extract_parameters(X, t, kernel, lagr_mult, C, threshold=1e-5):
    '''Computes the intercept from the support vector constraints.
    Inputs
       X:
                  predictors
       t:
                  targets
                  a kernel to be used
        lagr mult: the Lagrange multipliers obtained by solving the dual QP
       threshold: threshold for choosing support vectors
    Returns
       lagr_mult: lagrange multipliers for the support vectors
                  set of support vectors
       sv_labels: targets t_n for the support vectors
       intercept: computed intercept
    t = np.array([-1 if l == 0 else 1 for l in t])
    # ----- INSERT CODE -----
    sv_inds = (lagr_mult > threshold) & (lagr_mult < (C-threshold))</pre>
    svs, sv_labels, lagr_mult = X[sv_inds], t[sv_inds], lagr_mult[sv_inds]
```

```
# Fit SVM on training data
sols = fit_svm(X_train, t_train, rbf_kernel)

# Extract parameters
lagr, svs, sv_labels, intercept = extract_parameters(X_train, t_train, rbf_kernel, sols, 1.0)
```

```
pcost
               dcost
                                  pres
                                        dres
                           gap
0: -1.2225e+01 -2.4493e+02 8e+02 1e+00 4e-16
1: -9.2861e+00 -1.1119e+02 1e+02 1e-15 4e-16
2: -1.4797e+01 -2.4685e+01 1e+01 3e-15 4e-16
3: -1.5743e+01 -1.7893e+01 2e+00 1e-15 2e-16
4: -1.5986e+01 -1.6566e+01 6e-01 1e-15 1e-16
 5: -1.6091e+01 -1.6235e+01 1e-01 2e-15 2e-16
6: -1.6117e+01 -1.6144e+01 3e-02 2e-15 2e-16
7: -1.6123e+01 -1.6125e+01 3e-03 3e-15 2e-16
8: -1.6123e+01 -1.6123e+01 7e-05 1e-15 2e-16
9: -1.6123e+01 -1.6123e+01 2e-06 4e-16 1e-16
Optimal solution found.
```

Task 1.2

Having learnt an SVM, we can use the calculated parameters to make predictions on novel samples. - Implement a function svm_predict(X, kernel, lagr_mult, svs, sv_labels, intercept) . - Use this function with your Gaussian RBF kernel (Practical 3) and compute the test accuracy on the 2d dataset.

```
# Calculate test accuracy
y_train = svm_predict(X_train, rbf_kernel, lagr, svs, sv_labels, intercept)
y_test = svm_predict(X_test, rbf_kernel, lagr, svs, sv_labels, intercept)
print("The accuracy on the training set is", accuracy(t_train, y_train), "and on the test se
t", accuracy(t_test, y_test))
```

The accuracy on the training set is 0.9828571428571429 and on the test set 1.0

Task 1.3

- Instead of using the Gaussian RBF kernel, use the linear kernel (dot product) defined in Practical 3.
- Compare results on with both kernels with sklearn implementation (SVC)
- Visualize the predictions on the test set, the learned support vectors and the decision boundary for both kernels.

```
def linear_kernel(x_n, x_m):
    return np.matmul(x_n.T, x_m)
```

```
# Fit SVM with linear kernel and calculate the test accuracy
# Fit SVM on training data
sols_lk = fit_svm(X_train, t_train, linear_kernel)

# Extract parameters
lagr_lk, svs_lk, sv_labels_lk, intercept_lk = extract_parameters(X_train, t_train, linear_kernel, sols_lk, 1.0)
```

```
pcost dcost gap pres dres
0: -1.6799e+02 -3.7889e+02 9e+02 2e+00 2e-14
1: -1.2879e+02 -2.5279e+02 1e+02 5e-15 1e-14
2: -1.3986e+02 -1.4886e+02 9e+00 3e-15 9e-15
3: -1.4000e+02 -1.4009e+02 9e-02 8e-16 8e-15
4: -1.4000e+02 -1.4000e+02 9e-04 4e-15 9e-15
5: -1.4000e+02 -1.4000e+02 9e-06 2e-15 1e-14
Optimal solution found.
```

y_train_lk = svm_predict(X_train, linear_kernel, lagr_lk, svs_lk, sv_labels_lk, intercept_lk)
y_test_lk = svm_predict(X_test, linear_kernel, lagr_lk, svs_lk, sv_labels_lk, intercept_lk)
print("The accuracy on the training set is", accuracy(t_train, y_train_lk), "and on the test
set", accuracy(t_test, y_test_lk))

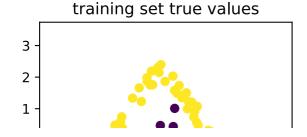
```
# Fit SVM using sklearn and calculate the test accuracy
from sklearn.svm import SVC
from sklearn.pipeline import make_pipeline
from sklearn.preprocessing import StandardScaler
clf = make_pipeline(StandardScaler(), SVC(gamma='auto'))
clf.fit(X_train, t_train)
```

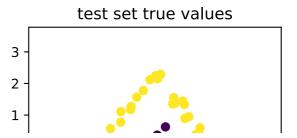
```
y_train_sk = clf.predict(X_train)
y_test_sk = clf.predict(X_test)
print("The accuracy on the training set is", accuracy(t_train, y_train_sk), "and on the test
set", accuracy(t_test, y_test_sk))
```

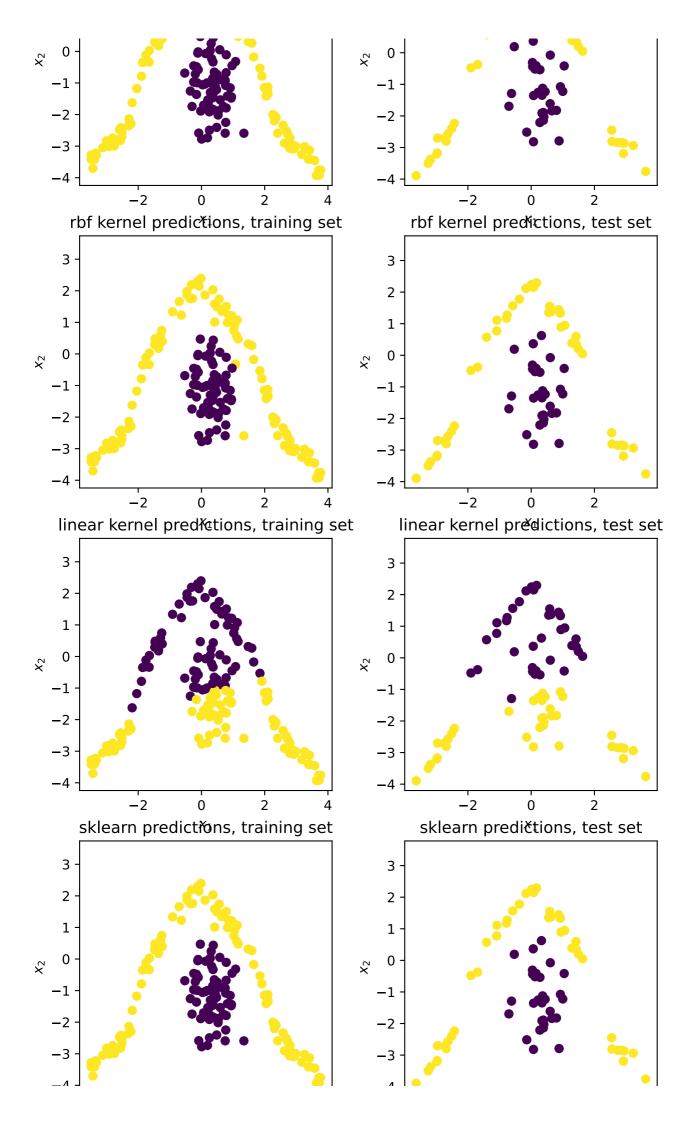
The accuracy on the training set is 0.9942857142857143 and on the test set 1.0

```
def plt2d(data, labels, title):
  plt.scatter(data[:,0], data[:,1], c = labels)
  plt.xlabel('$x_1$')
  plt.ylabel('$x_2$')
  plt.axis('square')
  plt.title(title)
```

```
# Visualize
# Visualize results
plt.figure(figsize=(8,16))
plt.subplot(4,2,1)
plt2d(X_train, t_train, title = 'training set true values')
plt.subplot(4,2,2)
plt2d(X_test, t_test, title = 'test set true values')
plt.subplot(4,2,3)
plt2d(X_train, y_train, title = 'rbf kernel predictions, training set')
plt.subplot(4,2,4)
plt2d(X_test, y_test, title = 'rbf kernel predictions, test set')
plt.subplot(4,2,5)
plt2d(X_train, y_train_lk, title = 'linear kernel predictions, training set')
plt.subplot(4,2,6)
plt2d(X_test, y_test_lk, title = 'linear kernel predictions, test set')
plt.subplot(4,2,7)
plt2d(X_train, y_train_sk, title = 'sklearn predictions, training set')
plt.subplot(4,2,8)
plt2d(X_test, y_test_sk, title = 'sklearn predictions, test set')
plt.show()
```

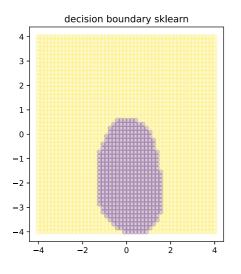


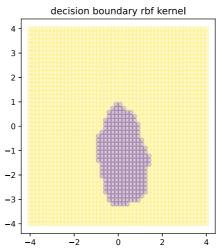


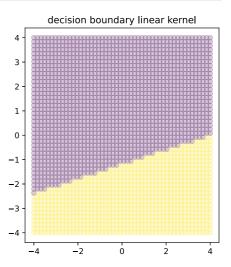


svg

```
def plot_decision_boundary(predict_fn):
    ''' Plot decision boundary.
    predict_fn: function handle to a predict function that takes
                as input a data point x and outputs the predicted
                label t
    grid = np.meshgrid(np.linspace(-4, 4, 50), np.linspace(-4, 4, 50))
    grid_x, grid_y = grid[0].flatten(), grid[1].flatten()
    grid_t = np.zeros(len(grid_x))
    for i, p in enumerate(zip(grid_x, grid_y)):
      tmp = np.array(p)
      tmp = np.reshape(tmp, (1, 2))
      grid_t[i] = predict_fn(tmp)
    plt.scatter(grid_x, grid_y, c=grid_t, alpha=0.2)
plt.figure(figsize = (15, 5))
plt.subplot(1, 3, 1)
predict_fn = lambda x: clf.predict(x)
plot_decision_boundary(predict_fn)
plt.title('decision boundary sklearn')
plt.subplot(1, 3, 2)
predict_fn = lambda x: svm_predict(x, rbf_kernel, lagr, svs, sv_labels, intercept)
plot decision boundary(predict fn)
plt.title('decision boundary rbf kernel')
plt.subplot(1, 3, 3)
predict_fn = lambda x: svm_predict(x, linear_kernel, lagr_lk, svs_lk, sv_labels_lk, intercept
_lk)
plot decision boundary(predict fn)
plt.title('decision boundary linear kernel')
plt.show()
```







Task 2: Decision Trees

Next, we will implement a simple decision tree classifier using the Wine dataset, one of the standard sklearn datasets.

We will use the Gini impurity as a criterion for splitting. It is defined for a set of labels as

$$G=\sum_{i=0}^C p(i)*(1-p(i))$$

Given labels l and split l_a and l_b , the weighted removed impurity can be computed by $G(l)-rac{|l_a|}{|l|}G(l_a)-rac{|l_b|}{|l|}G(l_b)$.

Here is a simple explanation of the Gini impurity that you may find useful: https://victorzhou.com/blog/giniimpurity/ (https://victorzhou.com/blog/gini-impurity/)

Task 2.1

- 1. Plot the distribution of the first feature of for each class of the wine dataset.
- 2. Implement a function gini impurity(t) that computes the Gini impurity for an array of labels t.
- 3. Calculate the removed Gini impurity for a split after 50 samples, i.e. between t[:50] and t[50:].

```
# Load Wine dataset and split into train+test set
from sklearn.datasets import load_wine
X, t = load_wine(return_X_y=True)
X_train, X_test, t_train, t_test = split_data(X = X, y = t)
```

```
# Plot distribution
import seaborn as sns
for i in np.unique(t):
    sns.distplot(X[(t==i)][:,1],kde=1,label='{}'.format(i))
plt.legend()
plt.show()
```

C:\Users\rene-\AppData\Local\Programs\Python\Python38\lib\site-packages\seaborn\distribution s.py:2557: FutureWarning: `distplot` is a deprecated function and will be removed in a future version. Please adapt your code to use either `displot` (a figure-level function with similar flexibility) or `histplot` (an axes-level function for histograms).

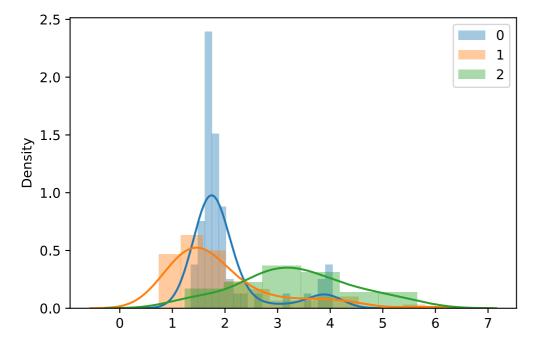
warnings.warn(msg, FutureWarning)

 $\label{libsite-packages} C:\Users\rene-\AppData\Local\Programs\Python\Python38\lib\site-packages\seaborn\distribution$ s.py:2557: FutureWarning: `distplot` is a deprecated function and will be removed in a future version. Please adapt your code to use either `displot` (a figure-level function with similar flexibility) or `histplot` (an axes-level function for histograms).

warnings.warn(msg, FutureWarning)

C:\Users\rene-\AppData\Local\Programs\Python\Python38\lib\site-packages\seaborn\distribution s.py:2557: FutureWarning: `distplot` is a deprecated function and will be removed in a future version. Please adapt your code to use either `displot` (a figure-level function with similar flexibility) or `histplot` (an axes-level function for histograms).

warnings.warn(msg, FutureWarning)



svg

```
# Compute Gini impurity
def gini_impurity(data):
    temp = 0
    for i in np.unique(data):
        temp += np.sum(data==i)/len(data) * (1- np.sum(data==i)/len(data))
    return temp
```

```
print('Gini of full dataset: ', gini_impurity(t))
print('Gini of split dataset: ', gini_impurity(t[50:]))
print('Gini of split dataset: ', gini_impurity(t[:50]))
```

```
Gini of full dataset: 0.6583133442747128
Gini of split dataset: 0.5467529296875
Gini of split dataset: 0.0
```

```
gini_before = gini_impurity(t)
gini_after = len(t[:50])/len(t) * gini_impurity(t[:50]) + len(t[50:])/len(t) * gini_impurity
(t[50:])
reduced_gini = gini_before - gini_after
print("The removed gini based on a split after 50 samples is:",reduced_gini)
print("\n")
print("gini before:", gini_before)
print("gini after:", gini_after)
```

```
The removed gini based on a split after 50 samples is: 0.2651426982072971

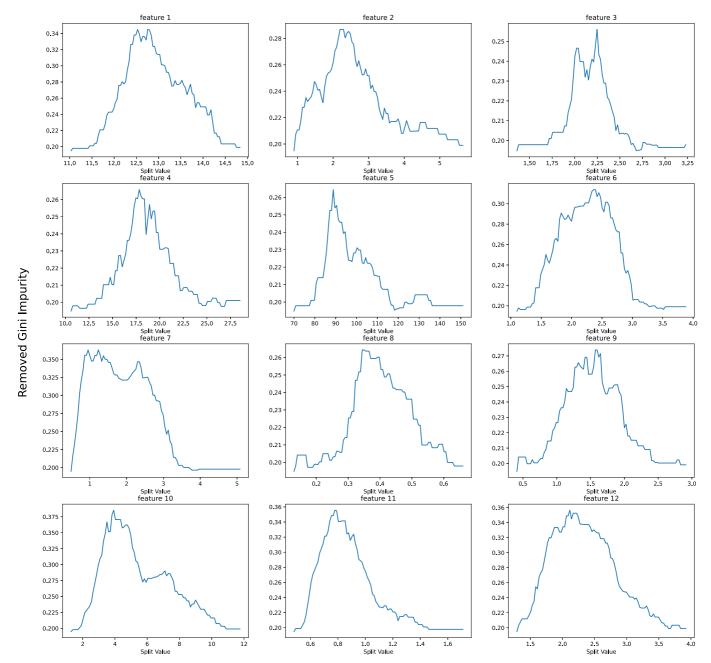
gini before: 0.6583133442747128
gini after: 0.3931706460674157
```

Split the data along the first 12 features and plot the removed Gini impurity at different values of this feature. Which is the optimal split?

```
# Vereinfacher durch pandas:
import pandas as pd
X = pd.DataFrame(X)
t = pd.Series(t)
```

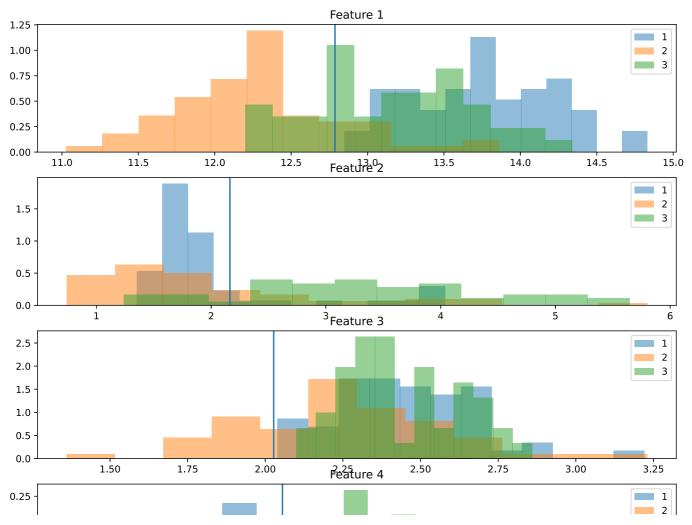
```
def splittest(wert, zeile, y):
    links, rechts = list(), list()
    n = len(zeile)
    for i in range(n):
        if zeile[i] < wert:
            links.append(int(y[i]))
        else:
            rechts.append(int(y[i]))
# For format reasons:
links, rechts = np.array(links), np.array(rechts)</pre>
```

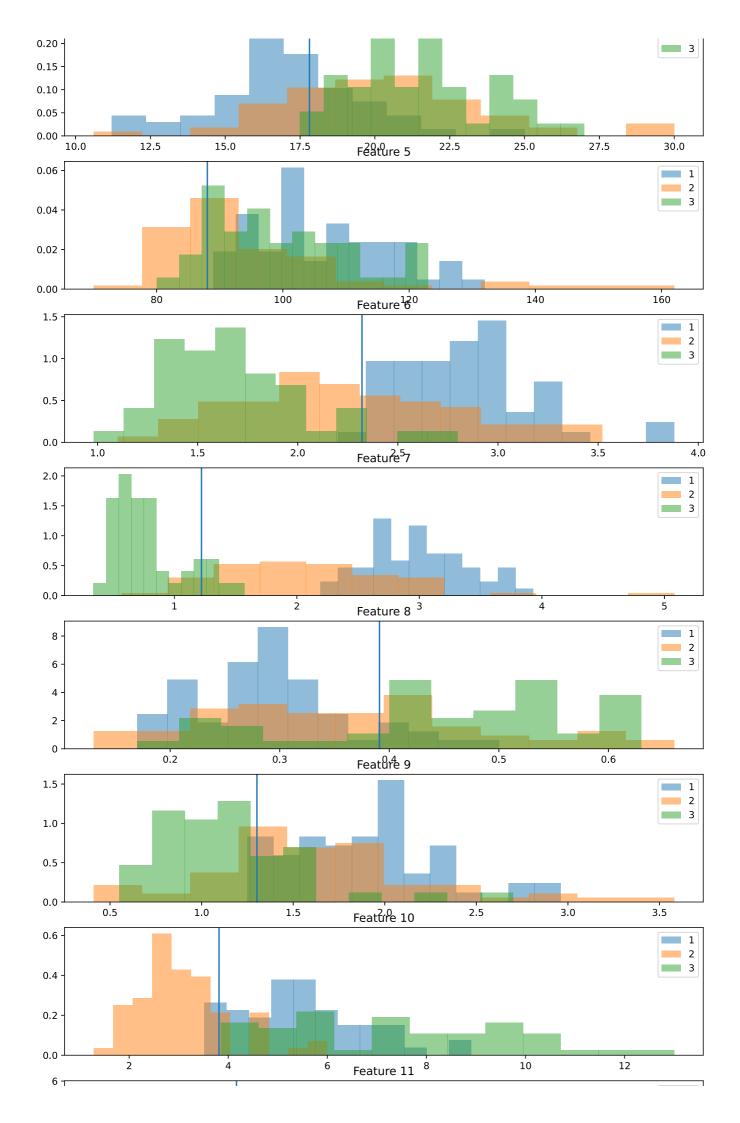
```
gini_reduc = np.ones((100,12))
thresholds = np.ones((100,12))
for jj in range(12):
   feat = X_train[:,jj]
   var_min = feat.min()
    var_max = feat.max()
    seq = np.linspace(var_min,var_max,100)
    thresholds[:,jj] = seq
    for ii in range(len(seq)):
        t_0 = t_train[feat < seq[ii]]</pre>
        t_1 = t_train[feat >= seq[ii]]
        gini_before = gini_impurity(t)
        gini_after = len(t_0)/len(t) * gini_impurity(t_0) + len(t_1)/len(t) * gini_impurity(t_0)
_1)
        gini_reduc[ii,jj] = gini_before - gini_after
fig, ax = plt.subplots(4,3,figsize = (20,20))
fig.text(0.07, .5, "Removed Gini Impurity", fontsize = 20, va='center', rotation='vertical')
count = 0
for jj in range(4):
    for ii in range(3):
        ax[jj,ii].plot(thresholds[:,count],gini_reduc[:,count])
        ax[jj,ii].set_title("feature " + str(count+1))
        ax[jj,ii].set_xlabel("Split Value")
        count += 1
```

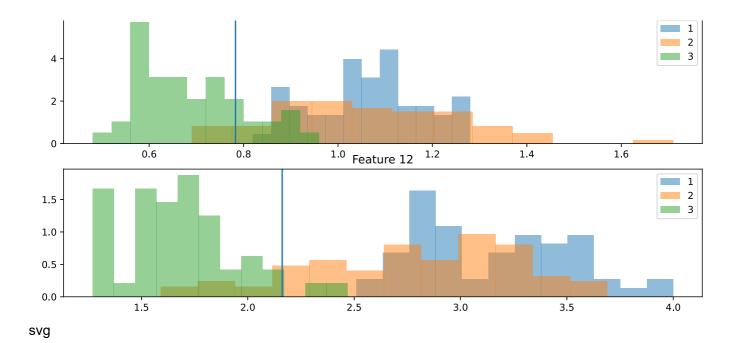


svg

```
fig, ax = plt.subplots(12, figsize = (12, 35))
for i in range(12):
    x_min, x_max = min(X.iloc[:, i]), max(X.iloc[:, i])
    x_range = np.linspace(start=x_min, stop=x_max, num=200)
   n, _ = X.shape
   bester_split, bester_gini = 0.0, 1.0
    for j in range(len(x_range)):
        links, rechts = splittest(wert = x_range[j], zeile=X.iloc[:, i], y=t) # Aufteilen
        links_gini, rechts_gini = gini_impurity(links), gini_impurity(rechts) #Gini für beide
Splits
        weighted_gini = (len(links)/n)*links_gini + (len(rechts)/n)*rechts_gini # Gewichteter
Gini
        if weighted_gini < bester_gini: # iteratives zuweisen</pre>
            bester_gini, bester_split = weighted_gini, x_range[j]
            bester_left, bester_right = links, rechts
    temp_ax = ax[i]
    feature1 = X.loc[t == 0, i]
    temp_ax.hist(feature1, alpha=0.5, bins=12, density=True, label = '1')
    feature2 = X.loc[t == 1, i]
    temp_ax.hist(feature2, alpha=0.5, bins=12, density=True, label = '2')
    feature3 = X.loc[t == 2, i]
    temp_ax.hist(feature3, alpha=0.5, bins=12, density=True, label = '3')
    temp_ax.axvline(x=bester_split)
    temp_ax.legend()
    name = 'Feature '
    name += str(i+1)
    temp_ax.set_title(name)
```







Task 2.3

- 1. Implement a function build_tree(X, t, depth) which recursively builds a tree. Use the classes Node and Leaf as a data structure to build your tree.
- 2. Implement a function <code>predict_tree(tree, x)</code> which makes a prediction for sample <code>x</code> . Obtain scores for the <code>wine dataset</code> and compare to <code>sklearn.tree.DecisionTree</code> .
- 3. Switch back to the synthetic 2d dataset from the beginning (kernel methods). Compute scores and visualize the decisions in a 2d grid.

```
class Node:
    def __init__(self, links, rechts, n_feat, thresholdold):
        self.links = links
        self.rechts = rechts
        self.n feat = n feat
        self.thresholdold = thresholdold
    # Funktion um Format ausgeben zu können
    def desc(self):
      return(self.n_feat, self.thresholdold)
    # Folgt ein Split/Child Node?
    def followingSplits(self):
      return True
    # Prediction Function:
    def predict_tree(self, x):
      print(self.followingSplits())
      if(self.followingSplits()):
        feat_n, threshold = self.desc()
        feat_x = x[feat_n]
        if(feat_x < threshold):</pre>
          return self.links.predict_tree(x)
          return self.rechts.predict tree(x)
      else:
        return self.desc()
```

```
class Leaf:
    def __init__(self, label):
        self.label = label

# braucht man erst für die plots wirklich

def desc(self):
    return(self.label)

# Um zu stoppen

def followingSplits(self):
    return False
```

```
def treeplot(node, x=[]):
    # cutoff points for plots
if node.followingSplits():
    x.append(node.desc())
    if node.links is not None:
        treeplot(node.links,x)
    if node.rechts is not None:
        treeplot(node.rechts,x)
```

```
def build_tree(X, t, depth, max_depth=3, n_labels=2):
    # wegen: Cannot cast array data from dtype('float64') to dtype('int64') according to the
 rule 'safe'
    if len(X) <=0:
      return
    if len(X) <= 1:
      return Leaf(np.argmax(np.bincount(t)))
    if depth == max_depth:
      return Leaf(np.argmax(np.bincount(t)))
    #start werte
    current_gini, temp_feat, current_split = 1, None , 0.0
    for i in range(X.shape[1]):
      x_{min}, x_{max} = min(X.iloc[:, i]), max(X.iloc[:, i])
      x_range = np.linspace(start=x_min, stop=x_max, num=200)
      n = X.shape[0]
      for j in range(len(x_range)):
          links, rechts = splittest(wert = x_range[j], zeile=X.iloc[:, i], y=t) # split colum
n into links and rechts
          links_gini, rechts_gini = gini_impurity(links), gini_impurity(rechts) # calculate
 gini coef for links and rechts
          weighted_gini = (len(links)/n)*links_gini + (len(rechts)/n)*rechts_gini # calculate
wighted gini
          if weighted_gini < current_gini: # save best gini</pre>
              current_gini, current_split = weighted_gini, x_range[j]
              temp_links, temp_rechts, temp_feat = links, rechts, i
    links_X, rechts_X = [], []
    for k in range(X.shape[0]): # fröhliches iterieren
      if X.iloc[k,temp_feat] < current_split:</pre>
        links_X.append(X.iloc[k,:])
      else:
        rechts_X.append(X.iloc[k,:])
    node = Node(None,None,temp_feat,current_split) # node class nutzen
    node.links = build tree(pd.DataFrame(np.array(links X)), temp links, depth+1, max depth,
 n labels)
    node.rechts = build_tree(pd.DataFrame(np.array(rechts_X)), temp_rechts, depth+1, max_dept
h, n_labels)
    return node
```

```
def predict_tree(node, x):
   if(node.followingSplits()):
      feat_n, threshold = node.desc()
      feat_x = x[feat_n]
      if(feat_x < threshold):
        return predict_tree(node.links, x)
      else:
        return predict_tree(node.rechts, x)
   else:
      return node.desc()</pre>
```

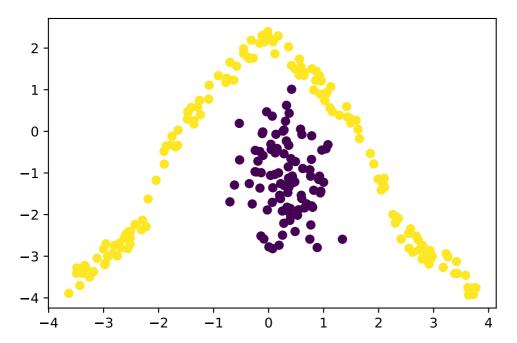
```
def split_data(X, y, frac=0.3, max_samples=None, seed=4711):
    if seed is not None:
        np.random.seed(seed)
    indices = np.random.permutation(len(X))
    indices = indices[:max_samples]
    indices_test, indices_train = indices[:int(len(indices)*frac)], indices[int(len(indices)*
frac):]
    assert (len(set(indices))) == len(indices)
    assert len(set(indices_test).intersection(indices_train)) == 0
   X_train, X_test = X[indices_train], X[indices_test]
   y_train, y_test = y[indices_train], y[indices_test]
    return X_train, X_test, y_train, y_test
X, t = load_wine(return_X_y=True)
X_train, X_test, t_train, t_test = split_data(X = X, y = t)
tree = build_tree(pd.DataFrame(X_train),t_train, 0, max_depth=3, n_labels=3)
count = 0
for i in range(len(X_test)):
 if(t_test[i] == predict_tree(tree, X_test[i])):
    count += 1
from sklearn import tree
clf = tree.DecisionTreeClassifier(max_depth=3)
clf = clf.fit(X_train, t_train)
predsk = clf.predict(X_test)
skcounter = 0
for i in range(len(predsk)):
 if(predsk[i] == t_test[i]):
    skcounter += 1
print("sklearn score:", skcounter/len(t_test))
print("Our score:", (count)/len(t_test))
sklearn score: 0.9245283018867925
Our score: 0.9245283018867925
X_2d, t_2d = np.load('data/nonlin_2d_data.npy')[:,:2], np.load('data/nonlin_2d_data.npy')[:,
X_train, X_test, y_train, y_test = split_data(X_2d, t_2d)
y_train = y_train.astype(int)
```

tree = build_tree(pd.DataFrame(X_train),y_train, 0, max_depth=3, n_labels=2)

```
correct = 0
for i in range(len(X_test)):
   if(y_test[i] == predict_tree(tree, X_test[i])):
      correct += 1
print(correct/len(y_test))
```

1.0

```
fig, ax = plt.subplots()
plt.scatter(X_2d[:,0], X_2d[:, 1], c=t_2d)
plt.show()
```

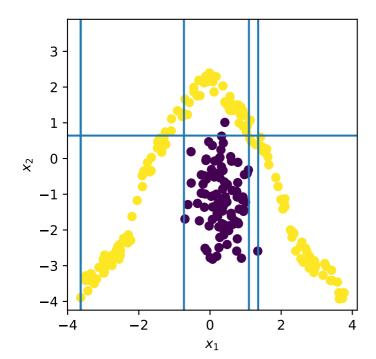


svg

```
cutoff = []
treeplot(tree, cutoff)
print(cutoff)
```

[(0, -0.7324396890211675), (0, -3.633659394608522), (0, -3.633659394608522), (0, 1.0942337613 46704), (1, 0.6419593869827454), (0, 1.35390440889765)]

```
fig, ax = plt.subplots()
plt.scatter(X_2d[:,0], X_2d[:, 1], c=t_2d)
plt.xlabel('$x_1$')
plt.ylabel('$x_2$')
plt.axis('square')
#ax.axhline(y=1)
for i in cutoff:
    if i[0]:
        ax.axhline(y=i[1])
else:
        ax.axvline(x=i[1])
```



svg