

CS273 Homework 3

Due: Friday October 25 2024 (11:59pm)

Instructions

This homework (and subsequent ones) will involve data analysis and reporting on methods and results using Python code. You will submit a **single PDF file** that contains everything to Gradescope. This includes any text you wish to include to describe your results, the complete code snippets of how you attempted each problem, any figures that were generated, and scans of any work on paper that you wish to include. It is important that you include enough detail that we know how you solved the problem, since otherwise we will be unable to grade it.

Your homeworks will be given to you as Jupyter notebooks containing the problem descriptions and some template code that will help you get started. You are encouraged to use these starter Jupyter notebooks to complete your assignment and to write your report. This will help you not only ensure that all of the code for the solutions is included, but also will provide an easy way to export your results to a PDF file (for example, doing *print preview* and *printing to pdf*). I recommend liberal use of Markdown cells to create headers for each problem and sub-problem, explaining your implementation/answers, and including any mathematical equations. For parts of the homework you do on paper, scan it in such that it is legible (there are a number of free Android/iOS scanning apps, if you do not have access to a scanner), and include it as an image in the Jupyter notebook.

Double check that all of your answers are legible on Gradescope, e.g. make sure any text you have written does not get cut off.

If you have any questions/concerns about using Jupyter notebooks, ask us on EdD. If you decide not to use Jupyter notebooks, but go with Microsoft Word or LaTeX to create your PDF file, make sure that all of the answers can be generated from the code snippets included in the document.

Summary of Assignment: 100 total points

- Problem 1: Logistic Regression (25 points)
 - Problem 1.1: Decision boundaries (10 points)
 - Problem 1.2: Gradient optimization (10 points)
 - Problem 1.3: Evaluation (5 points)
- Problem 2: Linear Support Vector Machines (15 points)
 - Problem 2.1: Fitting & Evaluation (8 points)
 - Problem 2.2: Decision boundary & margin (7 points)
- Problem 3: Feature Expansions (20 points)
 - Problem 3.1: Polynomial Features (10 points)
 - Problem 3.2: Using Regularization (10 points)
- Problem 4: Logistic Regression on MNIST Data (35 points)
 - Problem 4.1: Initial Training (10 points)
 - Problem 4.2: Regularization (10 points)
 - Problem 4.3: Interpreting the Weights (5 points)
 - Problem 4.4: Evaluating class probabilities (5 points)
 - Problem 4.4: Learning Curves (5 points)
- Statement of Collaboration (5 points)

Before we get started, let's import some libraries that you will make use of in this assignment. Make sure that you run the code cell below in order to import these libraries.

Important: In the code block below, we set `seed=1234` . This is to ensure your code has reproducible results and is important for grading. Do not change this. If you are not using the provided Jupyter notebook, make sure to also set the random seed as below.

Important: Do not change any codes we give you below, except for those waiting for you to complete. This is to ensure your code has reproducible results and is important for grading.

```
In [3]: import numpy as np
import matplotlib.pyplot as plt

from sklearn.datasets import fetch_openml          # common data set access
from sklearn.preprocessing import StandardScaler  # scaling transform
```

```

from sklearn.model_selection import train_test_split # validation tools
from sklearn.metrics import zero_one_loss
from sklearn.inspection import DecisionBoundaryDisplay

from sklearn.preprocessing import PolynomialFeatures
from sklearn.preprocessing import StandardScaler

from sklearn.linear_model import SGDClassifier      # Used in 2D data problems
from sklearn.linear_model import LogisticRegression # Used in MNIST data problem

import requests          # we'll use these for reading data from a url
from io import StringIO

import warnings
warnings.filterwarnings('ignore')

# Some keyword arguments for making nice looking decision plots.
plot_kwargs = {'cmap': 'jet',      # another option: viridis
               'response_method': 'predict',
               'plot_method': 'pcolormesh',
               'shading': 'auto',
               'alpha': 0.5,
               'grid_resolution': 100}

# Fix the random seed for reproducibility
# !! Important !! : do not change this
seed = 1234
np.random.seed(seed)

```

Binary Classification Dataset

First, let's load our Housing dataset from HW1. To start, we will extract a two-dimensional binary classification problem, which will allow us to visualize the problem, training, and resulting model.

```

In [5]: # Load the features and labels from an online text file
url = 'https://sli.ics.uci.edu/extras/cs178/data/nyc_housing.txt'
with requests.get(url) as link:

```

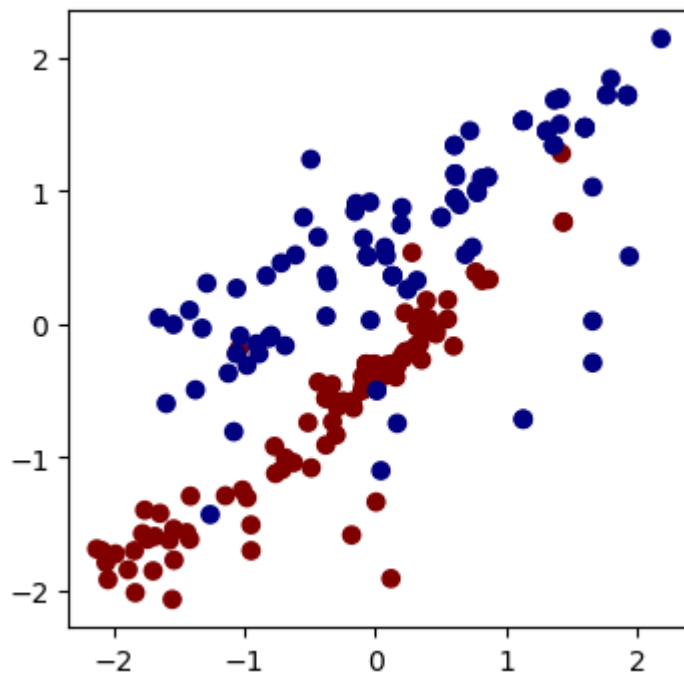
```

datafile = StringIO(link.text)
nych = np.genfromtxt(datafile,delimiter=',')
nych_X, nych_y = nych[:, :-1], nych[:, -1]

# Process the data to be only two classes and two real-valued & normalized features:
X, y = nych_X[nych_y<2, :2], nych_y[nych_y<2]
X -= X.mean(axis=0, keepdims=True) # remove mean
X /= X.std(axis=0, keepdims=True)  # & scale
y = 2*y - 1                       # classical binary: positive/negative

# Visualize the resulting dataset:
plt.figure(figsize=(4,4))
plt.scatter(X[:,0], X[:,1], c=y, cmap='jet');

```



Problem 1: Logistic Regression

The `scikit` package contains several implementations of logistic regression models for classification. In order to emphasize the similarities between different models, we will use the `SGDClassifier` object, which is a bit of a misnomer since SGD is an optimization technique, not a model. The object implements several types of linear classifiers, optimized using SGD or SGD-like training, depending on the loss function selected.

Problem 1.1: Decision Boundaries

First, let's build a linear classifier and manually set its parameters. Suppose that we initialize our linear classifier to make its predictions as,

$$\hat{y} = T(\theta_0 + \theta_1 x_1 + \theta_2 x_2)$$

with $[\theta_0, \theta_1, \theta_2] = [-2, 2, 1]$.

(a) What is the decision boundary of this classifier? (Answer in the form $x_2 = ax_1 + b$.)

Answer to (a): The decision boundary occurs when the argument of T is zero:

$$\theta_0 + \theta_1 x_1 + \theta_2 x_2 = 0.$$

Substitute the given values ($\theta_0 = -2$), ($\theta_1 = 2$), and ($\theta_2 = 1$):

$$-2 + 2x_1 + 1x_2 = 0.$$

Rearranging this equation to solve for (x_2), we get:

$$x_2 = -2x_1 + 2.$$

Thus, the decision boundary is:

$$x_2 = -2x_1 + 2.$$

Let's initialize the classifier and look at its decision function.

We will set the classifier to use the logistic negative log-likelihood surrogate loss (`loss=log_loss`); the other parameters prevent re-initializing the model later (`warm_start=True`) and set the stochastic gradient step size schedule (`learning_rate='adaptive'`) is

a simple backoff method, and initial step size `eta0=1e-3` is a small initial step size, so we can see the early progress).

(b) Add code below to plot your answer above on the decision function and verify that your answer matches `scikit` 's output:

```
In [10]: logreg = SGDClassifier(loss='log_loss', warm_start=True, learning_rate='adaptive', eta0 = 1e-3)
```

```
# Now let's initialize the model manually:
```

```
logreg.classes_ = np.unique(y)           # class IDs from the data
```

```
logreg.coef_ = np.array([[2.,1.]])
```

```
logreg.intercept_ = np.array([-2.])      # r(x) = 2*x1 + 1*x2 + (-2)
```

```
figure, axes = plt.subplots(1, 1, figsize=(4,4))
```

```
DecisionBoundaryDisplay.from_estimator(logreg, X, ax=axes, **plot_kwargs)
```

```
axes.scatter(X[:, 0], X[:, 1], c=y, edgecolor=None, s=12, cmap='jet')
```

```
### YOUR CODE STARTS HERE
```

```
# Generate x1 values within an appropriate range for the plot
```

```
x1_vals = np.linspace(-0.5, X[:, 0].max(), 100)
```

```
# Calculate the corresponding x2 values using the derived equation
```

```
x2_vals = -2 * x1_vals + 2
```

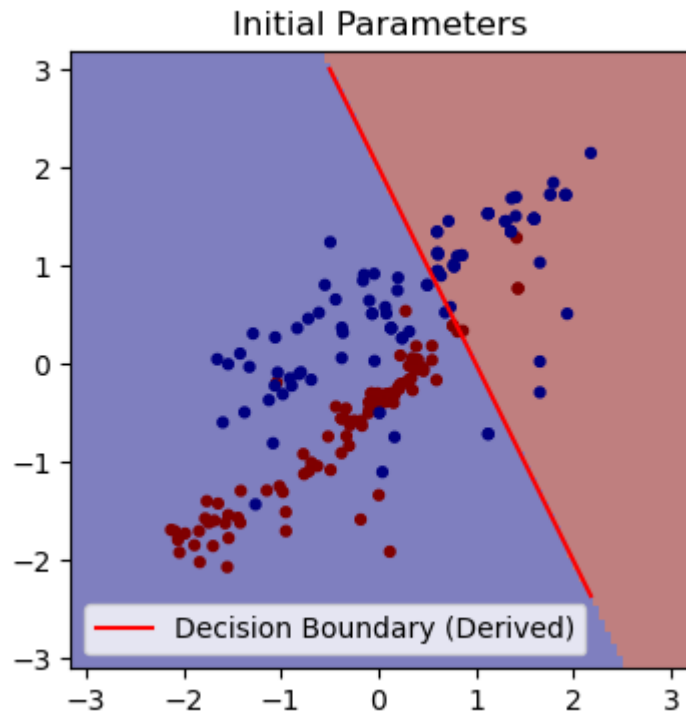
```
# Plot the line you derived in part 1 on the figure, in an appropriate range of values
```

```
axes.plot(x1_vals, x2_vals, 'r-', label='Decision Boundary (Derived)')
```

```
axes.legend()
```

```
### YOUR CODE ENDS HERE
```

```
axes.set_title(f'Initial Parameters');
```



Problem 1.2: Gradient Optimization

Start training your model using stochastic gradient descent, and looking at the classifier and decision boundary as you progress. For this part, we use `partial_fit`, a function that does a single epoch of stochastic gradient descent, and does not reset the internal state of the optimization loop (number of iterations, etc.), so that subsequent calls "pick up" right where the previous calls left off.

We'll initialize the model as before; then, train your model and visualize its current decision function (using `DecisionBoundaryDisplay`) after each of:

- 1 epoch
- 25 epochs
- 100 epochs
- 1000 epochs (final model)

Note that each call to `partial_fit` performs one epoch of SGD.

```
In [12]: np.random.seed(seed)

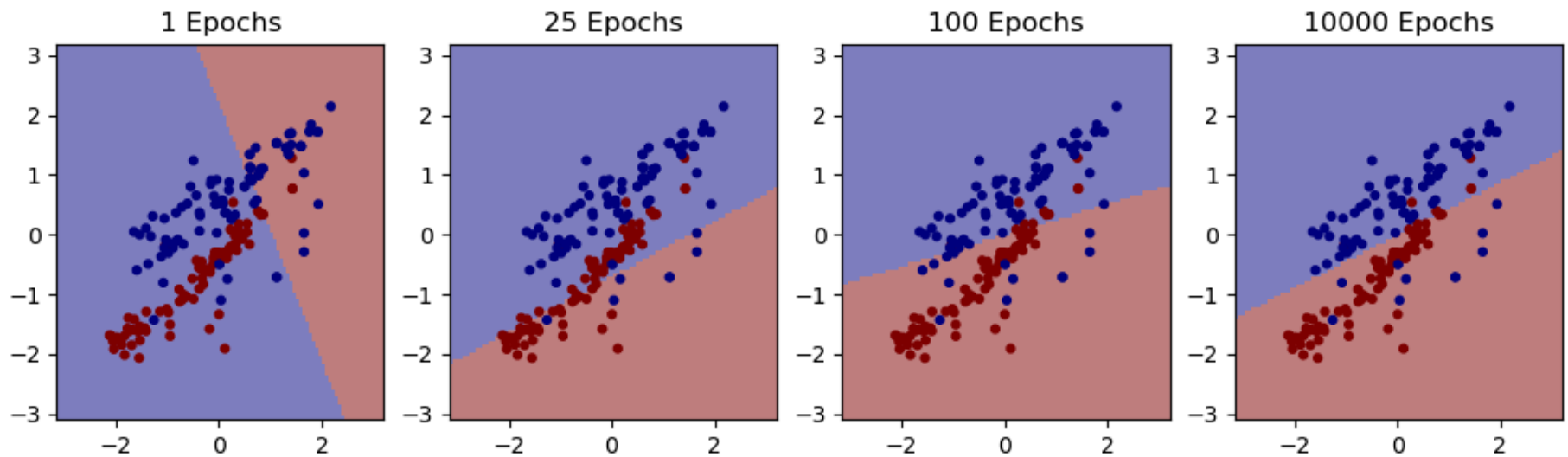
logreg = SGDClassifier(loss='log_loss', warm_start=True, learning_rate='adaptive', eta0 = 1e-3)
logreg.coef_ = np.array([[2.,1.]])
logreg.intercept_ = np.array([-2.]) #  $r(x) = 2x_1 + 1x_2 + (-2)$ 

plot_iters = [1,25,100,10000]
figure, axes = plt.subplots(1, 4, figsize=(12,3))

### YOUR CODE STARTS HERE
# Train the model using partial_fit and visualize decision boundaries at specified epochs
for i, (ax, epochs) in enumerate(zip(axes, plot_iters)):
    for j in range(epochs): # Train the model for the given number of epochs
        if j == 0: # On the first iteration, include the classes parameter
            logreg.partial_fit(X, y, classes=np.unique(y))
        else:
            logreg.partial_fit(X, y)

    # Plot the decision boundary after the specified number of epochs
    DecisionBoundaryDisplay.from_estimator(logreg, X, ax=ax, **plot_kwargs)
    ax.scatter(X[:, 0], X[:, 1], c=y, edgecolor=None, s=12, cmap='jet')
    ax.set_title(f'{epochs} Epochs')

### YOUR CODE ENDS HERE
```

Problem 1.3: Evaluation

Using your final model after training, display its learned linear coefficients and evaluate its (training) error rate.

Manually compute the linear response at the point $(x_1, x_2) = (-1, 0)$, and use the logistic function to evaluate the model's estimated probability that this point is in each class. (You can use the model's built-in function `predict_proba` to check your answer, if you like.)

```
In [14]: from sklearn.metrics import accuracy_score
from scipy.special import expit # Sigmoid function
# Display model parameters
print("Learned coefficients:", logreg.coef_)
print("Learned intercept:", logreg.intercept_)
# Evaluate model performance
y_pred = logreg.predict(X)
error_rate = 1 - accuracy_score(y, y_pred)
print("Training error rate:", error_rate)

# Manual computation: linear response and predicted probability at (-1,0)
theta_0 = logreg.intercept_
theta_1 = logreg.coef_[0][0]
```

```

theta_2 = logreg.coef_[0][1]
point = np.array([-1, 0])

# Compute the linear response
linear_response = theta_0 + theta_1 * point[0] + theta_2 * point[1]
print("Linear response at the point (-1, 0):", linear_response)

# Use the logistic function to compute the probability
probability = expit(linear_response)
print("Estimated probability of belonging to each class:", 1 - probability, probability)

```

```

Learned coefficients: [[ 1.56416414 -3.54684729]]
Learned intercept: [-0.02672695]
Training error rate: 0.09499999999999997
Linear response at the point (-1, 0): [-1.59089109]
Estimated probability of belonging to each class: [0.83074144] [0.16925856]

```

```

In [15]: logreg.predict_proba([[-1,0]]).round(2) # Evaluate on final model to check your answer

```

```

Out[15]: array([[0.83, 0.17]])

```

Problem 2: (Linear) Support Vector Machines

As we saw in lecture, a linear support vector machine optimizes the "margin" around the data. Our current data set is not linearly separable, so we will need to use a "Soft Margin" SVM. Soft-margin Linear SVMs are equivalent to a linear classifier trained using an L2-regularized hinge loss; so, we can implement the SVM using exactly the same `SGDClassifier` model, using the same learner (linear classifier) and an identical learning algorithm (stochastic gradient), but changing the loss function.

To make our model as "close" to a hard-margin SVM as possible, we set the L2 regularization to be very small. This also can make the optimization a bit slow, so we'll use a lot of iterations and turn off any early stopping criteria.

Problem 2.1: Training & Evaluation

Fit your model to the data, then print out its linear coefficients and the resulting (training) error rate:

```

In [18]: np.random.seed(seed)

```

```

learner = SGDClassifier(loss='hinge',          # hinge loss = primal linear SVM form
                        penalty='l2',alpha=1e-20, # small L2 regularization is "closest" to Hard SVM
                        learning_rate='adaptive',eta0=1e-3, # same optimization as before
                        tol=0.,max_iter=10000,n_iter_no_change=1000) # prevent any early stopping

### YOUR CODE STARTS HERE

# Train the model, display your parameters & evaluate its performance
# Train the model
learner.fit(X, y)

# Display the learned parameters
print("Learned coefficients (weights):", learner.coef_)
print("Learned intercept (bias):", learner.intercept_)

# Evaluate the model's performance on the training data
train_predictions = learner.predict(X)
training_accuracy = accuracy_score(y, train_predictions)
training_error_rate = 1 - training_accuracy

print("Training accuracy:", training_accuracy)
print("Training error rate:", training_error_rate)

### YOUR CODE ENDS HERE

```

```

Learned coefficients (weights): [[ 1.72112436 -2.63600577]]
Learned intercept (bias): [0.14569216]
Training accuracy: 0.925
Training error rate: 0.07499999999999996

```

Problem 2.2: Decision boundary & margins

Now, display the decision function learned by your linear SVM. In addition, on top of the decision boundary plot, display the SVM's margins, i.e.,

$$r(x) = \theta_0 + \theta_1 x_1 + \theta_2 x_2 = +1$$

and

$$r(x) = \theta_0 + \theta_1 x_1 + \theta_2 x_2 = -1$$

(Recall that the decision boundary, as you plotted earlier, is given by $r(x) = 0$.)

```
In [20]: figure, axes = plt.subplots(1, 1, figsize=(4,4))
DecisionBoundaryDisplay.from_estimator(learner, X, ax=axes, **plot_kwargs)
axes.scatter(X[:, 0], X[:, 1], c=y, edgecolor=None, s=12, cmap='jet')

### YOUR CODE STARTS HERE

# Draw (e.g. with dashed lines) the set of points that are on the +1 and -1 margins

# (Hint: this is almost the same as drawing the decision boundary earlier, except that
# you need to use your trained parameters, and solve  $r(x)=+1$  and  $r(x)=-1$  instead of  $r(x)=0$ .)
# Get the coefficients for the decision boundary equation
w = learner.coef_[0]
b = learner.intercept_[0]

# Plotting range
x_min, x_max = axes.get_xlim()
x_vals = np.linspace(x_min, x_max, 100)

# Compute the decision boundary:  $r(x) = 0 \rightarrow w_1x_1 + w_2x_2 + b = 0$ 
# Rearrange to solve for  $x_2$ :  $x_2 = -(w_1/w_2) * x_1 - b/w_2$ 
decision_boundary = -(w[0] / w[1]) * x_vals - b / w[1]

# Compute the +1 margin:  $r(x) = +1 \rightarrow w_1x_1 + w_2x_2 + b = 1$ 
margin_pos = -(w[0] / w[1]) * x_vals - (b - 1) / w[1]

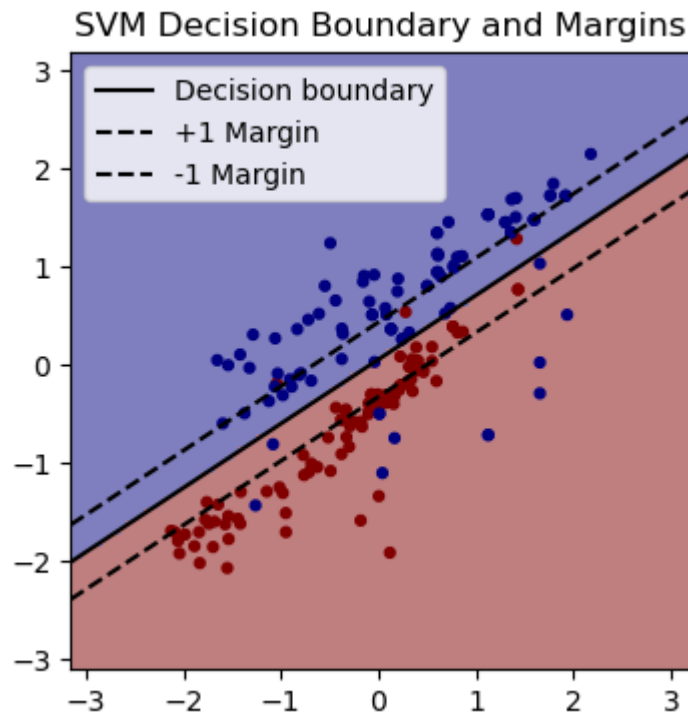
# Compute the -1 margin:  $r(x) = -1 \rightarrow w_1x_1 + w_2x_2 + b = -1$ 
margin_neg = -(w[0] / w[1]) * x_vals - (b + 1) / w[1]

# Plot the decision boundary
axes.plot(x_vals, decision_boundary, 'k-', label='Decision boundary')

# Plot the +1 and -1 margins as dashed lines
axes.plot(x_vals, margin_pos, 'k--', label='+1 Margin')
axes.plot(x_vals, margin_neg, 'k--', label='-1 Margin')

# Set title and legend
axes.set_title('SVM Decision Boundary and Margins')
```

```
axes.legend()  
  
plt.show()  
### YOUR CODE ENDS HERE
```



Problem 3: Feature Expansion

If we feel that our linear classifier is insufficiently flexible, one option is to provide it with more features. Just like in our linear regression models, additional features, such as polynomial features, make the resulting model more adaptable to the data.

In this problem, we will expand our features using `PolynomialFeatures`, and look at the resulting logistic regression model's decision function.

Note that, when creating new features, especially high-order polynomials, it is a good idea to scale the data after the feature transform. As in the HW2 solutions, the easiest way to expand the feature set and rescale the data is to use the `Pipeline` object in `sklearn`.

Adapt the code below to fit and display the decision function for degrees 1, 2, 5, and 20.

```
In [22]: from sklearn.pipeline import Pipeline
np.random.seed(seed)

degrees=[1,2,5,20]
figure,axes = plt.subplots(1,4,figsize=(12,3))

for i,d in enumerate(degrees):

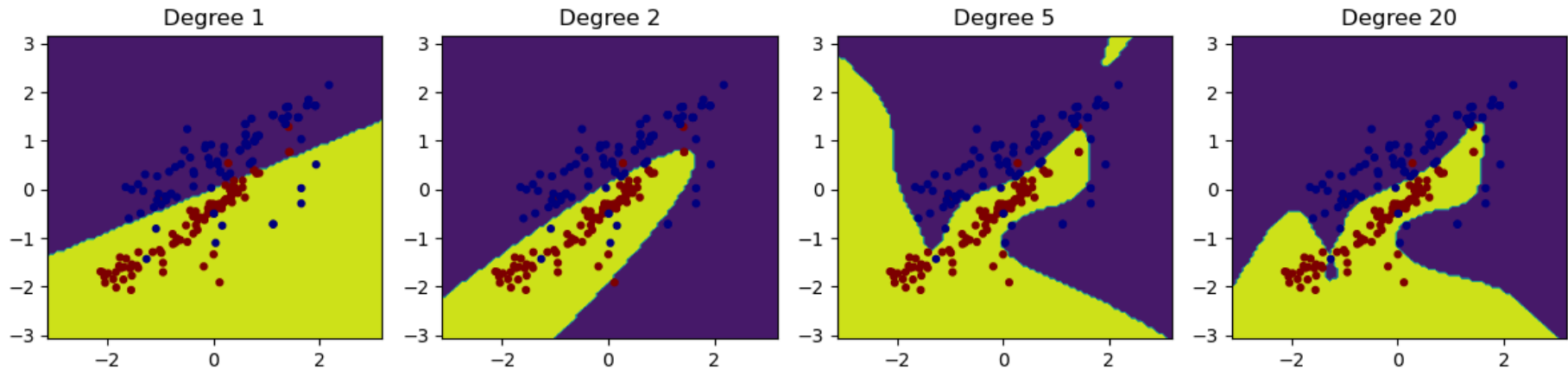
    # Each item in the pipeline is a pair, (name, transform); the end is (name, learner):
    learner = Pipeline( [('poly',PolynomialFeatures(degree=d)),
                        ('scale',StandardScaler()),
                        ('logreg',SGDClassifier(loss='log_loss',
                                                penalty='l2',alpha=1e-20,
                                                learning_rate='adaptive', eta0=1e-2,
                                                tol=0.,max_iter=100000,n_iter_no_change=1000))
                        ])

    ### YOUR CODE STARTS HERE

    # Fit the model
    learner.fit(X, y)

    # Display the resulting decision function and training data
    DecisionBoundaryDisplay.from_estimator(learner, X, ax=axes[i], response_method="predict")
    axes[i].scatter(X[:, 0], X[:, 1], c=y, edgecolor=None, s=12, cmap='jet')
    axes[i].set_title(f'Degree {d}')

plt.tight_layout()
plt.show()
    ### YOUR CODE ENDS HERE
```



Problem 3.2: Regularization

Our higher-order models are most likely overfitting (although we can't tell for sure, since we didn't save any data for validation). Let's re-learn the model using some regularization to see how it affects the resulting decision function.

Try increasing the L2 regularization to `1e-3`, `1e-1`, and `10` and display the resulting decision functions. Discuss how these compare to each other, and to the (nearly) unregularized version in the previous question.

```
In [24]: from sklearn.pipeline import Pipeline
np.random.seed(seed)

d = 20
alphas = [1e-3, 1e-1, 10.]
figure, axes = plt.subplots(1, 3, figsize=(10, 3))

for i, alpha in enumerate(alphas):
    # Each item in the pipeline is a pair, (name, transform); the end is (name, learner):
    learner = Pipeline([('poly', PolynomialFeatures(degree=d)),
                        ('scale', StandardScaler()),
                        ('logreg', SGDClassifier(loss='log_loss',
                                                penalty='l2', alpha=alpha,
                                                learning_rate='adaptive', eta0=1e-2,
                                                tol=0., max_iter=100000, n_iter_no_change=1000))])
```

```

    ])

    ### YOUR CODE STARTS HERE

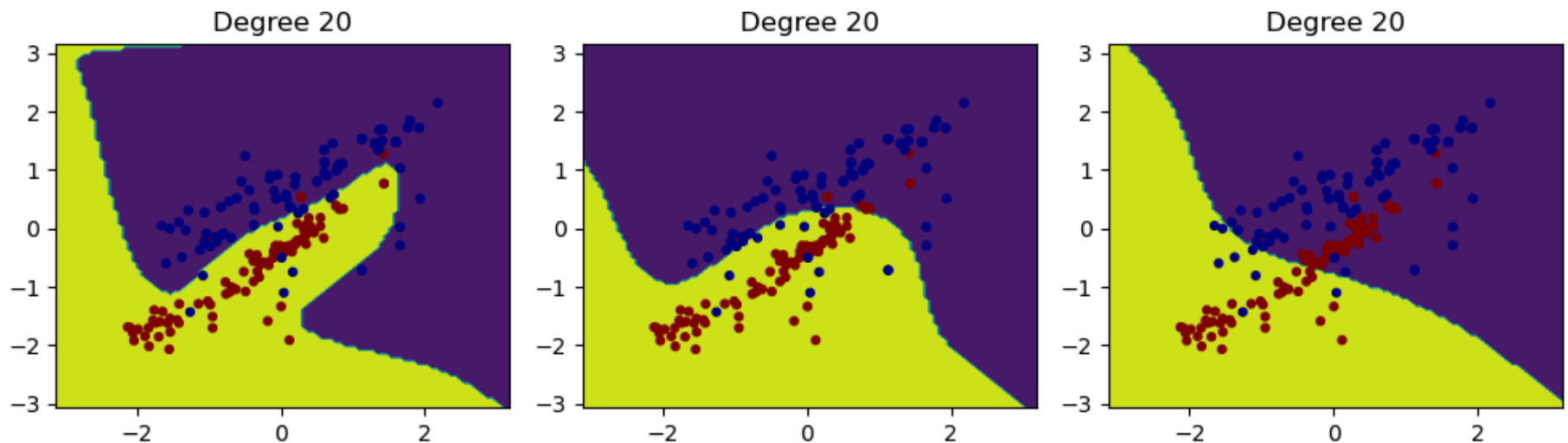
    # Fit the model
    learner.fit(X, y)

    # Display the resulting decision function and training data
    DecisionBoundaryDisplay.from_estimator(learner, X, ax=axes[i], response_method="predict")
    axes[i].scatter(X[:, 0], X[:, 1], c=y, edgecolor=None, s=12, cmap='jet')
    axes[i].set_title(f'Degree {d}')

plt.tight_layout()
plt.show()

    ### YOUR CODE ENDS HERE

```



Discuss

Lower regularization ($\alpha=1e-3$) may result in slight overfitting and significant fluctuations in decision boundaries. Moderate regularization ($\alpha=1e-1$) smooths decision boundaries and reduces overfitting of the model to the data. Strong regularization ($\alpha=10$) may lead to underfitting, with boundaries becoming too simple to capture patterns in the data well. Compared to

unnormalized versions, adding regularization helps smooth decision boundaries and avoid overfitting of the model to noise. As regularization increases, the model becomes more linear, reducing the influence of higher-order polynomials.

Problem 4: Logistic Regression on MNIST

Finally, let us now build a linear classifier (specifically, a logistic regression model) on a higher-dimensional, multi-class problem: the MNIST data set.

The MNIST dataset is an image dataset consisting of 70,000 hand-written digits (from 0 to 9), each of which is a 28x28 grayscale image. For each image, we also have a label, corresponding to which digit is written.

Problem 4.0: Setting up the Data

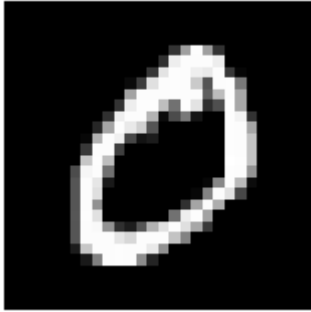
First, we'll load our dataset, split it into a training set and a testing set, and do some basic pre-processing. Here you are given code that does this for you, and you only need to run it.

```
In [28]: # Load the features and labels for the MNIST dataset
# This might take a minute to download the images.
X, y = fetch_openml('mnist_784', as_frame=False, return_X_y=True)

# Convert labels to integer data type
y = y.astype(int)
```

Each data point in the MNIST dataset is 768-dimensional, with each feature corresponding to a pixel intensity of a 28×28 scan of a digit. To visualize a data point, we can re-shape the feature vector into the shape of the image, and then display it using `imshow` :

```
In [30]: plt.figure(figsize=(2,2))
plt.imshow( X[1,:].reshape(28,28) , cmap='gray');
plt.axis('off');
```



As before, we will normalize the data before learning using the scikit-learn class `StandardScaler` to standardize both the training and testing features. Notice that we **only** fit the `StandardScaler` on the training data, and *not* the testing data.

```
In [140... X_tr, X_te, y_tr, y_te = train_test_split(X, y, test_size=0.1, random_state=seed, shuffle=True)

scaler = StandardScaler()
scaler.fit(X_tr)
X_tr = scaler.transform(X_tr)      # We can forget about the original values & work
X_te = scaler.transform(X_te)      # just with the transformed values from here
                                   # (This does make it harder to visualize a data point, though)
```

Problem 4.1: Initial Training (10 points)

For this part of the problem, you will train on **just** the first 10000 training data points, and compute the training and test error rates.

- Be sure to set the random seed with `random_state=seed` for consistency.
- Other than the random seed, just use the default values of the learner for this part.
- Here, the training error rate is defined on the first 10k data points (i.e., the points that were used for training the model)
- The test error rate is defined on the full test data from your split.

```
In [144... m_tr = 10000

X_tr_subset = X_tr[:m_tr, :]
y_tr_subset = y_tr[:m_tr]

# Construct a logistic regression classifier (random_state = seed)
```

```

learner_mnist = LogisticRegression(random_state=seed)

### YOUR CODE STARTS HERE ###

# Fit your model to the (small subset of the) training data
learner_mnist.fit(X_tr_subset, y_tr_subset)
# Compute the training error (on the small training subset) and testing error (on the test data)
y_train_pred = learner_mnist.predict(X_tr_subset)
y_test_pred = learner_mnist.predict(X_te)

training_error = 1 - accuracy_score(y_tr_subset, y_train_pred)
testing_error = 1 - accuracy_score(y_te, y_test_pred)

# Display the errors
print(f'Training Error: {training_error:.4f}')
print(f'Testing Error: {testing_error:.4f}')
### YOUR CODE ENDS HERE ###

```

Training Error: 0.0014

Testing Error: 0.1204

Your model should learn a set of linear coefficients for each of the 10 classes:

```

In [36]: print(f'Coefficients shape: {learner_mnist.coef_.shape}') # should be 10 x 768
         print(f'Intercepts shape: {learner_mnist.intercept_.shape}') # should be 10

```

Coefficients shape: (10, 784)

Intercepts shape: (10,)

Problem 4.2: Regularization (10 points)

Suspecting that we are overfitting to our limited data set, we decide to try to use regularization. (This should reduce our model's variance, and thus its tendency to overfit.) Try re-training your logistic regression model at various levels of regularization.

The `LogisticRegression` class in `sklearn` takes an "inverse regularization" parameter, `C` (effectively the same as the value R we saw in soft-margin Support Vector Machines). Re-train your model with values of $C \in \{.0001, .001, .01, .1, 1.0, 10.\}$ and compute the training and test error rates of each setting. Plot the training and test error rates together as a function of C (plot using `semilogx` for it to look nice) and state what value of C you would select and why.

```
In [38]: m_tr = 10000
C_vals = [.0001,.001,.01,.1,1.,10.];

### YOUR CODE STARTS HERE ###
train_errors = []
test_errors = []
for C in C_vals:

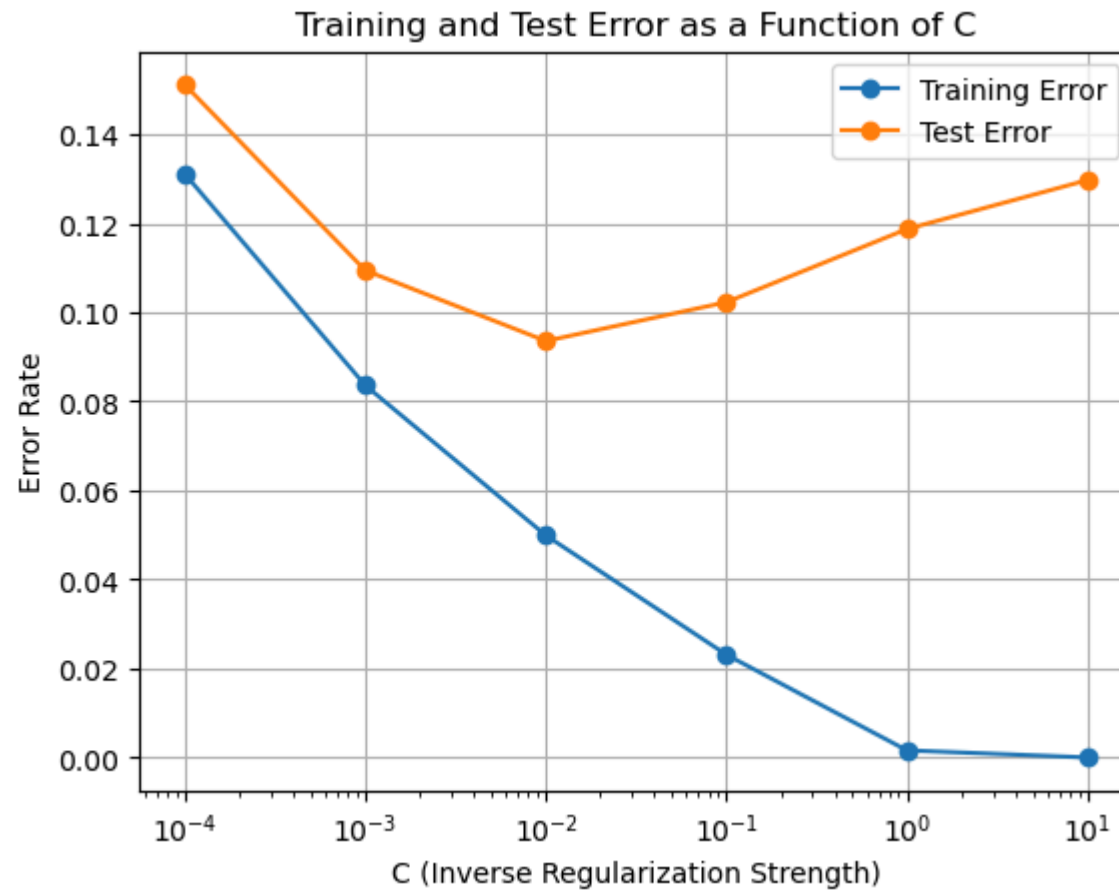
    learner_mnist = LogisticRegression(C=C, random_state=seed, max_iter=10000)
    learner_mnist.fit(X_tr_subset, y_tr_subset)

    # Compute the training error
    y_train_pred = learner_mnist.predict(X_tr_subset)
    train_error = 1 - accuracy_score(y_tr_subset, y_train_pred)
    train_errors.append(train_error)

    # Compute the test error
    y_test_pred = learner_mnist.predict(X_te)
    test_error = 1 - accuracy_score(y_te, y_test_pred)
    test_errors.append(test_error)

# Plot the training and testing error rates as a function of C
plt.figure()
plt.semilogx(C_vals, train_errors, label='Training Error', marker='o')
plt.semilogx(C_vals, test_errors, label='Test Error', marker='o')
plt.xlabel('C (Inverse Regularization Strength)')
plt.ylabel('Error Rate')
plt.title('Training and Test Error as a Function of C')
plt.legend()
plt.grid(True)
plt.show()

### YOUR CODE ENDS HERE ###
```



I will choose $C=10e-2$ since it has the lowest test error. When $C < 10e-2$ the model is underfitting while $C > 10e-2$ the model becomes overfitting.

Problem 4.3: Interpreting the weights (5 points)

Now that we have a model that we believe might perform well, let's try to understand what properties of the data it is using to make its predictions. Since our model is just using a linear combination of the input pixels, we can display the coefficient (slope) associated with each pixel, to see whether that pixel's being bright (high value) is positively associated with a given class, or is negatively associated with that class.

First, re-train your model using your selected value of C .

```
In [98]: ### YOUR CODE START HERE ###

# Re-train your model with your selected value of C
learner_mnist = LogisticRegression(C=.01, random_state=seed, max_iter=10000)
learner_mnist.fit(X_tr_subset, y_tr_subset)
### YOUR CODE ENDS HERE ###
```

```
Out[98]: 

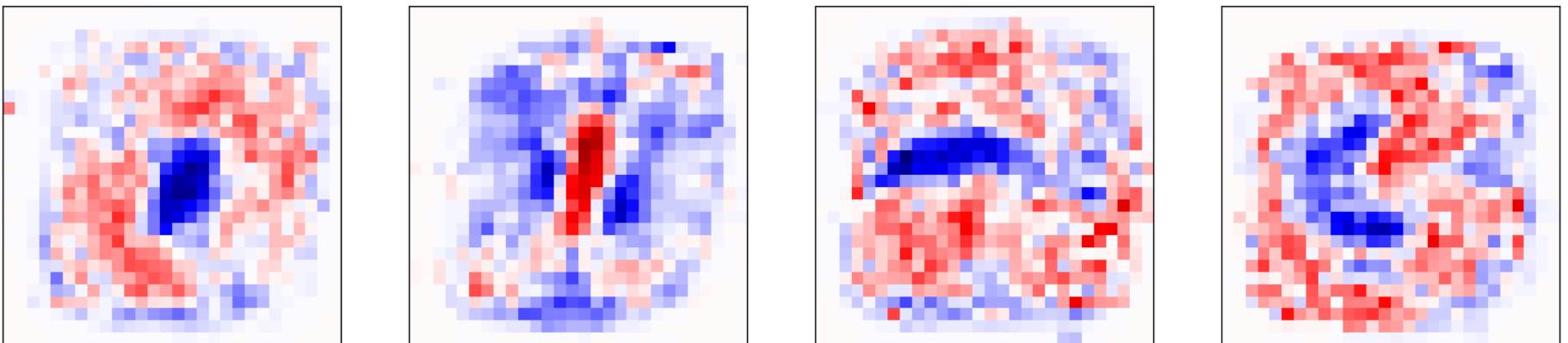
LogisticRegression
LogisticRegression(C=0.01, max_iter=10000, random_state=1234)


```

Run the provided code to display the coefficients of the first four classes' linear responses, re-shaped to the same size as the input image. (Here, red is positive, blue is negative, and white is zero.) Do the responses make sense? Discuss.

```
In [100... fig, ax = plt.subplots(1,4, figsize=(18,8))

mu = learner_mnist.coef_.mean(0).reshape(28,28)
for i in range(4):
    ax[i].imshow(learner_mnist.coef_[i,:].reshape(28,28)-mu,cmap='seismic',vmin=-.25,vmax=.25);
    ax[i].set_xticks([]); ax[i].set_yticks([]);
```



DISCUSS

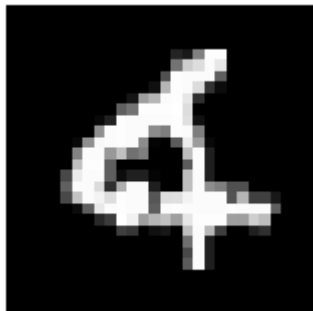
The responses can indeed be meaningful as they provide insight into which features (pixels) are more important for the classification of different categories. A positive coefficient (in red) means that an increase in the corresponding pixel value increases the likelihood of the input being classified into that category, while a negative coefficient (in blue) implies a decrease in likelihood. Thus the number 0,1,2,3 can be recognized in red in these four figures, respectively.

Problem 4.4

The multilogistic classifier uses the negative log-likelihood loss, just like the logistic classifier, but produces a predicted probability for each class based on that class's linear response.

In this problem, we'll consider a particular (somewhat ambiguous) data point:

```
In [103... idx = 14290
plt.figure(figsize=(2,2))
plt.imshow( X[idx,:].reshape(28,28) , cmap='gray');
plt.axis('off');
```



(a) Using your model parameters, **manually** compute the linear response values for each of the 10 classes on this data point. (You can do this easily using matrix multiplication and addition of arrays.)

```
In [106... # Extract the input data point (reshaped to 1D if needed)
X_input = X[idx, :].reshape(1, -1) # Reshape to 2D for matrix multiplication

# Compute the linear response for each class
linear_responses = np.dot(X_input, learner_mnist.coef_.T) + learner_mnist.intercept_
```

```
# Display the responses
print("Linear response values for each class:")
print(linear_responses)
```

Linear response values for each class:

```
[[ 222.84513787 -572.9254554  -22.51348394 -372.61970792  216.12551476
 -127.49720846  447.80608611 -418.72960749  231.62273054  395.88599395]]
```

(b) Use the multi-logit or `softmax` transformation to convert these responses into estimated class probabilities.

```
In [118... from scipy.special import softmax

# Suppose linear_responses contains the response values for each of the 10 classes
# linear_responses = np.array([...]) # Replace with your actual linear response values

# Use the softmax function to convert the linear responses to probabilities
softmax_probabilities = softmax(linear_responses)

# Display the estimated probabilities for each class
print("Estimated class probabilities using the softmax function:")
print(softmax_probabilities)
```

Estimated class probabilities using the softmax function:

```
[[1.99848796e-098 0.00000000e+000 5.53102977e-205 0.00000000e+000
 2.41216105e-101 1.40913873e-250 1.00000000e+000 0.00000000e+000
 1.29646739e-094 2.82742098e-023]]
```

(c) Do these probabilities make sense, given the observation? Discuss briefly.

These probabilities make sense. The estimated class probabilities obtained indicate a very strong prediction for class 6 (which often corresponds to the digit "4" in the MNIST dataset), with a probability very close to 1, while the probabilities for all other classes are extremely small or effectively zero.

Note: To check your answer, you can compare to the values given by the learner's built-in `predict_proba()` function:

```
In [128... learner_mnist.predict_proba(X[idx:idx+1,:]).round(2)
```

```
Out[128... array([[0., 0., 0., 0., 0., 0., 1., 0., 0., 0.]])
```


Problem 4.5: Learning Curves (10 points)

Another way to reduce overfitting is to increase the amount of data used for training the model (if possible). Build a logistic regression model, but with no regularization

- Train a logistic regression classifier (with the default settings in sklearn) using the first `m_tr` feature vectors in `X_tr`, where `m_tr = [100, 1000, 5000, 10000, 20000, 50000, 63000]`. You should use the `LogisticRegression` class from scikit-learn in your implementation. **Make sure to use the argument `random_state=seed` for reproducibility.**
- Create a plot of the training error and testing error for your logistic regression model as a function of the number of training data points. Be sure to include an x-label, y-label, and legend in your plot. Use a log-scale on the x-axis. Give a short (one or two sentences) description of what you see in your plot.
- Add a comment with your thoughts after the plot: although we ran out of data at 63k examples, can you tell how much additional data could help, with this model?

```
In [146... train_sizes = [100, 1000, 5000, 10000, 20000, 50000, 63000]
train_errors = []
test_errors = []

C = np.inf      # No regularization!

### YOUR CODE STARTS HERE ###

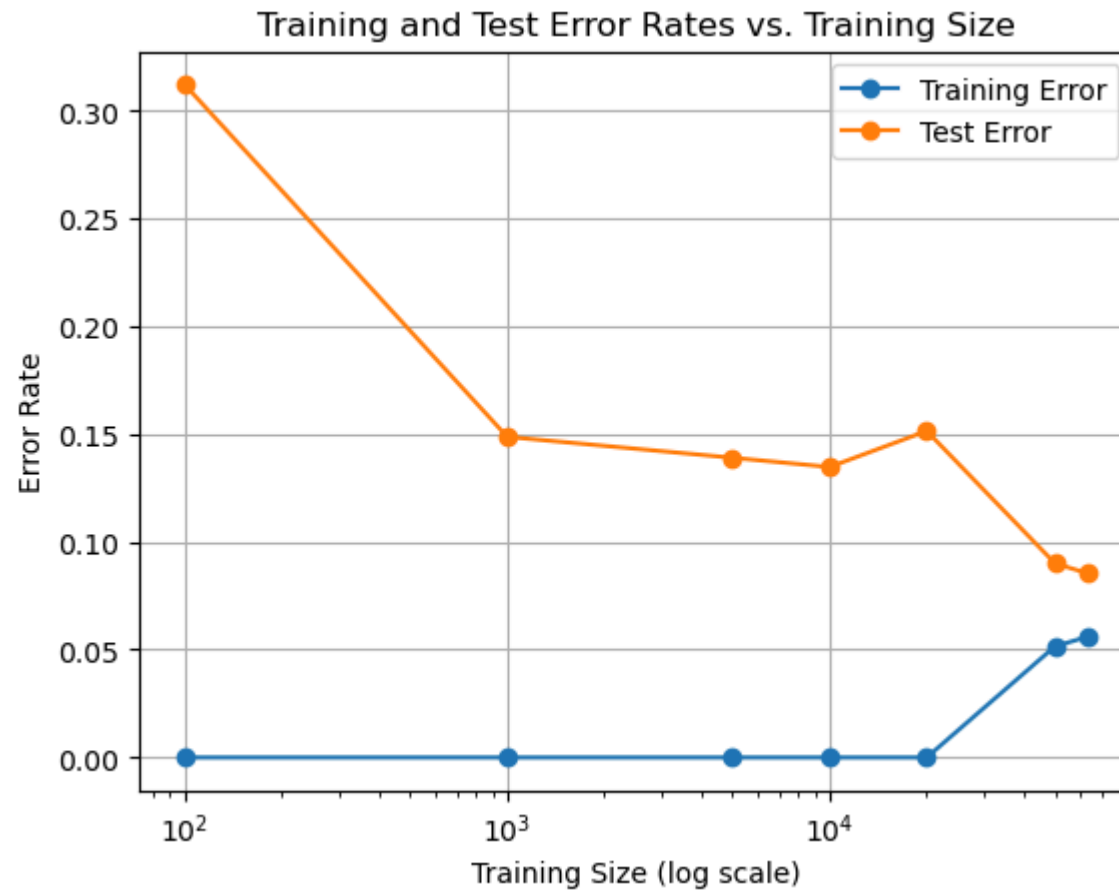
# Train a logistic regression model with each data size m and C=infinity
for m_tr in train_sizes:
    X_tr_subset = X_tr[:m_tr, :]
    y_tr_subset = y_tr[:m_tr]
    learner_mnist = LogisticRegression(C=np.inf, random_state=seed, max_iter=10000)
    learner_mnist.fit(X_tr_subset, y_tr_subset)

    # Compute the training and test error rates
    # Compute the training error
    y_train_pred = learner_mnist.predict(X_tr_subset)
    train_error = 1 - accuracy_score(y_tr_subset, y_train_pred)
    train_errors.append(train_error)
```

```
# Compute the test error
y_test_pred = learner_mnist.predict(X_te)
test_error = 1 - accuracy_score(y_te, y_test_pred)
test_errors.append(test_error)

# Plot the resulting performance as a function of m
plt.figure()
plt.plot(train_sizes, train_errors, label='Training Error', marker='o')
plt.plot(train_sizes, test_errors, label='Test Error', marker='o')
plt.xscale('log')
plt.xlabel('Training Size (log scale)')
plt.ylabel('Error Rate')
plt.title('Training and Test Error Rates vs. Training Size')
plt.legend()
plt.grid()
plt.show()

### YOUR CODE ENDS HERE ###
```



```
In [153... print("train error:", train_errors)
print("test error:", test_errors)
```

```
train error: [0.0, 0.0, 0.0, 0.0, 0.0, 0.05154000000000003, 0.055968253968254]
```

```
test error: [0.31200000000000006, 0.1487142857142857, 0.139, 0.13471428571428568, 0.15142857142857147, 0.08999999999999997, 0.08542857142857141]
```

COMMENT / DISCUSS

Comment: As the data size gradually increases, the degree of overfitting in the model will decrease. When the data size is greater than 10E4, the training error will be greater than 0. As the data size gradually increases, the testing error will gradually decrease, indicating

that the model's classification ability for new data is gradually improving.

Discuss: When additional training data is added, the test error continues to decrease, indicating that the model could potentially benefit from more data to further refine its understanding and generalization to unseen examples. If additional training examples were available, particularly those that capture more diverse patterns or edge cases within the dataset, it might help improve the model's performance, especially in reducing the test error.



Statement of Collaboration (5 points)

It is **mandatory** to include a Statement of Collaboration in each submission, with respect to the guidelines below. Include the names of everyone involved in the discussions (especially in-person ones), and what was discussed.

All students are required to follow the academic honesty guidelines posted on the course website. For programming assignments, in particular, I encourage the students to organize (perhaps using EdD) to discuss the task descriptions, requirements, bugs in my code, and the relevant technical content before they start working on it. However, you should not discuss the specific solutions, and, as a guiding principle, you are not allowed to take anything written or drawn away from these discussions (i.e. no photographs of the blackboard, written notes, referring to EdD, etc.). Especially after you have started working on the assignment, try to restrict the discussion to EdD as much as possible, so that there is no doubt as to the extent of your collaboration.

This assignment was completed independently by me.