

Assessment 3

SCHOOL OF ADVANCED TECHNOLOGY

INT10: Advanced Pattern Recognition

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1 Objectives

This assessment aims at evaluating students' ability to exploit the advanced pattern recognition knowledge, which is accumulated during lectures, and after-class study, to analyze, design, implement, develop, test and document the pattern classification methods with Neural Networks (MLP), Discriminant Function (MQDF), and Kernel Methods (Support Vector Machine) typically on the image data.

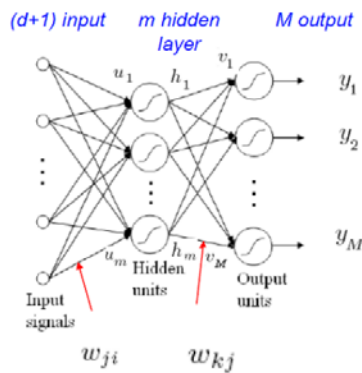
2 Materials

Pycharm

3 Methods, result and discussion

1 Implement Multilayer Perceptron (MLP) on MNIST

An artificial neural network is composed of many artificial neurons that are linked together according to a specific network architecture. The model transforms the inputs into meaningful outputs by iterating a combination of linear and nonlinear functions.



$$W = \arg \min_W \frac{1}{2} \sum_{n=1}^N \sum_{k=1}^M [y_k(x^n) - t_k^n]^2$$

$$y_k(x) = g[v_k(x)]$$

$$v_k(x) = \sum_{j=1}^m w_{kj} h_j + w_{k0}$$

$$h_j(x) = g[u_j(x)]$$

$$u_j(x) = \sum_{i=1}^d w_{ji} x_i + w_{j0}$$

Design an MLP model to achieve the classification task on MNIST. Show the training process and classification results.

Many factors will affect the performance, such as, the number of nodes in hidden layers, the

selection of the activation functions, the normalization method of the samples, the loss function and its regularization term, the learning rate and the stopping policy of training, etc.

- Load data set

train_set, valid_set and test_set are training set, validation set and test set respectively, organized in the form of tuple (x, y), x represents sample, y represents sample label

```
1. def load_data():
2.     f = gzip.open("mnist.pkl.gz", "rb")
3.     train_set, valid_set, test_set = pickle.load(f, encoding="latin1")
4.     f.close()
5.     return [train_set, valid_set, test_set]
```

```
1. datasets = load_data()
```

Result:

```
[(array([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        ...,
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.], dtype=float32), array([5, 0, 4, ..., 8, 4, 8], dtype=int64)), (array([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.], dtype=float32), array([3, 0, 6, ..., 5, 6, 8], dtype=int64)), (array([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.], dtype=float32), array([7, 2, 1, ..., 4, 5, 6], dtype=int64))]
```

- Define the loss function

For multi-class problems, the negative log-likelihood function is used as the loss function. $p_{y_given_x}$ is the prediction of the model. It is an $m \times 10$ two-dimensional array. Each row represents a sample, and each column represents the model prediction sample is 0~9 Probability. $Np.arange(y.shape[0])$ is a one-dimensional array, containing $[0, 1, \dots, m-1]$, y is the true label corresponding to the sample, is a one-dimensional array, $p_{y_given_x}[[0, 1, \dots, m-1][y[0], \dots, y[m-1]]]$ is equivalent to $p_{y_given_x}[0, y[0]]$, $p_{y_given_x}[1, y[1]]$, ..., $p_{y_given_x}[m-1, y[m-1]]$, that is, read

the predicted value of the model corresponding to the true category of each sample, we want this value bigger.

$$C = -\log y_r$$

$$y_r = \frac{e^{z_r}}{\sum_k e^{z_k}}$$

Derivation:

1、 $i=r$

$$\frac{\partial C}{\partial z_r} = \frac{\partial C}{\partial y_r} \frac{\partial y_r}{\partial z_r} = -\frac{1}{y_r} \left(\frac{e^{z_r}}{\sum_k e^{z_k}} - e^{z_r} \frac{e^{z_r}}{(\sum_k e^{z_k})^2} \right) = -\frac{1}{y_r} (y_r - y_r^2) = y_r - 1$$

2、 $i \neq r$

$$\frac{\partial C}{\partial z_i} = \frac{\partial C}{\partial y_r} \frac{\partial y_r}{\partial z_i} = -\frac{1}{y_r} \left(-e^{z_r} \frac{e^{z_i}}{(\sum_k e^{z_k})^2} \right) = -\frac{1}{y_r} (-y_r^2) = y_r - 0$$

Unify, have:

$$\frac{\partial C}{\partial z_i} = y_i - \hat{y}_i$$

Code:

```
1. # Define loss function
2. def loss_function(p_y_given_x, y):
3.     return -np.mean(np.log(p_y_given_x)[np.arange(y.shape[0]), y])
```

Result:

The screenshot shows a Jupyter Notebook with a file explorer on the left containing files like 'mnist_testdata.hdf5', 'mnist_traindata.hdf5', 'process_dataset.py', 't10k-images-idx3-ubyte.gz', 't10k-labels-idx1-ubyte.gz', 'test_softmaxRegression_mlp', and 'train-images-idx3-ubyte.gz'. The main area shows a code cell with the following Python code:

```
15 # Define loss function
16 def loss_function(p_y_given_x, y):
17     print("loss:")
18     print(-np.mean(np.log(p_y_given_x)[np.arange(y.shape[0]), y]))
19     return -np.mean(np.log(p_y_given_x)[np.arange(y.shape[0]), y])
```

Below the code cell, the output of the notebook is displayed in a terminal-like window:

```
load_data()
mlp_mnist_1hidden
I:\Anaconda\python.exe C:/Users/123/Desktop/MLP/mlp_mnist_1hidden.py
loss:
5.662540242887119
loss:
23.8429716693345
```

- Count the number of error samples

First, according to the predicted probability of the network, choose the one with the highest probability as the network predicted label

Code:

```
1. # Get the predicted label of each sample
```

```

2. def pred_num_label(p_y_given_x):
3.     y_pred = []
4.     for i in range(p_y_given_x.shape[0]):
5.         max_index = np.argmax(p_y_given_x[i] == np.max(p_y_given_x[i]))
6.         y_pred.append(max_index[0,0])
7.     # print(y_pred[:50])
8.     return np.array(y_pred)

```

Result:

```

19 # Get the predicted label of each sample
20 def pred_num_label(p_y_given_x):
21     y_pred = []
22     for i in range(p_y_given_x.shape[0]):
23         max_index = np.argmax(p_y_given_x[i] == np.max(p_y_given_x[i]))
24         y_pred.append(max_index[0,0])
25     print("predicted label of each sample:")

```

I:\Anaconda\python.exe C:/Users/123/Desktop/MLP/mlp_mnist_1hidden.py
predicted label of each sample:
[1 4 1 ... 1 1 1]

- Construct the sample real label matrix

In the original data set, the label of each sample corresponds to a scalar, which directly indicates which number the sample is. But in calculating the gradient, y needs to be expanded into a vector with the same dimension as the model output.

Code:

```

1. # Construct the sample real label matrix
2. def truth_label_matrix(y):
3.     matrix = np.zeros((len(y), 10))
4.     for i in np.arange(len(y)):
5.         matrix[i, y[i]] = 1
6.     return matrix

```

Result:

```

35 for i in np.arange(len(y)):
36     matrix[i, y[i]] = 1
37     print("sample real label matrix:")
38     truth_label_matrix()

```

sample real label matrix:
[[0. 0. 0. ... 0. 0. 1.]
 [0. 0. 0. ... 0. 0. 1.]
 [0. 0. 0. ... 0. 1. 0.]
 ...
 [0. 0. 0. ... 0. 0. 0.]

- sigmoid function and its derivative

Code:

```
1. # Sigmoid
2. def sigmoid(x):
3.     return 1.0 / (1 + np.exp(-x))
4. # Sigmoid derivative
5. def sigmoid_derivative(x):
6.     return sigmoid(x) * (1 - sigmoid(x))
```

Result:

The first screenshot shows the implementation of the sigmoid function in a Jupyter Notebook. The code defines a function `sigmoid(x)` that returns $1.0 / (1 + \exp(-x))$. The output of the function is displayed as a 3x3 matrix of values, ranging from approximately 0.03 to 0.87.

```
def sigmoid(x):
    print("Sigmoid")
    print(1.0 / (1 + np.exp(-x)))
    return 1.0 / (1 + np.exp(-x))
```

The second screenshot shows the implementation of the sigmoid derivative function in a Jupyter Notebook. The code defines a function `sigmoid_derivative(x)` that returns `sigmoid(x) * (1 - sigmoid(x))`. The output of the function is displayed as a 3x3 matrix of values, ranging from approximately 0.05 to 0.25.

```
# Sigmoid derivative
def sigmoid_derivative(x):
    print("Sigmoid derivative")
    print(sigmoid(x) * (1 - sigmoid(x)))
    return sigmoid(x) * (1 - sigmoid(x))
```

- Initialize weight w

Determine the number of hidden layers, the number of hidden layer neurons, and initialize the connection weight, the neural network model is determined

Code:

```
50
57 def hidden_layer(n_input, n_output, hidden_unit, rand):
58     temp = []
59     try:
60         if len(hidden_unit) == 1:
61             w_0 = np.asarray(rand.uniform(low=-np.sqrt(6. / (n_input + hidden_unit[0])),
62                                           high=np.sqrt(6. / (n_input + hidden_unit[0])),
63                                           size=(n_input, hidden_unit[0])))
64             w_0 *= 4
65             temp.append(w_0)
66             w_1 = np.asarray(rand.uniform(low=-np.sqrt(6. / (hidden_unit[0] + n_output)),
67                                           high=np.sqrt(6. / (hidden_unit[0] + n_output)),
68                                           size=(hidden_unit[0], n_output)))
69             w_1 *= 4
70             temp.append(w_1)
71         else:
72             w_0 = 4 * np.asarray(rand.uniform(low=-np.sqrt(6. / (n_input + hidden_unit[0])),
73                                               high=np.sqrt(6. / (n_input + hidden_unit[0])),
74                                               size=(n_input, hidden_unit[0])))
75             temp.append(w_0)
76             for i in range(len(hidden_unit) - 1):
77                 temp.append(4 * np.asarray(rand.uniform(low=-np.sqrt(6. / (hidden_unit[i] + hidden_unit[i + 1])),
78                                                         high=np.sqrt(6. / (hidden_unit[i] + hidden_unit[i + 1])),
79                                                         size=(hidden_unit[i], hidden_unit[i + 1]))))
80             temp.append(4 * np.asarray(rand.uniform(low=-np.sqrt(6. / (hidden_unit[-1] + n_output)),
81                                                     high=np.sqrt(6. / (hidden_unit[-1] + n_output)),
82                                                     size=(hidden_unit[-1], n_output))))
83         except:
84             print("Warning: No input")
85             pass
86     return temp
```

Result:

```
mnist.pkl          86         print("hidden layer")
mnist.pkl.gz       87         print(temp)
mnist_testdata.hdf5 88         return temp
mnist_traindata.hdf5 hidden_layer()

mlp_mnist_1hidden x
I:\Anaconda\python.exe C:/Users/123/Desktop/MLP/mlp_mnist_1hidden.py
hidden layer
[array([[ 0.10300014,  0.24935554,  0.14300967, ...,  0.26158428,
         0.22310819, -0.12770949],
       [ 0.11495814, -0.24748337,  0.01365309, ..., -0.12329714,
        -0.19550261, -0.22514715],
       [-0.13076902,  0.1730739 ,  0.11202465, ..., -0.0494331 ,
         0.06451429, -0.03003983],
```

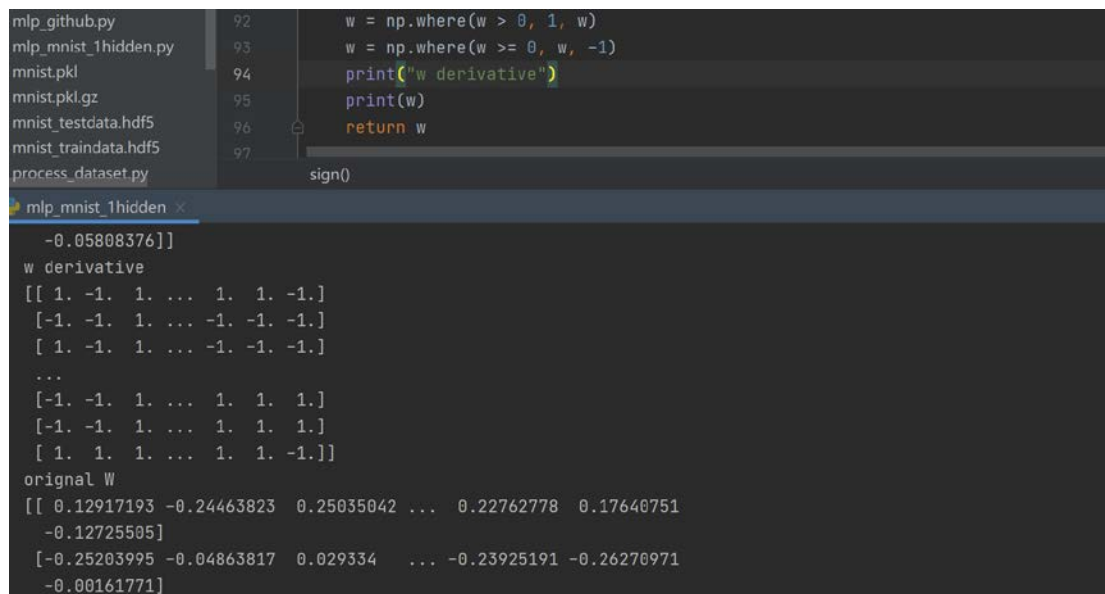
- Calculate the derivative of $|w|$

Calculate the derivative of $|w|$, needed when using L1 regularization

Code:

```
1. # Calculate the derivative of |w|, needed when using L1 regularization
2. def sign(w):
3.     w = np.where(w > 0, 1, w)
4.     w = np.where(w >= 0, w, -1)
5.     return w
```

Result:



```
mlp_github.py 92 w = np.where(w > 0, 1, w)
mlp_mnist_1hidden.py 93 w = np.where(w >= 0, w, -1)
mnist.pkl 94 print("w derivative")
mnist.pkl.gz 95 print(w)
mnist_testdata.hdf5 96 return w
mnist_traindata.hdf5 97
process_dataset.py sign()

mlp_mnist_1hidden x
-0.05808376]]
w derivative
[[ 1. -1.  1. ...  1.  1. -1.]
 [-1. -1.  1. ... -1. -1. -1.]
 [ 1. -1.  1. ... -1. -1. -1.]
 ...
 [-1. -1.  1. ...  1.  1.  1.]
 [-1. -1.  1. ...  1.  1.  1.]
 [ 1.  1.  1. ...  1.  1. -1.]]
original W
[[ 0.12917193 -0.24463823  0.25035042 ...  0.22762778  0.17640751
 -0.12725505]
 [-0.25203995 -0.04863817  0.029334 ... -0.23925191 -0.26270971
 -0.00161771]]
```

- Forward spread

Forward propagation, get model prediction value

Code:

```
1. # Forward spread
2. def forward(x, W):
3.     layer_input = [x] # input
4.     layer_output = [x] # output
5.
6.     # hidden layer
7.     for i in range(len(W) - 1):
8.         layer_input.append(np.dot(x, W[i]))
9.         temp_ = sigmoid(np.dot(x, W[i]))
10.        layer_output.append(temp_)
11.    # output layer
12.    layer_input.append(np.dot(temp_, W[-1]))
```

```

13.         layer_output.append((np.transpose(np.exp(np.dot(temp_, W[-
14.         1])))) / np.sum(np.exp(np.dot(temp_, W[-1])), axis=1)).T)
15.         return (layer_output, layer_input)

```

Result:

```

layer input
[array([[0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       ...,
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.]], dtype=float32), array([[ 2.3306077 ,  0.81466284, -2.14460153, ...,  0.20070811,
        -0.75865222,  1.17101196],
       [ 0.84820775,  1.78792717,  0.13623135, ...,  0.32425543,
        -1.34257206, -0.93801444]

```

```

layer output
[array([[0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       ...,
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.]], dtype=float32), array([[0.91138043, 0.69310224, 0.10483677, ..., 0.55000926, 0.31893896,
        0.76332788],
       [0.70019104, 0.85667295, 0.53400526, ..., 0.58036098, 0.2070874 ,

```

● Back propagation

Using the backpropagation algorithm, calculate the gradient of each layer connected W, because:

$$\frac{\partial C}{\partial w_{ij}^l} = \frac{\partial C}{\partial z_j^l} \frac{\partial z_j^l}{\partial w_{ij}^l}$$

The key is to find $\frac{\partial C}{\partial z_j^l}$ and make it equal to δ_j^l

➤ Output layer

$$\frac{\partial C}{\partial z_j^L} = \frac{\partial C}{\partial y_j^L} \frac{\partial y_j^L}{\partial z_j^L} = \nabla C(y^L) * \delta'(z_j^L)$$

➤ Un-output layer

$$\frac{\partial C}{\partial z_j^l} = \sum_k \frac{\partial C}{\partial z_k^{l+1}} \frac{\partial z_k^{l+1}}{\partial a_j^l} \frac{\partial a_j^l}{\partial z_j^l} = \sum_k \delta_k^{l+1} \omega_{jk}^{l+1} \delta'(z_j^l)$$

Code:

```

1.     # Back propagation
2.     def back(layer_output, layer_input, w, y, L1_reg, L2_reg):
3.         grad_w = [np.zeros(weight.shape) for weight in w]

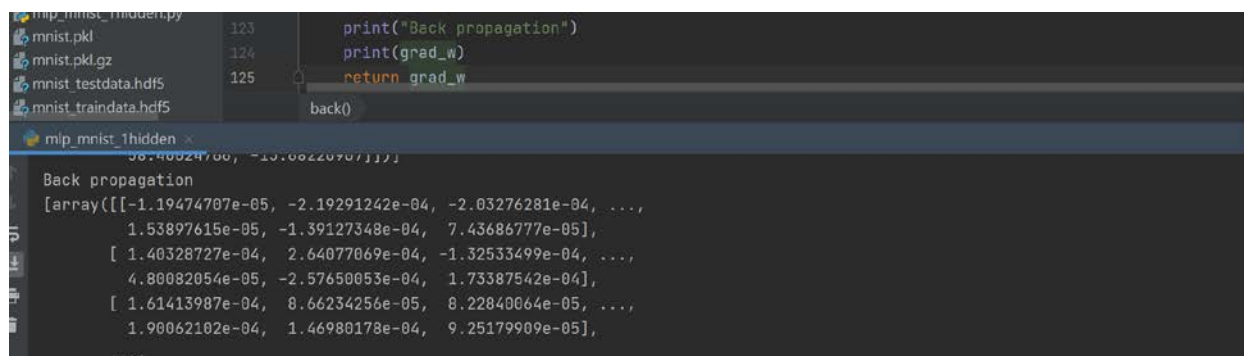
```

```

4.
5.     # Backpropagation
6.     delta = layer_output[-1] - y # last layer, m*10
7.     grad_w[-1] = np.dot(np.transpose(layer_output[-2]), delta)
8.     for l in range(2, len(w) + 1):
9.         delta = np.dot(delta, w[-
10.         l + 1].transpose()) * sigmoid_derivative(layer_input[-l]) # other layer
11.         grad_w[-l] = np.dot(np.transpose(layer_output[-
12.         l - 1]), delta) / y.shape[0] + L1_reg * sign(w[-l]) + L2_reg * w[
13.         -l]
14.     return grad_w

```

Result:



```

Back propagation
[array([[ -1.19474707e-05, -2.19291242e-04, -2.03276281e-04, ...,
         1.53897615e-05, -1.39127348e-04,  7.43686777e-05],
        [ 1.40328727e-04,  2.64077069e-04, -1.32533499e-04, ...,
         4.80082054e-05, -2.57650053e-04,  1.73387542e-04],
        [ 1.61413987e-04,  8.66234256e-05,  8.22840064e-05, ...,
         1.90062102e-04,  1.46980178e-04,  9.25179909e-05],
        ...,

```

After calculating the gradient, update w in the opposite direction of the gradient

$$\omega^{n+1} = \omega^n - \alpha \nabla C(\omega^n)$$

Code:

```

1.     # Parameter update
2.     def gradient_update(y_truth_matrix, w_current, layer_output, layer_input, alpha, L1_
3.     reg, L2_reg):
4.         if len(w_current) > 1:
5.             # Calculate the gradient using backpropagation
6.             grad_w = back(layer_output, layer_input, w_current, y_truth_matrix, L1_reg,
7.             L2_reg)
8.             # Parameter update
9.             new_w = [w - nw * alpha for w, nw in zip(w_current, grad_w)]
10.
11.     return new_w

```

Result:

```
mnist_testdata.hdf5 133 new_w = [w - nw * alpha for w, nw in zip(w_current, grad_w)]
mnist_traindata.hdf5 134 print("Parameter update")
process_dataset.py 135 print(new_w)
t10k-images-idx3-ubyte.gz 136 return new_w
t10k-labels-idx1-ubyte.gz 137
test_softmaxRegression_mlp gradient_update()

mlp_mnist_hidden x
Parameter update
[array([[ 0.25107862, -0.00288859, -0.24436217, ...,  0.07747148,
        -0.09842982, -0.25721582],
       [ 0.22615457, -0.11934802,  0.03936936, ..., -0.10211156,
        0.14714912,  0.04290086],
       [-0.16622887,  0.03450829, -0.14240778, ..., -0.16318103,
```

● Gradient descent

Code:

```
1. # Gradient descent
2. def gradient_descent(x, y, w_initial, alpha, L1_reg, L2_reg):
3.     w = w_initial
4.
5.     y_truth_matrix = truth_label_matrix(y)
6.
7.     # Forward propagation to get model prediction value
8.     layer_output, layer_input = forward(x, w)
9.     probabilities = layer_output[-1]
10.    if L1_reg != 0.00:
11.        L1 = sum([abs(w_i).sum() for w_i in w]) # L1 norm
12.    else:
13.        L1 = 0
14.    if L2_reg != 0.00:
15.        L2_sqr = sum([(w_i ** 2).sum() for w_i in w]) # L2 norm
16.    else:
17.        L2_sqr = 0
18.
19.    # Total loss = negative log likelihood cost + regularization penalty
20.    cost = loss_function(probabilities, y) + L1_reg * L1 + 0.5 * L2_reg * L2_sqr #
Error before weight w update
21.
22.    # Update parameters
23.    w = gradient_update(y_truth_matrix, w, layer_output, layer_input, alpha, L1_reg,
L2_reg)
24.
25.    return (w, cost)
```

Result:

```
mnist_testdata.hdf5 160 print("Gradient descent w:")
mnist_traindata.hdf5 161 print(w)
process_dataset.py 162 print("Gradient descent cost:")
t10k-images-idx3-ubyte.gz 163 print(cost)
t10k-labels-idx1-ubyte.gz 164 return (w, cost)
test_softmaxRegression_mlp gradient_descent()

mlp_mnist_1hidden x
0.99777242, 0.20002002]]]]
Gradient descent cost:
25.28182519217677
Gradient descent w:
[array([[ 0.18564151,  0.14414483,  0.05545726, ..., -0.05901251,
         -0.20489878,  0.11700711],
        [ 0.17715691,  0.00707062,  0.00777012, ...,  0.05004444,
```

- Validation and testing model

Code:

```
1. def model_validation(valid_set_x, valid_set_y, W):
2.     p_y_given_valid_x = forward(valid_set_x, W)[0][-1]
3.     valid_y_pred = pred_num_lable(p_y_given_valid_x)
4.     cost_valid = loss_function(p_y_given_valid_x, valid_set_y)
5.     error_num_valid = error_num(valid_y_pred, valid_set_y)
6.     return (cost_valid, error_num_valid)
7.
8. def model_test(test_set_x, test_set_y, W):
9.
10.    p_y_given_test_x = forward(test_set_x, W)[0][-1]
11.    test_y_pred = pred_num_lable(p_y_given_test_x)
12.    error_num_test = error_num(test_y_pred, test_set_y)
13.    test_precision = 1 - error_num_test / test_set_x.shape[0]
14.    return test_precision
```

- Model construction and optimization

Code:

```

176
177 def model_build(datasets, w_initial, alpha, epochs, threshold, batch_size, L1_reg, L2_reg):
178     W = w_initial
179
180     # Partition data set
181     train_set_x, train_set_y = datasets[0]
182     valid_set_x, valid_set_y = datasets[1]
183     test_set_x, test_set_y = datasets[2]
184
185     # Divide the data set into smaller batches
186     n_train_batches = train_set_x.shape[0] // batch_size
187
188     # train model
189     best_valid_cost = 0
190     # Gradient descent iterative validation_frequency times,
191     # verify the model performance once on the validation set
192     validation_frequency = 100
193     epoch = 0
194     done_looping = False
195     while (epoch < epochs) and (not done_looping):
196         epoch += 1
197         for batch_index in range(n_train_batches):
198             x = train_set_x[batch_index * batch_size: (batch_index + 1) * batch_size] # X per training
199             y = train_set_y[batch_index * batch_size: (batch_index + 1) * batch_size] # Corresponding y
200
201             W, cost_train = gradient_descent(x, y, W, alpha, L1_reg, L2_reg)

```

```

203     cost_valid, error_num_valid = model_validation(valid_set_x, valid_set_y, W)
204     this_validation_loss = error_num_valid / valid_set_x.shape[0] # error rate
205
206     num_iter = (epoch - 1) * n_train_batches + batch_index
207     if num_iter % validation_frequency == 0:
208         # Track performance changes on the validation set
209         result_temp = []
210         result_temp.append((1 - this_validation_loss) * 100)
211         print("After gradient descent iteration %d times, Accuracy is: %f%%"
212               % (num_iter, (1 - this_validation_loss) * 100))
213
214     # Determine whether to early stopping
215     if abs(cost_valid - best_valid_cost) < threshold:
216         done_looping = True
217         print("The error on the validation set no longer decreases, and the model training ends")
218         break
219     else:
220         best_valid_cost = cost_valid
221
222 if not done_looping:
223     print("The maximum number of epochs is reached, and the model training ends")
224
225 # model test
226 test_precision = model_test(test_set_x, test_set_y, W)
227 print("The accuracy of the model on the test set is: %f%%" % (test_precision * 100))

```

```

239
240 def mlp(datasets):
241     # Specify hyperparameters
242     alpha = 0.005 # learning rate
243     epochs = 1000 # epoch num
244     threshold = 0.00001 # Gradient descent early stop threshold
245     batch_size = 300 # batch size
246     L1_reg = 0.00 # L1 regularization parameters
247     L2_reg = 0.001 # L2 regularization parameters
248
249     # 初始化权重w
250     input_layer_num = 784 # input layer number
251     output_layer_num = 10 # output layer number
252     hidden_layer_num = [500] # hidden layer number
253     rand = np.random.RandomState(int(time.time()))
254     W = hidden_layer(input_layer_num, output_layer_num, hidden_layer_num, rand) # Initialize connection weight
255
256     # Model building and tuning
257     model_build(datasets, W, alpha, epochs, threshold, batch_size, L1_reg, L2_reg)
258

```

Result:

Train model

```

mlp_mnist_1hidden x
I:\Anaconda\python.exe C:/Users/123/Desktop/MLP/mlp_mnist_1hidden.py
After gradient descent iteration 0 times, Accuracy is: 16.300000%
After gradient descent iteration 100 times, Accuracy is: 65.250000%
After gradient descent iteration 200 times, Accuracy is: 88.310000%
After gradient descent iteration 300 times, Accuracy is: 77.190000%
After gradient descent iteration 400 times, Accuracy is: 89.310000%
After gradient descent iteration 500 times, Accuracy is: 87.650000%
After gradient descent iteration 600 times, Accuracy is: 71.620000%
After gradient descent iteration 700 times, Accuracy is: 90.840000%
After gradient descent iteration 800 times, Accuracy is: 89.680000%

```

- Final result

```

After gradient descent iteration 11500 times, Accuracy is: 94.450000%
After gradient descent iteration 11600 times, Accuracy is: 94.150000%
After gradient descent iteration 11700 times, Accuracy is: 92.440000%
The error on the validation set no longer decreases, and the model training ends
The accuracy of the model on the test set is: 94.030000%

```

2 Implement Modified Quadratic Discriminant Function (MQDF) on MNIST

The MQDF model is a classifier based on Bayesian decision theory. Its discriminant function is given as below:

$$\begin{aligned}
g_0(\mathbf{x}, \omega_i) &= -(\mathbf{x} - \mu_i)^T \Sigma_i^{-1} (\mathbf{x} - \mu_i) - \log |\Sigma_i| \\
&= -[\Phi_i^T (\mathbf{x} - \mu_i)]^T \Lambda_i^{-1} \Phi_i^T (\mathbf{x} - \mu_i) - \log |\Sigma_i| \\
&= -\sum_{j=1}^d \frac{1}{\lambda_{ij}} [(\mathbf{x} - \mu_i)^T \phi_{ij}]^2 - \sum_{j=1}^d \log \lambda_{ij}
\end{aligned}
\quad \Sigma_i = \Phi_i \Lambda_i \Phi_i^T$$

Design and implement a MQDF model to achieve the classification task on MNIST. Show the training process and classification results. The work in [1] can be used as reference for your work.

- Load data

```

10
11 def load_data(scaled=False):
12
13     print('Loading dataset...')
14     data_class_list = set()
15     train_x = []
16     train_y = []
17
18     with open('mnist_train.csv', 'r', encoding="utf-8") as f:
19         reader = csv.reader(f)
20         for row in reader:
21             if row == []:
22                 continue
23             train_y.append(row[-1])
24             data_class_list.add(int(row[-1]))
25             train_x.append(row[:-1])
26
27     test_x = []
28     test_y = []
29     with open('mnist_test.csv', 'r', encoding="utf-8") as f:
30         reader = csv.reader(f)
31         for row in reader:
32             if row == []:
33                 continue
34             test_y.append(row[-1])
35
36             data_class_list.add(int(row[-1]))
37             test_x.append(row[:-1])
38
39     class_num = len(data_class_list)
40     feature_num = len(train_x[0])
41
42     train_x = np.array(train_x, np.float64)
43     train_y = np.array(train_y, np.int)
44     test_x = np.array(test_x, np.float64)
45     test_y = np.array(test_y, np.int)
46
47     if scaled:
48         train_length = len(train_x)
49         x_whole = np.vstack((train_x, test_x))
50         from sklearn.preprocessing import scale
51         x_whole = scale(x_whole, axis=0, with_std=True)
52
53         train_x = x_whole[:train_length, :]
54         test_x = x_whole[train_length:, :]
55
56     print("Load data end!")
57     return class_num, feature_num, train_x, train_y, test_x, test_y

```


Result:

```

156 print("Load data")
157 print(class_num)
158 print(feature_num)
159 print(train_x)
160 print(train_y)
161 print(test_x)
162 print(test_y)
163
if __name__ == "__main__":

```

MQDF x

```

Load data
10
784
[[0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]
 ...
 [0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]]
[5 0 4 ... 5 6 8]
[[0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]
 [0. 0. 0. ... 0. 0. 0.]
 ...
 [0. 0. 0. ... 0. 0. 0.]

```

- K-L decomposition

Bayes decision theory is to judge the input sample into the category with the largest posterior probability (Maximum A Posterior, MAP). The quadratic discriminant function (QDF) is the specific manifestation of Bayes decision when the input data meets the Gaussian distribution.

According to Bayes' criterion, the posterior probability of the input sample x belonging to the category is:

$$P(\omega_i|x) = \frac{P(\omega_i)p(x|\omega_i)}{p(x)} \quad 1-1$$

The Bayes criterion can become:

$$x \in \omega_k = \arg \min g_{qdf}(x, \omega_i) \quad 1-2$$

among them :

$$g_{qdf}(x, \omega_i) = (x - \mu_i)^t \Sigma_i^{-1} (x - \mu_i) + \ln |\Sigma_i| \quad 1-3$$

Using K-L decomposition, the covariance matrix Σ_i can be decomposed into the following form:

$$\Sigma_i = P_i \Lambda_i P_i^t \quad 1-4$$

Among them $\Lambda_i = \text{diag}(\lambda_{i,1}, \lambda_{i,2}, \dots, \lambda_{i,d})$, $P_i = (p_{i,1}, p_{i,2}, \dots, p_{i,d})$, $\lambda_{i,j}$ are the j -th eigenvalue of Σ_i , and P_i is the corresponding eigenvector. Bringing 1-4 into equation 1-3, there are:

$$g_{qdf}(x, \omega_i) = \sum_{j=1}^d \frac{[(x - \mu_i)^t p_{i,j}]^2}{\lambda_{i,j}} + \sum_{j=1}^d \ln \lambda_{i,j}$$

1 - 5

Code:

```
def MQDF(class_num, train_x, train_y, k):
    """
    Build the MQDF model
    """

    feature_num_ = len(train_x[0]) # number of features
    assert(k < feature_num_ and k > 0)

    data = []
    train_length = len(train_x)
    for i in range(class_num):
        data.append(list())

    for i in range(train_length):
        class_index = int(train_y[i])
        data[class_index].append(train_x[i])

    mean = []
    cov_matrix = []
    prior = []

    for i in range(class_num):
        data[i] = np.matrix(data[i], dtype=np.float64)
        mean.append(data[i].mean(0).T)

        # np.cov treat each row as one feature, so data[i].T has to be transposed
        cov_matrix.append(np.matrix(np.cov(data[i].T)))
        prior.append(len(data[i]) * 1.0 / train_length)

    eigenvalue_list = [] # store the first largest k eigenvalue lists of each class
    eigenvector_list = [] # the first largest k eigenvector_lists, column-wise of each class
    delta = [0] * class_num # delta for each class
    for i in range(class_num):
        covariance = cov_matrix[i]
        eig_value, eig_vector = linalg.eigh(covariance)

        # sort the eigvalues
        index_ = eig_value.argsort()
        index_ = index_[::-1] # reverse the array
        eig_value = eig_value[index_]
        eig_vector = eig_vector[:, index_]

        eigenvector_list.append(eig_vector[:, 0:k])
        eigenvalue_list.append(eig_value[:k])

    # delta via ML estimation
    delta[i] = (covariance.trace() - sum(eigenvalue_list[i])) * 1.0 / (feature_num_ - k)

    return mean, eigenvalue_list, eigenvector_list, delta
```

Result:

```
K-L  
mean  
[matrix([[0.00000000e+00],  
          [0.00000000e+00],  
          [0.00000000e+00],  
          [0.00000000e+00],  
          [0.00000000e+00],  
          [0.00000000e+00]],
```

```
eigenvalue_list
[array([567256.85692849, 407965.73262991, 257402.56682999, 215674.95059997,
        130735.3079817 , 115421.02049092, 100191.43232102, 90892.99024019]), array([5120
        59299.20586441, 40307.29571767, 37179.55454379, 29503.89482626]), array([3969
        164761.36238806, 136638.48516282, 114256.66221255, 98658.61603268]), array([3644
        127818.9159466 , 103378.13114165, 85725.8767925 , 80396.16522126]), array([3171
        134334.87352031, 113788.60959256, 100070.24951814, 88969.95391862]), array([5175
        118182.69242695, 102815.01940639, 95206.85594221, 75191.48095084]), array([4854
        120310.044074 , 110001.51105054, 100010.00715444, 80002.44044725]), array([3010
```

```
eigenvector_list
[matrix([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        ...,
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.])), matrix([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.]
```

```
delta
[matrix([[1753.8439463]]), matrix([[497.17054524]]), matrix([[2260.74667863]]), matrix([[1929.4408
```

- MQDF predict

The parameters required by the QDF classifier can be obtained from the training data according to the maximum likelihood estimation. In comparison, the training is relatively simple and direct, and can obtain high accuracy, so it is widely used in practice. But when the input feature dimensionality is high and the training data is insufficient, there will be a curse of dimensionality phenomenon, which is manifested in the QDF classifier, which is an estimation error. It can be seen from formula (1-5) that for eigenvalues, the same estimation error has a much greater impact on the result when the eigenvalue is smaller than when the eigenvalue is larger.

$$g_2(x, \omega_i) = \frac{1}{h_i^2} \left\{ \|x - \mu_i\|^2 - \sum_{j=1}^k \left(1 - \frac{h_i^2}{\lambda_i} \right) [(x - \mu_i)^t p_{ij}]^2 + \sum_{j=1}^k \ln \lambda_i + (d - k) \ln h_i^2 \right\}$$

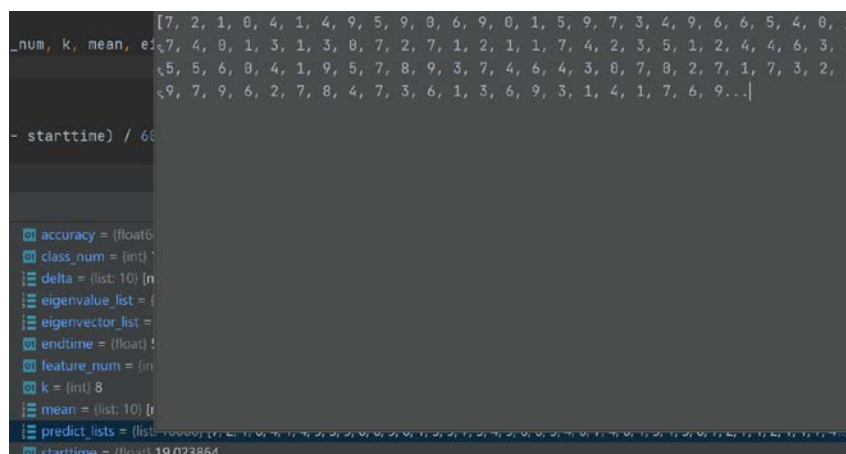
Code:

```

1. # formula
2.     for i in range(class_num):
3.
4.         minus = np.linalg.norm(x.reshape((d,)) - mean[i].reshape((d,))) ** 2
5.         matrix_minus = [0] * d
6.         for j in range(k):
7.             matrix_minus[j] = (((x - mean[i]).T * eigenvector_list[i][:, j])[0,0]
8.         ])**2
9.
10.        g = 0
11.        for j in range(k):
12.
13.            g += (matrix_minus[j] * 1.0 / eigenvalue_list[i][j])
14.
15.            g += ((minus - sum(matrix_minus)) / delta[i])
16.
17.            for j in range(k):
18.
19.                g += math.log(eigenvalue_list[i][j])
20.
21.            g += ((d - k) * math.log(delta[i]))
22.
23.        if g < min_posteriori:
24.            min_posteriori = g
25.            prediction = i

```

Result:



The screenshot shows a Jupyter Notebook interface. At the top, there is a large array of data, likely a feature matrix, with dimensions (10, 10). Below this, there is a list of parameters and variables defined in the code cell:

```

- starttime) / 60
-
-
- accuracy = (float64)
- class_num = (int) 10
- delta = (list: 10) [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
- eigenvector_list = (list: 10) [[0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0], ...]
- eigenvalue_list = (list: 10) [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
- endtime = (float64)
- feature_num = (int) 10
- k = (int) 8
- mean = (list: 10) [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
- predict_lists = (list: 10) [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
- starttime = (float64) 19.023864

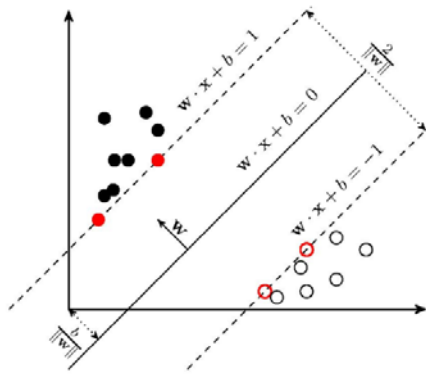
```

- Accuracy

```
Code running: 0.60min
Final correct: 94.38
```

3 Implement Support Vector Machine (SVM) on MNIST

A Support Vector Machine constructs a hyperplane or set of hyperplanes in a high- or infinite-dimensional space, which can be used for classification, regression, or other tasks.



$$\max. W(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1, j=1}^n \alpha_i \alpha_j y_i y_j K(\mathbf{x}_i, \mathbf{x}_j)$$

$$b = 1 - \frac{1}{|i : 0 < \alpha_i < C|} \sum_{i: 0 < \alpha_i < C} \sum_{j=1}^s \alpha_j y_j k(\mathbf{x}_j, \mathbf{x}_i)$$

$$C \geq \alpha_i \geq 0, \sum_{i=1}^n \alpha_i y_i = 0$$

$$\text{Polynomial Kernel: } k(\mathbf{x}_1, \mathbf{x}_2) = (\mathbf{x}_1^T \mathbf{x}_2 + 1)^d$$

$$\text{RBF Kernel: } k(\mathbf{x}_1, \mathbf{x}_2) = \exp(-||\mathbf{x}_1 - \mathbf{x}_2||^2 / 2s^2)$$

Github is a good choice for searching codes (<https://github.com>). Download the python code available in [https://github.com/prashantkh19/MNIST SVM](https://github.com/prashantkh19/MNIST_SVM). Adjust and modify the codes to implement two SVM models with Polynomial Kernel and RBF Kernel respectively on MNIST. Show the training process and classification results.

A. Principle

Describe sklearn definition and parameters of SVM.

Here we choose SVC, which means using SVM for classification. The specific use is here. About the introduction of the parameters in the code:

➤ `C = 1.0`

$$L(w, b, \xi, \alpha, \beta) := \frac{1}{2} w^T w + C \sum_{i=1}^m \xi_i + \sum_{i=1}^m \alpha_i (1 - \xi_i - y_i (w^T \phi(x_i) + b)) + \sum_{i=1}^m \beta_i (-\xi_i)$$

It can be seen that C is used to balance structural risk and experience risk.

➤ `kernel='rbf'`

Specifies the kernel type to be used in the algorithm.

It must be one of 'linear', 'poly', 'rbf', 'sigmoid', 'precomputed' or a callable.

If none is given, 'rbf' will be used. If a callable is given it is used to pre-compute the kernel matrix from data matrices; that matrix should be an array of shape ``(n_samples, n_samples)``.

➤ `gamma='scale'`

Kernel coefficient for 'rbf', 'poly' and 'sigmoid'.

if ``gamma='scale'`` (default) is passed then it uses:

$$\gamma = \frac{1}{(n_features * X.var())}$$

if 'auto', uses :

$$\frac{1}{n_features}$$

B. Code and Result

- Load data

Code:

```
1. def load_data():
2.     """
3.     Return pattern recognition data containing tuples of training data, verification
   data, and test data
4.     The training data contains 50,000 pictures, and the test data and verification d
   ata only contain 10,000 pictures
5.     """
6.     f = gzip.open('mnist.pkl.gz', 'rb')
7.     training_data, validation_data, test_data = pickle.load(f, encoding='bytes')
8.     f.close()
9.     print("load dataset successful")
10.    return (training_data, validation_data, test_data)
```

Result:

```
mnist_traindata.hdf5
process_dataset.py
t10k-images-idx3-ubyte.gz
t10k-labels-idx1-ubyte.gz
test_softmaxRegression_mlp
train-images-idx3-ubyte.gz
train-labels-idx1-ubyte.gz
weights_bias.hdf5

14     f = gzip.open('mnist.pkl.gz', 'rb')
15     training_data, validation_data, test_data = pickle.load(f, encoding='bytes')
16     f.close()
17     print("load dataset successful")
18     print(training_data)
19     return (training_data, validation_data, test_data)
20
load_data()

SVM_PI
I:\Anaconda\python.exe C:/Users/123/Desktop/SVM/SVM_PI.py
Load dataset successful
(array([[0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       ...,
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.],
       [0., 0., 0., ..., 0., 0., 0.]], dtype=float32), array([5, 0, 4, ..., 8, 4, 8], dtype=int64))
```

● SVM

Code:

```
1. # Pass the parameters of the training model
2.     clf = svm.SVC(C=100.0, kernel='rbf', gamma=0.03)
3.
4.     # model train
5.     clf.fit(training_data[0], training_data[1])
6.
7.     # test
8.     predictions = [int(a) for a in clf.predict(test_data[0])]
9.     num_correct = sum(int(a == y) for a, y in zip(predictions, test_data[1]))
```

Result:

```
I:\Anaconda\python.exe C:/Users/123/Desktop/SVM/SVM_RBF.py
start SVM
start load data
load dataset successful
num of correct
9848
SVM end
9848 of 10000 test values correct.
The accuracy rate is 0.9848.
Code running10.17min
```

- Change kernel

Code:

```
1. # Pass the parameters of the training model
2. clf = svm.SVC(C=100.0, kernel='poly', gamma=0.03)
```

Result:

4 Analysis and Comparison

Analyze and compare the advantages and disadvantages of these three models in classification from different aspects, such as effectiveness, efficiency, complexity. Your analysis and conclusion should be well justified theoretically and/or empirically.

	<i>time</i>	<i>accuracy</i>	<i>CPU utilization(i5-1021U)</i>
<i>MLP</i>	<i>42min</i>	<i>94.86</i>	<i>95%</i>
<i>MQDF</i>	<i>1.36min</i>	<i>95.53</i>	<i>30%</i>
<i>SVM RBF</i>	<i>7.96min</i>	<i>98.48</i>	<i>30%</i>
<i>SVM poly</i>	<i>2.46min</i>	<i>97.78</i>	<i>29%</i>

The MLP algorithm requires a large number of back propagation and gradient descent processes, so the gradient descent threshold becomes the main parameter that determines the time complexity of the entire algorithm. We can see that the MLP algorithm consumes a lot of time and occupies a lot of CPU because it has a large number of calculation processes and the overall space complexity is very high.

Because MQDF is a Bayesian classification algorithm, it is a kind of lazy algorithm. It does not need train and only uses calculation accuracy. Therefore, MQDF has good time complexity, runs fast, and has accuracy very high.

For SVM, the support vector is the training result of the SVM, and it is the support vector that plays a decisive role in the SVM classification decision. In essence, it avoids the traditional process from induction to deduction, realizes efficient "transduction inference" from training samples to

forecast samples, and greatly simplifies the usual classification and regression problems. The final decision function of SVM is determined by only a few support vectors, and the complexity of the calculation depends on the number of support vectors, not the dimensionality of the sample space, which avoids the "dimension disaster" in a sense. Therefore, the time consumption corresponding to different kernels is also different.

4 Comparison and improvement

A. Processing of data sets

i. PKL

Because in the program, directly importing the IDX FILE FORMAT format database is more troublesome, so the database should be preprocessed in advance and processed as Numpy format and transferred to B format.

- .pkl data file: In Python, the Pickle module converts any Python object into a system byte, similar to the JSON format, but not readable by humans.
- One-hot encoding: Use n bits to express n states, the correct state is expressed by 1, and the others are all 0. For example, 2 is represented as [0,0,1,0,0,0,0,0,0]
- """Read into the MNIST data set

Parameters

normalize: Normalize the pixel value of the image to 0.0~1.0

one_hot_label:

When one_hot_label is True, the label is returned as a one-hot array

One-hot array refers to an array like [0,0,1,0,0,0,0,0,0]

flatten: whether to expand the image into a one-dimensional array

Returns

(Training image, training label), (test image, test label)

"""

Code and Result:



```
1. # Record data set location
2. url_base = 'http://yann.lecun.com/exdb/mnist/'
3. key_file = {
4.     'train_img': 'train-images-idx3-ubyte.gz',
5.     'train_label': 'train-labels-idx1-ubyte.gz',
6.     'test_img': 't10k-images-idx3-ubyte.gz',
7.     'test_label': 't10k-labels-idx1-ubyte.gz'
8. }
```

Result:

```
[2]: # Record data set Location
url_base = 'http://yann.lecun.com/exdb/mnist/'
key_file = {
    'train_img': 'train-images-idx3-ubyte.gz',
    'train_label': 'train-labels-idx1-ubyte.gz',
    'test_img': 't10k-images-idx3-ubyte.gz',
    'test_label': 't10k-labels-idx1-ubyte.gz'
}

[4]: display(key_file)

{'train_img': 'train-images-idx3-ubyte.gz',
 'train_label': 'train-labels-idx1-ubyte.gz',
 'test_img': 't10k-images-idx3-ubyte.gz',
 'test_label': 't10k-labels-idx1-ubyte.gz'}
```



```
1. def _change_one_hot_label(X):
2.     T = np.zeros((X.size, 10))
3.     for idx, row in enumerate(T):
4.         row[X[idx]] = 1
5.
6.     return T
```

```

1. def load_mnist(normalize=True, flatten=True, one_hot_label=False):
2.
3.     if not os.path.exists(save_file):
4.         init_mnist()
5.
6.     with open(save_file, 'rb') as f:
7.         dataset = pickle.load(f)
8.
9.     if normalize:
10.        for key in ('train_img', 'test_img'):
11.            dataset[key] = dataset[key].astype(np.float32)
12.            dataset[key] /= 255.0
13.
14.    if one_hot_label:
15.        dataset['train_label'] = _change_one_hot_label(dataset['train_label'])
16.        dataset['test_label'] = _change_one_hot_label(dataset['test_label'])
17.
18.    if not flatten:
19.        for key in ('train_img', 'test_img'):
20.            dataset[key] = dataset[key].reshape(-1, 1, 28, 28)
21.
22.    return (dataset['train_img'], dataset['train_label']), (dataset['test_img'], dataset['test_label'])

```

Result:

```

Done
Converting train-labels-idx1-ubyte.gz to NumPy Array ...
Done
Converting t10k-images-idx3-ubyte.gz to NumPy Array ...
Done
Converting t10k-labels-idx1-ubyte.gz to NumPy Array ...
Done
Creating pickle file ...
Done!

```



ii. CVS

Code:

```
1. def convert(imgf, labelf, outf, n):
2.     f = open(imgf, "rb")
3.     o = open(outf, "w")
4.     l = open(labelf, "rb")
5.
6.     f.read(16)
7.     l.read(8)
8.     images = []
9.
10.    for i in range(n):
11.        image = [ord(l.read(1))]
12.        for j in range(28*28):
13.            image.append(ord(f.read(1)))
14.        images.append(image)
15.
16.    for image in images:
17.        o.write(",".join(str(pix) for pix in image)+"\n")
18.    f.close()
19.    o.close()
20.    l.close()
21.
22.    convert("MNIST/train-images.idx3-ubyte", "MNIST/train-labels.idx1-ubyte",
23.            "mnist_train.csv", 60000)
24.    convert("MNIST/t10k-images.idx3-ubyte", "MNIST/t10k-labels.idx1-ubyte",
25.            "mnist_test.csv", 10000)
26.
27.    print("Convert Finished!")
```

Result:



mnist_test.csv



mnist_train.csv

B. MLP parameter improvements

We have set certain hyperparameters in the program, we can change them to get the best prediction effect.

```
def mlp(datasets):
    # Specify hyperparameters
    alpha = 0.005      # learning rate
    epochs = 1000      # epoch num
    threshold = 0.00001 # Gradient descent early stop threshold
    batch_size = 300   # batch size
    L1_reg = 0.00      # L1 regularization parameters
    L2_reg = 0.001     # L2 regularization parameters
```

<i>alpha</i>	<i>0.001</i>	<i>0.005</i>	<i>0.01</i>	<i>0.015</i>	<i>0.02</i>	<i>0.05</i>	<i>0.1</i>
<i>accuracy</i>	<i>90.91</i>	<i>94.03</i>	<i>94.30</i>	<i>94.86</i>	<i>94.72</i>	<i>92.74</i>	<i>92.37</i>

<i>Activate function</i>	<i>sigmoid</i>	<i>tanh</i>
<i>accuracy</i>	<i>94.86</i>	<i>94.61</i>
<i>Time</i>	<i>42min</i>	<i>29min</i>

C. MQDF parameter improvements

<i>K</i>	<i>2</i>	<i>4</i>	<i>6</i>	<i>8</i>	<i>10</i>	<i>12</i>	<i>14</i>
<i>accuracy</i>	<i>89.11</i>	<i>92.27</i>	<i>93.61</i>	<i>94.38</i>	<i>94.82</i>	<i>95.07</i>	<i>94.88</i>

D. SVM parameter improvements

Ensure that other parameters remain unchanged

<i>gamma</i>	<i>0.01</i>	<i>0.02</i>	<i>0.03</i>	<i>0.04</i>	<i>0.05</i>	<i>0.06</i>
<i>accuracy</i>	<i>0.9823</i>	<i>0.9847</i>	<i>0.9848</i>	<i>0.984</i>	<i>0.9828</i>	<i>0.9804</i>

5 Conclusion

We can see that the accuracy of the three strategies is already ideal, but the time consumed by the three strategies and the size of the model are very different, depending on the internal algorithms corresponding to the different strategies. Moreover, under a certain strategy, adjusting the parameters can often achieve a relatively good predictive value.

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