```
In [1]:
```

```
import pandas as pd
from sklearn.datasets import load_iris
from sklearn.datasets import load_wine
# from sklearn.neural_network import MLPClassifier
from sklearn.tree import DecisionTreeClassifier
# from sklearn.neighbors import KNeighborsClassifier
from sklearn.base import BaseEstimator
from sklearn.metrics import accuracy_score, recall_score, precision_score, fl_score
from sklearn.model_selection import train_test_split
import matplotlib.pyplot as plt
from sklearn.ensemble import BaggingClassifier, AdaBoostClassifier, RandomForestClassifier
import numpy as np
import numpy as np
import warnings
warnings.filterwarnings('ignore')
```

Ładowanie danych

```
In [2]:
```

```
iris = load iris(as frame=True)
wine = load_wine(as_frame=True)
glass = \{\}
glass['data'] = pd.read csv('glass.data', index col=0, header=None, names=[
    'Id', 'RI', 'Na', 'Mg',
    'Al', 'Si', 'K', 'Ca',
    'Ba', 'Fe', 'target'
])
glass['target'] = glass['data']['target']
glass['data'] = glass['data'].drop(columns=['target'])
seeds = {}
seeds['data'] = pd.read csv('seeds dataset.txt', sep='\t', header=None, names=[
    'area',
    'perimeter',
    'compactness',
    'length of kernel',
    'width_of_kernel',
    'asymmetry_coefficient',
    'length of kernel groove',
    'target'
])
seeds['target'] = seeds['data']['target']
seeds['data'] = seeds['data'].drop(columns=['target'])
```

In [3]:

```
random_state = 42
def get_estimator():
    return DecisionTreeClassifier(max_depth=1, random_state=random_state)
```

In [4]:

```
def test_model(model: BaseEstimator, X_train, X_test, y_train, y_test):
    model.fit(X_train, y_train)
    y_pred = model.predict(X_test)
    print('Accuracy:', accuracy_score(y_test, y_pred))
    print('Recall:', recall_score(y_test, y_pred, average='weighted'))
    print('Precision:', precision_score(y_test, y_pred, average='weighted'))
    print('F1:', f1_score(y_test, y_pred, average='weighted'))
```

In [5]:

```
def split_data(data):
    return train_test_split(data['data'], data['target'], test_size=0.2, random_state=ra
ndom_state)
```

```
In [6]:
```

```
data = split_data(iris)
test_model(get_estimator(), *data)
data = split_data(wine)
test_model(get_estimator(), *data)
data = split_data(glass)
test_model(get_estimator(), *data)
data = split_data(seeds)
test_model(get_estimator(), *data)
```

Recall: 0.63333333333333333 Precision: 0.46833333333333334 F1: 0.5195402298850574 Precision: 0.5671296296296295 F1: 0.5951417004048583 Accuracy: 0.5116279069767442 Recall: 0.5116279069767442 Precision: 0.3162790697674419 F1: 0.37209302325581395 Accuracy: 0.5952380952380952 Recall: 0.5952380952380952 Precision: 0.4362244897959183 F1: 0.48107448107448103

In [7]:

```
def plot bagging(dataset):
   X_train, X_test, y_train, y_test = split_data(dataset)
   fig, ax = plt.subplots(3, sharey=True)
   acc = []
   recall = []
   prec = []
   f1 = []
   n = stimators = list(range(5, 15))
   for i in n estimators:
       model = BaggingClassifier(base_estimator=get_estimator(), n_estimators=i,
                                  random state=random state).fit(X train, y train)
       y pred = model.predict(X test)
       acc.append(accuracy_score(y_test, y_pred))
       recall.append(recall_score(y_test, y_pred, average='weighted'))
       prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1_score(y_test, y_pred, average='weighted'))
   ax[0].plot(n_estimators, acc, label='accuracy')
   ax[0].plot(n_estimators, recall, label='recall')
   ax[0].plot(n_estimators, prec, label='precision')
   ax[0].plot(n_estimators, f1, label='f1')
   ax[0].set xlabel('estimators')
   acc = []
   recall = []
   prec = []
   f1 = []
   \max \text{ samples} = \text{list(np.arange(0.1, 1.1, 0.1))}
   for i in max samples:
       model = BaggingClassifier(base estimator=get estimator(), max samples=i,
                                  random state=random state).fit(X train, y train)
       y pred = model.predict(X test)
       acc.append(accuracy_score(y_test, y_pred))
       recall.append(recall_score(y_test, y_pred, average='weighted'))
       prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1_score(y_test, y_pred, average='weighted'))
   ax[1].plot(max_samples, acc, label='accuracy')
   ax[1].plot(max_samples, recall, label='recall')
   ax[1].plot(max_samples, prec, label='precision')
   ax[1].plot(max samples, f1, label='f1')
   ax[1].set xlabel('samples (fraction of dataset)')
```

```
acc = []
recall = []
prec = []
f1 = []
max features = list(range(1, len(dataset['data'].columns)+1))
for i in max features:
    model = BaggingClassifier(base estimator=get estimator(), max features=i,
                              random state=random state).fit(X train, y train)
    y pred = model.predict(X test)
    acc.append(accuracy score(y test, y pred))
    recall.append(recall score(y test, y pred, average='weighted'))
    prec.append(precision score(y test, y pred, average='weighted'))
    f1.append(f1_score(y_test, y_pred, average='weighted'))
ax[2].plot(max_features, acc, label='accuracy')
ax[2].plot(max features, recall, label='recall')
ax[2].plot(max_features, prec, label='precision')
ax[2].plot(max features, f1, label='f1')
ax[2].set xlabel('features')
ax[0].set ylabel('score')
ax[0].legend()
fig.set size inches (19.2, 14.4)
fig.suptitle('Bagging', fontsize=16)
print('Without bootstrap')
model = BaggingClassifier(base estimator=get estimator(), bootstrap=False,
                              random state=random state)
test model (model, X train, X test, y train, y test)
print('\nWith bootstrap')
model = BaggingClassifier(base estimator=get estimator(), bootstrap=True,
                              random state=random state)
test model (model, X train, X test, y train, y test)
```

In [8]:

```
def plot boost(dataset):
   X_train, X_test, y_train, y_test = split_data(dataset)
    fig, ax = plt.subplots(2, sharey=True)
   acc = []
   recall = []
    prec = []
    f1 = []
    n = stimators = list(range(25, 75, 5))
    for i in n estimators:
        model = AdaBoostClassifier(base estimator=get estimator(), n estimators=i,
                                  random state=random state).fit(X train, y train)
        y pred = model.predict(X_test)
        acc.append(accuracy score(y test, y pred))
        recall.append(recall_score(y_test, y_pred, average='weighted'))
        prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1 score(y test, y pred, average='weighted'))
    ax[0].plot(n estimators, acc, label='accuracy')
    ax[0].plot(n estimators, recall, label='recall')
    ax[0].plot(n estimators, prec, label='precision')
    ax[0].plot(n estimators, f1, label='f1')
    ax[0].set xlabel('estimators')
    acc = []
   recall = []
    prec = []
    f1 = []
    learning_rate = list(np.arange(0.5, 1.5, 0.1))
    for i in learning rate:
       model = AdaBoostClassifier(base estimator=get estimator(), learning rate=i,
                                  random state=random state).fit(X train, y train)
        y pred = model.predict(X test)
        acc.append(accuracy score(y test, y pred))
        recall.append(recall_score(y_test, y_pred, average='weighted'))
        prec.append(precision score(y test, y pred, average='weighted'))
```

```
f1.append(f1_score(y_test, y_pred, average='weighted'))
ax[1].plot(learning_rate, acc, label='accuracy')
ax[1].plot(learning_rate, recall, label='recall')
ax[1].plot(learning_rate, prec, label='precision')
ax[1].plot(learning_rate, f1, label='f1')
ax[1].set_xlabel('learning_rate')

ax[0].set_ylabel('score')
ax[0].legend()
fig.set_size_inches(19.2, 9.6)
fig.suptitle('Boosting', fontsize=16)
```

In [9]:

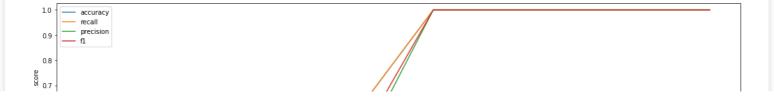
```
def plot forest(dataset):
    X_train, X_test, y_train, y_test = split data(dataset)
    fig, ax = plt.subplots(4, sharey=True)
    acc = []
   recall = []
   prec = []
   f1 = []
   max_samples = list(np.arange(0.1, 1.0, 0.1))
   for i in max samples:
        model = RandomForestClassifier(max samples=i,
                                  random state=random state).fit(X train, y train)
        y pred = model.predict(X test)
        acc.append(accuracy_score(y_test, y_pred))
        recall.append(recall_score(y_test, y_pred, average='weighted'))
        prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1_score(y_test, y_pred, average='weighted'))
    ax[0].plot(max samples, acc, label='accuracy')
    ax[0].plot(max samples, recall, label='recall')
    ax[0].plot(max samples, prec, label='precision')
    ax[0].plot(max samples, f1, label='f1')
   ax[0].set xlabel('samples (fraction of dataset)')
   acc = []
   recall = []
   prec = []
   f1 = []
   max features = list(range(1, len(dataset['data'].columns)+1))
    for i in max features:
        model = RandomForestClassifier(max features=i,
                                  random state=random state).fit(X train, y train)
        y pred = model.predict(X test)
        acc.append(accuracy_score(y_test, y_pred))
        recall.append(recall_score(y_test, y_pred, average='weighted'))
        prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1_score(y_test, y_pred, average='weighted'))
    ax[1].plot(max_features, acc, label='accuracy')
    ax[1].plot(max_features, recall, label='recall')
    ax[1].plot(max features, prec, label='precision')
    ax[1].plot(max features, f1, label='f1')
    ax[1].set xlabel('features')
    acc = []
    recall = []
   prec = []
    f1 = []
    n = stimators = list(range(50, 150, 10))
    for i in n estimators:
        model = RandomForestClassifier(n estimators=i,
                                  random state=random state).fit(X train, y train)
        y pred = model.predict(X test)
        acc.append(accuracy_score(y_test, y_pred))
        recall.append(recall_score(y_test, y_pred, average='weighted'))
        prec.append(precision_score(y_test, y_pred, average='weighted'))
        f1.append(f1_score(y_test, y_pred, average='weighted'))
    ax[2].plot(n_estimators, acc, label='accuracy')
    ax[2].plot(n estimators, recall, label='recall')
```

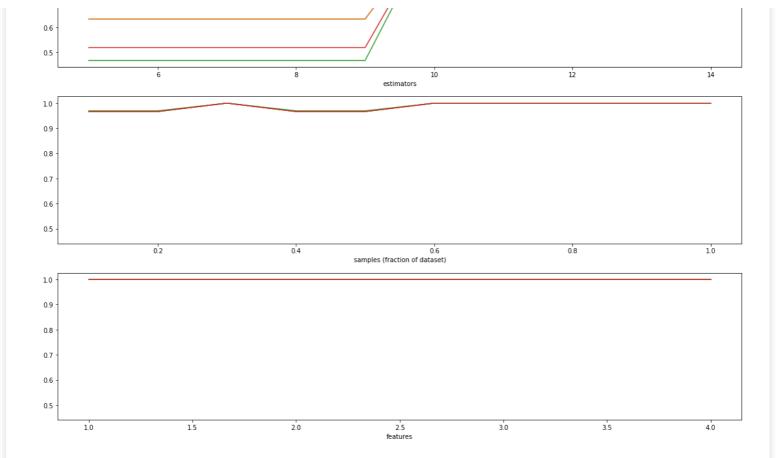
```
ax[2].plot(n_estimators, prec, label='precision')
ax[2].plot(n_estimators, f1, label='f1')
ax[2].set xlabel('trees')
acc = []
recall = []
prec = []
f1 = []
max depth = list(range(5, 15))
for i in max depth:
    model = RandomForestClassifier(max depth=i,
                              random state=random state).fit(X train, y train)
    y pred = model.predict(X test)
    acc.append(accuracy score(y test, y pred))
    recall.append(recall_score(y_test, y_pred, average='weighted'))
    prec.append(precision_score(y_test, y_pred, average='weighted'))
    f1.append(f1_score(y_test, y_pred, average='weighted'))
ax[3].plot(max_depth, acc, label='accuracy')
ax[3].plot(max_depth, recall, label='recall')
ax[3].plot(max_depth, prec, label='precision')
ax[3].plot(max depth, f1, label='f1')
ax[3].set_xlabel('max depth')
ax[0].set ylabel('score')
ax[0].legend()
fig.set size inches (19.2, 19.2)
fig.suptitle('Random Forest', fontsize=16)
```

Iris

```
In [10]:
dataset = iris
In [11]:
data = split data(dataset)
test model(get estimator(), *data)
Accuracy: 0.63333333333333333
Recall: 0.633333333333333333
Precision: 0.4683333333333334
F1: 0.5195402298850574
In [12]:
plot bagging(dataset)
Without bootstrap
Accuracy: 0.63333333333333333
Recall: 0.633333333333333333
Precision: 0.4683333333333334
F1: 0.5195402298850574
With bootstrap
Accuracy: 1.0
Recall: 1.0
Precision: 1.0
F1: 1.0
```

Bagging

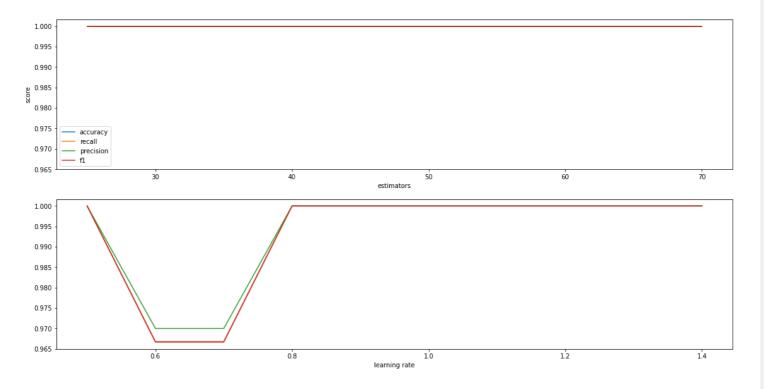




In [13]:

plot_boost(dataset)

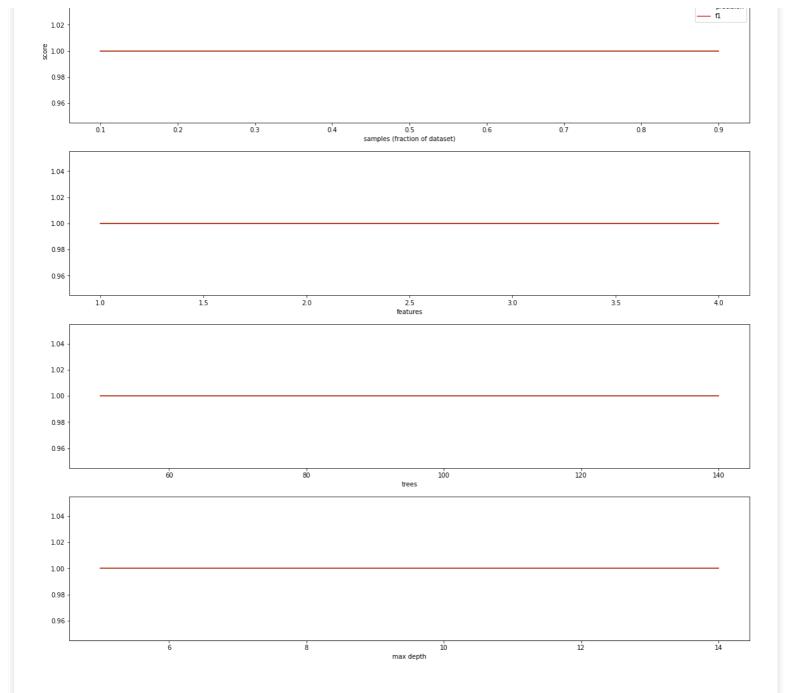
Boosting



In [14]:

plot_forest(dataset)

Random Forest



Wine

```
In [15]:
```

dataset = wine

In [16]:

data = split_data(dataset)
test_model(get_estimator(), *data)

F1: 0.5951417004048583

In [17]:

```
plot_bagging(dataset)
```

Without bootstrap

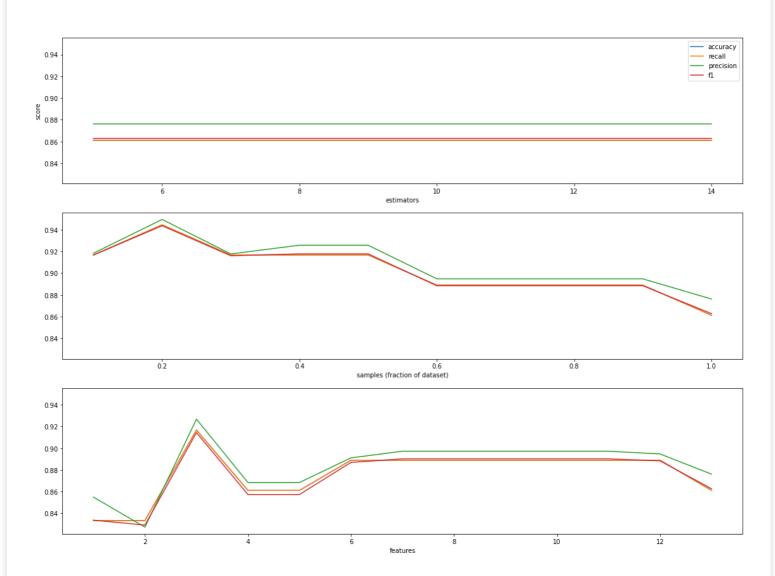
F1: 0.5951417004048583

with bootstrap

Accuracy: 0.8611111111111112 Recall: 0.861111111111112 Precision: 0.8760893246187363

F1: 0.8626321110192079

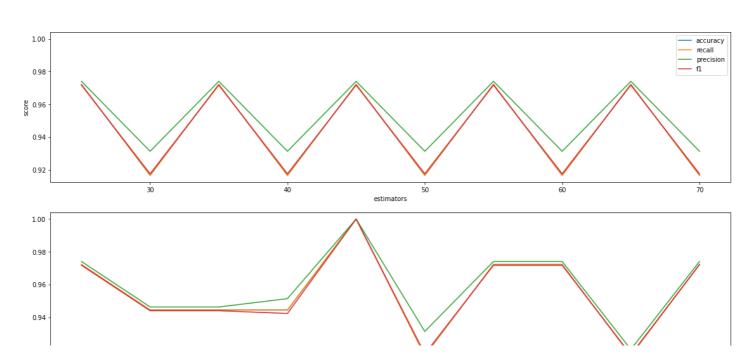
Bagging



In [18]:

plot_boost(dataset)

Boosting

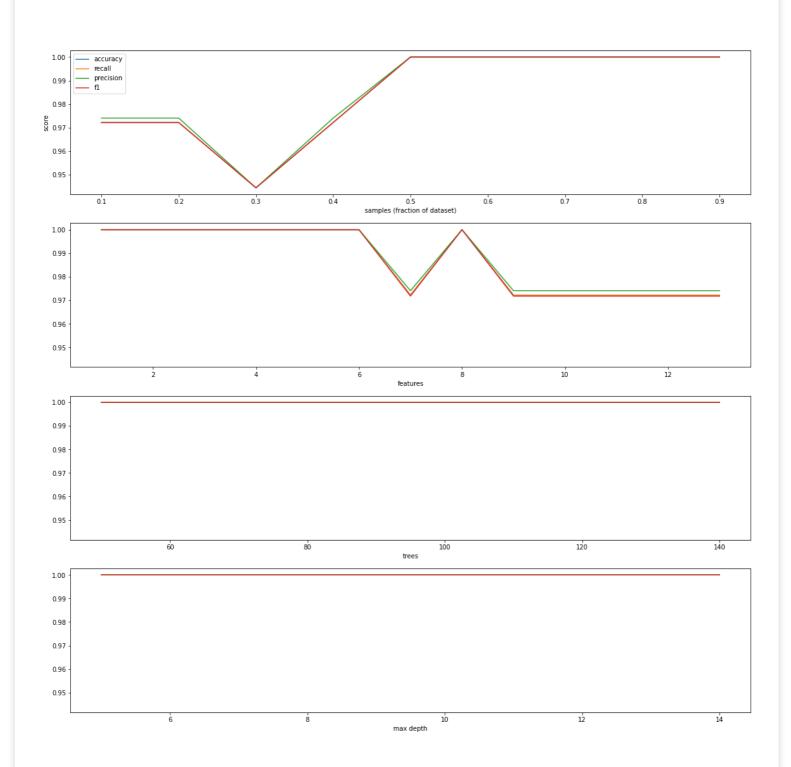


```
0.92 0.6 0.8 10 12 14
```

In [19]:

```
plot_forest(dataset)
```

Random Forest



Glass

```
In [20]:
```

```
dataset = glass
```

In [21]:

```
data = split_data(dataset)
test_model(get_estimator(), *data)
```

Accuracy: 0.5116279069767442 Recall: 0.5116279069767442 Precision: 0.3162790697674419

F1: 0.37209302325581395

In [22]:

plot bagging(dataset)

Without bootstrap

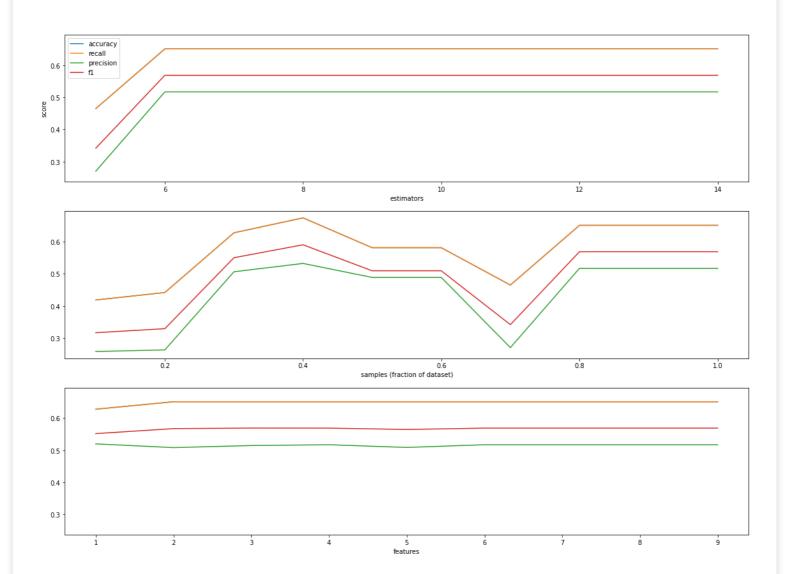
Accuracy: 0.5116279069767442
Recall: 0.5116279069767442
Precision: 0.3162790697674419
F1: 0.37209302325581395

With bootstrap

Accuracy: 0.6511627906976745 Recall: 0.6511627906976745 Precision: 0.5172888616891065

F1: 0.5689922480620155

Bagging

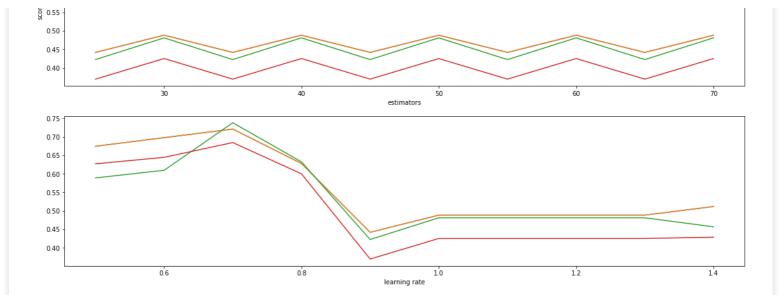


In [23]:

plot_boost(dataset)

Boosting

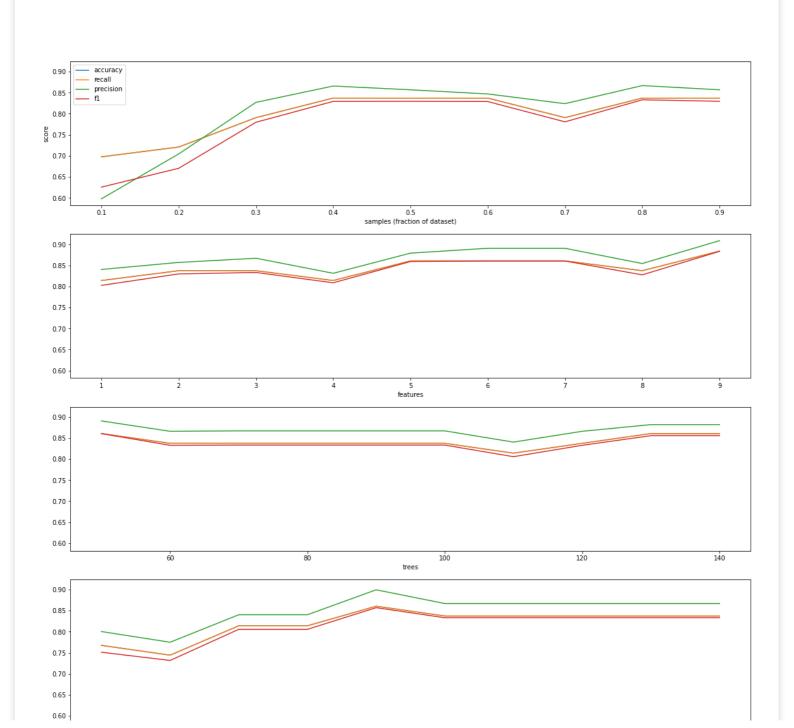




In [24]:

plot_forest(dataset)

Random Forest



6 8 10 12 14 max depth

Seeds

In [25]:

dataset = seeds

In [26]:

data = split_data(dataset)
test_model(get_estimator(), *data)

Accuracy: 0.5952380952380952 Recall: 0.5952380952380952 Precision: 0.4362244897959183

F1: 0.48107448107448103

In [27]:

plot bagging(dataset)

Without bootstrap

Accuracy: 0.5952380952380952 Recall: 0.5952380952380952 Precision: 0.4362244897959183

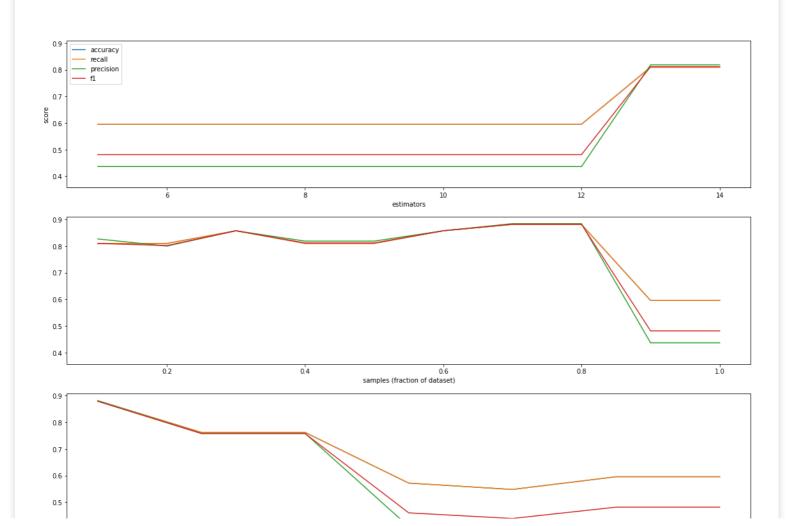
F1: 0.48107448107448103

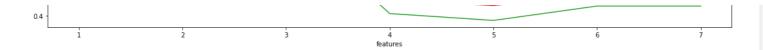
With bootstrap

Accuracy: 0.5952380952380952 Recall: 0.5952380952380952 Precision: 0.4362244897959183

F1: 0.48107448107448103

Bagging

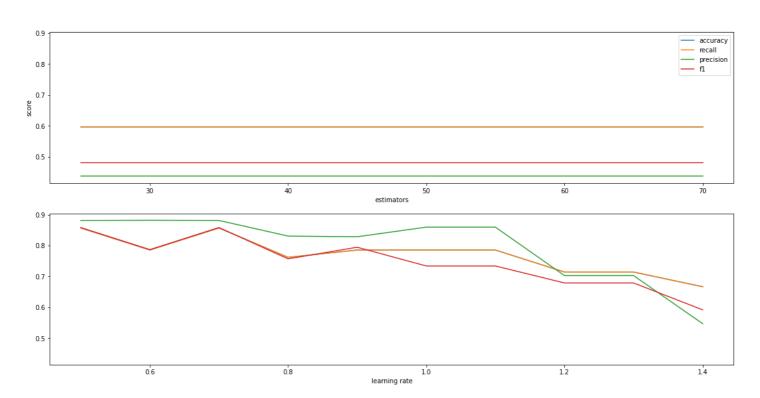




In [28]:

plot_boost(dataset)

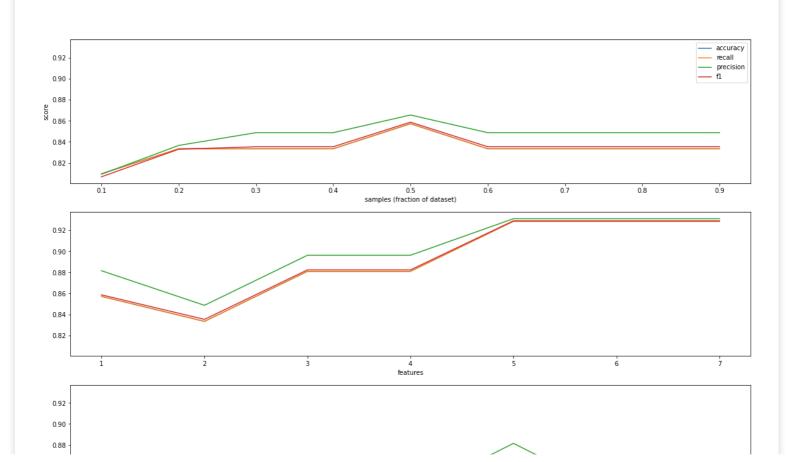
Boosting

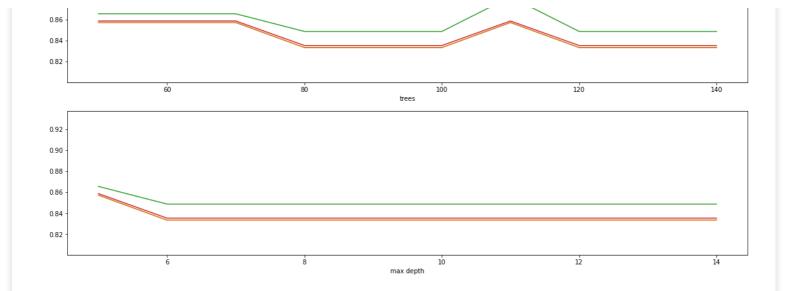


In [29]:

plot_forest(dataset)

Random Forest





Obserwacje i wnioski

Dla wszystkich 3 zbiorów najlepsze wyniki osiągnął algorytm random forest, ale też najdłużej działał.

Algorytm bagging z reguły działał lepiej dla większej liczby estymatorów. Wpływ liczby próbek i atrybutów zależał od zbioru.

W przypadku AdaBoost wpływ liczby estymatorów znacząco różni się dla wszystkich 3 zbiorów: w Wine działa lepiej dla nieparzystej liczby estymatorów, w Glass dla parzystej, a w Seeds nie ma wpływu na jakość działania. Dla 2 z 3 zbiorów jakość spadała dla zbyt dużego learning rate.

Dla RandomForest jakość predykcji rośnie wraz z ilością próbek dla 2 z 3 zbiorów. Podobnie z ilością atrybutów. Ilość drzew w lesie nie ma wielkiego wpływu na wyniki, a zwiększenie maksymalnej wysokości drzewa poprawia wyniki.

Sposób działania powyższych metod nie wyklucza możliwości użycia walidacji krzyżowej do zmniejszenia zagrożenia przeuczeniem, chociaż te metody same dzielą dane na mniejsze zbiory, co po dodaniu podziału z walidacji krzyżowej mogło by stworzyć przy małych zestawach danych wejściowych. Dodatkowo dla metod z bootstrapem, gdzie próbki pobierane są ze zwracaniem mniejszy zbiór danych może zwiększyć zagrożenie przeuczeniem.