In [1]: import numpy as np

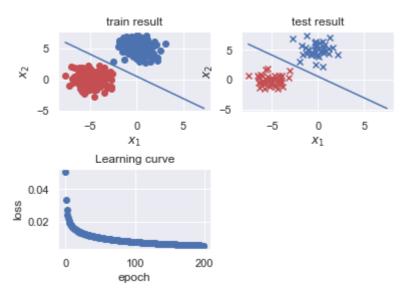
```
from drawer import *
        from data gen import *
        from sklearn.model selection import train test split
        import seaborn as sns
        sns.set theme()
        SEED=0
In [2]: ## define the function
In [3]: def sigmoid(s):
            return 1/(1+np.exp(-s))
        class CrossEntropy():
            "use the sigmoid function as the activate function"
            def backward(self, X, y, w):
                s=-y*(X@w)
                h=sigmoid(s)
                return (-y*X.T)@h
            def call (self, X, y, w):
                N=y.shape[0]
                return 1/N*np.sum(-np.log(sigmoid(y*(X@w))))
        class SGD():
            def init (self,learning rate=0.1):
                self.learning rate=learning rate
            # def zero grad(self):
            # self.grad[:]=0
            def step(self, w, grad):
                w = w - self.learning rate*grad
                return w
In [4]: class Logistic Regression():
            def init (self, data, loss fn, optimizer):
                self.data=(np.hstack((np.ones((data[1].shape[0],1)),data[0])),da
        ta[1])
                self.w=np.random.randn(3)
                self.loss fn=loss fn
                self.optimizer=optimizer
            def predict(self, X, is test=True):
                ""测试时做增广, fit里用的话不再做增广(初始化的时候已经做过了)""
                if is test:
                    X=np.hstack((np.ones((X.shape[0],1)),X))
                prob=sigmoid(X@self.w)
                return np.sign(prob-0.5),prob
            def validate(self, data):
                X=data[0]
                y=data[1]
                h , =np.sign(self.predict(X))
                mistake indices = np.where(h!=y)[0]
                return (X.shape[0]-len(mistake indices))/X.shape[0]
            def fit(self,epoch,batch size):
```

```
loss history={}
                for e in range(epoch):
                    batch num=self.data[1].shape[0]//batch size
                     for i in range(batch num):
                         try:
                             batch data=(self.data[0][i*batch size:(i+1)*batch si
        ze,:],self.data[1][i*batch size:(i+1)*batch size])
                         except:
                             batch data=(self.data[0][i*batch size:,:],self.data[
        1][i*batch size:])
                            batch size=self.data[1].shape[0]%batch size
                         # print(batch data)
                         pred , _ = self.predict(batch_data[0],is_test=False)
                         loss=self.loss fn(batch data[0],batch data[1],self.w)
                         # self.optimizer.zero grad()
                         grad=0
                         grad=self.loss fn.backward(batch data[0],batch data[1],s
        elf.w)
                         self.w=self.optimizer.step(self.w,grad)
                    loss history[e]=loss
                return loss history
In [5]: data=data generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
        X_train, X_test, y_train, y_test=train_test_split(data[0], data[1], train_siz
        e=0.8, test size=0.2, random state=SEED)
In [6]: def algorithm(learning rate=0.01,epoch=200,batch size=40,fig title=None)
             # sqd method
            loss fn=CrossEntropy()
            optimizer=SGD(learning rate=learning rate)
            model=Logistic Regression((X train,y train),loss fn,optimizer)
            loss history=model.fit(epoch,batch size)
            f=plt.figure()
            data=(X train, y train)
            plt.subplot(221)
            draw(data, model.w, marker='o')
            plt.title('train result')
            print(str(fig title),'train accuracy:', model.validate(data))
            data=(X test,y test)
            plt.subplot(222)
            draw(data, model.w, marker='x')
            plt.title('test result')
            print(str(fig title), 'test accuracy:', model.validate(data))
            # learning curve
            plt.subplot(2,2,3)
            plt.scatter(x=loss history.keys(),y=loss history.values())
            plt.xlabel('epoch')
            plt.ylabel('loss')
            plt.title('Learning curve')
            plt.subplots adjust(hspace=0.7)
            f.suptitle(fig title)
```

train accuracy: 1.0

algorithm(fig title='')

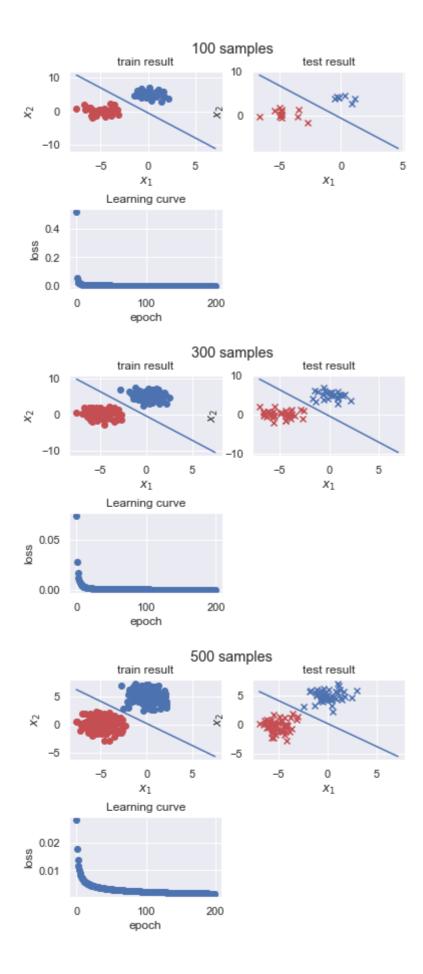
test accuracy: 1.0



展开实验

测试样本数量影响

```
In [7]: data=data generator([-5,0],np.eye(2),[0,5],np.eye(2),100,seed=SEED)
        X train, X test, y train, y test=train test split(data[0], data[1], train siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig_title='100 samples')
        data=data generator([-5,0],np.eye(2),[0,5],np.eye(2),300,seed=SEED)
        X_train, X_test, y_train, y_test=train_test_split(data[0], data[1], train_siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig title='300 samples')
        data=data generator([-5,0],np.eye(2),[0,5],np.eye(2),500,seed=SEED)
        X_train, X_test, y_train, y_test=train_test_split(data[0], data[1], train_siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig title='500 samples')
        100 samples train accuracy: 1.0
        100 samples test accuracy: 1.0
        300 samples train accuracy: 1.0
        300 samples test accuracy: 1.0
        500 samples train accuracy: 1.0
        500 samples test accuracy: 1.0
```



结论 可见数据量较大时, 较少的epoch即可获得较低的loss, 只用进行较少的迭代便可收敛

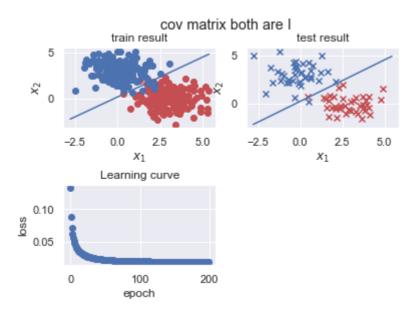
测试样本分布影响

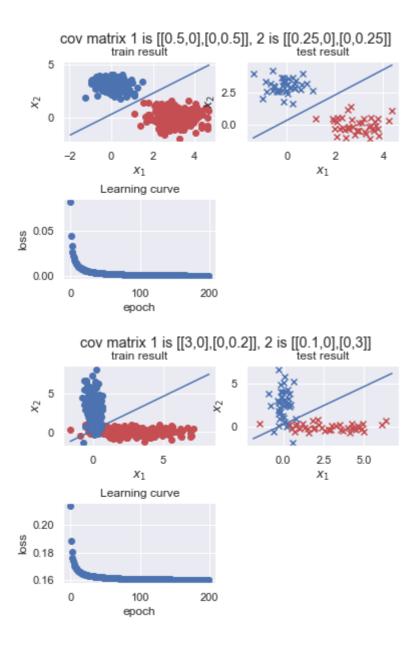
```
In [8]: data=data generator([3,0],np.eye(2),[0,3],np.eye(2),400,seed=SEED)
        X train, X test, y train, y test=train test split(data[0], data[1], train siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig title='cov matrix both are I')
        print('\n')
        data=data generator([3,0],[[0.5,0],[0,0.5]],[0,3],[[0.25,0],[0,0.25]],40
        0, seed=SEED)
        X train, X test, y train, y test=train test split(data[0], data[1], train siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig title='cov matrix 1 is [[0.5,0],[0,0.5]], 2 is [[0.25,0],[0,0.5]]
        0,0.25]]')
        print('\n')
        data=data generator([3,0],[[3,0],[0,0.2]],[0,3],[[0.1,0],[0,3]],400,seed
        =SEED)
        X_train, X_test, y_train, y_test=train_test_split(data[0], data[1], train_siz
        e=0.8, test size=0.2, random state=SEED)
        algorithm(fig title='cov matrix 1 is [[3,0],[0,0.2]], 2 is [[0.1,0],[0,3]
        ]]')
        print('\n')
```

cov matrix both are I train accuracy: 0.96875 cov matrix both are I test accuracy: 0.975

cov matrix 1 is [[0.5,0],[0,0.5]], 2 is [[0.25,0],[0,0.25]] train accura cy: 1.0 cov matrix 1 is [[0.5,0],[0,0.5]], 2 is [[0.25,0],[0,0.25]] test accurac y: 1.0

cov matrix 1 is [[3,0],[0,0.2]], 2 is [[0.1,0],[0,3]] train accuracy: 0.959375 cov matrix 1 is [[3,0],[0,0.2]], 2 is [[0.1,0],[0,3]] test accuracy: 0.9375





可见, 当样本较为分散时, 结果较差, 而当样本聚集程度较高, 方差比较小时, 结果较好

测试学习率的影响

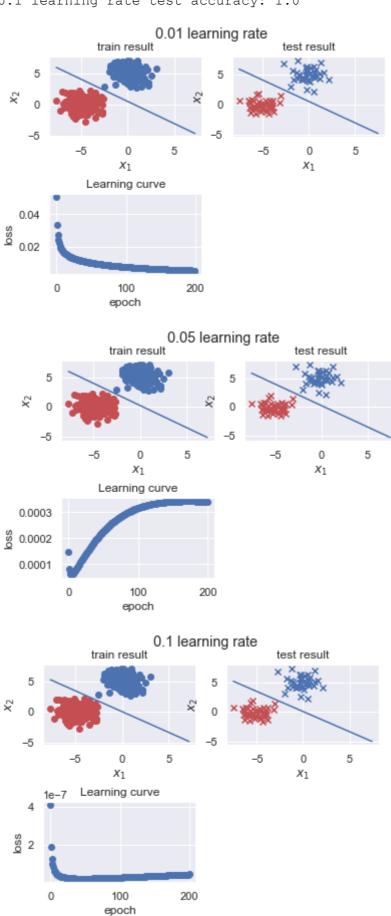
```
In [9]: data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(learning_rate=0.01,fig_title='0.01 learning rate')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(learning_rate=0.05,fig_title='0.05 learning rate')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(learning_rate=0.1,fig_title='0.1 learning_rate')
```

0.01 learning rate train accuracy: 1.0

0.01 learning rate test accuracy: 1.0 0.05 learning rate train accuracy: 1.0 0.05 learning rate test accuracy: 1.0 0.1 learning rate train accuracy: 1.0 0.1 learning rate test accuracy: 1.0



可见,适当调大学习率可以使损失函数收敛的更快,但如果没有调节好的话可能反而会使得损失函数变大

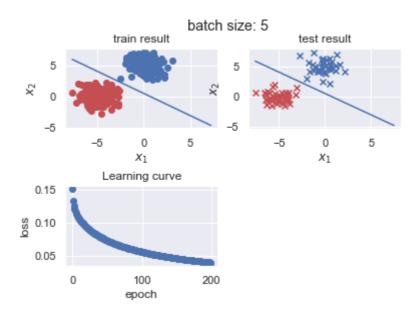
测试batch size的影响

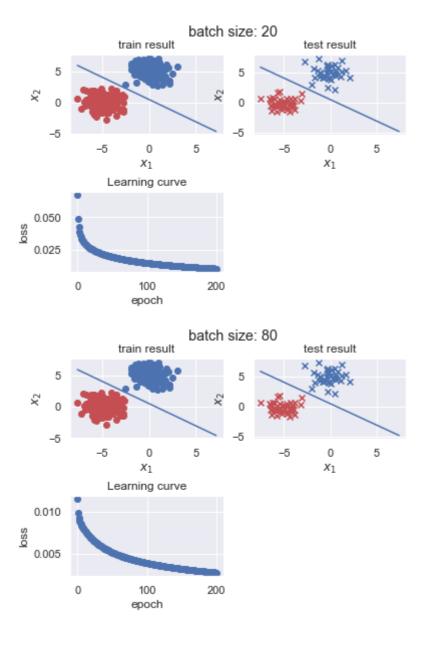
```
In [10]: data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(batch_size=5,fig_title='batch size: 5')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(batch_size=20,fig_title='batch size: 20')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(batch_size=80,fig_title='batch size: 80')
```

batch size: 5 train accuracy: 1.0 batch size: 5 test accuracy: 1.0 batch size: 20 train accuracy: 1.0 batch size: 20 test accuracy: 1.0 batch size: 80 train accuracy: 1.0 batch size: 80 test accuracy: 1.0





可见, 较大的batch size会使得结果收敛更快, 在较少的epoch数量后就得到较为理想的结果

测试epoch的影响

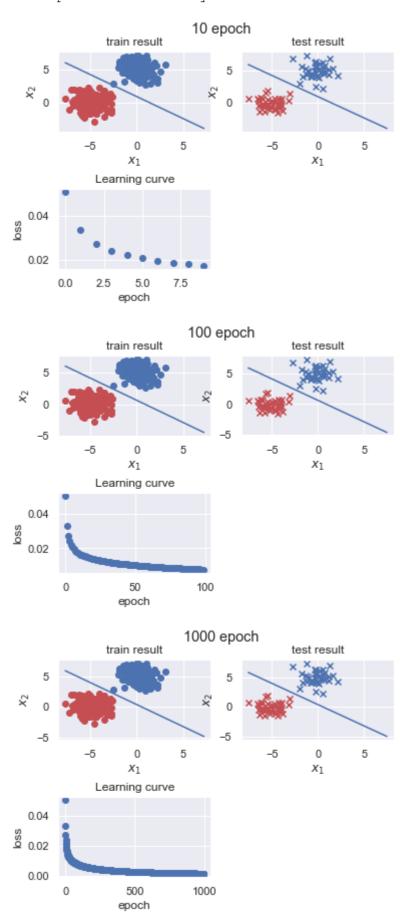
```
In [11]: data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(epoch=10,fig_title='10 epoch')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(epoch=100,fig_title='100 epoch')

data=data_generator([-5,0],np.eye(2),[0,5],np.eye(2),400,seed=SEED)
    X_train,X_test,y_train,y_test=train_test_split(data[0],data[1],train_siz
    e=0.8,test_size=0.2,random_state=SEED)
    algorithm(epoch=1000,fig_title='1000 epoch')
```

10 epoch train accuracy: 1.0

10 epoch test accuracy: 1.0 100 epoch train accuracy: 1.0 100 epoch test accuracy: 1.0 1000 epoch train accuracy: 1.0 1000 epoch test accuracy: 1.0



可以看出,对于当前这种样本分布,样本具有较为明显的线性分界线时,很少的epoch即可以达到收敛状态,并不需要进行很多次的迭代